



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 23, 2024 – 01:57 AM EDT

PDB ID : 4V8A
Title : The structure of thermorubin in complex with the 70S ribosome from *Thermus thermophilus*.
Authors : Bulkley, D.; Johnson, F.A.; Steitz, T.A.
Deposited on : 2011-12-05
Resolution : 3.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

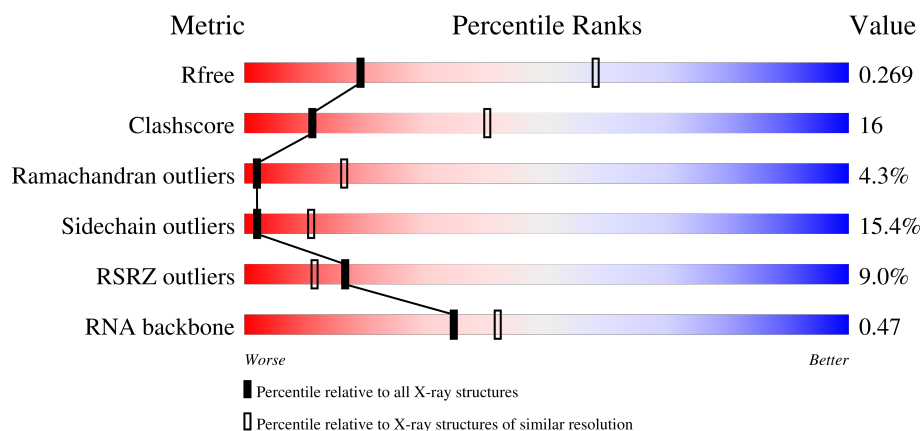
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	1370 (3.20-3.20)
Clashscore	180529	1497 (3.20-3.20)
Ramachandran outliers	177936	1479 (3.20-3.20)
Sidechain outliers	177891	1478 (3.20-3.20)
RSRZ outliers	164620	1371 (3.20-3.20)
RNA backbone	3690	1111 (3.50-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	AA	2915	
1	BA	2915	
2	AB	122	
2	BB	122	

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Mol	Chain	Length	Quality of chain
3	AD	276	
3	BD	276	
4	AE	206	
4	BE	206	
5	AF	205	
5	BF	205	
6	AG	182	
6	BG	182	
7	AH	180	
7	BH	180	
8	AI	148	
8	BI	148	
9	AN	140	
9	BN	140	
10	AO	122	
10	BO	122	
11	AP	150	
11	BP	150	
12	AQ	141	
12	BQ	141	
13	AR	118	
13	BR	118	
14	AS	112	
14	BS	112	
15	AT	146	

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Mol	Chain	Length	Quality of chain
15	BT	146	
16	AU	118	
16	BU	118	
17	AV	101	
17	BV	101	
18	AW	113	
18	BW	113	
19	AX	96	
19	BX	96	
20	AY	110	
20	BY	110	
21	AZ	206	
21	BZ	206	
22	A0	85	
22	B0	85	
23	A1	98	
23	B1	98	
24	A2	72	
24	B2	72	
25	A3	60	
25	B3	60	
26	A4	71	
26	B4	71	
27	A5	60	
27	B5	60	

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Mol	Chain	Length	Quality of chain
28	A6	54	
28	B6	54	
29	A7	49	
29	B7	49	
30	A8	65	
30	B8	65	
31	CA	1521	
31	DA	1521	
32	CB	256	
32	DB	256	
33	CC	239	
33	DC	239	
34	CD	209	
34	DD	209	
35	CE	162	
35	DE	162	
36	CF	101	
36	DF	101	
37	CG	156	
37	DG	156	
38	CH	138	
38	DH	138	
39	CI	128	
39	DI	128	
40	CJ	105	

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Mol	Chain	Length	Quality of chain
40	DJ	105	
41	CK	129	
41	DK	129	
42	CL	132	
42	DL	132	
43	CM	126	
43	DM	126	
44	CN	61	
44	DN	61	
45	CO	89	
45	DO	89	
46	CP	88	
46	DP	88	
47	CQ	105	
47	DQ	105	
48	CR	88	
48	DR	88	
49	CS	93	
49	DS	93	
50	CT	106	
50	DT	106	
51	CU	27	
51	DU	27	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
52	T8B	AA	3001	-	-	X	-
52	T8B	BA	3001	-	-	X	-

2 Entry composition

There are 53 unique types of molecules in this entry. The entry contains 279316 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a RNA chain called 23S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AA	2827	Total	C	N	O	P	0	0	0
			60900	27102	11403	19569	2826			
1	BA	2827	Total	C	N	O	P	0	0	0
			60900	27102	11403	19569	2826			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AA	272A	G	U	conflict	GB AP008226.1
BA	272A	G	U	conflict	GB AP008226.1

- Molecule 2 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	AB	120	Total	C	N	O	P	0	0	0
			2574	1146	476	833	119			
2	BB	120	Total	C	N	O	P	0	0	0
			2574	1146	476	833	119			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AB	120	G	A	conflict	GB AP008226.1
BB	120	G	A	conflict	GB AP008226.1

- Molecule 3 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	AD	275	Total	C	N	O	S	0	0	0
			2136	1349	423	361	3			
3	BD	275	Total	C	N	O	S	0	0	0
			2136	1349	423	361	3			

- Molecule 4 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	AE	204	Total	C	N	O	S	0	0	0
			1555	982	297	270	6			
4	BE	204	Total	C	N	O	S	0	0	0
			1555	982	297	270	6			

- Molecule 5 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	AF	203	Total	C	N	O	S	0	0	1
			1576	1005	297	272	2			
5	BF	203	Total	C	N	O	S	0	0	1
			1576	1005	297	272	2			

- Molecule 6 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	AG	181	Total	C	N	O	S	0	0	0
			1368	879	242	244	3			
6	BG	181	Total	C	N	O	S	0	0	0
			1368	879	242	244	3			

- Molecule 7 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	AH	174	Total	C	N	O	S	0	0	0
			1317	837	243	236	1			
7	BH	174	Total	C	N	O	S	0	0	0
			1317	837	243	236	1			

- Molecule 8 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	AI	145	Total	C	N	O	S	0	0	0
			1046	674	180	191	1			
8	BI	145	Total	C	N	O	S	0	0	0
			1046	674	180	191	1			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AI	110	GLU	ASP	conflict	UNP Q5SLQ1

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Chain	Residue	Modelled	Actual	Comment	Reference
BI	110	GLU	ASP	conflict	UNP Q5SLQ1

- Molecule 9 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	AN	140	Total	C	N	O	S	0	0	0
			1112	717	207	184	4			
9	BN	140	Total	C	N	O	S	0	0	0
			1112	717	207	184	4			

- Molecule 10 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	AO	122	Total	C	N	O	S	0	0	0
			923	583	168	168	4			
10	BO	122	Total	C	N	O	S	0	0	0
			923	583	168	168	4			

- Molecule 11 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	AP	147	Total	C	N	O	S	0	0	0
			1119	695	227	194	3			
11	BP	147	Total	C	N	O	S	0	0	0
			1119	695	227	194	3			

- Molecule 12 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	AQ	141	Total	C	N	O	S	0	0	0
			1122	715	212	188	7			
12	BQ	141	Total	C	N	O	S	0	0	0
			1122	715	212	188	7			

- Molecule 13 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	AR	118	Total	C	N	O	S	0	0	0
			968	604	203	160	1			
13	BR	118	Total	C	N	O	S	0	0	0
			968	604	203	160	1			

- Molecule 14 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
14	AS	110	Total	C	N	O	0	0	0
			865	544	172	149			
14	BS	110	Total	C	N	O	0	0	0
			865	544	172	149			

- Molecule 15 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	AT	131	Total	C	N	O	S	0	0	0
			1063	666	213	183	1			
15	BT	131	Total	C	N	O	S	0	0	0
			1063	666	213	183	1			

- Molecule 16 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	AU	116	Total	C	N	O	S	0	0	0
			959	608	201	149	1			
16	BU	116	Total	C	N	O	S	0	0	0
			959	608	201	149	1			

- Molecule 17 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	AV	101	Total	C	N	O	S	0	0	0
			771	495	140	135	1			
17	BV	101	Total	C	N	O	S	0	0	0
			771	495	140	135	1			

- Molecule 18 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
18	AW	112	Total	C	N	O	S	0	0	0
			881	554	172	153	2			
18	BW	112	Total	C	N	O	S	0	0	0
			881	554	172	153	2			

- Molecule 19 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	AX	95	Total	C	N	O	S	0	0	0
			742	483	134	124	1			
19	BX	95	Total	C	N	O	S	0	0	0
			742	483	134	124	1			

- Molecule 20 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	AY	107	Total	C	N	O	S	0	0	0
			785	503	145	131	6			
20	BY	107	Total	C	N	O	S	0	0	0
			785	503	145	131	6			

- Molecule 21 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
21	AZ	198	Total	C	N	O	S	0	0	0
			1522	972	269	279	2			
21	BZ	198	Total	C	N	O	S	0	0	0
			1522	972	269	279	2			

- Molecule 22 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	A0	76	Total	C	N	O	S	0	0	0
			594	368	125	100	1			
22	B0	76	Total	C	N	O	S	0	0	0
			594	368	125	100	1			

- Molecule 23 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
23	A1	97	Total	C	N	O	S	0	0	0
			745	469	144	131	1			
23	B1	97	Total	C	N	O	S	0	0	0
			745	469	144	131	1			

- Molecule 24 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	A2	70	Total	C	N	O	S	0	0	0
			588	365	118	103	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	B2	70	Total	C	N	O	S	0	0	0
			588	365	118	103	2			

- Molecule 25 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	A3	59	Total	C	N	O	S	0	0	0
			458	293	87	78				
25	B3	59	Total	C	N	O	S	0	0	0
			458	293	87	78				

- Molecule 26 is a protein called 50S ribosomal protein L31.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	A4	46	Total	C	N	O	S	0	0	0
			349	223	57	64	5			
26	B4	46	Total	C	N	O	S	0	0	0
			349	223	57	64	5			

- Molecule 27 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
27	A5	59	Total	C	N	O	S	0	0	0
			455	286	90	74	5			
27	B5	59	Total	C	N	O	S	0	0	0
			455	286	90	74	5			

- Molecule 28 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	A6	53	Total	C	N	O	S	0	0	0
			449	278	90	77	4			
28	B6	53	Total	C	N	O	S	0	0	0
			449	278	90	77	4			

- Molecule 29 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	A7	48	Total	C	N	O	S	0	0	0
			418	257	104	55	2			
29	B7	48	Total	C	N	O	S	0	0	0
			418	257	104	55	2			

- Molecule 30 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	A8	64	Total	C	N	O	S	0	0	0
			509	326	99	82	2			
30	B8	64	Total	C	N	O	S	0	0	0
			509	326	99	82	2			

- Molecule 31 is a RNA chain called 16S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
31	CA	1498	Total	C	N	O	P	0	0	0
			32208	14334	5974	10402	1498			
31	DA	1498	Total	C	N	O	P	0	0	0
			32208	14334	5974	10402	1498			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
CA	?	-	U	deletion	GB AP008226.1
DA	?	-	U	deletion	GB AP008226.1

- Molecule 32 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
32	CB	229	Total	C	N	O	S	0	0	0
			1777	1134	318	320	5			
32	DB	229	Total	C	N	O	S	0	0	0
			1777	1134	318	320	5			

- Molecule 33 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
33	CC	206	Total	C	N	O	S	0	0	0
			1450	906	279	264	1			
33	DC	206	Total	C	N	O	S	0	0	0
			1450	906	279	264	1			

- Molecule 34 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	CD	208	Total	C	N	O	S	0	0	0
			1520	960	283	272	5			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	DD	208	Total	C	N	O	S	0	0	0
			1520	960	283	272	5			

- Molecule 35 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
35	CE	148	Total	C	N	O	S	0	0	0
			1105	699	204	198	4			
35	DE	148	Total	C	N	O	S	0	0	0
			1105	699	204	198	4			

- Molecule 36 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
36	CF	100	Total	C	N	O	S	0	0	0
			781	495	137	146	3			
36	DF	100	Total	C	N	O	S	0	0	0
			781	495	137	146	3			

- Molecule 37 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
37	CG	155	Total	C	N	O	S	0	0	0
			1167	727	224	210	6			
37	DG	155	Total	C	N	O	S	0	0	0
			1167	727	224	210	6			

- Molecule 38 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
38	CH	138	Total	C	N	O	S	0	0	0
			1045	665	188	190	2			
38	DH	138	Total	C	N	O	S	0	0	0
			1045	665	188	190	2			

- Molecule 39 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
39	CI	125	Total	C	N	O	0	0	0
			852	533	163	156			
39	DI	125	Total	C	N	O	0	0	0
			852	533	163	156			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
CI	58	ARG	HIS	conflict	UNP P80374
DI	58	ARG	HIS	conflict	UNP P80374

- Molecule 40 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
40	CJ	96	Total	C	N	O	0	0	0
			659	408	131	120			
40	DJ	96	Total	C	N	O	0	0	0
			659	408	131	120			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
CJ	75	LEU	ILE	conflict	UNP Q5SHN7
DJ	75	LEU	ILE	conflict	UNP Q5SHN7

- Molecule 41 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
41	CK	114	Total	C	N	O	S	0	0	0
			828	516	155	154	3			
41	DK	114	Total	C	N	O	S	0	0	0
			828	516	155	154	3			

- Molecule 42 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	CL	122	Total	C	N	O	S	0	0	0
			909	570	179	159	1			
42	DL	122	Total	C	N	O	S	0	0	0
			909	570	179	159	1			

- Molecule 43 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
43	CM	114	Total	C	N	O	S	0	0	0
			801	494	164	142	1			
43	DM	114	Total	C	N	O	S	0	0	0
			801	494	164	142	1			

- Molecule 44 is a protein called 30S ribosomal protein S14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
44	CN	60	Total	C	N	O	S	0	0	0
			478	303	99	72	4			
44	DN	60	Total	C	N	O	S	0	0	0
			478	303	99	72	4			

- Molecule 45 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
45	CO	88	Total	C	N	O	S	0	0	0
			724	453	143	126	2			
45	DO	88	Total	C	N	O	S	0	0	0
			724	453	143	126	2			

- Molecule 46 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
46	CP	82	Total	C	N	O	S	0	0	0
			651	416	123	111	1			
46	DP	82	Total	C	N	O	S	0	0	0
			651	416	123	111	1			

- Molecule 47 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
47	CQ	99	Total	C	N	O	S	0	0	0
			823	528	151	142	2			
47	DQ	99	Total	C	N	O	S	0	0	0
			823	528	151	142	2			

- Molecule 48 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
48	CR	68	Total	C	N	O	0	0	0
			514	329	98	87			
48	DR	68	Total	C	N	O	0	0	0
			514	329	98	87			

- Molecule 49 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
49	CS	78	Total	C	N	O	S	0	0	0
			544	342	105	95	2			
49	DS	78	Total	C	N	O	S	0	0	0
			544	342	105	95	2			

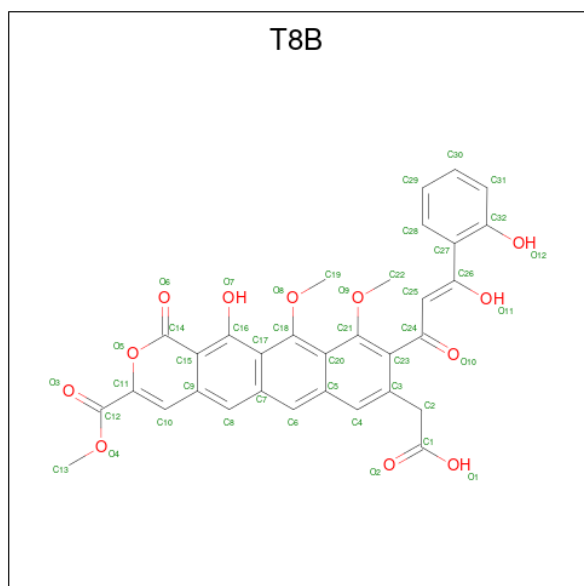
- Molecule 50 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
50	CT	96	Total	C	N	O	S	0	0	0
			708	435	151	120	2			
50	DT	96	Total	C	N	O	S	0	0	0
			708	435	151	120	2			

- Molecule 51 is a protein called 30S ribosomal protein THX.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
51	CU	23	Total	C	N	O	0	0	0
			199	122	48	29			
51	DU	23	Total	C	N	O	0	0	0
			199	122	48	29			

- Molecule 52 is Thermorubin (three-letter code: T8B) (formula: $C_{32}H_{24}O_{12}$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
52	AA	1	Total	C	O	0	0
			44	32	12		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
52	BA	1	Total	C	O	0	0
			44	32	12		

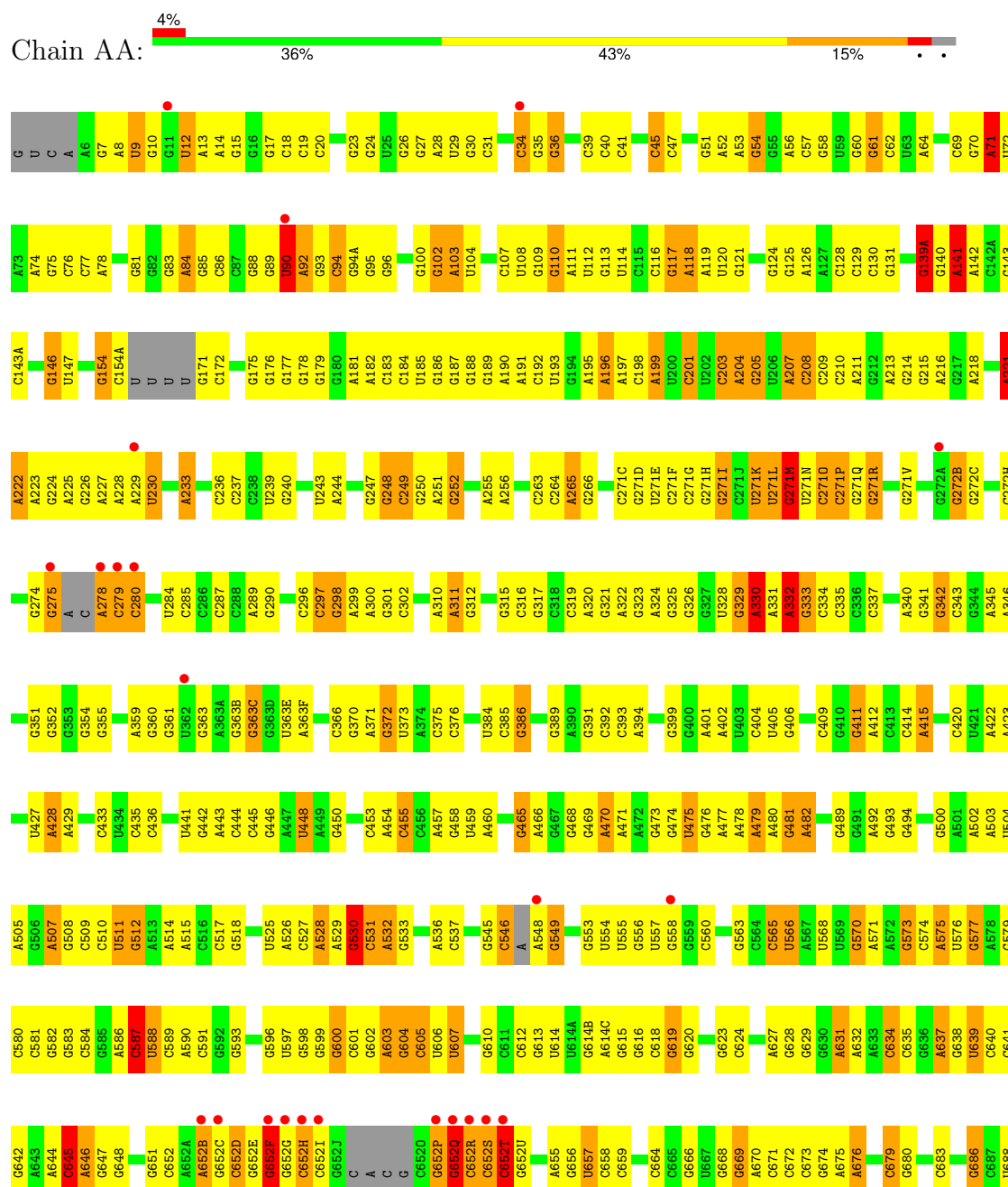
- Molecule 53 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
53	AA	2	Total	Mg	0	0
			2	2		
53	BA	2	Total	Mg	0	0
			2	2		

3 Residue-property plots

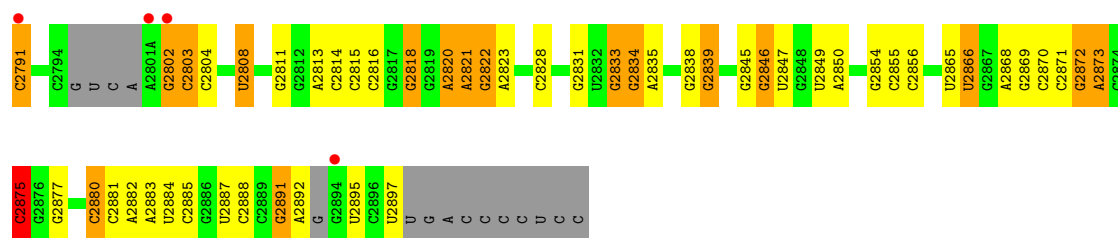
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: 23S ribosomal RNA

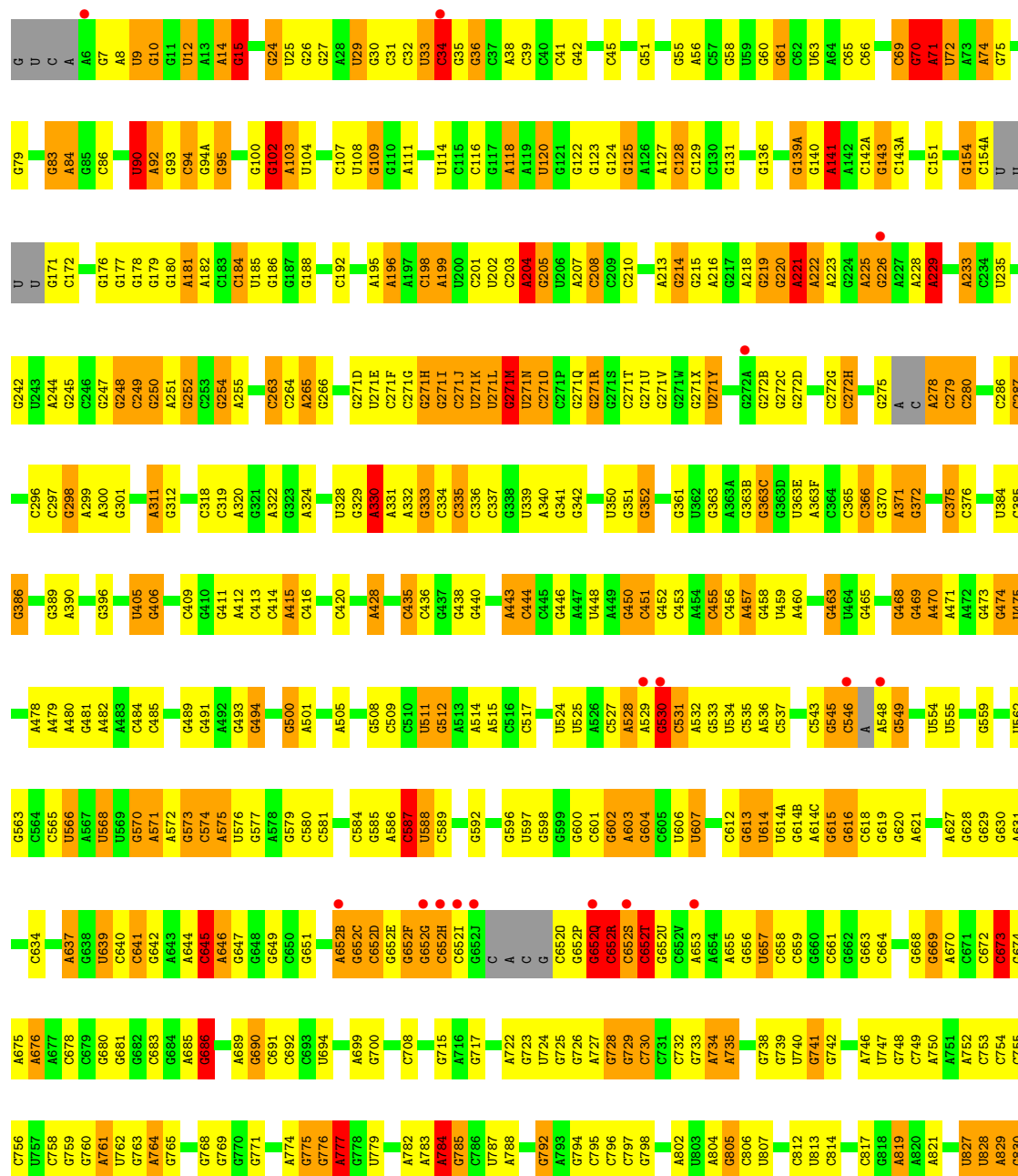


C1636	A1637	U1540	A1472	U1396	U1323	A1246	A1177	U	A1045	C975	C902	G836	G769	A659
A1638	U1639	A1542	G1473	C1399	C1327	A1247	C1178	G	A1046	G975A	C903	C837	G770	G830
C1640	A1545	A1545	G1478	C1403	G1328	U1248	C1179	G1107	A1047	C976	C904	C838	G771	C693
C1644	C1546	C1546	G1479	C1404	U1329	U1249	C1180	C1109	A1048	A774	U905	C839	A774	
G1645	C1547	C1547	U1481	U1405	C1330	C1251	G1183	C1110	C1049	G906	U907	C840	G775	G696
C1646	C1548	C1548	G1482	U1406	A1331	G1252	C1184	A1111	A1050	C908	C908	C841	G776	C697
G1647	G1484	G1484	G1484	C1407	G1332	A1253	C1185	A1112	C1051	A909	A910	C842	A777	C698
C1648	A1485	C1408	G1485	C1408	U1335	C1257	G1187	U1113	A	A983	A911	G845	G778	A699
C1650	A1486	C1409	A1486	C1409	G1336	C1257	U1188	G1114	G	C985	G912	G848	U779	G700
G1651	G1487	G1410	G1487	G1410	A1336	C1258	A1189	G1115	G	C986	G912	G849	G780	G701
A1652	G1488	C1411	G1488	C1411	G1337	G1259	G1190	G1116	A	A987	U913	A849	A781	G702
A1653	U1489	A1412	U1489	A1412	G1338	G1260	G1191	C1117	G	G988	C914	C850	A782	
A1654	A1490	G1413	A1490	G1413	G1339	C1261	G1192	C1118	G	C990	C915	G854	A783	A705
A1655	G1491	C1416	G1491	C1416	U1340	U1265	G1193	C1119	G	A990	A916	G855	A784	A706
A1656	C1492	C1417	C1492	C1417	U1341	G1266	U1198	G1120	U	C991	A917	G856	A785	A707
A1570	A1494	G1418	A1494	G1418	A1342	U1267	U1199	G1121	U	C992	A918	C857	G786	C708
A1571	A1495	G1419	A1495	G1419	G1343	U1270	U1205	G1122	G	C993	A919	C858	A787	U709
A1496	U1420	G1421	U1420	G1421	G1344	C1270	A1206	C1123	C	C994	G920	U858	A788	
U1497	G1422	G1423	G1422	G1423		G1271	G1203	G1124	U	C995	G921	U859		G715
G1500	G1423		G1500		A1354	A1274	U1204	G1125	U	A996	U922	U860	G792	A716
C1501	G1427	A1427	C1501	A1427	G1355	A1275	G1206	U1130	A	A1001	G928	U861	A792	A717
G1502	C1428	C1428	U1503	C1428	U1356	A1278		U1131	A	G1002	G932	A861	A793	A718
U1503	G1429	G1429	C1504	G1429	G1357	U1282	A1210	A1132	G	G1003	A933	A862	G794	C719
C1505	C1430	C1430	C1505	C1430	U1358	U1283	G1211	U1133	A	C1004	G938	A863	C795	C720
C1506	U1431	A1359	C1506	A1359	A1359	G1284	G1212	G1138	G	C1005	G942	A864	G796	C721
A1507	C1432	A1360	A1507	C1432	A1360	G1285	A1213	G1139	C	C1006	A941	A870	A804	G726
A1508	U1433		A1508	U1433		G1286	A1214	C1140	C	C1007	U942	G876	G805	A727
A1509	A1434		C1509	A1434		A1287	G1215	U1141	A	C1008	U943	G877	G806	G728
A1509A			A1509A		A1366	U1288	G1216	U1142	C	G944	U944	G875	U807	G729
U1514	C1437	C1437	U1514	C1437	A1367	U1288	A1220	A1143	U	C	G946	G876	C731	C730
G1515	U1438	U1438	G1515	U1438	G1368	U1292	C1221	G1144	C	U1012	G947	U877	C812	C732
C1516	C1440	C1440	C1516	C1440	G1369	U1292	C1221A		U	C1013	G948	A878	U813	C733
U1517	G1441	G1441	U1517	G1441	C1370	C1293	C1222	A1148	U	C1014	G949	G879	C814	
G1518	G1442	G1442	G1518	G1442	U1371	C1293	G1223		A	C1015	G950	G880	C815	A734
G1519	A1445	A1445	G1519	A1445	A1372	C1297	A1226	C1153	A	C1016	C951	G881	C816	A735
G1520	G1448	G1448	G1520	G1448	G1374	C1298	G1227	A1154	A	C1017	C952	G882	C817	C736
G1525	A1449	A1449	G1525	A1449	C1375	U1300	G1228	A1155	A	U1018	G953	G883	G818	
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G1527	G1455	G1455	G1527	G1455	G1377	A1308	G1230	U1159	G	A1021	U958	C885	A820	U747
A1528	G1459	G1459	A1528	G1459	A1378	A1309	G1231	G1160	U	G1022	A959	C886	A821	U748
C1529	A1460	A1460	C1529	A1460	G1380	G1310	G1232	G1161	C	G1025	C961	C887	U822	C749
C1530	C1461	C1461	C1530	C1461	G1381	G1311	G1235	U1165	G	U1026	G962	C888	G823	A752
C1531	G1462	G1462	C1531	G1462	C1382	U1312	G1236	C1166	A	A1027	U963	C889	U824	C753
C1532	C1463	C1463	C1532	C1463	C1383	U1313	A1237	U1167	U	A1028	G966	C890	U825	C754
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U	C1466	C1466	U	C1466	U1316	U1316	U1240	G	C	G1039	G969	A896	G830	A761
C	C1467	C1467	C	C1467	C1317	G1318	A1241	G1171	U	G1042	C970	C897	G831	
G1537	G1470	G1470	G1537	G1470	C1387	G1319	A1242	G1173	C	G1043	C971	C898	G832	A764
G1538	A1471	A1471	G1538	A1471	A1393	C1320	G1243	A1174	C	C1044	G974	A899	U833	G765
G1539			G1539		A1395		G1245	G1176	C			A901	A835	G768

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G2727	A2639	A2639	A2561	G2488	C2416	U2348	A2286	G2206	G2141	U2074	G1998	A1914	U1905	C1711
U2728	G2640	G2640	U2562	G2489	G2419	G2349	A2287	G2207	C2142	U2075	C1999	A1915	G1814	C1712
G2729	G2641	G2641	U2563	G2490	G2420	G2350	A2288	A2208	U2143	C2078	C2007	A1916	G1815	U1720
G2730	G2645	G2645	U2564	U2493	G2421	A2352	G2289	G2219	C2145	U2079		U1917	G1816	U1721
G2731	G2646	G2646	U2565	G2494	A2422	G2353	U2291	G2223	C2146	G2080	G2010	A1918	U1817	G1722
G2732	G2667	G2667	U2566	G2495	G2423	G2354	C2292	G2224	G2147	C2081	U2011	A1919	G1818	A1722
A2733	U2647	U2647	C2568	C2496	G2424	C2355	C2293	A2225	G2148	A2082	G2012	C1920	A1819	U1730
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G2735			C2498	U2357	G2426	U2357	U2295	A2227	U2150	C2084	A2014	A1927	A1741	G1741
G2736			C2499	G2358	G2427	G2358	C2297	A2228	G2151	G2085	A2015	G1929	G1742	G1742
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G2738			C2502	A2360	G2429	A2361	A2298		G2153		U2017	G1931	G1826	G1746
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A2749							A2310	U2243	G2165		G2029		G1839	G1763
A2750							A2311	U2244	C2166	C1943	A2030	U1946	G1840	G1764
G2751							U2312		U2167	C1944	A2031	G1947	C1844	C1765
G2752							G2313	A2247	G2168	G1948	G2032	G1948	U1766	U1766
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G2755							G2315	U2249	C2170	U2109	G2034		A1848	A1848
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A2757							G2317	G2251	U2172	C2111		U1963	G1850	C1774
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G2759							G2319	G2253	C2174	U2113	C2040	G1965	A1854	U1779
G2760							A2320		C2175	A2114	U2041	C1965	A1780	A1780
G2761							G2321	G2256	A2176	G2115	A2042	A1966	C1761	C1761
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A2765							G2324	C2261	C2179	U2118	G2049	A1969	A1784	A1784
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G2769							A2328		G2184	U2122		G1973	A1877	A1788
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A2776							G2330	A2269	G2186	G2124	G2056	G1975	A1789	A1789
G2777							G2331		C2187	G2125	A2057	U1976	C1882	C1790
A2778							U2332	U2272	G2188	A2126	A2058	A1977	G1883	A1791
							A2333	A2273	G2189	G2127			A1884	G1792
A2781							G2334	A2274	U2190	U2130	A2060	G1980	C1793	C1793
G2782							G2335	A2275	G2191	G2131	G2061	A1981	G1888	U1794
G2783							A2336	C2276	G2192	U2132	C2062	C1982	A2062	C1795
C2784							G2337	G2277	G2193	U2133	C2063	A1889	U1796	U1796
G2785							G2338			A2134	C2064	A1986	G1891	C1797
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A2713								G2280	C2200	C2136	C2067	U1991	C1800	C1800
G2714							C2342	G2281	C2201	U2068	U2068	G1992	G1801	C1801
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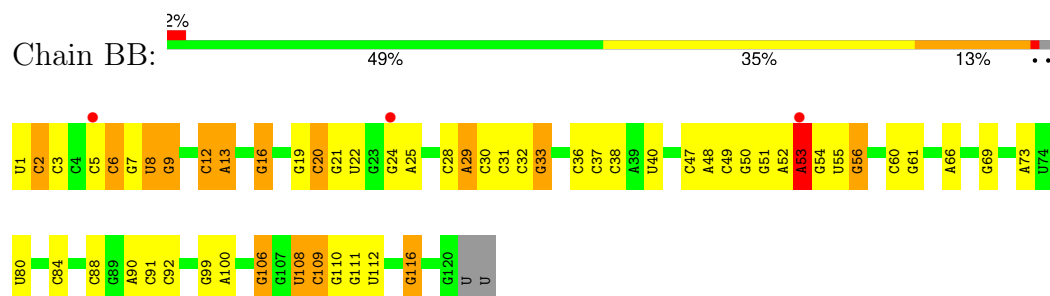


• Molecule 1: 23S ribosomal RNA

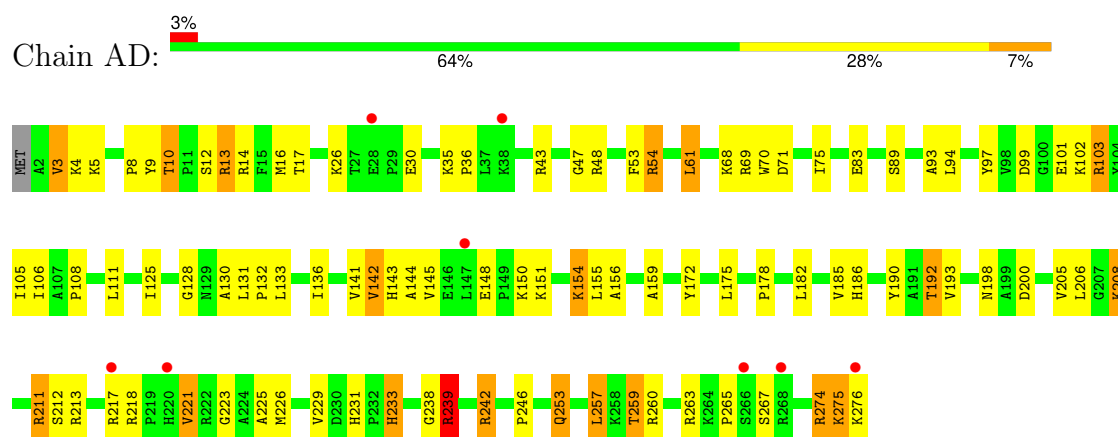




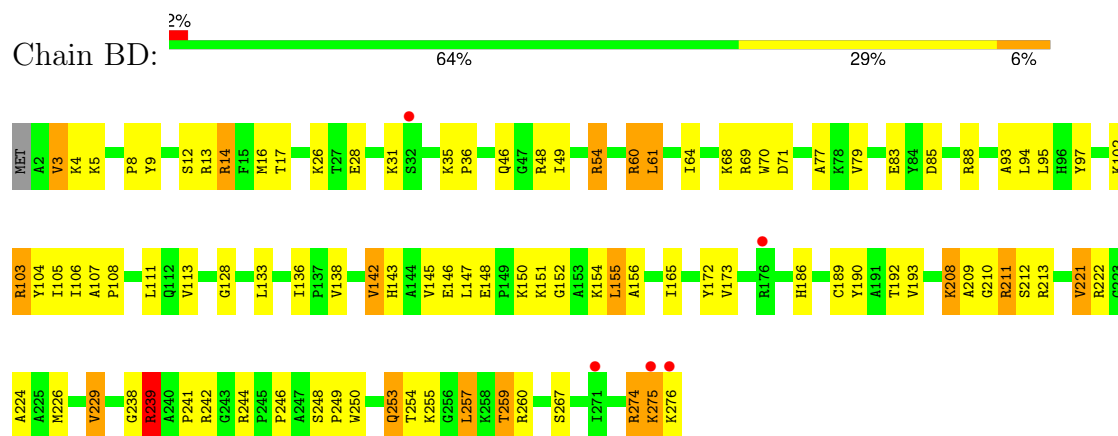
- Molecule 2: 5S ribosomal RNA



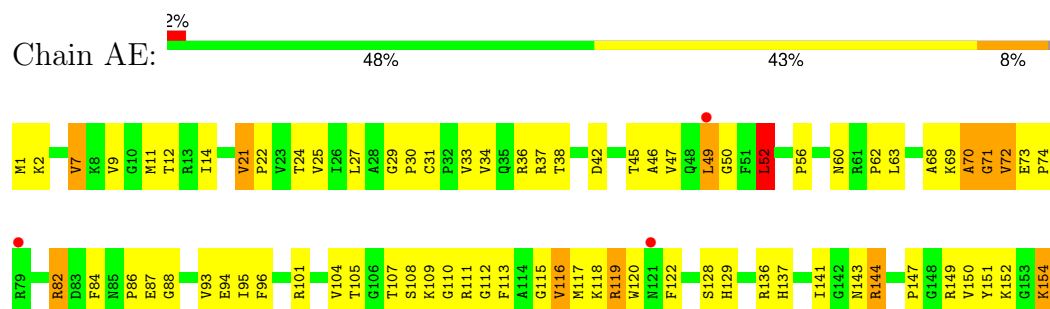
- Molecule 3: 50S ribosomal protein L2



- Molecule 3: 50S ribosomal protein L2

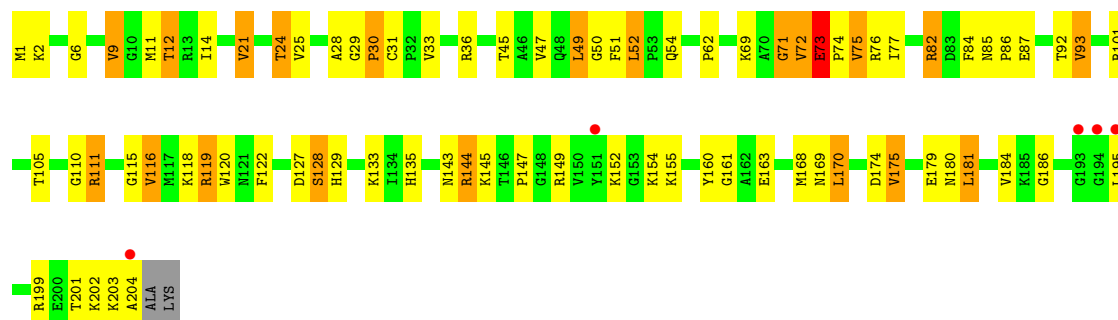


- Molecule 4: 50S ribosomal protein L3

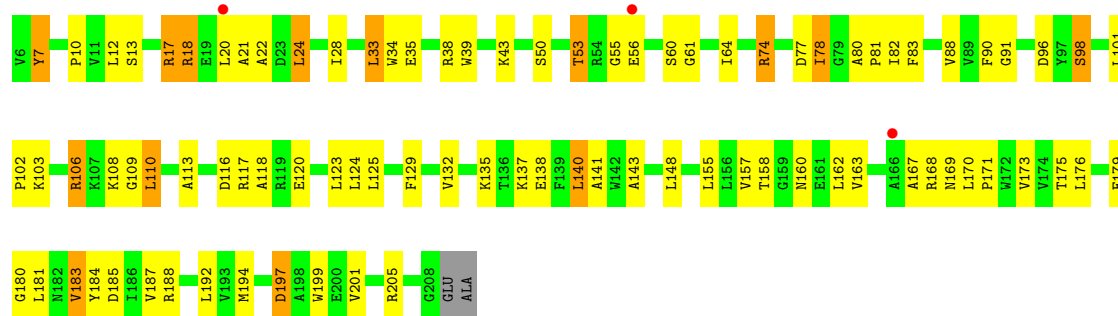




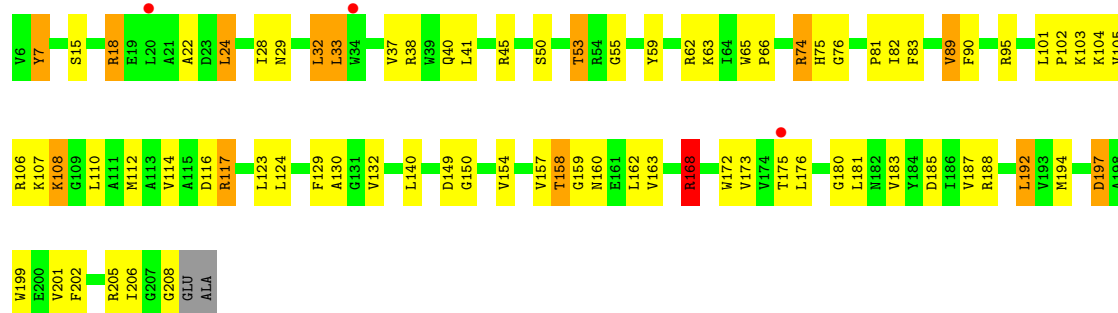
• Molecule 4: 50S ribosomal protein L3



• Molecule 5: 50S ribosomal protein L4

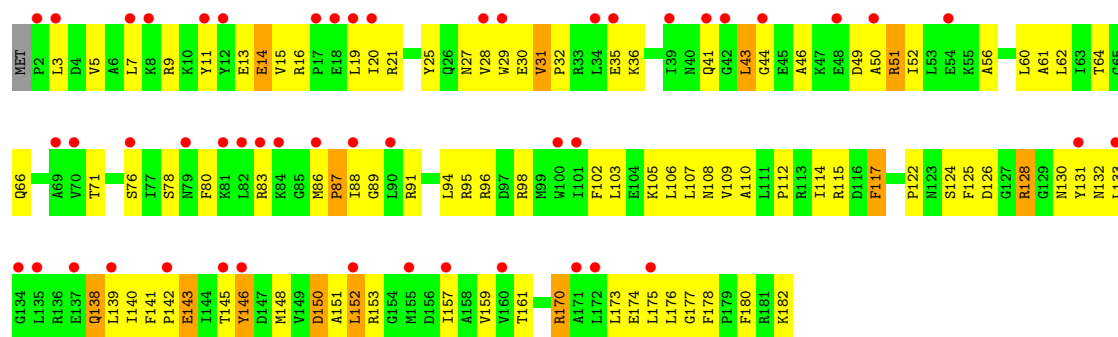


• Molecule 5: 50S ribosomal protein L4

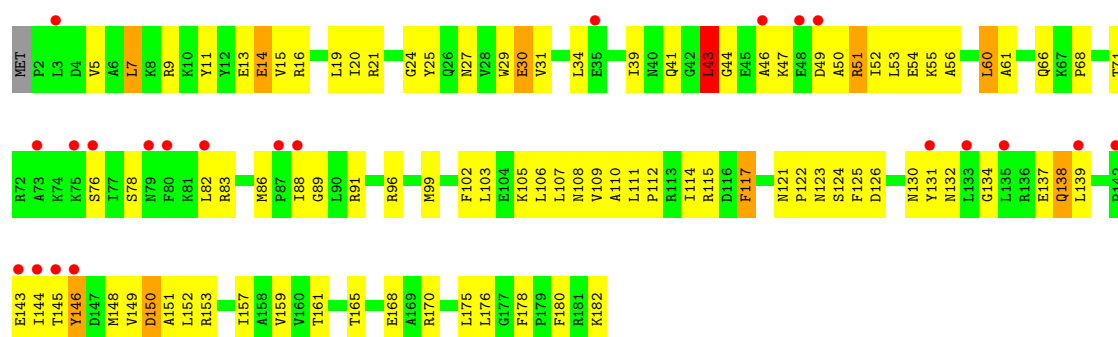


• Molecule 6: 50S ribosomal protein L5

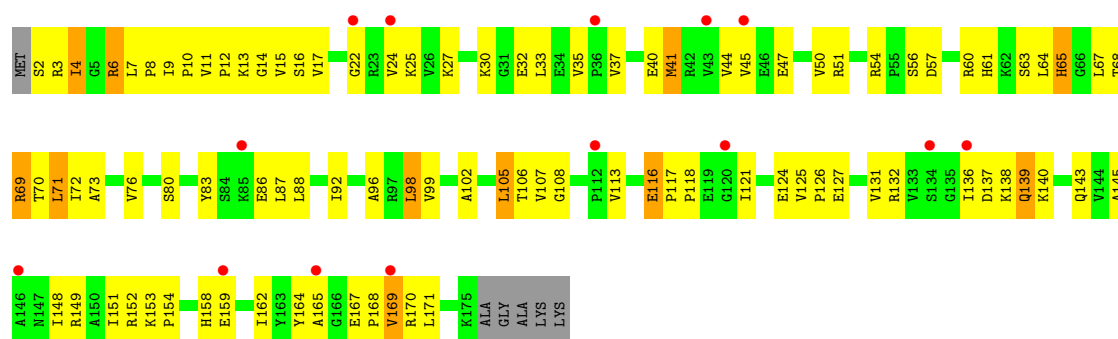
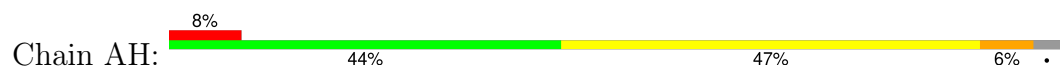




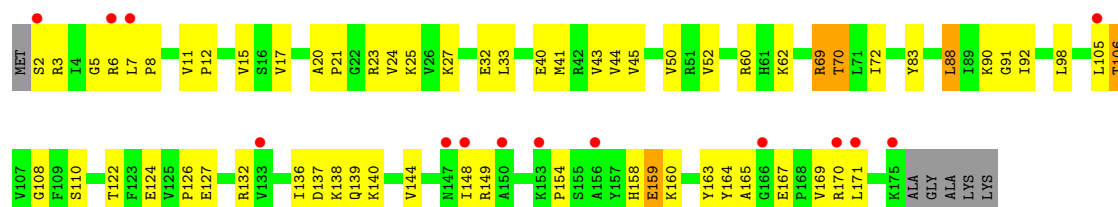
• Molecule 6: 50S ribosomal protein L5



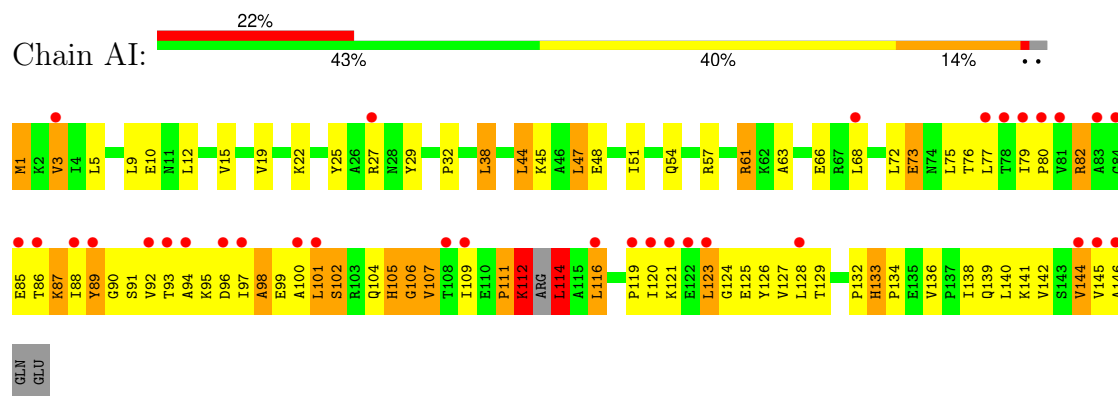
• Molecule 7: 50S ribosomal protein L6



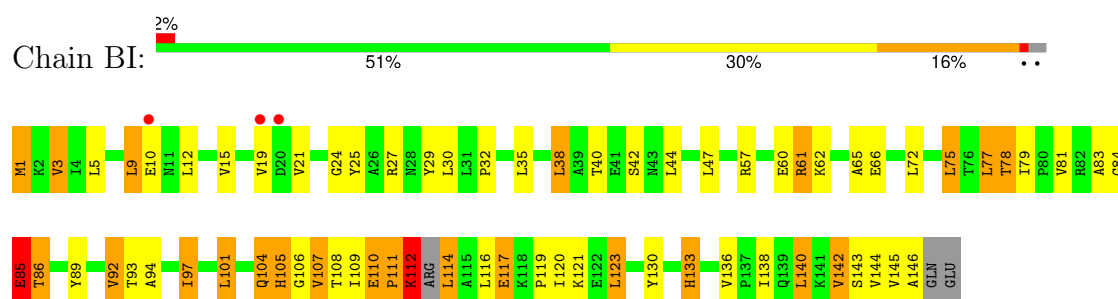
• Molecule 7: 50S ribosomal protein L6



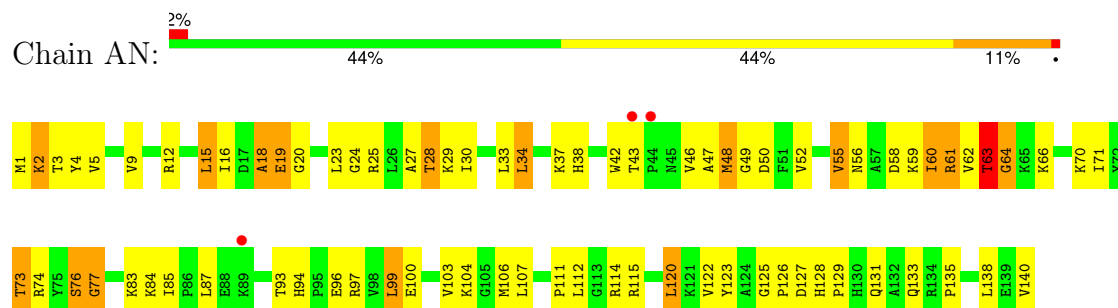
- Molecule 8: 50S ribosomal protein L9



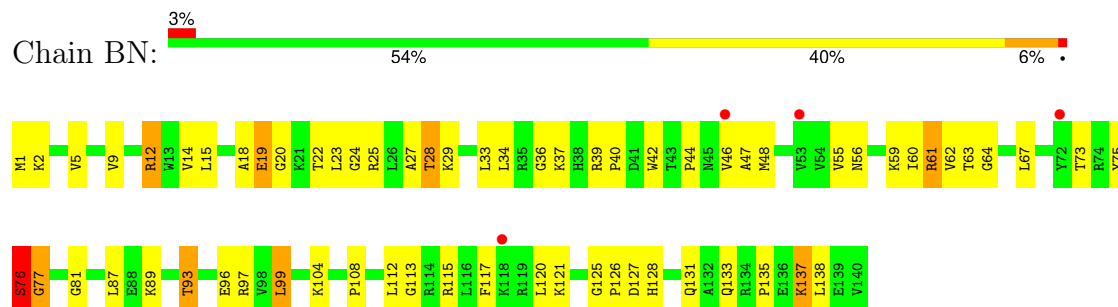
- Molecule 8: 50S ribosomal protein L9



- Molecule 9: 50S ribosomal protein L13

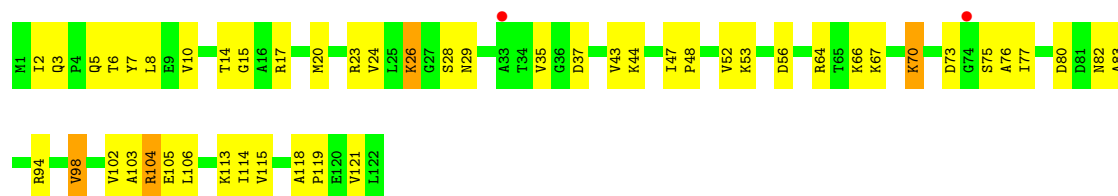


- Molecule 9: 50S ribosomal protein L13

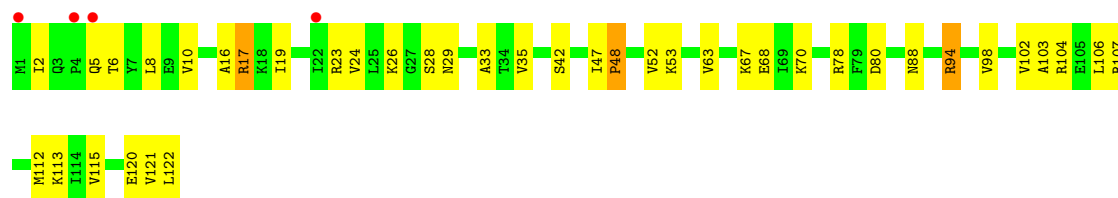


- Molecule 10: 50S ribosomal protein L14

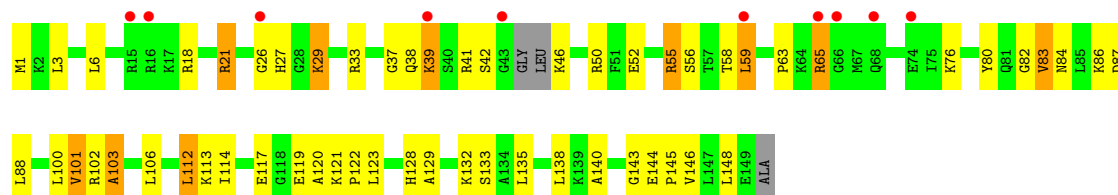




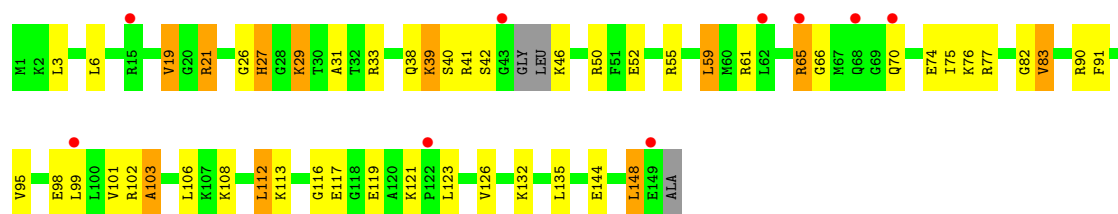
• Molecule 10: 50S ribosomal protein L14



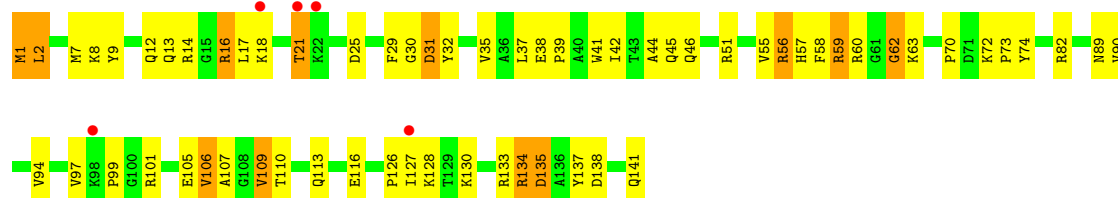
• Molecule 11: 50S ribosomal protein L15



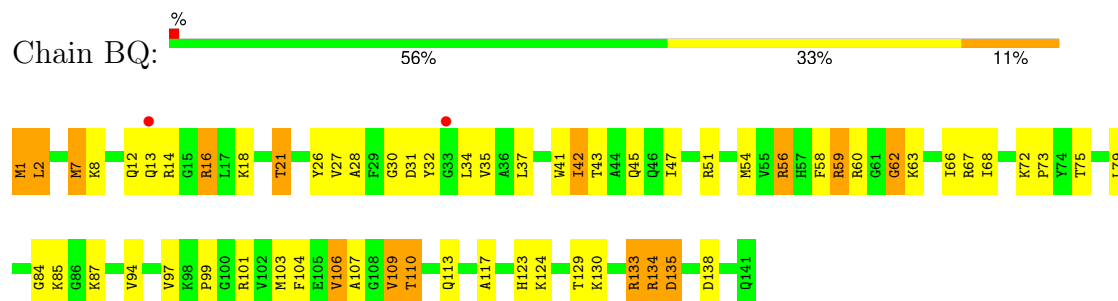
• Molecule 11: 50S ribosomal protein L15



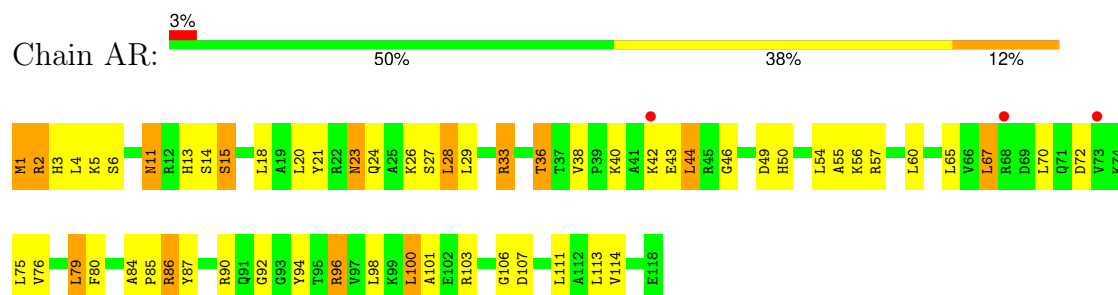
• Molecule 12: 50S ribosomal protein L16



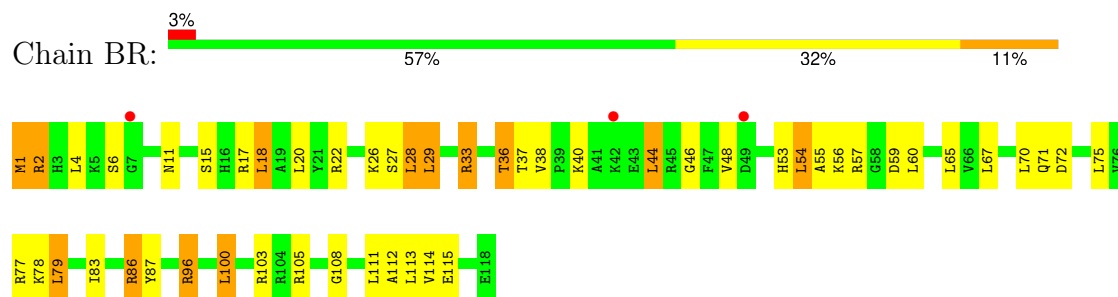
- Molecule 12: 50S ribosomal protein L16



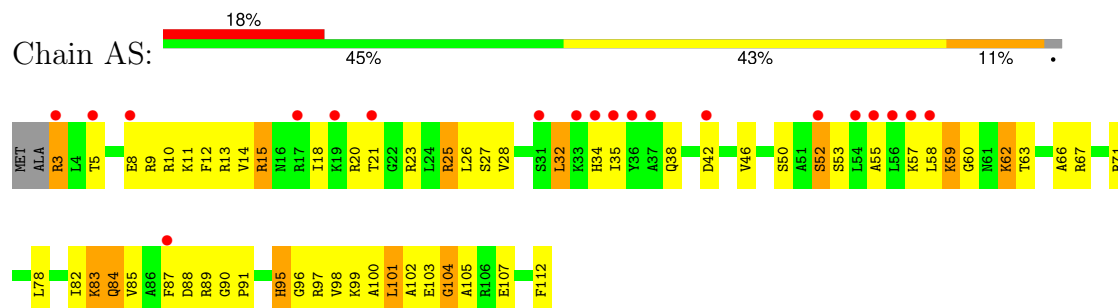
- Molecule 13: 50S ribosomal protein L17



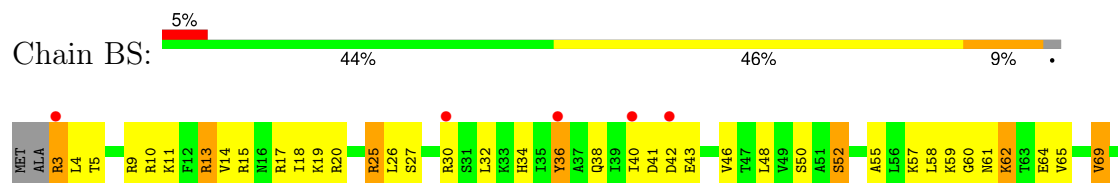
- Molecule 13: 50S ribosomal protein L17

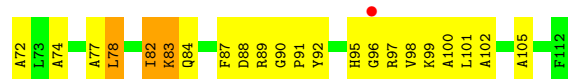


- Molecule 14: 50S ribosomal protein L18



- Molecule 14: 50S ribosomal protein L18





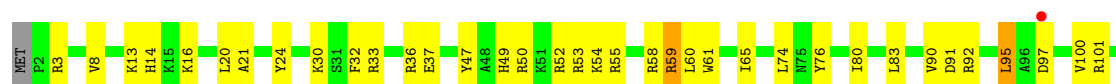
- Molecule 15: 50S ribosomal protein L19



- Molecule 15: 50S ribosomal protein L19



- Molecule 16: 50S ribosomal protein L20



- Molecule 16: 50S ribosomal protein L20

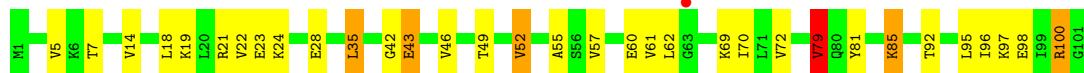


- Molecule 17: 50S ribosomal protein L21

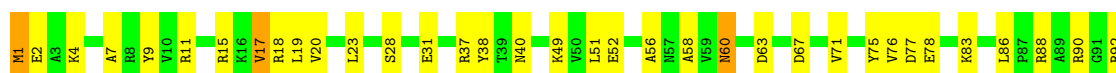




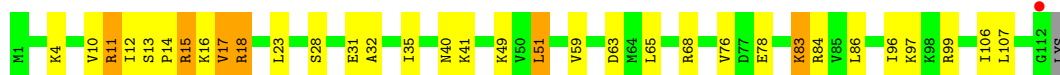
- Molecule 17: 50S ribosomal protein L21



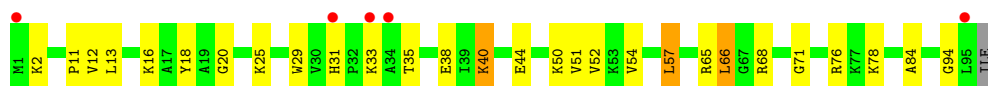
- Molecule 18: 50S ribosomal protein L22



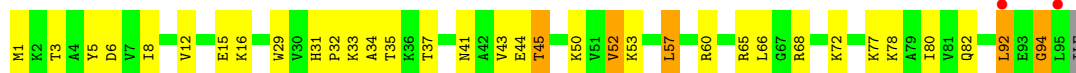
- Molecule 18: 50S ribosomal protein L22



- Molecule 19: 50S ribosomal protein L23



- Molecule 19: 50S ribosomal protein L23

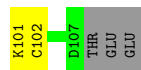


- Molecule 20: 50S ribosomal protein L24

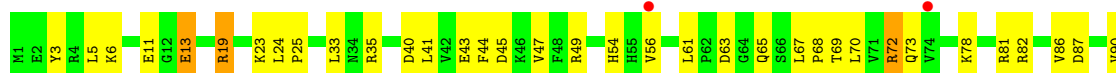




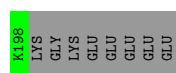
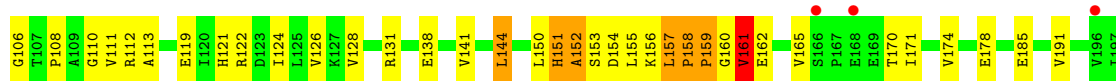
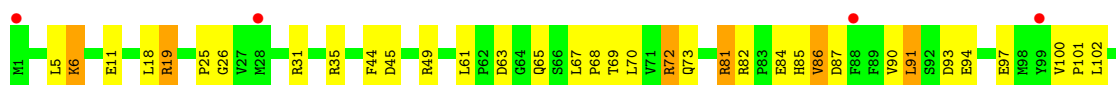
- Molecule 20: 50S ribosomal protein L24



- Molecule 21: 50S ribosomal protein L25



- Molecule 21: 50S ribosomal protein L25



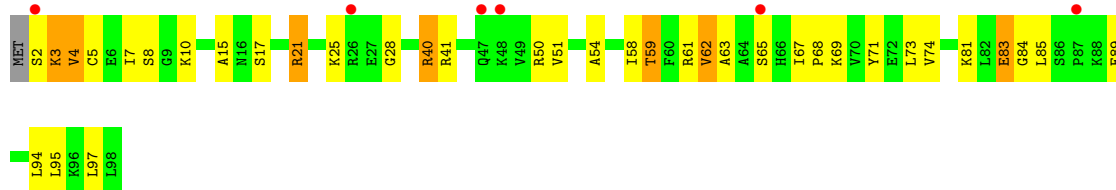
- Molecule 22: 50S ribosomal protein L27



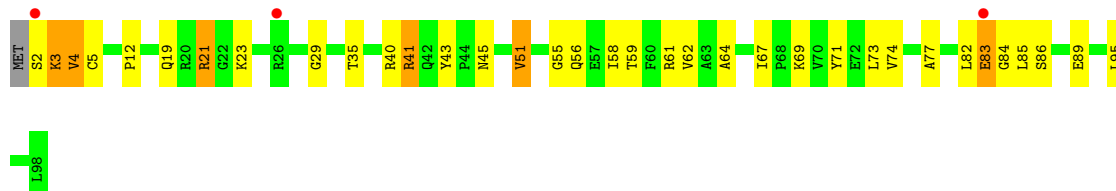
- Molecule 22: 50S ribosomal protein L27



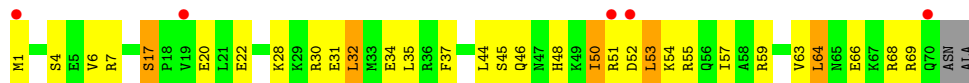
- Molecule 23: 50S ribosomal protein L28



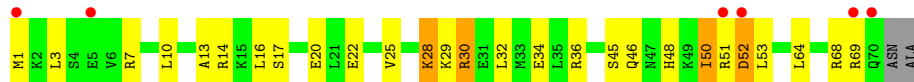
- Molecule 23: 50S ribosomal protein L28



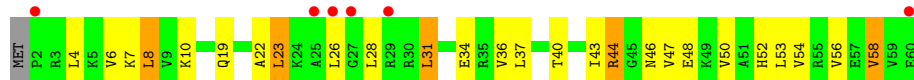
- Molecule 24: 50S ribosomal protein L29



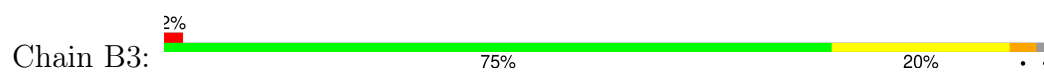
- Molecule 24: 50S ribosomal protein L29



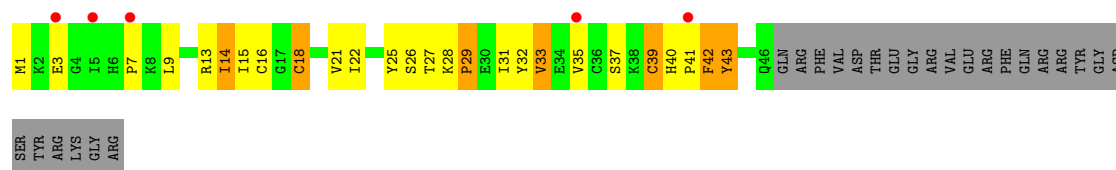
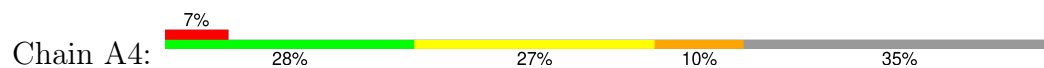
- Molecule 25: 50S ribosomal protein L30



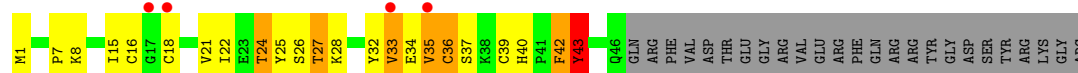
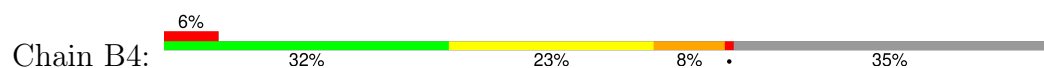
- Molecule 25: 50S ribosomal protein L30



- Molecule 26: 50S ribosomal protein L31



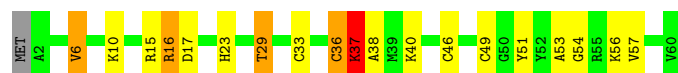
- Molecule 26: 50S ribosomal protein L31



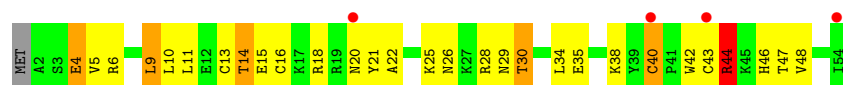
- Molecule 27: 50S ribosomal protein L32



- Molecule 27: 50S ribosomal protein L32



- Molecule 28: 50S ribosomal protein L33

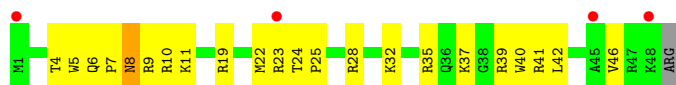


- Molecule 28: 50S ribosomal protein L33

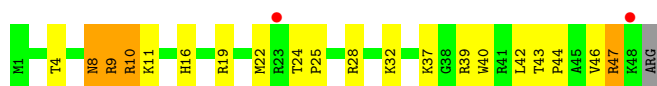




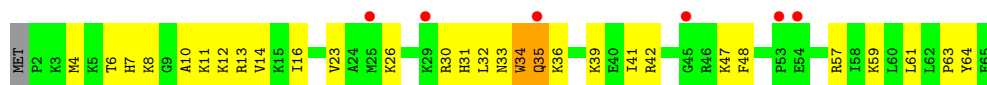
- Molecule 29: 50S ribosomal protein L34



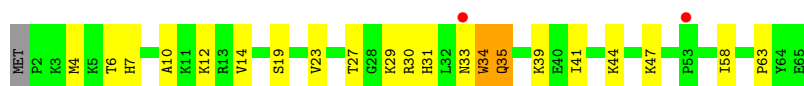
- Molecule 29: 50S ribosomal protein L34



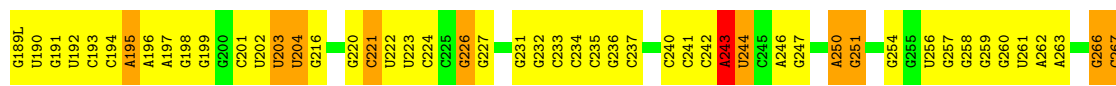
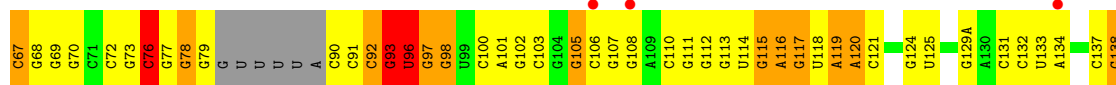
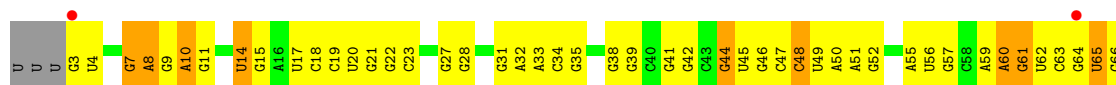
- Molecule 30: 50S ribosomal protein L35

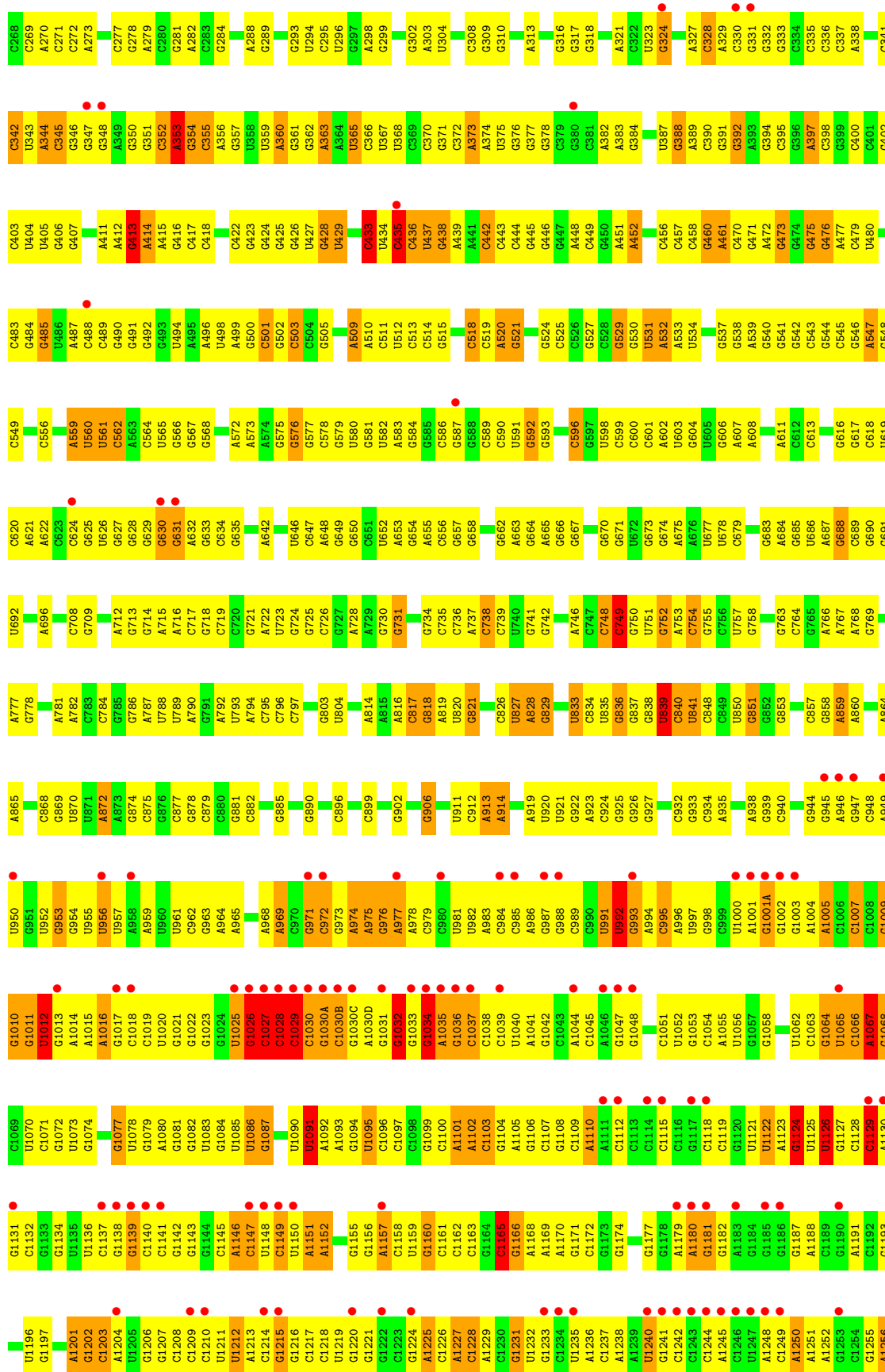


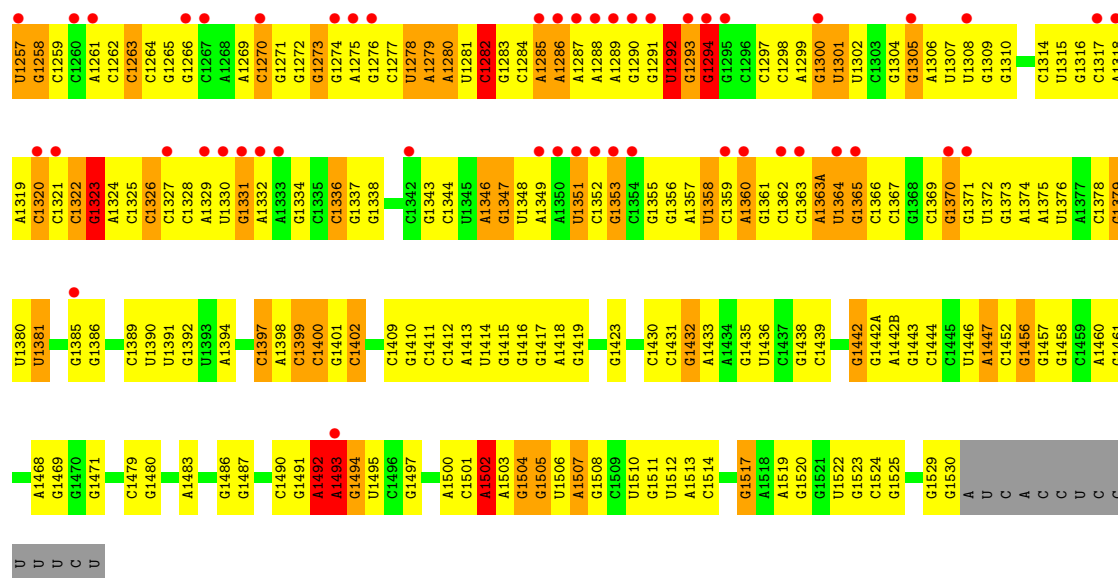
- Molecule 30: 50S ribosomal protein L35



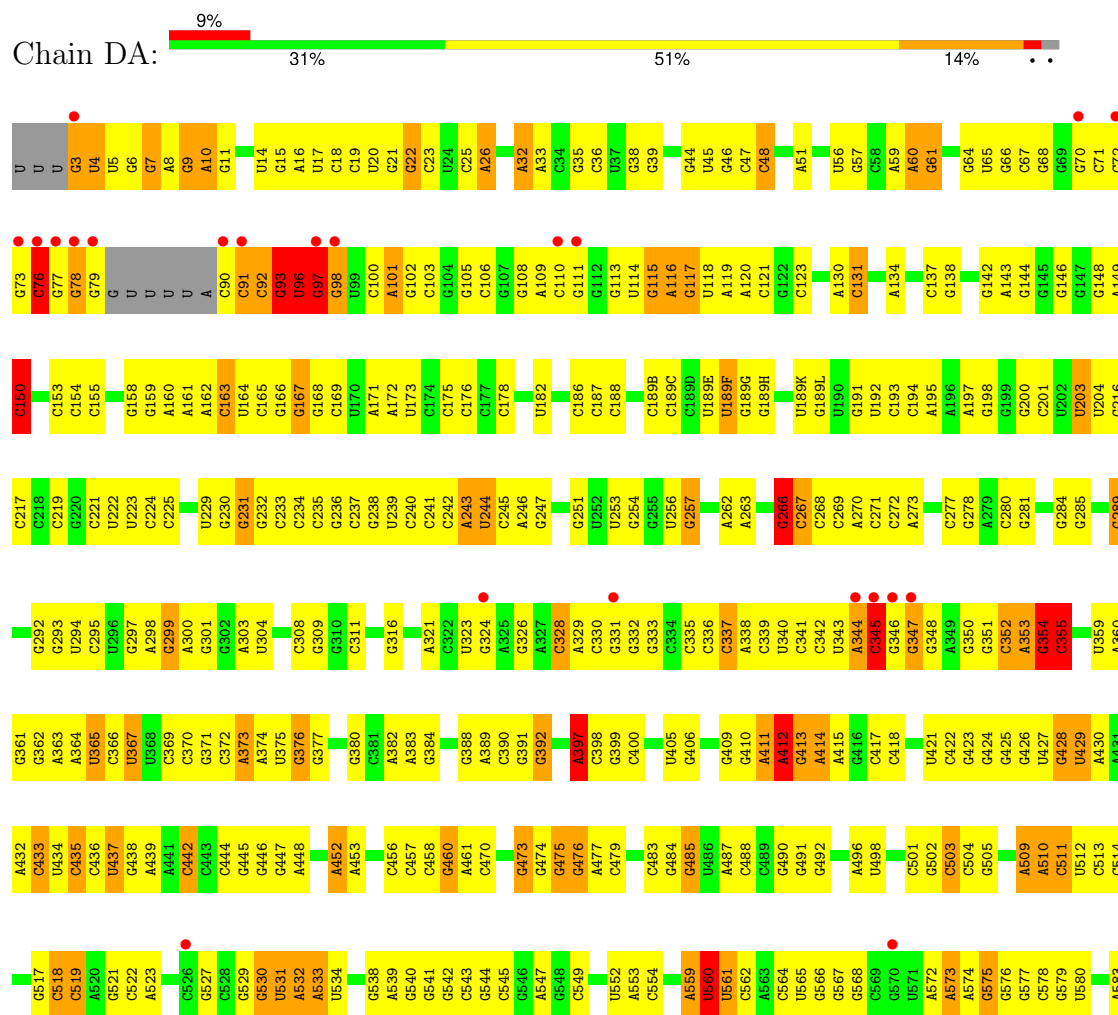
- Molecule 31: 16S ribosomal RNA

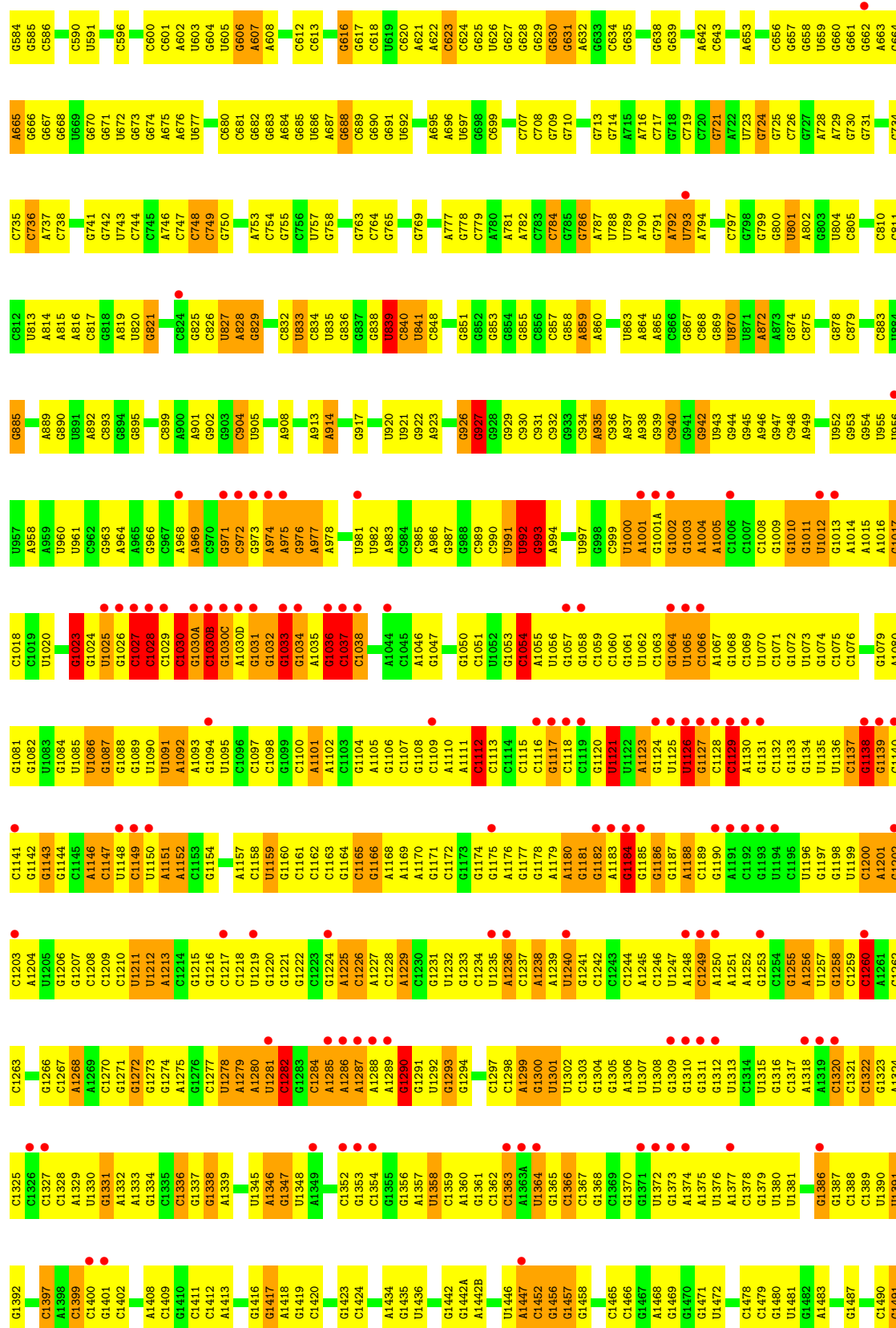






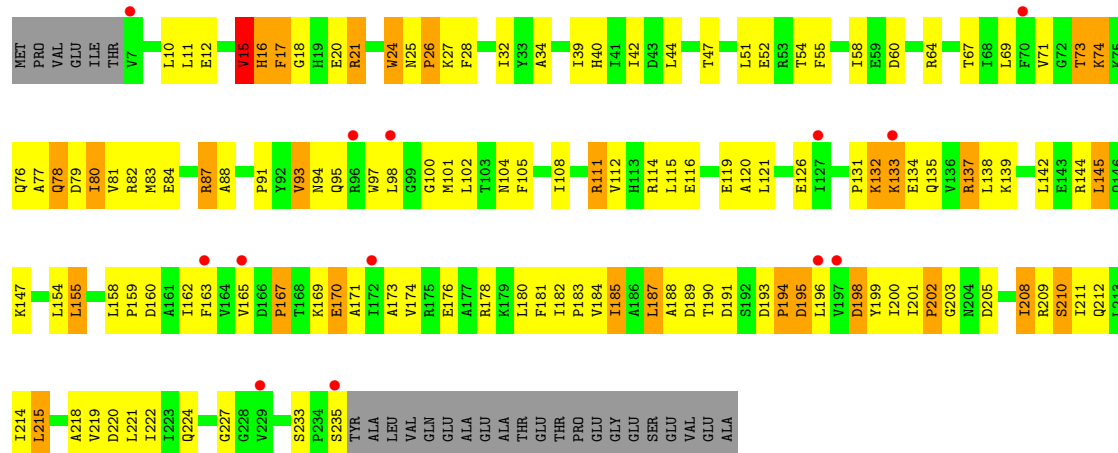
- Molecule 31: 16S ribosomal RNA



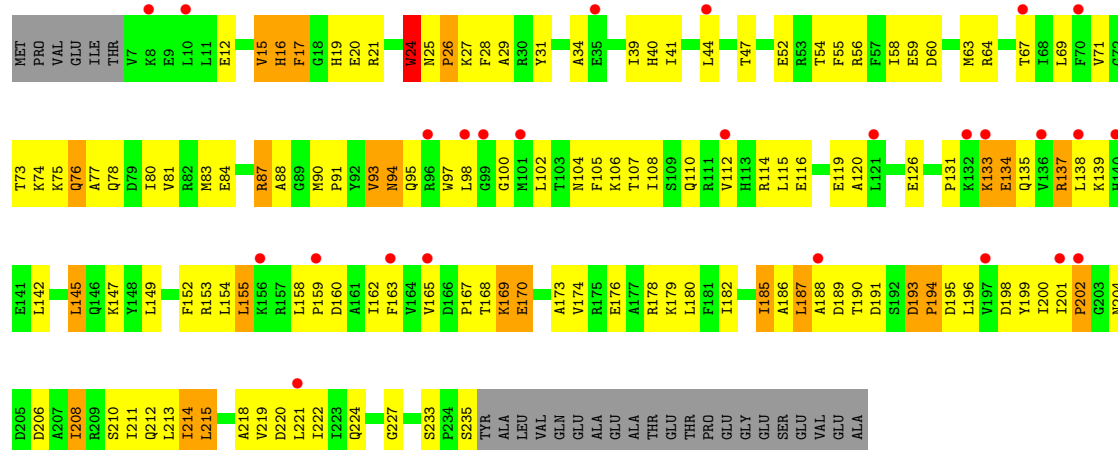
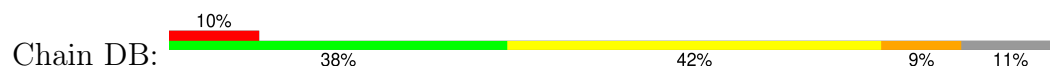




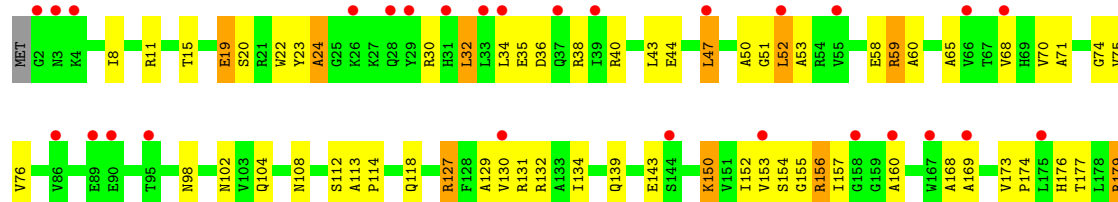
• Molecule 32: 30S ribosomal protein S2

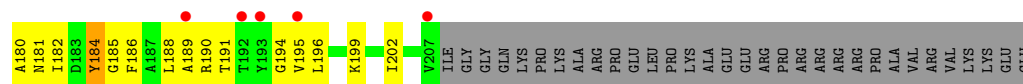


• Molecule 32: 30S ribosomal protein S2

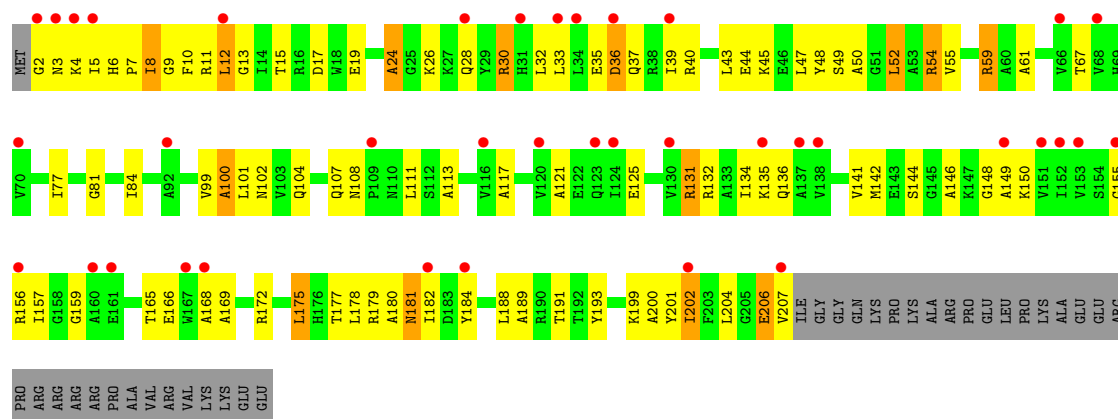


• Molecule 33: 30S ribosomal protein S3

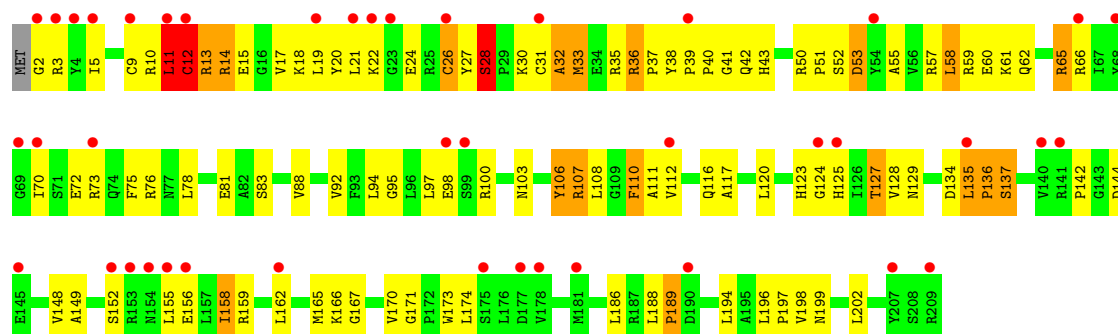




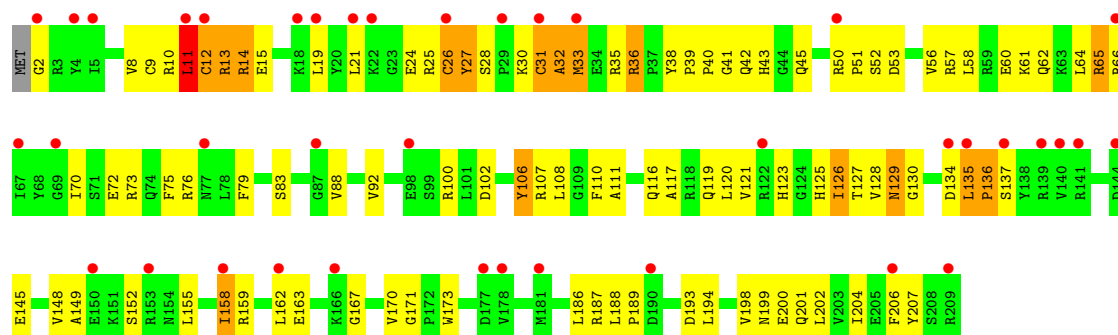
• Molecule 33: 30S ribosomal protein S3



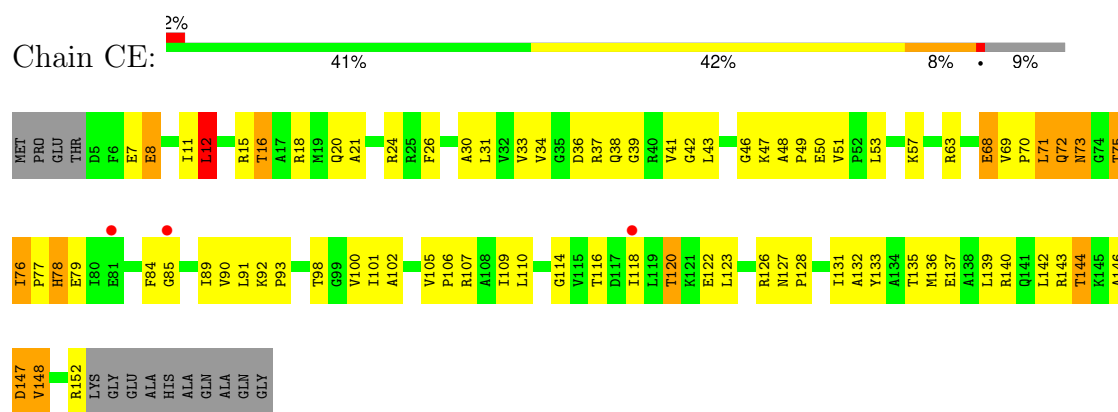
• Molecule 34: 30S ribosomal protein S4



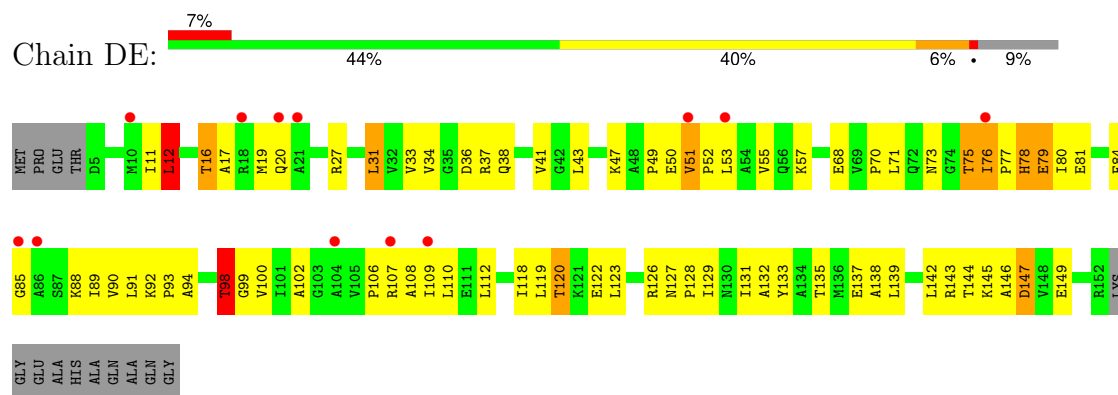
• Molecule 34: 30S ribosomal protein S4



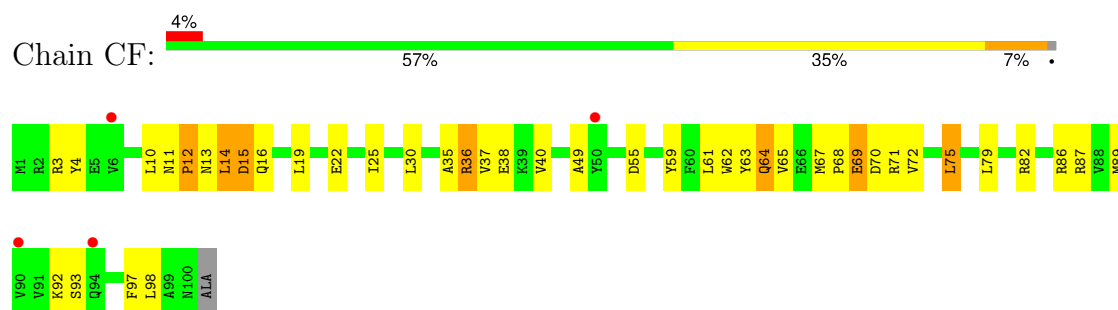
• Molecule 35: 30S ribosomal protein S5



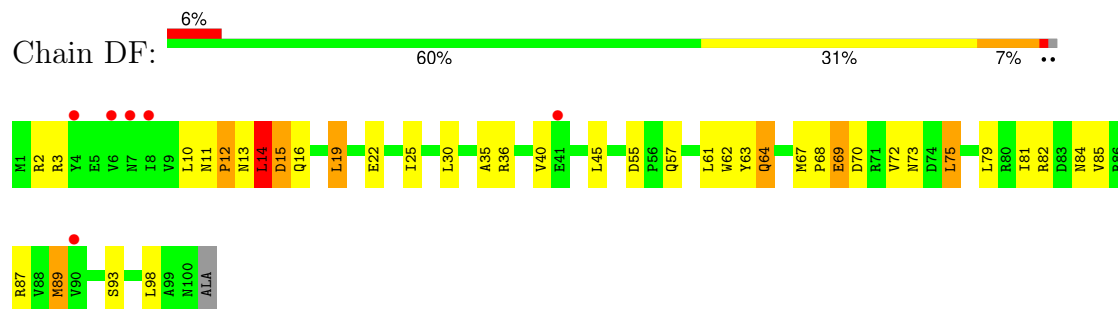
• Molecule 35: 30S ribosomal protein S5



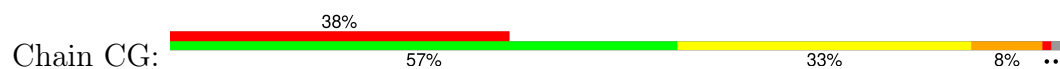
• Molecule 36: 30S ribosomal protein S6

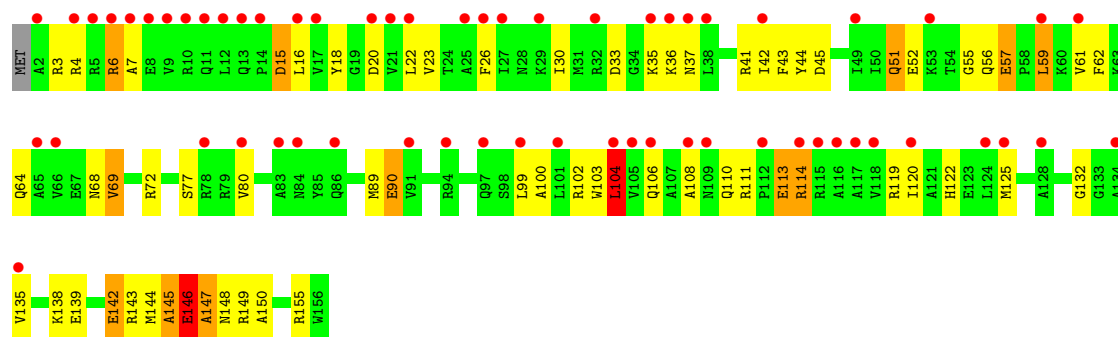


• Molecule 36: 30S ribosomal protein S6

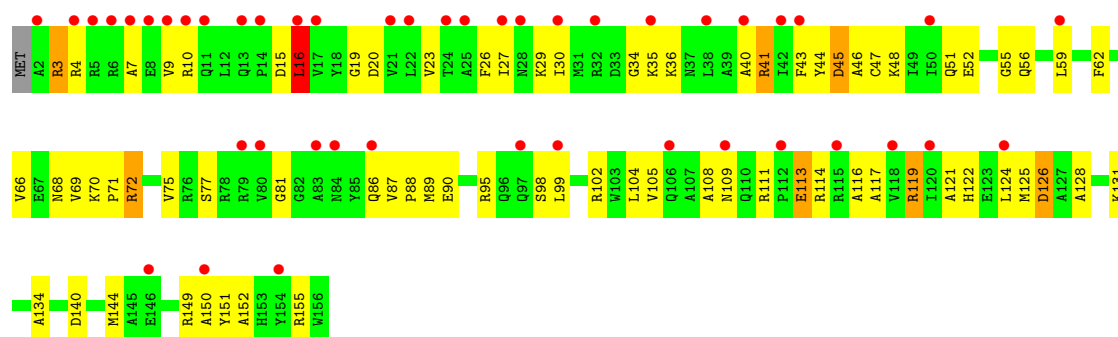


• Molecule 37: 30S ribosomal protein S7

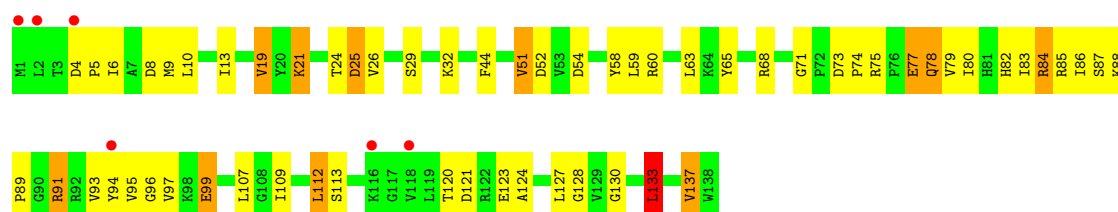




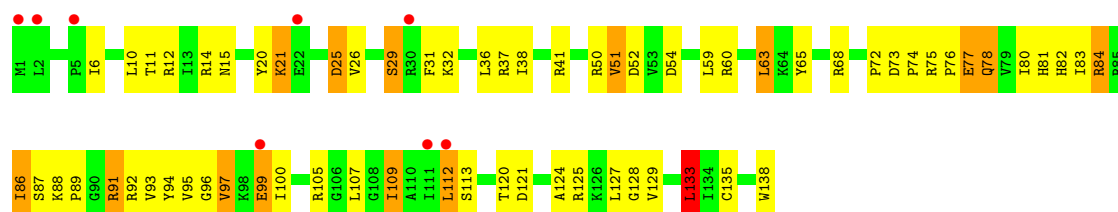
- Molecule 37: 30S ribosomal protein S7



- Molecule 38: 30S ribosomal protein S8

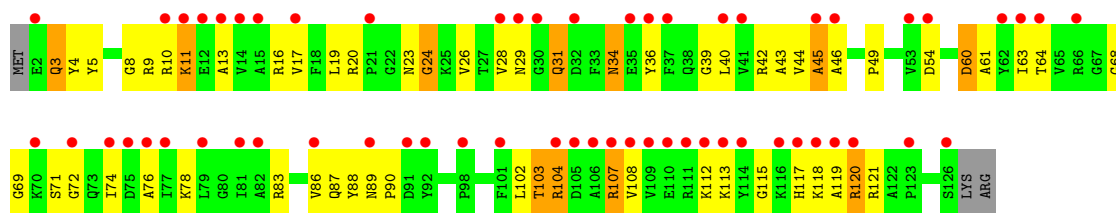


- Molecule 38: 30S ribosomal protein S8

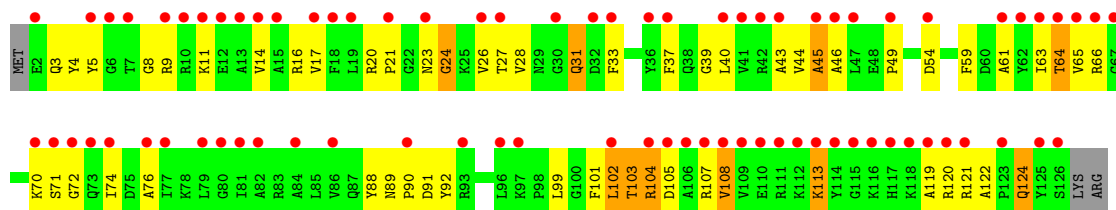


- Molecule 39: 30S ribosomal protein S9

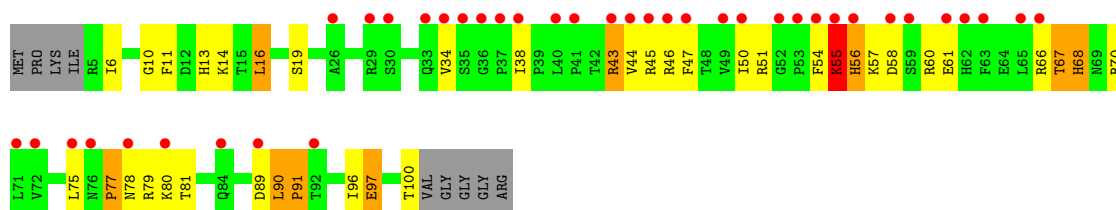
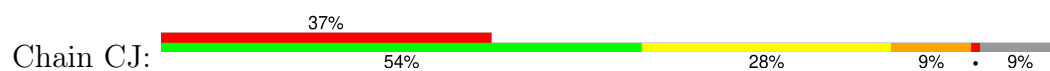




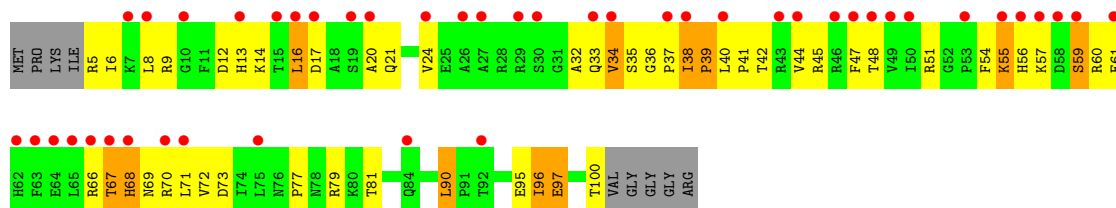
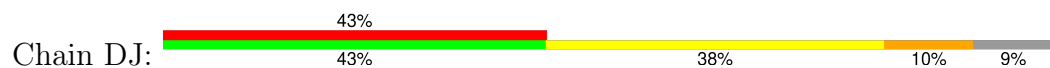
• Molecule 39: 30S ribosomal protein S9



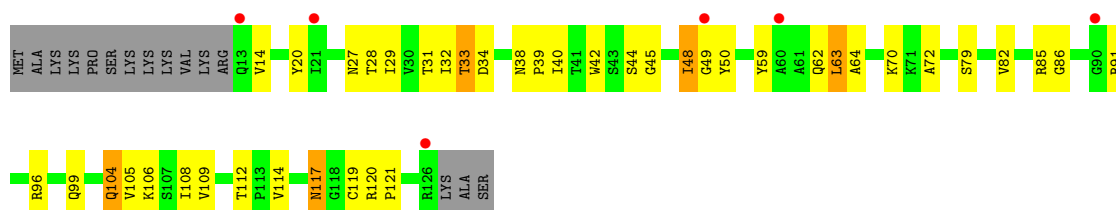
• Molecule 40: 30S ribosomal protein S10



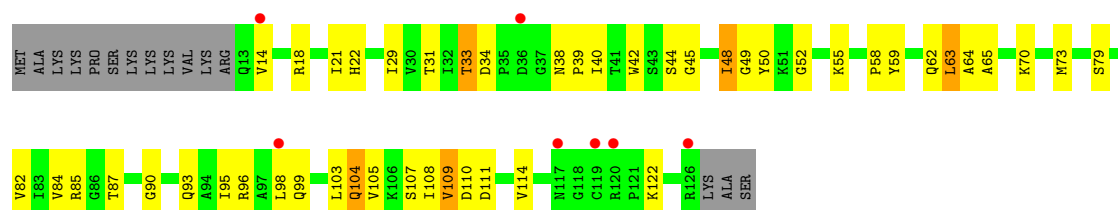
• Molecule 40: 30S ribosomal protein S10



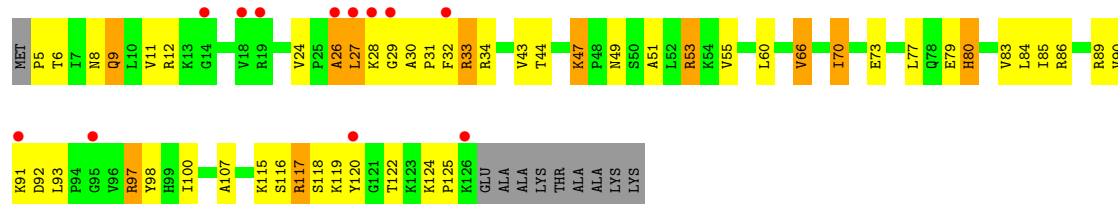
• Molecule 41: 30S ribosomal protein S11



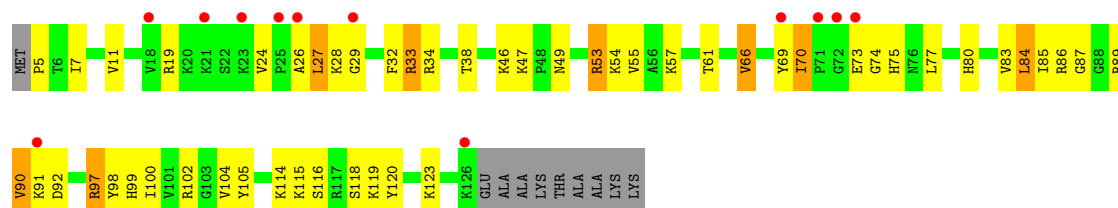
- Molecule 41: 30S ribosomal protein S11



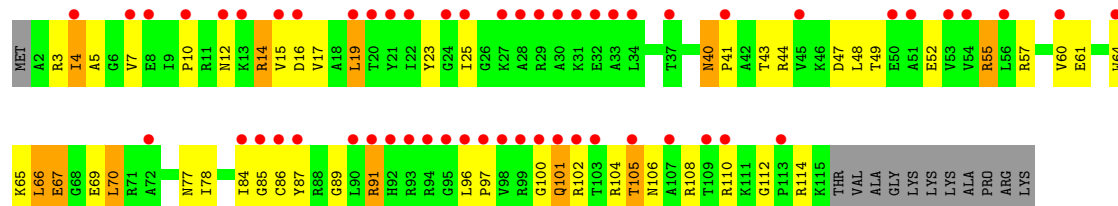
- Molecule 42: 30S ribosomal protein S12



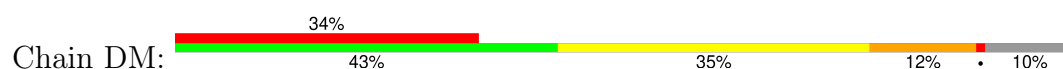
- Molecule 42: 30S ribosomal protein S12



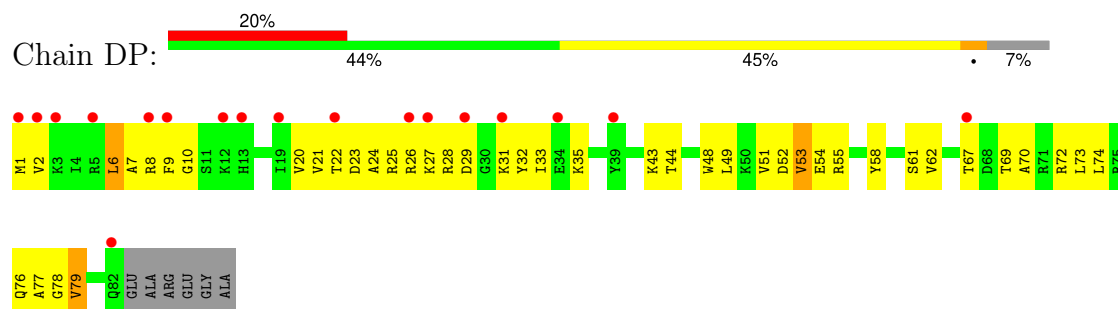
- Molecule 43: 30S ribosomal protein S13



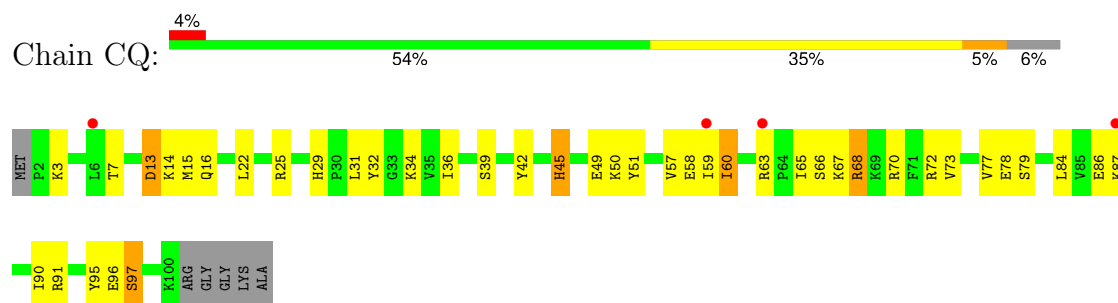
- Molecule 43: 30S ribosomal protein S13



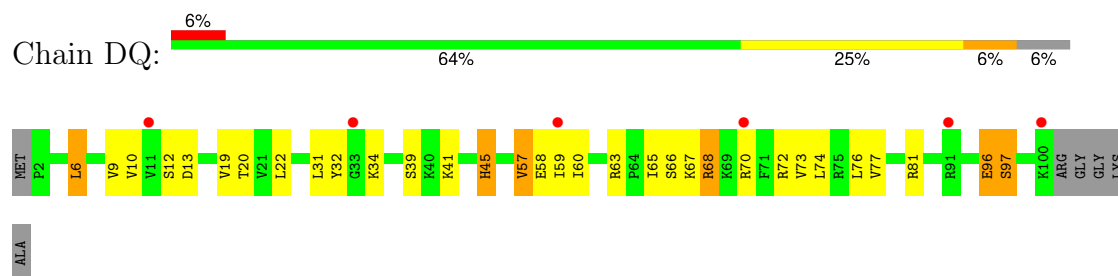
• Molecule 46: 30S ribosomal protein S16



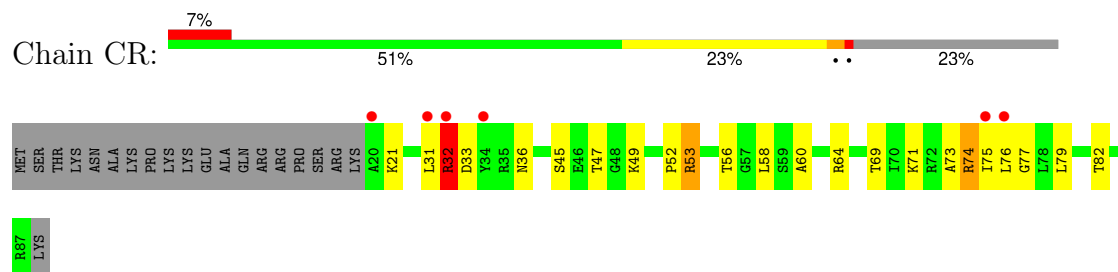
• Molecule 47: 30S ribosomal protein S17



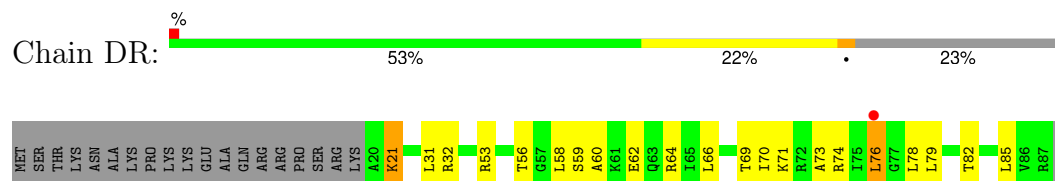
• Molecule 47: 30S ribosomal protein S17



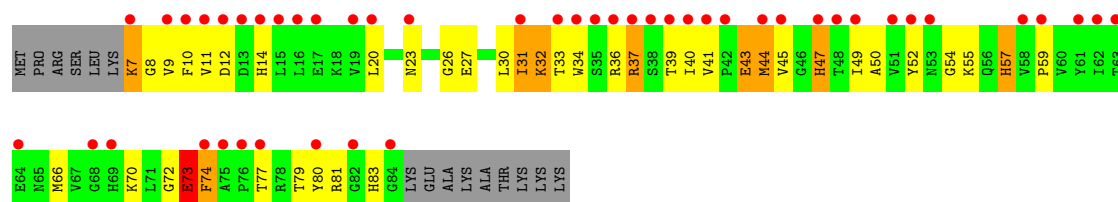
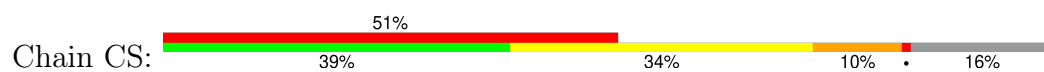
• Molecule 48: 30S ribosomal protein S18



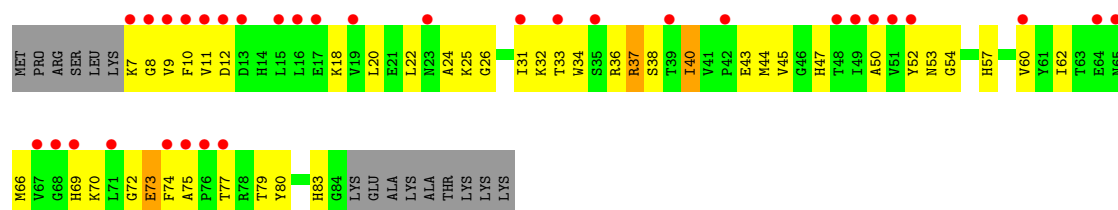
• Molecule 48: 30S ribosomal protein S18



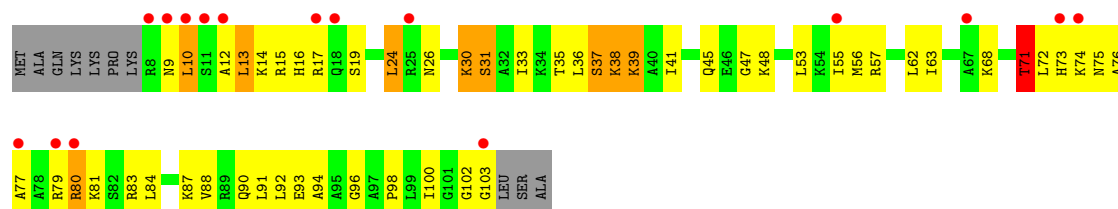
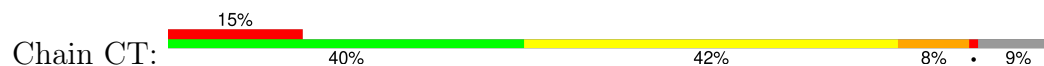
• Molecule 49: 30S ribosomal protein S19



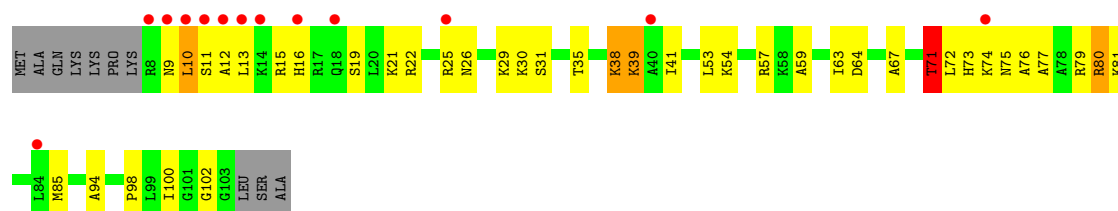
- Molecule 49: 30S ribosomal protein S19



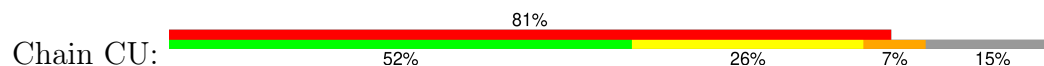
- Molecule 50: 30S ribosomal protein S20



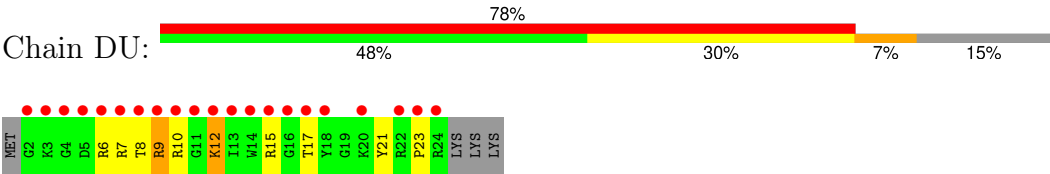
- Molecule 50: 30S ribosomal protein S20



- Molecule 51: 30S ribosomal protein THX



- Molecule 51: 30S ribosomal protein THX



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	209.37Å 445.46Å 619.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.29 – 3.20 49.29 – 3.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.29-3.20) 99.8 (49.29-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 2.91Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.228 , 0.273 0.226 , 0.269	Depositor DCC
R_{free} test set	47287 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	75.4	Xtriage
Anisotropy	0.180	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 60.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	279316	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.63% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MG, T8B

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	AA	0.82	17/68203 (0.0%)	1.30	624/106459 (0.6%)
1	BA	1.20	70/68203 (0.1%)	1.37	800/106459 (0.8%)
2	AB	0.94	0/2879	1.25	21/4492 (0.5%)
2	BB	0.88	0/2879	1.26	21/4492 (0.5%)
3	AD	0.54	0/2186	0.75	2/2944 (0.1%)
3	BD	0.67	0/2186	0.81	4/2944 (0.1%)
4	AE	0.55	0/1588	0.76	0/2145
4	BE	0.72	0/1588	0.83	1/2145 (0.0%)
5	AF	0.51	0/1609	0.70	0/2177
5	BF	0.73	0/1609	0.77	0/2177
6	AG	0.61	0/1393	0.66	0/1892
6	BG	0.46	0/1393	0.64	0/1892
7	AH	0.58	0/1343	0.68	1/1820 (0.1%)
7	BH	0.59	0/1343	0.70	0/1820
8	AI	0.63	1/1061 (0.1%)	0.78	0/1451
8	BI	0.50	0/1061	0.74	0/1451
9	AN	0.52	0/1139	0.72	0/1538
9	BN	0.74	0/1139	0.78	0/1538
10	AO	0.50	0/933	0.72	1/1257 (0.1%)
10	BO	0.67	0/933	0.74	0/1257
11	AP	0.50	0/1135	0.75	1/1510 (0.1%)
11	BP	0.64	0/1135	0.81	2/1510 (0.1%)
12	AQ	0.53	0/1143	0.74	0/1527
12	BQ	0.64	0/1143	0.74	0/1527
13	AR	0.51	0/982	0.74	0/1312
13	BR	0.69	0/982	0.82	1/1312 (0.1%)
14	AS	0.64	0/875	0.79	0/1168
14	BS	0.53	0/875	0.79	1/1168 (0.1%)
15	AT	0.52	0/1077	0.73	0/1444
15	BT	0.61	0/1077	0.79	1/1444 (0.1%)
16	AU	0.56	0/977	0.69	0/1301
16	BU	0.88	1/977 (0.1%)	0.81	1/1301 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
17	AV	0.58	0/782	0.67	0/1049
17	BV	0.70	0/782	0.77	0/1049
18	AW	0.56	0/891	0.75	0/1197
18	BW	0.82	0/891	0.80	0/1197
19	AX	0.55	0/756	0.77	1/1016 (0.1%)
19	BX	0.66	0/756	0.76	1/1016 (0.1%)
20	AY	0.50	0/798	0.77	0/1073
20	BY	0.61	0/798	0.80	1/1073 (0.1%)
21	AZ	0.57	0/1555	0.68	0/2118
21	BZ	0.49	0/1555	0.71	0/2118
22	A0	0.50	0/602	0.69	0/804
22	B0	0.66	0/602	0.77	0/804
23	A1	0.51	0/752	0.72	0/1003
23	B1	0.62	0/752	0.76	0/1003
24	A2	0.59	0/590	0.68	0/781
24	B2	0.60	0/590	0.74	0/781
25	A3	0.45	0/463	0.69	0/623
25	B3	0.65	0/463	0.74	0/623
26	A4	0.65	0/358	0.74	0/487
26	B4	0.56	0/358	0.74	1/487 (0.2%)
27	A5	0.67	1/469 (0.2%)	0.83	2/634 (0.3%)
27	B5	0.79	0/469	0.88	0/634
28	A6	0.59	0/456	0.70	0/609
28	B6	0.68	0/456	0.74	0/609
29	A7	0.57	0/426	0.75	0/561
29	B7	0.78	0/426	0.84	0/561
30	A8	0.50	0/516	0.73	0/679
30	B8	0.70	0/516	0.82	0/679
31	CA	0.80	10/36054 (0.0%)	1.18	176/56272 (0.3%)
31	DA	0.77	7/36054 (0.0%)	1.19	211/56272 (0.4%)
32	CB	0.51	0/1811	0.69	0/2452
32	DB	0.56	0/1811	0.69	0/2452
33	CC	0.56	0/1474	0.65	0/2003
33	DC	0.58	0/1474	0.65	0/2003
34	CD	0.53	0/1550	0.72	3/2106 (0.1%)
34	DD	0.84	2/1550 (0.1%)	0.78	4/2106 (0.2%)
35	CE	0.49	0/1121	0.70	1/1517 (0.1%)
35	DE	0.52	0/1121	0.72	1/1517 (0.1%)
36	CF	0.49	0/794	0.64	0/1082
36	DF	0.49	0/794	0.67	1/1082 (0.1%)
37	CG	0.57	0/1186	0.65	0/1603
37	DG	0.56	0/1186	0.62	0/1603
38	CH	0.44	0/1065	0.67	0/1445

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
38	DH	0.44	0/1065	0.64	0/1445
39	CI	0.62	0/867	0.71	0/1180
39	DI	0.62	0/867	0.69	0/1180
40	CJ	0.60	0/672	0.74	1/919 (0.1%)
40	DJ	0.60	0/672	0.70	1/919 (0.1%)
41	CK	0.47	0/843	0.71	0/1144
41	DK	0.47	0/843	0.67	0/1144
42	CL	0.44	0/925	0.67	0/1251
42	DL	0.46	0/925	0.69	0/1251
43	CM	0.67	0/811	0.72	0/1103
43	DM	0.63	0/811	0.73	1/1103 (0.1%)
44	CN	0.60	0/487	0.68	0/649
44	DN	0.59	0/487	0.74	0/649
45	CO	0.49	0/735	0.64	0/981
45	DO	0.47	0/735	0.61	0/981
46	CP	0.51	0/667	0.70	0/905
46	DP	0.43	0/667	0.65	0/905
47	CQ	0.46	0/836	0.68	0/1117
47	DQ	0.47	0/836	0.66	0/1117
48	CR	0.43	0/519	0.64	0/699
48	DR	0.50	0/519	0.67	0/699
49	CS	0.69	0/558	0.88	1/759 (0.1%)
49	DS	0.76	1/558 (0.2%)	0.87	3/759 (0.4%)
50	CT	0.47	0/710	0.72	0/940
50	DT	0.42	0/710	0.68	0/940
51	CU	0.64	0/203	0.67	0/266
51	DU	0.59	0/203	0.70	0/266
All	All	0.86	110/303650 (0.0%)	1.16	1892/454928 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
26	A4	0	1
34	CD	0	1
34	DD	0	1
42	CL	0	1
42	DL	0	1
All	All	0	5

All (110) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
34	DD	12	CYS	CB-SG	19.94	2.16	1.82
34	DD	26	CYS	CB-SG	15.79	2.09	1.82
31	DA	1492	A	C6-N6	-12.21	1.24	1.33
31	CA	1492	A	C2-N3	12.12	1.44	1.33
31	CA	1493	A	N9-C4	-11.96	1.30	1.37
1	BA	1142(A)	A	N9-C4	-10.62	1.31	1.37
31	DA	1492	A	C2-N3	10.06	1.42	1.33
1	BA	1332	G	N9-C4	-9.83	1.30	1.38
31	CA	1492	A	N3-C4	8.93	1.40	1.34
31	CA	1492	A	C5-C4	8.89	1.45	1.38
1	BA	676	A	N9-C4	-8.36	1.32	1.37
1	BA	330	A	N9-C4	-8.36	1.32	1.37
1	BA	2028	U	C4-O4	8.30	1.30	1.23
31	DA	1493	A	N7-C5	8.26	1.44	1.39
1	AA	528	A	N9-C4	-8.17	1.32	1.37
1	AA	1913	A	N9-C4	8.15	1.42	1.37
49	DS	7	LYS	CB-CG	7.98	1.74	1.52
1	BA	1602	U	C4-O4	7.93	1.29	1.23
1	BA	2249	U	C4-O4	7.78	1.29	1.23
1	BA	1021	A	N9-C4	-7.48	1.33	1.37
1	BA	676	A	N3-C4	-7.31	1.30	1.34
1	BA	1332	G	N3-C4	-7.17	1.30	1.35
1	BA	1210	A	C5-C6	-7.01	1.34	1.41
1	BA	528	A	N3-C4	-7.01	1.30	1.34
1	BA	983	A	N9-C4	-6.96	1.33	1.37
31	CA	1493	A	C8-N7	-6.81	1.26	1.31
1	AA	1142(A)	A	N9-C4	-6.78	1.33	1.37
1	BA	945	A	N9-C4	-6.71	1.33	1.37
31	CA	1492	A	N9-C4	6.60	1.41	1.37
1	BA	450	G	C6-O6	6.59	1.30	1.24
31	DA	1493	A	P-OP2	-6.56	1.37	1.49
1	BA	71	A	N9-C4	-6.54	1.33	1.37
1	AA	1045	A	N9-C4	6.51	1.41	1.37
1	BA	1332	G	C5-C6	-6.46	1.35	1.42
1	BA	777	A	N3-C4	-6.42	1.30	1.34
1	BA	2287	A	N9-C4	-6.28	1.34	1.37
1	BA	750	A	C6-N1	-6.27	1.31	1.35
1	AA	2685	G	C6-O6	6.22	1.29	1.24
27	A5	49	CYS	CB-SG	-6.14	1.71	1.82
1	BA	1045	A	N9-C4	6.13	1.41	1.37
8	AI	89	TYR	CD2-CE2	-6.11	1.30	1.39
1	AA	1332	G	N9-C4	-6.08	1.33	1.38
1	BA	2013	A	N9-C4	-6.08	1.34	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	1913	A	N7-C5	-6.05	1.35	1.39
1	AA	2119	A	N9-C4	6.04	1.41	1.37
1	AA	528	A	N3-C4	-6.04	1.31	1.34
1	BA	1268	A	C6-N1	-6.01	1.31	1.35
1	AA	1698	A	N9-C4	-5.97	1.34	1.37
1	AA	1913	A	C2-N3	5.92	1.38	1.33
1	BA	71	A	C5-C6	-5.91	1.35	1.41
1	BA	1186	G	C8-N7	-5.89	1.27	1.30
31	DA	1492	A	N9-C4	5.78	1.41	1.37
1	BA	2764	A	N3-C4	-5.74	1.31	1.34
1	AA	2249	U	C4-O4	5.71	1.28	1.23
1	BA	1308	A	C6-N1	-5.71	1.31	1.35
1	BA	527	C	N3-C4	-5.70	1.29	1.33
1	BA	2176	A	N9-C4	5.69	1.41	1.37
31	CA	1492	A	C5-C6	5.68	1.46	1.41
1	BA	1187	G	C6-O6	5.67	1.29	1.24
31	CA	1492	A	N7-C5	-5.67	1.35	1.39
1	AA	1914	C	N1-C2	5.61	1.45	1.40
1	BA	2424	C	N3-C4	-5.60	1.30	1.33
1	BA	1274	A	N9-C4	-5.59	1.34	1.37
1	BA	2590	A	N9-C4	-5.59	1.34	1.37
1	BA	2778	A	P-O5'	-5.59	1.54	1.59
31	CA	1492	A	N9-C8	5.57	1.42	1.37
1	BA	1785	A	N7-C5	-5.52	1.35	1.39
1	BA	1671	U	C4-O4	5.51	1.28	1.23
1	BA	559	G	C5-C4	-5.51	1.34	1.38
1	BA	1045	A	N3-C4	5.47	1.38	1.34
31	DA	1492	A	N1-C2	5.46	1.39	1.34
1	BA	210	C	N1-C6	-5.42	1.33	1.37
1	BA	1022	G	N9-C4	-5.42	1.33	1.38
1	BA	2451	A	C6-N1	-5.41	1.31	1.35
1	BA	2613	U	C2-N3	-5.39	1.33	1.37
1	BA	1142(A)	A	N3-C4	-5.37	1.31	1.34
1	AA	2790	A	N9-C4	5.37	1.41	1.37
31	CA	171	A	N9-C4	5.36	1.41	1.37
1	BA	735	A	N3-C4	-5.35	1.31	1.34
1	BA	1332	G	C2-N3	-5.32	1.28	1.32
1	BA	2576	G	C8-N7	-5.32	1.27	1.30
1	BA	531	C	N1-C6	-5.31	1.33	1.37
1	BA	192	C	N1-C6	-5.30	1.33	1.37
1	BA	2247	A	N9-C4	-5.27	1.34	1.37
1	BA	2685	G	C6-O6	5.27	1.28	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	BA	1779	U	N3-C4	-5.26	1.33	1.38
1	AA	1915	U	O3'-P	-5.26	1.54	1.61
1	BA	460	A	C6-N1	-5.26	1.31	1.35
1	BA	1257	C	N3-C4	-5.25	1.30	1.33
31	DA	1255	G	C6-N1	5.25	1.43	1.39
1	AA	227	A	N9-C4	5.25	1.41	1.37
1	BA	1913	A	N3-C4	-5.20	1.31	1.34
1	BA	1847	A	N9-C4	5.20	1.41	1.37
1	BA	2015	A	N7-C5	-5.19	1.36	1.39
1	BA	2873	A	N3-C4	-5.18	1.31	1.34
1	AA	1915	U	C3'-O3'	-5.17	1.34	1.42
1	BA	1308	A	N7-C5	-5.12	1.36	1.39
1	BA	469	G	C5-C4	-5.11	1.34	1.38
1	BA	2057	A	N3-C4	-5.11	1.31	1.34
1	BA	1333	C	N1-C6	-5.10	1.34	1.37
1	BA	2580	U	N1-C2	-5.10	1.33	1.38
1	BA	1256	G	N9-C8	-5.09	1.34	1.37
1	BA	71	A	N7-C5	-5.08	1.36	1.39
1	BA	1254	A	N3-C4	-5.07	1.31	1.34
1	BA	2497	A	P-O5'	-5.07	1.54	1.59
1	BA	1204	A	N9-C4	-5.04	1.34	1.37
1	BA	575	A	N7-C5	-5.04	1.36	1.39
1	BA	2311	A	N9-C4	5.04	1.40	1.37
16	BU	69	CYS	CB-SG	-5.03	1.73	1.81
1	BA	2451	A	N3-C4	-5.01	1.31	1.34

All (1892) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
31	DA	1492	A	C6-N1-C2	-28.45	101.53	118.60
31	DA	1492	A	C5-C6-N1	26.41	130.90	117.70
31	CA	1492	A	C8-N9-C4	-21.05	97.38	105.80
1	BA	1332	G	C2-N3-C4	-19.65	102.08	111.90
1	BA	1332	G	N3-C4-C5	17.99	137.59	128.60
1	AA	1913	A	C4-C5-C6	17.05	125.53	117.00
1	BA	1332	G	N3-C4-N9	-16.77	115.94	126.00
31	CA	1493	A	C8-N9-C4	15.82	112.13	105.80
1	BA	1779	U	C5-C6-N1	-14.96	115.22	122.70
31	CA	1492	A	N7-C8-N9	14.60	121.10	113.80
1	AA	1913	A	C6-N1-C2	-14.27	110.04	118.60
1	BA	2249	U	N3-C4-C5	-14.14	106.12	114.60
31	DA	1492	A	N3-C4-C5	-14.09	116.94	126.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	2028	U	N3-C4-C5	-13.81	106.31	114.60
1	BA	1332	G	C5-N7-C8	-13.80	97.40	104.30
31	DA	1492	A	C8-N9-C4	-13.58	100.37	105.80
1	AA	1332	G	C2-N3-C4	-13.55	105.12	111.90
31	CA	1492	A	N3-C4-C5	-13.27	117.51	126.80
1	AA	1913	A	N3-C4-C5	-13.21	117.55	126.80
1	BA	330	A	C2-N3-C4	-13.06	104.07	110.60
31	DA	1492	A	N1-C2-N3	13.04	135.82	129.30
1	BA	1602	U	N3-C4-C5	-13.02	106.79	114.60
31	CA	1493	A	N9-C4-C5	-12.86	100.66	105.80
1	BA	450	G	C5-C6-N1	-12.42	105.29	111.50
1	BA	1021	A	C2-N3-C4	-12.36	104.42	110.60
49	CS	8	GLY	N-CA-C	12.03	143.18	113.10
1	BA	1142(A)	A	C2-N3-C4	-11.84	104.68	110.60
1	BA	450	G	C4-C5-N7	-11.78	106.09	110.80
31	DA	1493	A	C8-N9-C4	11.69	110.48	105.80
31	CA	1493	A	N3-C4-C5	11.61	134.93	126.80
31	DA	1030	C	N1-C2-O2	11.60	125.86	118.90
31	DA	1492	A	N7-C8-N9	11.59	119.60	113.80
1	BA	1187	G	C5-C6-N1	-11.38	105.81	111.50
1	BA	450	G	N9-C4-C5	11.15	109.86	105.40
1	AA	2685	G	C5-C6-N1	-11.11	105.95	111.50
49	DS	8	GLY	N-CA-C	11.11	140.87	113.10
1	BA	1915	U	C4'-C3'-C2'	-11.10	91.50	102.60
1	BA	141	A	C5-N7-C8	-11.03	98.39	103.90
1	BA	1332	G	N1-C6-O6	11.02	126.51	119.90
1	BA	528	A	C8-N9-C4	-11.02	101.39	105.80
1	AA	1779	U	C5-C6-N1	-10.94	117.23	122.70
1	BA	570	G	C4-C5-N7	-10.92	106.43	110.80
31	CA	1493	A	C6-N1-C2	10.90	125.14	118.60
1	AA	2137	C	C6-N1-C2	-10.68	116.03	120.30
31	CA	1126	U	N1-C2-O2	10.62	130.23	122.80
1	BA	2028	U	N3-C4-O4	10.57	126.80	119.40
1	AA	1332	G	N3-C4-N9	-10.56	119.66	126.00
1	BA	1671	U	N3-C4-O4	10.55	126.78	119.40
1	BA	676	A	C2-N3-C4	-10.53	105.33	110.60
1	BA	141	A	N7-C8-N9	10.49	119.04	113.80
1	BA	945	A	N1-C6-N6	10.47	124.88	118.60
1	AA	912	C	C6-N1-C2	-10.45	116.12	120.30
1	BA	566	U	C6-N1-C2	10.43	127.26	121.00
31	CA	1493	A	N1-C2-N3	-10.42	124.09	129.30
31	DA	1493	A	N9-C4-C5	-10.37	101.65	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	1828	G	C5-C6-O6	10.35	134.81	128.60
31	CA	1492	A	C6-N1-C2	-10.34	112.40	118.60
1	BA	2685	G	C5-C6-N1	-10.32	106.34	111.50
1	BA	2685	G	N3-C2-N2	-10.31	112.69	119.90
2	AB	30	C	C6-N1-C2	-10.29	116.18	120.30
1	BA	71	A	C5-N7-C8	-10.29	98.76	103.90
1	BA	1332	G	N3-C2-N2	-10.26	112.72	119.90
1	AA	1913	A	N1-C2-N3	10.24	134.42	129.30
1	BA	2249	U	C6-N1-C2	-10.22	114.87	121.00
1	BA	2544	G	N1-C6-O6	10.20	126.02	119.90
1	BA	456	C	C6-N1-C2	10.12	124.35	120.30
1	BA	570	G	C5-C6-O6	10.12	134.67	128.60
1	AA	1913	A	N3-C4-N9	10.07	135.46	127.40
1	BA	298	G	C5-N7-C8	-10.07	99.26	104.30
1	AA	2828	C	C6-N1-C2	10.05	124.32	120.30
1	BA	141	A	N1-C6-N6	10.01	124.61	118.60
1	BA	1332	G	N7-C8-N9	9.98	118.09	113.10
1	AA	528	A	C2-N3-C4	-9.97	105.62	110.60
1	AA	1332	G	N3-C4-C5	9.92	133.56	128.60
1	BA	568	U	N3-C4-C5	-9.91	108.65	114.60
1	BA	450	G	N3-C2-N2	-9.89	112.98	119.90
31	DA	899	C	C6-N1-C2	9.86	124.25	120.30
1	BA	528	A	N7-C8-N9	9.85	118.73	113.80
1	BA	528	A	C5-N7-C8	-9.83	98.99	103.90
1	AA	1963	U	C2-N1-C1'	9.78	129.44	117.70
1	AA	676	A	N7-C8-N9	9.73	118.67	113.80
31	DA	1492	A	N1-C6-N6	-9.72	112.77	118.60
1	BA	1210	A	N1-C6-N6	9.72	124.43	118.60
1	BA	856	C	C6-N1-C2	-9.71	116.41	120.30
31	CA	1492	A	C5-N7-C8	-9.70	99.05	103.90
31	DA	1492	A	N3-C4-N9	9.68	135.15	127.40
31	DA	1492	A	C5-N7-C8	-9.61	99.10	103.90
31	CA	1126	U	N3-C2-O2	-9.60	115.48	122.20
1	AA	1698	A	C2-N3-C4	-9.59	105.81	110.60
31	DA	1492	A	C4-N9-C1'	9.57	143.53	126.30
1	BA	1142(A)	A	N3-C4-C5	9.53	133.47	126.80
1	BA	450	G	C8-N9-C4	-9.51	102.59	106.40
1	BA	298	G	C4-C5-N7	9.44	114.57	110.80
1	BA	945	A	C5-N7-C8	-9.40	99.20	103.90
1	BA	12	U	N3-C2-O2	-9.39	115.63	122.20
1	BA	179	G	N1-C6-O6	9.38	125.53	119.90
1	BA	528	A	C2-N3-C4	-9.37	105.91	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1022	G	N3-C2-N2	-9.34	113.36	119.90
1	AA	1840	G	N1-C6-O6	9.33	125.50	119.90
1	BA	1671	U	N3-C4-C5	-9.30	109.02	114.60
1	AA	1913	A	C6-C5-N7	-9.29	125.80	132.30
1	AA	1828	G	C4-C5-N7	-9.26	107.10	110.80
31	CA	992	U	C2-N1-C1'	9.23	128.77	117.70
31	DA	1492	A	C5-C6-N6	-9.22	116.33	123.70
1	AA	450	G	C5-C6-N1	-9.17	106.92	111.50
1	BA	1332	G	C4-C5-N7	9.16	114.46	110.80
1	AA	2439	A	N1-C6-N6	9.11	124.07	118.60
1	BA	676	A	C5-N7-C8	-9.08	99.36	103.90
1	BA	652(H)	C	C5-C6-N1	9.07	125.53	121.00
1	BA	1187	G	N3-C2-N2	-9.06	113.56	119.90
1	AA	12	U	N1-C2-O2	9.05	129.14	122.80
1	BA	2137	C	C6-N1-C2	-9.03	116.69	120.30
1	BA	933	A	C5-N7-C8	-9.01	99.39	103.90
1	AA	2447	G	N1-C6-O6	8.99	125.30	119.90
31	CA	839	U	N1-C2-O2	8.97	129.08	122.80
1	BA	587	C	C6-N1-C2	-8.96	116.72	120.30
1	BA	1779	U	C2-N1-C1'	-8.95	106.97	117.70
1	BA	271(M)	G	N3-C4-N9	8.91	131.34	126.00
1	AA	1828	G	C5-C6-O6	8.90	133.94	128.60
1	AA	1210	A	N7-C8-N9	8.86	118.23	113.80
1	AA	1828	G	N9-C4-C5	8.86	108.94	105.40
31	DA	992	U	C2-N1-C1'	8.82	128.28	117.70
1	AA	330	A	C2-N3-C4	-8.79	106.20	110.60
31	CA	839	U	N3-C2-O2	-8.79	116.05	122.20
1	BA	1203	G	C5-C6-O6	8.78	133.87	128.60
31	CA	1492	A	C5-C6-N1	8.76	122.08	117.70
31	DA	839	U	N3-C2-O2	-8.74	116.08	122.20
1	BA	2489	G	N1-C6-O6	8.72	125.13	119.90
31	CA	1492	A	C6-C5-N7	-8.71	126.20	132.30
1	AA	2473	U	N1-C2-O2	8.66	128.86	122.80
1	BA	528	A	N1-C2-N3	8.66	133.63	129.30
1	BA	659	C	C6-N1-C2	8.66	123.76	120.30
1	BA	1372	U	C5-C6-N1	8.65	127.03	122.70
1	BA	1204	A	C2-N3-C4	-8.62	106.29	110.60
1	AA	2293	C	C6-N1-C2	8.62	123.75	120.30
1	BA	141	A	C6-C5-N7	-8.59	126.28	132.30
1	AA	1142(A)	A	C2-N3-C4	-8.59	106.30	110.60
1	AA	2685	G	N3-C4-N9	-8.59	120.85	126.00
1	BA	208	C	C6-N1-C2	8.57	123.73	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	570	G	N9-C4-C5	8.56	108.83	105.40
1	AA	201	C	C6-N1-C2	8.56	123.72	120.30
1	BA	1256	G	C8-N9-C1'	-8.56	115.88	127.00
31	CA	1017	G	C5-C6-O6	8.54	133.72	128.60
1	BA	1779	U	C4-C5-C6	8.52	124.81	119.70
31	CA	1502	A	N1-C2-N3	8.52	133.56	129.30
1	AA	893	C	C2-N1-C1'	8.51	128.16	118.80
31	DA	1030	C	C2-N3-C4	8.51	124.16	119.90
1	BA	1699	G	N9-C4-C5	8.51	108.80	105.40
31	CA	1027	C	C5-C4-N4	8.49	126.14	120.20
1	BA	2032	G	C5-N7-C8	-8.48	100.06	104.30
1	BA	271(M)	G	N3-C4-C5	-8.47	124.36	128.60
1	BA	456	C	N3-C2-O2	8.46	127.82	121.90
1	BA	893	C	C2-N1-C1'	8.46	128.11	118.80
1	BA	1784	A	N7-C8-N9	-8.46	109.57	113.80
31	DA	1282	C	C2-N1-C1'	8.45	128.10	118.80
1	AA	618	C	C6-N1-C2	8.40	123.66	120.30
1	AA	2218	U	C2-N1-C1'	8.39	127.77	117.70
34	DD	12	CYS	CA-CB-SG	8.39	129.10	114.00
31	DA	1030	C	C5-C6-N1	8.37	125.19	121.00
1	BA	734	A	N1-C6-N6	8.35	123.61	118.60
31	CA	1492	A	N3-C4-N9	8.34	134.07	127.40
1	BA	1313	U	C2-N1-C1'	8.34	127.70	117.70
31	DA	1492	A	C6-C5-N7	-8.32	126.47	132.30
1	AA	2619	C	C6-N1-C2	8.32	123.63	120.30
1	BA	2206	G	C4-N9-C1'	-8.32	115.69	126.50
1	AA	1210	A	C8-N9-C4	-8.29	102.48	105.80
1	AA	2061	G	N1-C6-O6	8.27	124.86	119.90
1	AA	12	U	N3-C2-O2	-8.26	116.42	122.20
1	BA	2286	A	N1-C6-N6	8.26	123.56	118.60
1	AA	1914	C	N1-C2-O2	8.26	123.85	118.90
1	BA	673	C	N1-C2-O2	8.25	123.85	118.90
1	BA	71	A	C2-N3-C4	-8.24	106.48	110.60
31	DA	299	G	C5-C6-N1	-8.24	107.38	111.50
1	AA	676	A	C5-N7-C8	-8.23	99.78	103.90
31	CA	1524	C	C6-N1-C2	8.21	123.58	120.30
1	BA	2685	G	N9-C4-C5	8.19	108.68	105.40
1	AA	676	A	C8-N9-C4	-8.18	102.53	105.80
1	AA	1210	A	C5-N7-C8	-8.18	99.81	103.90
1	AA	2866	U	C2-N1-C1'	8.16	127.49	117.70
1	AA	2131	G	N7-C8-N9	8.15	117.18	113.10
49	DS	7	LYS	CB-CA-C	8.14	126.68	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	2544	G	C5-C6-O6	-8.12	123.72	128.60
1	AA	1332	G	N1-C2-N3	8.12	128.77	123.90
31	DA	1493	A	C8-N9-C1'	-8.12	113.09	127.70
1	BA	1049	C	C6-N1-C2	-8.09	117.06	120.30
1	BA	330	A	N3-C4-C5	8.08	132.46	126.80
1	BA	945	A	C6-C5-N7	-8.06	126.66	132.30
1	BA	570	G	C5-C6-N1	-8.05	107.47	111.50
1	AA	179	G	N1-C6-O6	8.05	124.73	119.90
1	BA	141	A	C4-C5-N7	8.02	114.71	110.70
1	BA	798	G	C8-N9-C4	8.02	109.61	106.40
1	AA	1021	A	C2-N3-C4	-8.02	106.59	110.60
1	AA	2137	C	N3-C2-O2	-8.02	116.29	121.90
2	AB	115	G	C8-N9-C4	8.01	109.61	106.40
1	BA	945	A	C4-C5-N7	8.00	114.70	110.70
1	AA	645	C	N1-C2-O2	7.99	123.70	118.90
31	DA	1030(B)	C	C2-N1-C1'	7.98	127.58	118.80
31	CA	1493	A	C4-C5-C6	-7.98	113.01	117.00
1	AA	2206	G	N3-C4-C5	7.98	132.59	128.60
1	AA	450	G	N3-C2-N2	-7.97	114.32	119.90
1	AA	1109	C	C4-C5-C6	7.96	121.38	117.40
1	BA	2022	U	C5-C4-O4	-7.96	121.12	125.90
1	BA	566	U	N1-C2-N3	-7.96	110.13	114.90
1	AA	208	C	C6-N1-C2	7.94	123.48	120.30
1	AA	1022	G	N3-C4-N9	-7.94	121.24	126.00
31	CA	1149	C	C2-N3-C4	7.94	123.87	119.90
1	BA	1142(A)	A	N3-C4-N9	-7.94	121.05	127.40
1	BA	1021	A	N1-C2-N3	7.93	133.26	129.30
1	AA	2249	U	N3-C4-C5	-7.92	109.85	114.60
1	BA	14	A	C8-N9-C4	7.92	108.97	105.80
1	AA	2685	G	N3-C2-N2	-7.92	114.36	119.90
1	BA	2419	U	N3-C4-C5	-7.91	109.86	114.60
1	BA	2419	U	N3-C4-O4	7.91	124.94	119.40
1	BA	676	A	N1-C2-N3	7.89	133.25	129.30
1	BA	2473	U	N3-C2-O2	-7.88	116.68	122.20
1	BA	676	A	N7-C8-N9	7.88	117.74	113.80
31	CA	1492	A	C2-N3-C4	7.88	114.54	110.60
1	BA	1332	G	C5-C6-N1	-7.87	107.56	111.50
1	BA	1784	A	C8-N9-C4	7.87	108.95	105.80
1	AA	915	C	C6-N1-C2	-7.87	117.15	120.30
1	BA	71	A	N7-C8-N9	7.85	117.73	113.80
1	BA	2249	U	C4-C5-C6	7.85	124.41	119.70
1	AA	1882	C	C2-N1-C1'	7.85	127.44	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
31	CA	1126	U	C2-N1-C1'	7.85	127.12	117.70
1	BA	2249	U	N3-C4-O4	7.85	124.89	119.40
31	DA	1126	U	C2-N1-C1'	7.84	127.11	117.70
31	DA	76	C	N1-C2-O2	7.83	123.60	118.90
1	BA	893	C	C6-N1-C1'	-7.83	111.40	120.80
1	BA	1210	A	C5-N7-C8	-7.83	99.99	103.90
1	BA	1963	U	C2-N1-C1'	7.83	127.09	117.70
1	AA	915	C	N3-C2-O2	-7.82	116.42	121.90
31	CA	1492	A	C4-C5-C6	7.82	120.91	117.00
1	AA	1698	A	C5-N7-C8	-7.81	99.99	103.90
1	BA	1913	A	O4'-C1'-N9	-7.81	101.95	108.20
31	CA	1034	G	C5-C6-O6	7.80	133.28	128.60
1	BA	204	A	C6-N1-C2	-7.80	113.92	118.60
1	BA	298	G	N7-C8-N9	7.79	117.00	113.10
1	BA	749	C	C6-N1-C2	7.79	123.42	120.30
1	AA	840	C	C6-N1-C2	7.79	123.42	120.30
1	BA	2287	A	C2-N3-C4	-7.78	106.71	110.60
1	BA	1791	A	N1-C6-N6	7.77	123.26	118.60
1	BA	885	C	N1-C2-O2	7.76	123.56	118.90
1	BA	1256	G	C4-N9-C1'	7.76	136.59	126.50
1	AA	2200	C	C2-N1-C1'	7.76	127.33	118.80
1	BA	139(A)	G	N7-C8-N9	7.74	116.97	113.10
1	BA	71	A	C4-C5-N7	7.73	114.56	110.70
1	AA	528	A	N3-C4-N9	-7.72	121.22	127.40
31	CA	992	U	N1-C2-O2	7.72	128.20	122.80
1	BA	2419	U	C6-N1-C2	-7.71	116.37	121.00
1	BA	2489	G	C5-C6-O6	-7.71	123.98	128.60
1	BA	2028	U	C6-N1-C2	-7.70	116.38	121.00
31	CA	1037	C	C5-C6-N1	7.68	124.84	121.00
1	BA	2785	C	C6-N1-C2	-7.68	117.23	120.30
31	DA	992	U	N1-C2-O2	7.67	128.17	122.80
1	BA	1022	G	N3-C2-N2	-7.67	114.53	119.90
1	AA	530	G	N1-C6-O6	-7.66	115.30	119.90
1	AA	1333	C	C2-N1-C1'	7.66	127.23	118.80
1	BA	2473	U	N1-C2-O2	7.63	128.14	122.80
1	AA	2439	A	C6-C5-N7	-7.62	126.96	132.30
1	BA	60	G	N9-C4-C5	-7.62	102.35	105.40
1	BA	139(A)	G	C8-N9-C4	-7.59	103.36	106.40
31	DA	839	U	C2-N1-C1'	7.59	126.81	117.70
31	DA	1492	A	C4-C5-C6	7.56	120.78	117.00
1	AA	2791	C	C6-N1-C2	-7.56	117.28	120.30
1	AA	570	G	C5-C6-N1	-7.55	107.72	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	792	G	N3-C4-N9	7.54	130.53	126.00
1	BA	124	G	C8-N9-C4	7.54	109.42	106.40
1	BA	933	A	C4-C5-N7	7.53	114.46	110.70
1	BA	141	A	C8-N9-C4	-7.52	102.79	105.80
31	DA	1126	U	N1-C2-O2	7.52	128.06	122.80
1	BA	2249	U	C2-N3-C4	7.51	131.50	127.00
1	BA	2028	U	C2-N3-C4	7.49	131.50	127.00
1	BA	1784	A	C5-N7-C8	7.49	107.65	103.90
1	AA	1328	G	N3-C4-N9	7.48	130.49	126.00
1	BA	12	U	N1-C2-O2	7.48	128.04	122.80
1	BA	452	G	N3-C4-C5	-7.46	124.87	128.60
1	AA	652(S)	C	N1-C2-O2	7.46	123.38	118.90
1	AA	1493	C	C2-N1-C1'	7.46	127.01	118.80
1	BA	1383	C	N3-C2-O2	7.46	127.12	121.90
1	AA	1022	G	N9-C4-C5	7.46	108.38	105.40
1	AA	2361	A	N9-C4-C5	-7.45	102.82	105.80
1	AA	1333	C	C6-N1-C1'	-7.45	111.86	120.80
31	DA	96	U	N1-C2-O2	7.45	128.01	122.80
1	BA	1698	A	C2-N3-C4	-7.43	106.88	110.60
1	BA	2134	A	N1-C6-N6	-7.43	114.14	118.60
1	BA	2685	G	C4-C5-N7	-7.43	107.83	110.80
31	DA	1027	C	C5-C4-N4	7.43	125.40	120.20
1	BA	2589	A	C8-N9-C4	7.42	108.77	105.80
1	AA	208	C	N3-C4-C5	7.42	124.87	121.90
31	DA	1493	A	N3-C4-N9	7.41	133.33	127.40
1	AA	2473	U	N3-C2-O2	-7.40	117.02	122.20
1	AA	2061	G	C8-N9-C4	7.37	109.35	106.40
1	BA	1263	U	C2-N1-C1'	7.37	126.55	117.70
31	CA	1149	C	N1-C2-O2	7.37	123.32	118.90
1	AA	1840	G	C6-C5-N7	-7.36	125.99	130.40
1	AA	1913	A	C5-N7-C8	7.35	107.57	103.90
31	DA	1149	C	C6-N1-C2	-7.35	117.36	120.30
1	AA	565	C	C6-N1-C2	7.34	123.24	120.30
1	BA	409	C	C6-N1-C2	7.34	123.24	120.30
1	BA	318	C	C6-N1-C2	-7.34	117.36	120.30
34	DD	31	CYS	CA-CB-SG	-7.33	100.80	114.00
1	BA	1332	G	N1-C2-N3	7.33	128.29	123.90
1	BA	613	G	C5-C6-O6	-7.32	124.21	128.60
31	CA	1091	U	N3-C2-O2	-7.32	117.07	122.20
1	AA	2473	U	C2-N1-C1'	7.32	126.49	117.70
1	AA	2218	U	C6-N1-C1'	-7.32	110.95	121.20
1	AA	2447	G	C5-C6-O6	-7.30	124.22	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	BE	127	ASP	CB-CG-OD1	-7.30	111.73	118.30
31	DA	1282	C	N3-C2-O2	-7.30	116.79	121.90
1	AA	1671	U	N3-C4-O4	7.29	124.50	119.40
1	BA	2447	G	N1-C6-O6	7.28	124.27	119.90
1	AA	1049	C	C6-N1-C2	-7.28	117.39	120.30
1	AA	1266	G	C8-N9-C4	7.28	109.31	106.40
1	AA	1828	G	C8-N9-C4	-7.28	103.49	106.40
1	BA	1982	C	C6-N1-C2	-7.27	117.39	120.30
31	CA	1087	G	N3-C4-C5	-7.27	124.96	128.60
1	AA	885	C	N1-C2-O2	7.26	123.26	118.90
1	BA	1772	G	N1-C6-O6	-7.26	115.54	119.90
31	DA	736	C	C5-C6-N1	7.26	124.63	121.00
1	AA	733	G	N3-C4-C5	7.26	132.23	128.60
1	BA	2447	G	C5-C6-O6	-7.26	124.25	128.60
31	CA	1165	C	C6-N1-C2	-7.26	117.40	120.30
31	DA	1282	C	N1-C2-O2	7.25	123.25	118.90
1	AA	1974	C	C6-N1-C2	7.24	123.20	120.30
1	AA	566	U	C6-N1-C2	7.23	125.34	121.00
31	DA	530	G	C4-N9-C1'	7.23	135.90	126.50
1	BA	933	A	N7-C8-N9	7.22	117.41	113.80
1	BA	90	U	C5-C6-N1	7.22	126.31	122.70
1	BA	1602	U	C4-C5-C6	7.22	124.03	119.70
1	BA	1699	G	C5-C6-O6	7.22	132.93	128.60
1	BA	2183	C	C2-N3-C4	7.21	123.51	119.90
31	DA	93	G	N3-C4-N9	7.21	130.33	126.00
3	BD	239	ARG	N-CA-C	-7.20	91.55	111.00
34	DD	12	CYS	N-CA-C	-7.20	91.56	111.00
1	AA	196	A	N1-C6-N6	7.20	122.92	118.60
1	AA	203	C	N3-C4-C5	7.19	124.78	121.90
31	CA	1493	A	C5-C6-N1	-7.19	114.11	117.70
1	BA	2036	C	C6-N1-C2	-7.18	117.43	120.30
1	AA	652(H)	C	C5-C6-N1	7.17	124.59	121.00
1	BA	1231	G	N3-C4-C5	7.17	132.18	128.60
1	AA	2035	G	C4-N9-C1'	-7.16	117.19	126.50
1	BA	1699	G	C4-C5-N7	-7.16	107.94	110.80
31	DA	1492	A	C8-N9-C1'	-7.15	114.83	127.70
1	BA	2218	U	C2-N1-C1'	7.14	126.27	117.70
1	BA	847	U	C2-N1-C1'	-7.13	109.14	117.70
1	BA	2894	G	C6-C5-N7	-7.13	126.12	130.40
1	BA	723	G	N3-C4-N9	7.13	130.28	126.00
1	BA	1142(A)	A	C5-C6-N1	-7.13	114.14	117.70
1	BA	1656	C	N3-C4-C5	-7.12	119.05	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	2206	G	C8-N9-C1'	7.12	136.25	127.00
1	BA	512	G	N3-C4-C5	7.11	132.16	128.60
31	CA	1017	G	C6-N1-C2	7.10	129.36	125.10
1	BA	71	A	N1-C6-N6	7.10	122.86	118.60
1	BA	1951	U	N3-C4-C5	-7.10	110.34	114.60
31	DA	1123	A	N1-C6-N6	-7.10	114.34	118.60
1	BA	645	C	N1-C2-O2	7.09	123.16	118.90
1	AA	2828	C	C5-C6-N1	-7.08	117.46	121.00
1	AA	1204	A	N1-C6-N6	7.08	122.85	118.60
1	BA	1506	C	C6-N1-C2	-7.07	117.47	120.30
1	BA	546	C	N1-C2-O2	7.07	123.14	118.90
1	BA	2104	G	C5-C6-O6	7.07	132.84	128.60
31	CA	839	U	C2-N1-C1'	7.06	126.17	117.70
1	BA	1332	G	C8-N9-C4	-7.06	103.58	106.40
1	BA	1290	C	N3-C4-C5	-7.05	119.08	121.90
31	DA	1397	C	C2-N1-C1'	7.05	126.56	118.80
1	BA	1324	G	N1-C6-O6	7.05	124.13	119.90
2	BB	36	C	C6-N1-C2	-7.05	117.48	120.30
1	AA	1123	C	C6-N1-C2	7.05	123.12	120.30
1	AA	1531	C	C2-N1-C1'	7.04	126.55	118.80
1	BA	2028	U	C5-C6-N1	7.04	126.22	122.70
1	BA	2473	U	C2-N1-C1'	7.04	126.15	117.70
1	BA	1225	G	N3-C4-N9	-7.04	121.78	126.00
1	AA	92	A	C8-N9-C4	-7.03	102.99	105.80
1	BA	587	C	C5-C6-N1	7.03	124.52	121.00
1	AA	528	A	N3-C4-C5	7.03	131.72	126.80
1	AA	1187	G	C8-N9-C4	-7.03	103.59	106.40
1	BA	1326	U	C5-C6-N1	-7.03	119.18	122.70
1	AA	1986	A	C8-N9-C4	7.03	108.61	105.80
1	BA	2251	G	N3-C4-N9	7.03	130.22	126.00
1	BA	1047	G	N3-C4-N9	7.01	130.21	126.00
1	BA	500	G	N9-C4-C5	7.01	108.20	105.40
1	BA	652(S)	C	N1-C2-O2	7.01	123.10	118.90
31	CA	1124	G	C5-C6-O6	7.00	132.80	128.60
1	AA	139(A)	G	C4-C5-N7	7.00	113.60	110.80
1	BA	1613	G	C5-C6-O6	6.99	132.79	128.60
1	BA	1359	A	N1-C2-N3	-6.99	125.81	129.30
31	DA	117	G	N3-C4-N9	6.99	130.19	126.00
1	AA	530	G	N9-C4-C5	6.98	108.19	105.40
1	AA	2241	A	C8-N9-C4	6.98	108.59	105.80
1	BA	330	A	N1-C2-N3	6.98	132.79	129.30
1	BA	2685	G	C8-N9-C4	-6.98	103.61	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	450	G	C4-C5-C6	6.97	122.98	118.80
1	AA	2361	A	N1-C6-N6	6.97	122.78	118.60
31	DA	96	U	N3-C2-O2	-6.97	117.32	122.20
1	AA	450	G	C4-C5-N7	-6.96	108.02	110.80
1	AA	298	G	N1-C6-O6	6.96	124.07	119.90
31	DA	839	U	N1-C2-O2	6.96	127.67	122.80
1	AA	928	G	N1-C6-O6	6.96	124.07	119.90
1	AA	1963	U	C6-N1-C1'	-6.95	111.48	121.20
1	BA	1827	C	N3-C2-O2	-6.95	117.04	121.90
1	BA	1210	A	C4-C5-N7	6.94	114.17	110.70
1	AA	141	A	N1-C6-N6	6.94	122.76	118.60
1	BA	1021	A	C5-C6-N1	-6.93	114.23	117.70
1	BA	1602	U	C6-N1-C2	-6.93	116.84	121.00
1	BA	2286	A	C6-C5-N7	-6.93	127.45	132.30
31	DA	1397	C	C6-N1-C2	-6.93	117.53	120.30
1	BA	1617	C	N1-C2-O2	-6.93	114.74	118.90
31	CA	1030(B)	C	C2-N1-C1'	6.93	126.42	118.80
1	AA	570	G	C4-C5-C6	6.92	122.95	118.80
2	AB	104	U	C6-N1-C2	6.92	125.15	121.00
1	BA	2591	C	C6-N1-C2	-6.92	117.53	120.30
31	CA	460	G	N7-C8-N9	6.91	116.56	113.10
1	AA	92	A	N7-C8-N9	6.90	117.25	113.80
31	DA	992	U	N3-C2-O2	-6.90	117.37	122.20
31	CA	226	G	C8-N9-C4	6.89	109.16	106.40
1	BA	90	U	N1-C2-N3	-6.88	110.77	114.90
1	AA	526	A	C8-N9-C4	-6.88	103.05	105.80
1	BA	12	U	C2-N1-C1'	6.88	125.95	117.70
1	AA	530	G	C8-N9-C4	-6.87	103.65	106.40
1	AA	893	C	C6-N1-C1'	-6.87	112.55	120.80
1	BA	2286	A	C5-N7-C8	-6.87	100.46	103.90
1	AA	568	U	N3-C4-C5	-6.87	110.48	114.60
1	AA	566	U	N3-C4-C5	6.87	118.72	114.60
1	BA	570	G	C5-N7-C8	6.87	107.73	104.30
1	BA	199	A	C8-N9-C4	6.86	108.54	105.80
1	BA	2544	G	C6-C5-N7	-6.86	126.28	130.40
1	AA	2104	G	C4-N9-C1'	6.86	135.41	126.50
1	BA	568	U	N3-C4-O4	6.85	124.20	119.40
31	CA	896	C	C6-N1-C2	6.85	123.04	120.30
1	BA	333	G	C4-N9-C1'	6.84	135.39	126.50
1	BA	2193	G	C5-C6-O6	-6.84	124.50	128.60
31	DA	117	G	C6-C5-N7	-6.84	126.30	130.40
1	AA	1698	A	N1-C6-N6	6.83	122.70	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	613	G	N1-C6-O6	6.83	124.00	119.90
1	AA	2104	G	C5-C6-O6	6.82	132.69	128.60
1	AA	2685	G	N9-C4-C5	6.82	108.13	105.40
1	AA	12	U	C2-N1-C1'	6.82	125.89	117.70
1	BA	298	G	N1-C6-O6	6.82	123.99	119.90
1	BA	2441	C	N3-C2-O2	-6.82	117.13	121.90
31	DA	1400	C	C6-N1-C2	6.82	123.03	120.30
1	BA	761	A	C5-N7-C8	-6.82	100.49	103.90
1	BA	2318	G	N7-C8-N9	6.81	116.51	113.10
1	AA	1602	U	N3-C4-C5	-6.81	110.51	114.60
31	DA	1028	C	C5-C6-N1	6.81	124.41	121.00
31	DA	150	C	C5-C6-N1	6.81	124.41	121.00
1	BA	2059	A	C8-N9-C4	6.81	108.52	105.80
1	BA	71	A	C6-C5-N7	-6.79	127.54	132.30
31	CA	1030	C	N1-C2-O2	6.79	122.98	118.90
1	AA	528	A	C5-N7-C8	-6.79	100.50	103.90
2	AB	64	C	C6-N1-C2	6.79	123.02	120.30
1	BA	500	G	C4-C5-N7	-6.78	108.09	110.80
1	BA	1210	A	C6-C5-N7	-6.77	127.56	132.30
1	AA	114	U	C2-N1-C1'	6.77	125.83	117.70
1	BA	546	C	C2-N1-C1'	6.77	126.24	118.80
1	BA	141	A	C5-C6-N6	-6.76	118.29	123.70
31	CA	1029	C	C2-N1-C1'	-6.76	111.36	118.80
1	AA	1382	G	N3-C4-C5	6.75	131.98	128.60
1	AA	1131	G	C4-N9-C1'	-6.74	117.74	126.50
1	AA	1899	G	N3-C2-N2	-6.74	115.18	119.90
1	BA	2056	G	N3-C2-N2	-6.74	115.18	119.90
1	BA	2590	A	C8-N9-C4	6.73	108.49	105.80
1	BA	2105	C	C6-N1-C2	-6.73	117.61	120.30
1	BA	2700	C	C6-N1-C2	6.73	122.99	120.30
1	BA	2505	G	C5-C6-O6	6.72	132.63	128.60
31	CA	1087	G	C8-N9-C4	-6.72	103.71	106.40
1	AA	2137	C	C2-N1-C1'	6.72	126.19	118.80
1	AA	1913	A	C8-N9-C4	-6.71	103.12	105.80
1	AA	2891	G	C5-C6-O6	-6.71	124.58	128.60
1	AA	1210	A	N1-C6-N6	6.70	122.62	118.60
1	AA	566	U	C5-C6-N1	-6.70	119.35	122.70
1	BA	2299	G	N3-C4-N9	-6.70	121.98	126.00
1	BA	1263	U	C6-N1-C1'	-6.70	111.82	121.20
1	BA	681	G	C8-N9-C4	6.69	109.08	106.40
1	AA	1914	C	C4-C5-C6	6.69	120.75	117.40
1	AA	1109	C	N3-C4-C5	-6.69	119.22	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2035	G	C8-N9-C1'	6.68	135.69	127.00
1	BA	2053	G	C5-C6-O6	-6.68	124.59	128.60
1	BA	1298	C	N1-C2-O2	6.68	122.91	118.90
1	BA	1638	C	C6-N1-C2	6.68	122.97	120.30
1	AA	2606	C	C6-N1-C2	6.68	122.97	120.30
1	AA	1602	U	N3-C4-O4	6.67	124.07	119.40
1	AA	2822	G	C8-N9-C4	6.67	109.07	106.40
1	BA	2056	G	C5-C6-O6	-6.67	124.60	128.60
1	BA	271(M)	G	C4-N9-C1'	6.66	135.16	126.50
1	AA	761	A	C5-N7-C8	-6.66	100.57	103.90
1	BA	2489	G	N9-C4-C5	-6.66	102.74	105.40
31	DA	438	G	N3-C4-C5	-6.66	125.27	128.60
1	BA	645	C	N3-C2-O2	-6.66	117.24	121.90
1	BA	2185	C	C2-N3-C4	6.66	123.23	119.90
1	AA	1204	A	C5-N7-C8	-6.65	100.57	103.90
1	AA	1698	A	C4-C5-N7	6.65	114.03	110.70
1	AA	2597	G	C8-N9-C4	6.65	109.06	106.40
31	DA	1030	C	N3-C2-O2	-6.64	117.25	121.90
1	BA	1022	G	N3-C4-N9	-6.64	122.02	126.00
1	BA	1840	G	N1-C6-O6	6.64	123.88	119.90
1	BA	180	G	N3-C4-N9	-6.63	122.02	126.00
1	BA	456	C	N1-C2-O2	-6.63	114.92	118.90
1	BA	1915	U	O4'-C1'-N1	6.63	113.50	108.20
2	BB	5	C	C6-N1-C2	6.63	122.95	120.30
31	CA	299	G	C5-C6-N1	-6.62	108.19	111.50
1	AA	1107	G	C4-N9-C1'	6.62	135.10	126.50
1	BA	1979	C	C6-N1-C2	-6.62	117.65	120.30
1	AA	1332	G	C5-N7-C8	-6.62	100.99	104.30
1	AA	2035	G	N3-C4-N9	-6.62	122.03	126.00
1	AA	1045	A	C2-N3-C4	6.61	113.91	110.60
31	CA	190	U	C5-C6-N1	6.61	126.01	122.70
1	BA	109	G	C5-C6-O6	6.61	132.57	128.60
1	AA	271(M)	G	N3-C4-N9	6.61	129.96	126.00
1	AA	645	C	C2-N1-C1'	6.60	126.06	118.80
31	CA	442	C	C5-C6-N1	6.60	124.30	121.00
31	DA	1126	U	N3-C2-O2	-6.60	117.58	122.20
1	BA	204	A	C5-C6-N1	6.60	121.00	117.70
1	AA	2137	C	N1-C2-O2	6.60	122.86	118.90
1	BA	2689	U	C2-N3-C4	-6.60	123.04	127.00
1	BA	614	U	N3-C2-O2	-6.59	117.58	122.20
1	AA	2439	A	C4-C5-N7	6.59	113.99	110.70
1	BA	60	G	C4-C5-N7	6.59	113.43	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	113	G	N3-C4-C5	6.58	131.89	128.60
1	BA	878	A	C8-N9-C4	-6.58	103.17	105.80
1	AA	512	G	N1-C6-O6	-6.58	115.95	119.90
1	AA	2131	G	C8-N9-C4	-6.58	103.77	106.40
1	BA	139(A)	G	C5-N7-C8	-6.58	101.01	104.30
1	BA	1827	C	C6-N1-C2	-6.58	117.67	120.30
2	BB	1	U	N1-C2-O2	6.57	127.40	122.80
1	AA	736	C	C6-N1-C2	6.57	122.93	120.30
31	CA	96	U	N3-C2-O2	-6.57	117.61	122.20
1	AA	34	C	N3-C4-C5	-6.56	119.27	121.90
1	BA	435	C	N1-C2-O2	6.55	122.83	118.90
31	CA	76	C	C2-N1-C1'	6.55	126.01	118.80
1	AA	2497	A	N9-C4-C5	-6.55	103.18	105.80
1	BA	2251	G	N3-C4-C5	-6.55	125.33	128.60
1	AA	1611	C	N3-C4-C5	6.54	124.52	121.90
1	BA	14	A	N9-C4-C5	-6.54	103.18	105.80
1	AA	124	G	N3-C4-C5	6.54	131.87	128.60
1	BA	452	G	N1-C6-O6	-6.54	115.98	119.90
1	BA	2137	C	C2-N1-C1'	6.54	125.99	118.80
1	AA	2318	G	C4-N9-C1'	6.53	134.99	126.50
1	BA	450	G	C4-C5-C6	6.53	122.72	118.80
1	AA	746	A	C8-N9-C4	6.53	108.41	105.80
31	DA	1272	G	C8-N9-C1'	-6.53	118.51	127.00
1	BA	1372	U	C5-C4-O4	-6.53	121.98	125.90
31	DA	1030	C	C6-N1-C2	-6.53	117.69	120.30
1	AA	1531	C	C6-N1-C2	-6.53	117.69	120.30
1	BA	2791	C	C2-N1-C1'	6.52	125.98	118.80
1	BA	1021	A	C5-N7-C8	-6.52	100.64	103.90
1	BA	1772	G	C5-C6-O6	6.52	132.51	128.60
1	AA	1022	G	C4-C5-N7	-6.52	108.19	110.80
2	AB	104	U	C5-C6-N1	-6.52	119.44	122.70
1	BA	330	A	C5-N7-C8	-6.52	100.64	103.90
31	CA	906	G	N3-C4-C5	-6.52	125.34	128.60
1	BA	525	U	C6-N1-C2	-6.52	117.09	121.00
1	AA	1393	A	C8-N9-C4	6.51	108.41	105.80
1	BA	512	G	C4-N9-C1'	-6.51	118.04	126.50
1	BA	271(M)	G	C8-N9-C1'	-6.50	118.54	127.00
31	DA	117	G	N9-C4-C5	-6.50	102.80	105.40
1	BA	915	C	N3-C2-O2	-6.50	117.35	121.90
1	AA	298	G	C4-C5-N7	6.50	113.40	110.80
31	DA	1272	G	C4-N9-C1'	6.49	134.94	126.50
1	AA	2420	C	C5-C4-N4	-6.49	115.66	120.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	143	G	N3-C4-C5	6.49	131.84	128.60
1	BA	1955	U	C5-C6-N1	-6.48	119.46	122.70
1	BA	452	G	C2-N3-C4	6.47	115.14	111.90
1	AA	1393	A	N9-C4-C5	-6.47	103.21	105.80
1	AA	1899	G	C8-N9-C4	-6.47	103.81	106.40
1	BA	198	C	C2-N1-C1'	6.47	125.92	118.80
1	BA	2419	U	C5-C6-N1	6.46	125.93	122.70
1	BA	1699	G	C8-N9-C4	-6.46	103.81	106.40
1	BA	1047	G	N3-C4-C5	-6.46	125.37	128.60
1	AA	928	G	C5-C6-O6	-6.45	124.73	128.60
1	AA	1962	C	N1-C2-O2	6.45	122.77	118.90
1	BA	2609	U	N3-C2-O2	6.45	126.72	122.20
1	BA	143	G	C4-N9-C1'	-6.45	118.11	126.50
1	AA	1142(A)	A	N3-C4-N9	-6.45	122.24	127.40
1	AA	1107	G	N3-C4-C5	-6.45	125.38	128.60
1	AA	2544	G	N1-C6-O6	6.45	123.77	119.90
1	BA	527	C	N3-C4-N4	-6.45	113.49	118.00
1	BA	1210	A	C5-C6-N6	-6.45	118.54	123.70
1	AA	190	A	N1-C2-N3	-6.45	126.08	129.30
1	BA	2187	G	C4-C5-N7	6.45	113.38	110.80
1	BA	2590	A	C4-C5-C6	-6.45	113.78	117.00
1	BA	2673	G	N1-C6-O6	6.44	123.77	119.90
1	AA	1246	A	C8-N9-C4	6.44	108.38	105.80
1	AA	1204	A	C2-N3-C4	-6.43	107.38	110.60
31	CA	1397	C	C2-N1-C1'	6.43	125.88	118.80
1	BA	690	G	C8-N9-C4	6.43	108.97	106.40
1	BA	2056	G	N1-C6-O6	6.43	123.75	119.90
31	CA	96	U	N1-C2-O2	6.43	127.30	122.80
1	BA	1982	C	N3-C4-C5	-6.42	119.33	121.90
31	DA	885	G	C8-N9-C4	6.42	108.97	106.40
19	AX	57	LEU	CA-CB-CG	6.42	130.07	115.30
31	DA	1492	A	C2-N3-C4	6.42	113.81	110.60
1	AA	1050	A	C8-N9-C4	-6.41	103.23	105.80
19	BX	57	LEU	CA-CB-CG	6.41	130.03	115.30
31	DA	1038	C	N3-C2-O2	-6.40	117.42	121.90
1	BA	652(H)	C	N1-C2-O2	6.40	122.74	118.90
1	BA	865	C	C6-N1-C2	6.40	122.86	120.30
31	DA	1027	C	N3-C4-N4	-6.40	113.52	118.00
1	AA	2263	C	C6-N1-C2	6.40	122.86	120.30
31	DA	503	C	C6-N1-C2	-6.40	117.74	120.30
31	DA	1030	C	C2-N1-C1'	6.39	125.83	118.80
1	AA	1313	U	C2-N1-C1'	6.39	125.37	117.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1962	C	C2-N1-C1'	6.39	125.83	118.80
1	BA	686	G	N3-C2-N2	6.39	124.37	119.90
31	CA	1030(B)	C	N1-C2-O2	6.39	122.73	118.90
34	CD	26	CYS	CA-CB-SG	6.38	125.49	114.00
31	DA	1420	C	C6-N1-C2	-6.38	117.75	120.30
2	AB	1	U	C2-N1-C1'	6.38	125.35	117.70
1	BA	978	G	C8-N9-C4	6.38	108.95	106.40
1	AA	271(M)	G	N3-C4-C5	-6.38	125.41	128.60
1	AA	2014	A	N1-C6-N6	6.37	122.42	118.60
1	BA	1021	A	N1-C6-N6	6.37	122.42	118.60
1	AA	2137	C	N3-C4-C5	-6.37	119.35	121.90
31	DA	90	C	C6-N1-C2	-6.37	117.75	120.30
31	CA	1493	A	N1-C6-N6	6.36	122.42	118.60
1	BA	741	G	C5-C6-O6	-6.36	124.78	128.60
1	BA	1913	A	C3'-C2'-C1'	-6.36	96.41	101.50
1	BA	2318	G	C8-N9-C4	-6.36	103.86	106.40
31	CA	76	C	C6-N1-C2	-6.36	117.76	120.30
1	BA	928	G	N1-C6-O6	6.36	123.71	119.90
1	AA	568	U	N1-C2-O2	-6.35	118.35	122.80
1	BA	827	U	C6-N1-C2	6.35	124.81	121.00
1	BA	2218	U	C6-N1-C1'	-6.35	112.31	121.20
1	AA	1992	G	C8-N9-C4	-6.35	103.86	106.40
1	BA	1772	G	C4-C5-N7	-6.35	108.26	110.80
1	AA	141	A	C5-N7-C8	-6.34	100.73	103.90
1	BA	1840	G	C4-C5-C6	6.34	122.61	118.80
31	CA	76	C	C5-C6-N1	6.34	124.17	121.00
1	AA	1664	A	C8-N9-C4	-6.34	103.26	105.80
1	BA	1653	G	N3-C4-C5	-6.34	125.43	128.60
1	BA	1997	G	C4-N9-C1'	-6.34	118.26	126.50
1	BA	2593	U	N3-C4-C5	-6.34	110.80	114.60
31	DA	71	C	N1-C2-O2	6.34	122.70	118.90
1	BA	1290	C	C6-N1-C2	-6.33	117.77	120.30
1	AA	2877	G	C8-N9-C4	6.33	108.93	106.40
1	BA	1840	G	C4-N9-C1'	6.33	134.73	126.50
1	AA	2497	A	C8-N9-C4	6.33	108.33	105.80
31	CA	992	U	C5-C6-N1	6.33	125.86	122.70
1	AA	1962	C	C5-C6-N1	6.32	124.16	121.00
1	BA	885	C	C2-N1-C1'	6.32	125.75	118.80
1	AA	507	A	C8-N9-C4	6.32	108.33	105.80
1	BA	2286	A	C4-C5-N7	6.32	113.86	110.70
1	AA	1966	A	C8-N9-C4	6.32	108.33	105.80
1	BA	678	C	C6-N1-C2	6.32	122.83	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2024	G	N3-C4-C5	6.31	131.76	128.60
31	CA	342	C	C6-N1-C2	-6.31	117.77	120.30
1	AA	1840	G	C5-C6-N1	-6.31	108.34	111.50
1	AA	465	G	N3-C4-C5	-6.31	125.44	128.60
1	AA	2456	C	C6-N1-C2	6.31	122.82	120.30
1	BA	2273	A	C8-N9-C4	-6.31	103.28	105.80
1	AA	1531	C	C5-C6-N1	6.31	124.15	121.00
1	AA	652(Q)	G	N3-C4-N9	6.31	129.78	126.00
1	AA	1963	U	C5-C6-N1	6.30	125.85	122.70
1	BA	661	C	C6-N1-C2	-6.30	117.78	120.30
1	AA	1204	A	C4-C5-N7	6.30	113.85	110.70
2	AB	54	G	C8-N9-C4	-6.29	103.88	106.40
1	AA	2419	U	N3-C4-C5	-6.29	110.83	114.60
1	BA	512	G	C4-C5-C6	-6.29	115.03	118.80
1	AA	494	G	C8-N9-C4	6.29	108.91	106.40
1	BA	446	G	N9-C4-C5	-6.29	102.89	105.40
1	BA	500	G	C5-C6-O6	6.28	132.37	128.60
1	BA	1187	G	N1-C6-O6	6.28	123.67	119.90
1	BA	1633	G	C5-C6-N1	-6.28	108.36	111.50
1	AA	399	G	C8-N9-C4	6.28	108.91	106.40
1	AA	2791	C	C2-N1-C1'	6.28	125.70	118.80
1	AA	530	G	C5-C6-O6	6.27	132.36	128.60
1	BA	1313	U	N3-C2-O2	-6.27	117.81	122.20
31	DA	530	G	C8-N9-C1'	-6.27	118.85	127.00
1	AA	2334	G	N9-C4-C5	-6.27	102.89	105.40
1	AA	2411	A	C8-N9-C4	6.27	108.31	105.80
1	BA	2827	C	C4-C5-C6	-6.26	114.27	117.40
31	DA	175	C	C6-N1-C2	-6.26	117.80	120.30
1	BA	2894	G	N7-C8-N9	6.26	116.23	113.10
1	BA	1951	U	C6-N1-C2	-6.25	117.25	121.00
1	BA	2053	G	N1-C6-O6	6.25	123.65	119.90
1	BA	645	C	C6-N1-C2	-6.25	117.80	120.30
1	AA	870	A	C8-N9-C4	6.25	108.30	105.80
1	AA	845	G	C4-N9-C1'	6.25	134.62	126.50
1	AA	1216	G	N1-C6-O6	6.25	123.65	119.90
1	BA	226	G	C5-C6-O6	-6.24	124.85	128.60
1	AA	1531	C	N1-C2-O2	6.24	122.64	118.90
1	BA	390	A	C8-N9-C4	6.24	108.30	105.80
1	BA	686	G	N9-C4-C5	-6.24	102.91	105.40
1	BA	2481	G	C5-C6-O6	-6.24	124.86	128.60
1	BA	2318	G	C4-N9-C1'	6.24	134.61	126.50
34	DD	26	CYS	CA-CB-SG	6.24	125.23	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	265	A	N1-C6-N6	6.24	122.34	118.60
1	BA	568	U	N1-C2-O2	-6.23	118.44	122.80
1	BA	584	C	N1-C2-O2	-6.23	115.16	118.90
1	BA	956	G	C5-C6-N1	-6.23	108.39	111.50
1	BA	1602	U	N3-C4-O4	6.23	123.76	119.40
31	DA	150	C	C6-N1-C2	-6.23	117.81	120.30
31	CA	1245	A	C5-C6-N6	-6.23	118.72	123.70
1	AA	2036	C	C6-N1-C2	-6.22	117.81	120.30
1	BA	673	C	N3-C2-O2	-6.22	117.55	121.90
1	AA	129	C	C5-C6-N1	-6.22	117.89	121.00
1	BA	1828	G	N1-C6-O6	-6.22	116.17	119.90
1	AA	2681	C	C6-N1-C2	6.22	122.79	120.30
1	AA	2253	G	N9-C4-C5	-6.22	102.91	105.40
1	BA	2482	G	C4-N9-C1'	6.22	134.58	126.50
1	BA	2137	C	N1-C2-O2	6.21	122.63	118.90
1	AA	2249	U	C6-N1-C2	-6.21	117.27	121.00
1	AA	1855	G	C8-N9-C4	6.21	108.89	106.40
31	CA	433	C	N3-C2-O2	-6.21	117.55	121.90
1	BA	60	G	C8-N9-C4	6.21	108.88	106.40
1	BA	1819	A	N1-C6-N6	-6.21	114.88	118.60
1	AA	71	A	C5-N7-C8	-6.20	100.80	103.90
2	AB	1	U	N1-C2-O2	6.20	127.14	122.80
1	AA	933	A	N1-C6-N6	6.20	122.32	118.60
1	AA	600	G	C8-N9-C4	6.19	108.88	106.40
1	BA	1275	A	N1-C6-N6	6.19	122.32	118.60
1	AA	827	U	N3-C2-O2	6.19	126.53	122.20
1	BA	652(Q)	G	N3-C4-N9	6.19	129.71	126.00
31	DA	699	C	C6-N1-C2	-6.19	117.82	120.30
31	CA	1034	G	C6-N1-C2	6.19	128.81	125.10
1	BA	568	U	C4-C5-C6	6.19	123.41	119.70
1	AA	1779	U	C4-C5-C6	6.18	123.41	119.70
1	BA	2429	G	N3-C4-C5	-6.18	125.51	128.60
1	AA	645	C	N3-C2-O2	-6.18	117.57	121.90
1	AA	512	G	N9-C4-C5	6.18	107.87	105.40
1	BA	142(A)	C	N1-C2-O2	-6.18	115.19	118.90
1	BA	2593	U	C6-N1-C2	-6.18	117.29	121.00
1	AA	1844	C	C6-N1-C2	-6.17	117.83	120.30
1	BA	1840	G	C6-C5-N7	-6.17	126.70	130.40
31	DA	1027	C	C6-N1-C1'	6.17	128.21	120.80
1	AA	90	U	C5-C6-N1	6.17	125.78	122.70
1	AA	679	C	N3-C2-O2	6.17	126.22	121.90
1	BA	1243	G	N3-C4-C5	6.17	131.68	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	2685	G	N3-C4-N9	-6.17	122.30	126.00
1	AA	2206	G	C8-N9-C4	6.16	108.86	106.40
1	AA	676	A	C6-C5-N7	-6.16	127.99	132.30
1	BA	1791	A	C6-C5-N7	-6.16	127.99	132.30
1	AA	1142(A)	A	C5-C6-N1	-6.15	114.62	117.70
1	AA	945	A	C5-N7-C8	-6.15	100.82	103.90
34	CD	12	CYS	CA-CB-SG	6.15	125.07	114.00
1	BA	202	U	N1-C2-O2	6.15	127.11	122.80
1	AA	124	G	C2-N3-C4	-6.15	108.83	111.90
31	DA	242	C	N1-C2-O2	-6.15	115.21	118.90
1	BA	1779	U	C5-C4-O4	6.14	129.59	125.90
1	AA	2685	G	C4-C5-N7	-6.14	108.34	110.80
1	BA	29	U	C5-C4-O4	-6.14	122.22	125.90
1	BA	265	A	C2-N3-C4	-6.14	107.53	110.60
31	CA	992	U	C6-N1-C1'	-6.14	112.61	121.20
31	CA	925	G	C5-C6-O6	-6.14	124.92	128.60
1	BA	652(R)	C	N1-C2-O2	6.13	122.58	118.90
1	BA	2022	U	N1-C2-N3	-6.13	111.22	114.90
1	AA	1647	G	N3-C4-N9	-6.13	122.32	126.00
1	BA	1204	A	C6-C5-N7	-6.13	128.01	132.30
1	AA	928	G	C6-C5-N7	-6.13	126.72	130.40
31	CA	1293	G	N3-C4-C5	-6.13	125.54	128.60
1	BA	1343	G	C4-N9-C1'	6.12	134.46	126.50
1	BA	2894	G	N1-C6-O6	6.12	123.57	119.90
1	BA	512	G	N3-C4-N9	-6.12	122.33	126.00
1	BA	1203	G	N1-C6-O6	-6.12	116.23	119.90
31	CA	1294	G	C8-N9-C1'	6.12	134.95	127.00
27	A5	58	LEU	CA-CB-CG	6.12	129.37	115.30
1	BA	468	G	C8-N9-C4	6.12	108.85	106.40
1	BA	263	C	C6-N1-C2	6.11	122.75	120.30
31	CA	1007	C	C5-C6-N1	6.11	124.06	121.00
1	BA	180	G	N3-C4-C5	6.11	131.65	128.60
1	BA	1882	C	C2-N1-C1'	6.11	125.52	118.80
1	BA	574	C	N1-C2-O2	-6.10	115.24	118.90
1	AA	129	C	C6-N1-C2	6.10	122.74	120.30
1	AA	948	G	N1-C6-O6	6.10	123.56	119.90
1	BA	450	G	N1-C2-N2	6.10	121.69	116.20
1	AA	885	C	C2-N1-C1'	6.09	125.50	118.80
31	DA	345	C	C2-N1-C1'	-6.09	112.10	118.80
1	AA	330	A	N1-C2-N3	6.09	132.34	129.30
1	BA	33	U	C2-N1-C1'	-6.09	110.40	117.70
1	BA	2866	U	C4-C5-C6	6.09	123.35	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	546	C	C6-N1-C2	-6.08	117.87	120.30
1	BA	729	G	N3-C2-N2	-6.08	115.64	119.90
1	BA	1699	G	N1-C6-O6	-6.08	116.25	119.90
1	AA	1004	C	C6-N1-C2	6.08	122.73	120.30
31	CA	1067	A	C2-N3-C4	6.08	113.64	110.60
1	AA	1216	G	C5-C6-O6	-6.07	124.96	128.60
2	AB	24	G	N3-C4-N9	6.07	129.64	126.00
31	DA	1282	C	C6-N1-C2	-6.07	117.87	120.30
1	BA	500	G	N1-C6-O6	-6.07	116.26	119.90
31	CA	1432	G	N1-C6-O6	6.07	123.54	119.90
31	DA	76	C	C5-C6-N1	6.07	124.03	121.00
1	AA	1616	A	N1-C6-N6	6.07	122.24	118.60
1	AA	1672	C	C6-N1-C2	6.07	122.73	120.30
1	AA	297	C	C6-N1-C2	-6.07	117.87	120.30
1	BA	676	A	C4-C5-N7	6.07	113.73	110.70
1	AA	1154	G	N3-C4-N9	6.06	129.64	126.00
1	AA	265	A	N7-C8-N9	6.06	116.83	113.80
1	BA	2896	C	C2-N1-C1'	6.06	125.46	118.80
15	BT	118	ARG	NE-CZ-NH1	6.06	123.33	120.30
1	AA	2721	A	C8-N9-C4	6.05	108.22	105.80
31	DA	1028	C	C6-N1-C2	-6.05	117.88	120.30
1	BA	795	C	N1-C2-O2	6.04	122.53	118.90
31	DA	1030(B)	C	C6-N1-C1'	-6.04	113.55	120.80
31	DA	412	A	C8-N9-C4	6.04	108.22	105.80
1	BA	151	C	C5-C6-N1	-6.04	117.98	121.00
1	BA	2206	G	N3-C4-C5	6.04	131.62	128.60
31	CA	1017	G	C5-C6-N1	-6.04	108.48	111.50
1	BA	723	G	C8-N9-C1'	-6.03	119.16	127.00
1	AA	1986	A	N7-C8-N9	-6.03	110.78	113.80
1	BA	90	U	C2-N3-C4	6.03	130.62	127.00
31	CA	1102	A	C8-N9-C4	-6.03	103.39	105.80
31	DA	1027	C	C2-N1-C1'	-6.03	112.17	118.80
1	BA	1201	C	C6-N1-C2	6.03	122.71	120.30
1	AA	933	A	N7-C8-N9	6.03	116.81	113.80
1	AA	2464	C	C6-N1-C2	6.03	122.71	120.30
1	BA	1982	C	C2-N1-C1'	6.03	125.43	118.80
31	DA	952	U	C2-N3-C4	6.03	130.62	127.00
1	AA	570	G	C4-N9-C1'	6.02	134.33	126.50
1	BA	584	C	N3-C2-O2	6.02	126.11	121.90
1	BA	333	G	C8-N9-C1'	-6.02	119.17	127.00
1	BA	2028	U	C4-C5-C6	6.02	123.31	119.70
1	AA	856	C	C6-N1-C2	-6.01	117.90	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	723	G	C4-N9-C1'	6.01	134.32	126.50
36	DF	19	LEU	CA-CB-CG	6.01	129.12	115.30
31	CA	1102	A	C6-N1-C2	-6.01	114.99	118.60
1	AA	1187	G	N7-C8-N9	6.01	116.10	113.10
1	AA	2206	G	C4-N9-C1'	-6.01	118.69	126.50
31	DA	178	C	C6-N1-C2	-6.00	117.90	120.30
2	BB	6	C	C6-N1-C2	6.00	122.70	120.30
1	BA	179	G	C5-C6-N1	-6.00	108.50	111.50
1	BA	2680	C	C6-N1-C2	6.00	122.70	120.30
31	CA	174	C	C6-N1-C2	-5.99	117.90	120.30
1	BA	2621	A	C8-N9-C4	5.99	108.20	105.80
1	BA	188	G	C8-N9-C4	5.99	108.80	106.40
1	AA	1266	G	N7-C8-N9	-5.99	110.11	113.10
1	BA	715	G	C6-C5-N7	-5.98	126.81	130.40
1	BA	1324	G	C5-C6-N1	-5.98	108.51	111.50
1	AA	1432	C	C6-N1-C2	5.98	122.69	120.30
1	BA	652(G)	G	N3-C4-N9	5.98	129.59	126.00
1	BA	2137	C	C5-C6-N1	5.98	123.99	121.00
31	DA	670	G	N3-C4-C5	-5.98	125.61	128.60
1	AA	1882	C	C6-N1-C2	-5.97	117.91	120.30
1	BA	912	C	C6-N1-C2	-5.97	117.91	120.30
1	BA	2685	G	C4-C5-C6	5.97	122.39	118.80
1	AA	1142(A)	A	C8-N9-C4	-5.97	103.41	105.80
1	BA	1187	G	C4-C5-C6	5.97	122.38	118.80
1	BA	729	G	N1-C2-N2	5.97	121.57	116.20
1	AA	213	A	C8-N9-C4	5.97	108.19	105.80
1	AA	945	A	N1-C6-N6	5.97	122.18	118.60
31	DA	1138	G	N3-C4-N9	5.97	129.58	126.00
1	BA	944	G	C4-N9-C1'	5.96	134.25	126.50
1	BA	2299	G	N3-C4-C5	5.96	131.58	128.60
1	BA	1324	G	C8-N9-C4	5.96	108.78	106.40
31	DA	76	C	C2-N1-C1'	5.96	125.36	118.80
1	AA	679	C	C6-N1-C2	5.96	122.68	120.30
1	BA	229	A	C8-N9-C4	-5.96	103.42	105.80
1	BA	1326	U	C6-N1-C2	5.96	124.57	121.00
1	BA	1698	A	C5-C6-N1	-5.95	114.72	117.70
1	AA	512	G	C8-N9-C1'	5.95	134.74	127.00
1	BA	2187	G	C5-C6-O6	-5.95	125.03	128.60
1	BA	1110	G	N9-C4-C5	5.95	107.78	105.40
31	CA	906	G	N3-C4-N9	5.95	129.57	126.00
1	BA	463	G	N1-C6-O6	-5.95	116.33	119.90
1	BA	2489	G	C6-C5-N7	-5.95	126.83	130.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2361	A	C4-C5-N7	5.95	113.67	110.70
1	BA	1488	G	C8-N9-C4	-5.95	104.02	106.40
31	CA	90	C	C6-N1-C2	-5.95	117.92	120.30
1	BA	1743	C	C6-N1-C2	-5.94	117.92	120.30
1	BA	2689	U	C5-C6-N1	-5.94	119.73	122.70
1	AA	1915	U	C2-N1-C1'	5.94	124.83	117.70
1	AA	1493	C	C6-N1-C1'	-5.94	113.67	120.80
1	BA	1109	C	C4-C5-C6	5.94	120.37	117.40
31	DA	943	U	C5-C4-O4	5.94	129.46	125.90
1	AA	912	C	N3-C2-O2	-5.94	117.75	121.90
1	BA	570	G	C8-N9-C4	-5.93	104.03	106.40
1	BA	1493	C	C2-N1-C1'	5.93	125.33	118.80
1	BA	570	G	C4-C5-C6	5.93	122.36	118.80
31	CA	992	U	N3-C2-O2	-5.93	118.05	122.20
1	AA	389	G	N9-C4-C5	-5.93	103.03	105.40
1	BA	639	U	C5-C4-O4	5.93	129.46	125.90
1	BA	915	C	N1-C2-O2	5.93	122.46	118.90
1	BA	143	G	N3-C4-N9	-5.93	122.44	126.00
1	BA	641	C	C6-N1-C2	-5.92	117.93	120.30
1	BA	2059	A	N7-C8-N9	-5.92	110.84	113.80
1	BA	2343	C	C6-N1-C2	-5.92	117.93	120.30
31	CA	1010	G	N7-C8-N9	5.92	116.06	113.10
31	CA	1397	C	C6-N1-C2	-5.92	117.93	120.30
31	DA	117	G	C8-N9-C1'	-5.92	119.30	127.00
1	AA	141	A	C6-C5-N7	-5.92	128.15	132.30
1	BA	652(T)	C	N1-C2-O2	5.92	122.45	118.90
1	BA	1142(A)	A	C5-N7-C8	-5.92	100.94	103.90
1	BA	1602	U	C5-C4-O4	5.92	129.45	125.90
1	AA	802	A	N9-C4-C5	-5.92	103.43	105.80
1	AA	298	G	C5-C6-O6	-5.92	125.05	128.60
1	AA	912	C	N3-C4-C5	-5.91	119.53	121.90
2	AB	114	C	C5-C6-N1	-5.91	118.04	121.00
1	BA	330	A	N3-C4-N9	-5.91	122.67	127.40
1	BA	945	A	N7-C8-N9	5.91	116.76	113.80
1	BA	1210	A	C2-N3-C4	-5.91	107.64	110.60
1	BA	2581	G	C5-C6-O6	5.91	132.15	128.60
1	BA	265	A	C5-N7-C8	-5.91	100.95	103.90
1	BA	1602	U	C2-N3-C4	5.91	130.54	127.00
1	BA	221	A	C8-N9-C4	-5.91	103.44	105.80
31	DA	1494	G	N3-C4-C5	-5.91	125.65	128.60
1	BA	2137	C	N3-C2-O2	-5.90	117.77	121.90
31	CA	529	G	N1-C6-O6	5.90	123.44	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	1183	G	N1-C6-O6	5.90	123.44	119.90
1	BA	2685	G	N1-C6-O6	5.90	123.44	119.90
1	AA	2440	C	N1-C2-O2	5.90	122.44	118.90
1	BA	761	A	C4-C5-C6	-5.90	114.05	117.00
31	DA	266	G	C4-C5-N7	5.90	113.16	110.80
1	AA	465	G	C4-C5-C6	5.89	122.34	118.80
1	AA	733	G	N3-C4-N9	-5.89	122.47	126.00
1	AA	2439	A	C5-N7-C8	-5.89	100.96	103.90
1	AA	582	G	N1-C6-O6	5.89	123.43	119.90
2	AB	31	C	N1-C2-O2	5.89	122.43	118.90
1	AA	2306	C	C5-C6-N1	5.88	123.94	121.00
1	AA	2318	G	N7-C8-N9	5.88	116.04	113.10
31	CA	1294	G	N9-C4-C5	5.88	107.75	105.40
31	DA	354	G	N3-C4-N9	5.88	129.53	126.00
31	DA	606	G	N3-C4-C5	-5.88	125.66	128.60
1	AA	1107	G	N3-C4-N9	5.88	129.53	126.00
1	AA	527	C	N3-C4-N4	-5.88	113.89	118.00
1	AA	1204	A	N7-C8-N9	5.88	116.74	113.80
1	BA	1049	C	C5-C6-N1	5.87	123.94	121.00
31	CA	365	U	C5-C6-N1	-5.87	119.76	122.70
1	BA	1959	G	C5-C6-O6	5.87	132.12	128.60
1	AA	141	A	C4-C5-N7	5.87	113.63	110.70
1	AA	2318	G	C8-N9-C4	-5.87	104.05	106.40
1	BA	1021	A	N3-C4-C5	5.86	130.90	126.80
1	BA	2251	G	C4-N9-C1'	5.86	134.12	126.50
1	AA	1333	C	C5-C4-N4	-5.86	116.10	120.20
1	AA	837	C	C6-N1-C2	-5.86	117.95	120.30
1	AA	2757	A	N7-C8-N9	5.86	116.73	113.80
1	BA	1314	C	C6-N1-C1'	-5.86	113.77	120.80
1	AA	1671	U	C5-C4-O4	-5.86	122.39	125.90
1	AA	1920	C	C6-N1-C2	-5.86	117.96	120.30
1	AA	2559	C	N3-C4-C5	5.86	124.24	121.90
1	BA	1187	G	C4-C5-N7	-5.85	108.46	110.80
1	BA	2574	G	C4-C5-N7	5.85	113.14	110.80
31	DA	1313	U	C5-C4-O4	-5.85	122.39	125.90
1	BA	527	C	C5-C4-N4	5.85	124.30	120.20
31	DA	76	C	C4-C5-C6	-5.85	114.48	117.40
1	AA	2514	U	C6-N1-C2	5.85	124.51	121.00
31	DA	1126	U	C5-C6-N1	5.85	125.62	122.70
1	BA	659	C	N3-C4-C5	5.84	124.24	121.90
1	AA	2617	C	C6-N1-C2	5.84	122.64	120.30
1	AA	1115	G	C4-N9-C1'	-5.83	118.92	126.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2036	C	C5-C6-N1	5.83	123.92	121.00
1	BA	1195	G	N3-C2-N2	-5.83	115.82	119.90
1	AA	733	G	C5-N7-C8	-5.83	101.39	104.30
1	AA	1914	C	N3-C2-O2	-5.83	117.82	121.90
1	BA	450	G	N3-C4-N9	-5.82	122.51	126.00
31	DA	1290	G	N3-C4-C5	-5.82	125.69	128.60
1	AA	110	G	C8-N9-C4	5.82	108.73	106.40
1	BA	2429	G	C8-N9-C4	-5.82	104.07	106.40
1	AA	271(M)	G	C4-N9-C1'	5.81	134.06	126.50
1	AA	1607	C	C6-N1-C2	5.80	122.62	120.30
1	AA	2544	G	N3-C2-N2	-5.80	115.84	119.90
1	BA	886	C	C2-N1-C1'	5.80	125.19	118.80
31	DA	1030(B)	C	N1-C2-O2	5.80	122.38	118.90
1	AA	1187	G	C5-C6-N1	-5.80	108.60	111.50
31	CA	1502	A	C2-N3-C4	-5.80	107.70	110.60
1	BA	204	A	N3-C4-C5	-5.80	122.74	126.80
1	BA	1022	G	C4-N9-C1'	-5.80	118.97	126.50
31	CA	1087	G	N3-C4-N9	5.80	129.48	126.00
1	BA	2540	C	C6-N1-C2	5.79	122.62	120.30
1	BA	226	G	N1-C6-O6	5.79	123.38	119.90
1	BA	1314	C	C2-N1-C1'	5.79	125.17	118.80
31	CA	460	G	C4-N9-C1'	5.79	134.03	126.50
31	DA	1184	G	C4-N9-C1'	-5.79	118.97	126.50
1	BA	500	G	C6-C5-N7	5.79	133.88	130.40
1	BA	1426	G	C4-C5-N7	5.79	113.12	110.80
31	DA	754	C	C2-N1-C1'	5.79	125.17	118.80
1	AA	2018	G	C5-C6-O6	-5.79	125.13	128.60
1	AA	666	G	N9-C4-C5	-5.78	103.09	105.40
31	DA	1184	G	C8-N9-C1'	5.78	134.52	127.00
1	AA	1005	C	C5-C6-N1	-5.78	118.11	121.00
1	BA	92	A	C8-N9-C4	-5.78	103.49	105.80
1	BA	2689	U	N1-C2-N3	5.78	118.37	114.90
31	DA	355	C	C6-N1-C2	-5.77	117.99	120.30
1	AA	652(H)	C	N1-C2-O2	5.77	122.36	118.90
31	DA	1378	C	C5-C6-N1	5.77	123.89	121.00
1	AA	1698	A	C6-C5-N7	-5.77	128.26	132.30
1	BA	271(J)	C	N1-C2-O2	5.77	122.36	118.90
1	BA	1963	U	C6-N1-C1'	-5.77	113.12	121.20
1	BA	465	G	C8-N9-C4	-5.77	104.09	106.40
1	AA	2306	C	C2-N1-C1'	5.76	125.14	118.80
1	BA	2251	G	C8-N9-C1'	-5.76	119.50	127.00
31	DA	1260	C	C6-N1-C2	-5.76	118.00	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1653	G	C4-N9-C1'	5.76	133.99	126.50
1	AA	2593	U	N3-C4-C5	-5.76	111.14	114.60
31	CA	96	U	C2-N1-C1'	5.76	124.61	117.70
31	CA	1150	U	C5-C4-O4	5.76	129.36	125.90
31	DA	245	C	C2-N1-C1'	-5.76	112.47	118.80
31	CA	1026	G	C8-N9-C4	-5.76	104.10	106.40
31	DA	293	G	N1-C6-O6	5.76	123.35	119.90
31	CA	1123	A	C5-C6-N6	5.75	128.30	123.70
31	DA	1037	C	C5-C6-N1	5.75	123.88	121.00
31	DA	1038	C	C6-N1-C2	-5.75	118.00	120.30
31	DA	438	G	N3-C4-N9	5.75	129.45	126.00
1	BA	1184	G	N1-C6-O6	5.75	123.35	119.90
1	BA	1204	A	C4-C5-N7	5.74	113.57	110.70
1	BA	2375	G	C8-N9-C4	5.74	108.70	106.40
1	AA	2382	G	N3-C4-N9	5.74	129.44	126.00
31	CA	956	U	C5-C6-N1	5.74	125.57	122.70
31	DA	1036	G	C4-N9-C1'	5.74	133.96	126.50
31	DA	1391	U	C5-C4-O4	5.74	129.34	125.90
1	AA	298	G	C5-N7-C8	-5.74	101.43	104.30
2	BB	80	U	C5-C4-O4	5.74	129.34	125.90
31	CA	93	G	N9-C4-C5	-5.73	103.11	105.40
1	BA	883	G	C5-C6-O6	-5.73	125.16	128.60
31	DA	1121	U	C5-C6-N1	5.73	125.56	122.70
31	DA	93	G	N9-C4-C5	-5.73	103.11	105.40
1	AA	696	G	C8-N9-C4	5.73	108.69	106.40
1	BA	1203	G	N9-C4-C5	5.72	107.69	105.40
1	BA	1383	C	N1-C2-O2	-5.72	115.47	118.90
31	CA	1087	G	C4-N9-C1'	5.72	133.94	126.50
31	DA	48	C	C2-N1-C1'	5.72	125.10	118.80
3	BD	274	ARG	C-N-CA	5.72	136.01	121.70
1	BA	2585	U	N1-C2-O2	5.72	126.81	122.80
1	AA	1193	G	C8-N9-C4	5.72	108.69	106.40
2	BB	106	G	N3-C2-N2	-5.72	115.90	119.90
31	DA	992	U	C6-N1-C1'	-5.71	113.20	121.20
1	BA	2281	C	C6-N1-C2	5.71	122.58	120.30
1	BA	1116	C	C2-N1-C1'	5.71	125.08	118.80
31	CA	1245	A	N1-C6-N6	5.71	122.03	118.60
31	DA	1290	G	C8-N9-C4	-5.71	104.12	106.40
1	BA	2187	G	N1-C6-O6	5.71	123.32	119.90
31	CA	435	C	C5-C6-N1	5.71	123.85	121.00
1	AA	1210	A	C4-C5-N7	5.70	113.55	110.70
1	BA	204	A	N9-C4-C5	5.70	108.08	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	1184	G	N3-C2-N2	-5.70	115.91	119.90
1	AA	635	C	C6-N1-C2	-5.70	118.02	120.30
1	BA	74	A	N1-C6-N6	-5.70	115.18	118.60
1	BA	723	G	N3-C4-C5	-5.70	125.75	128.60
1	BA	1936	A	N1-C6-N6	5.70	122.02	118.60
1	AA	512	G	C4-N9-C1'	-5.69	119.10	126.50
1	AA	2440	C	N3-C2-O2	-5.69	117.92	121.90
1	BA	196	A	N1-C6-N6	5.69	122.02	118.60
2	AB	19	G	N3-C4-N9	5.69	129.41	126.00
1	BA	298	G	N3-C4-C5	5.69	131.44	128.60
31	DA	1366	C	C2-N3-C4	5.69	122.74	119.90
1	AA	1142(A)	A	C5-N7-C8	-5.69	101.06	103.90
1	AA	1448	G	N1-C6-O6	5.69	123.31	119.90
1	AA	128	C	C2-N1-C1'	5.68	125.05	118.80
1	AA	2586	C	C5-C6-N1	5.68	123.84	121.00
1	BA	1992	G	N1-C6-O6	-5.68	116.49	119.90
1	AA	1408	C	N1-C2-O2	-5.68	115.49	118.90
1	BA	2351	G	C8-N9-C1'	-5.68	119.62	127.00
1	AA	1399	C	C6-N1-C2	5.68	122.57	120.30
2	AB	116	G	C8-N9-C4	5.68	108.67	106.40
1	BA	136	G	C8-N9-C4	5.68	108.67	106.40
1	BA	883	G	N3-C4-N9	5.68	129.41	126.00
20	BY	91	GLU	N-CA-C	-5.67	95.69	111.00
1	BA	2689	U	N3-C4-O4	-5.67	115.43	119.40
31	DA	316	G	N1-C6-O6	5.67	123.30	119.90
31	DA	899	C	N3-C4-C5	5.67	124.17	121.90
31	DA	904	C	C6-N1-C2	5.67	122.57	120.30
1	BA	435	C	N3-C2-O2	-5.67	117.93	121.90
31	CA	1290	G	N3-C4-N9	5.67	129.40	126.00
1	AA	12	U	C5-C6-N1	5.67	125.53	122.70
1	AA	1005	C	C6-N1-C2	5.67	122.57	120.30
1	BA	893	C	N1-C2-O2	5.67	122.30	118.90
1	AA	1261	C	C6-N1-C2	5.66	122.57	120.30
1	AA	2667	C	C6-N1-C2	-5.66	118.04	120.30
31	CA	1012	U	C2-N3-C4	-5.66	123.60	127.00
31	DA	1279	A	C8-N9-C4	-5.66	103.54	105.80
1	AA	1644	C	N1-C2-O2	5.66	122.29	118.90
31	CA	1087	G	N7-C8-N9	5.66	115.93	113.10
1	AA	2712	U	C2-N1-C1'	-5.65	110.92	117.70
1	BA	1948	G	C8-N9-C4	-5.65	104.14	106.40
1	BA	1957	C	C6-N1-C2	5.65	122.56	120.30
1	AA	2207	G	C8-N9-C1'	-5.65	119.66	127.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	948	G	N1-C6-O6	5.65	123.29	119.90
2	AB	54	G	N7-C8-N9	5.65	115.92	113.10
1	BA	219	G	N3-C4-C5	-5.65	125.78	128.60
1	BA	685	A	C8-N9-C4	-5.65	103.54	105.80
31	DA	365	U	C5-C6-N1	-5.65	119.88	122.70
1	BA	265	A	N1-C6-N6	5.65	121.99	118.60
31	DA	1502	A	N1-C2-N3	5.65	132.12	129.30
1	AA	265	A	C6-C5-N7	-5.64	128.35	132.30
1	AA	2378	A	N1-C6-N6	5.64	121.98	118.60
31	DA	530	G	N3-C4-N9	5.64	129.39	126.00
1	AA	1193	G	N9-C4-C5	-5.64	103.14	105.40
1	BA	1045	A	C2-N3-C4	5.64	113.42	110.60
31	DA	1282	C	C6-N1-C1'	-5.64	114.03	120.80
31	DA	337	C	C6-N1-C2	-5.64	118.04	120.30
1	AA	205	G	C8-N9-C4	5.64	108.66	106.40
1	AA	1112	G	C4-N9-C1'	-5.64	119.17	126.50
1	AA	2875	C	C6-N1-C2	5.64	122.55	120.30
31	CA	1027	C	N3-C4-N4	-5.64	114.06	118.00
1	BA	652(T)	C	C2-N3-C4	5.63	122.72	119.90
1	BA	2742	C	C6-N1-C2	5.63	122.55	120.30
1	AA	1955	U	C5-C6-N1	-5.63	119.88	122.70
1	AA	2334	G	C8-N9-C4	5.63	108.65	106.40
1	BA	2464	C	N3-C2-O2	5.63	125.84	121.90
1	AA	1351	C	C5-C6-N1	-5.63	118.19	121.00
31	DA	1509	C	C6-N1-C2	5.63	122.55	120.30
2	AB	115	G	N7-C8-N9	-5.63	110.29	113.10
1	BA	34	C	N1-C2-O2	5.63	122.28	118.90
1	BA	1372	U	N3-C4-O4	5.63	123.34	119.40
1	BA	1510	G	N9-C4-C5	-5.63	103.15	105.40
31	CA	1397	C	C5-C6-N1	5.63	123.81	121.00
1	AA	230	U	C2-N1-C1'	5.62	124.45	117.70
1	AA	460	A	N9-C4-C5	-5.62	103.55	105.80
1	AA	1827	C	N3-C2-O2	-5.62	117.96	121.90
11	BP	39	LYS	N-CA-C	-5.62	95.82	111.00
1	AA	2062	A	N1-C6-N6	5.62	121.97	118.60
1	AA	130	C	N3-C4-C5	5.62	124.15	121.90
1	AA	874	G	C4-N9-C1'	-5.62	119.20	126.50
1	AA	512	G	C5-C6-O6	5.62	131.97	128.60
1	AA	2629	A	N1-C6-N6	5.62	121.97	118.60
1	BA	2347	C	N1-C2-O2	5.62	122.27	118.90
1	BA	2078	C	C6-N1-C2	-5.62	118.05	120.30
31	CA	503	C	C6-N1-C2	-5.62	118.05	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2079	U	N1-C2-O2	-5.61	118.87	122.80
1	BA	975	C	N3-C4-C5	5.61	124.15	121.90
1	BA	2030	A	C8-N9-C4	5.61	108.05	105.80
1	BA	2164	C	C4-C5-C6	5.61	120.21	117.40
11	BP	26	GLY	C-N-CA	5.61	135.73	121.70
1	AA	1168	G	N3-C4-N9	5.61	129.37	126.00
1	BA	1108	U	N3-C2-O2	-5.61	118.27	122.20
1	BA	1840	G	C8-N9-C1'	-5.61	119.71	127.00
1	BA	265	A	C6-C5-N7	-5.60	128.38	132.30
1	BA	375	C	C6-N1-C2	5.60	122.54	120.30
1	BA	446	G	C4-C5-N7	5.60	113.04	110.80
1	BA	2581	G	C8-N9-C4	-5.60	104.16	106.40
1	AA	652(H)	C	C2-N1-C1'	5.60	124.96	118.80
1	AA	679	C	C2-N1-C1'	-5.60	112.64	118.80
1	AA	280	C	C2-N1-C1'	5.60	124.96	118.80
1	AA	769	G	C8-N9-C4	5.60	108.64	106.40
1	BA	1992	G	N3-C4-C5	-5.60	125.80	128.60
31	DA	1279	A	N7-C8-N9	5.60	116.60	113.80
31	DA	1249	C	C5-C6-N1	5.60	123.80	121.00
1	BA	199	A	N7-C8-N9	-5.59	111.00	113.80
2	BB	16	G	C4-C5-N7	5.59	113.04	110.80
2	BB	16	G	C5-C6-O6	-5.59	125.24	128.60
1	BA	12	U	C6-N1-C2	-5.59	117.64	121.00
1	BA	1359	A	C6-N1-C2	5.59	121.95	118.60
1	BA	2030	A	N1-C6-N6	5.59	121.95	118.60
1	BA	298	G	C5-C6-O6	-5.59	125.25	128.60
1	BA	71	A	C8-N9-C4	-5.58	103.57	105.80
1	BA	1936	A	N9-C4-C5	-5.58	103.57	105.80
31	DA	560	U	C2-N1-C1'	5.58	124.40	117.70
1	AA	1286	A	N9-C4-C5	5.58	108.03	105.80
1	BA	511	U	C5-C6-N1	-5.58	119.91	122.70
1	BA	1359	A	C4-C5-C6	-5.58	114.21	117.00
1	BA	1828	G	C4-C5-N7	-5.58	108.57	110.80
31	CA	1323	G	N3-C4-C5	-5.58	125.81	128.60
31	CA	1030	C	C2-N3-C4	5.58	122.69	119.90
31	DA	1493	A	N1-C6-N6	5.58	121.95	118.60
1	AA	2599	G	C8-N9-C4	5.58	108.63	106.40
1	AA	2597	G	N7-C8-N9	-5.57	110.31	113.10
1	BA	1898	U	C5-C4-O4	5.57	129.24	125.90
1	BA	870	A	C8-N9-C4	5.57	108.03	105.80
1	AA	141	A	C2-N3-C4	-5.57	107.82	110.60
1	BA	2271	G	N3-C4-N9	5.56	129.34	126.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	912	C	C2-N1-C1'	5.56	124.92	118.80
2	AB	19	G	C4-N9-C1'	5.56	133.73	126.50
2	AB	113	G	C8-N9-C4	5.56	108.62	106.40
1	BA	676	A	C6-C5-N7	-5.56	128.41	132.30
1	BA	1840	G	C5-C6-N1	-5.56	108.72	111.50
31	DA	117	G	C4-N9-C1'	5.56	133.73	126.50
1	BA	525	U	N3-C4-C5	-5.56	111.27	114.60
1	BA	15	G	N3-C2-N2	-5.55	116.01	119.90
1	BA	2351	G	C4-N9-C1'	5.55	133.72	126.50
1	BA	2362	G	C8-N9-C4	5.55	108.62	106.40
31	CA	1010	G	C8-N9-C4	-5.55	104.18	106.40
1	AA	745	G	N1-C6-O6	5.55	123.23	119.90
1	AA	1359	A	N1-C2-N3	-5.55	126.52	129.30
1	AA	2104	G	C8-N9-C1'	-5.55	119.78	127.00
1	AA	34	C	C6-N1-C2	-5.55	118.08	120.30
1	AA	1142(A)	A	N1-C2-N3	5.55	132.07	129.30
31	CA	1112	C	C6-N1-C2	-5.55	118.08	120.30
31	DA	1088	G	N3-C4-C5	-5.55	125.83	128.60
1	AA	2321	G	C4-N9-C1'	5.54	133.71	126.50
1	AA	2678	C	C6-N1-C2	5.54	122.52	120.30
31	CA	1123	A	N3-C4-N9	-5.54	122.97	127.40
1	BA	1915	U	O4'-C4'-C3'	-5.54	98.46	104.00
1	BA	639	U	N3-C4-C5	-5.54	111.28	114.60
1	AA	1828	G	N1-C6-O6	-5.53	116.58	119.90
1	BA	90	U	N3-C4-O4	5.53	123.27	119.40
1	BA	2491	U	C5-C4-O4	-5.53	122.58	125.90
1	BA	1204	A	C5-N7-C8	-5.53	101.13	103.90
1	BA	2626	C	N3-C4-C5	5.53	124.11	121.90
1	AA	512	G	C6-C5-N7	5.53	133.72	130.40
1	BA	2096	U	N1-C2-O2	5.53	126.67	122.80
31	CA	1231	G	C8-N9-C4	-5.53	104.19	106.40
1	BA	1203	G	C8-N9-C4	-5.53	104.19	106.40
1	AA	2435	A	C8-N9-C4	-5.53	103.59	105.80
31	CA	1058	G	N1-C6-O6	5.53	123.22	119.90
1	BA	1680	U	N1-C2-N3	5.53	118.22	114.90
1	BA	2121	G	C4-C5-N7	-5.52	108.59	110.80
1	BA	2609	U	N1-C2-O2	-5.52	118.93	122.80
1	BA	2827	C	N3-C4-C5	5.52	124.11	121.90
31	CA	995	C	N1-C2-O2	5.52	122.21	118.90
1	BA	932	G	N3-C4-N9	-5.52	122.69	126.00
1	AA	196	A	N9-C4-C5	-5.52	103.59	105.80
1	AA	2056	G	C4-C5-N7	5.52	113.01	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	2552	U	N1-C2-O2	-5.52	118.94	122.80
1	AA	1663	C	C6-N1-C2	5.51	122.51	120.30
1	AA	2791	C	N1-C2-O2	5.51	122.21	118.90
31	CA	360	A	C8-N9-C4	-5.51	103.59	105.80
1	AA	389	G	C5-C6-O6	-5.51	125.29	128.60
1	BA	1938	A	N1-C6-N6	5.51	121.91	118.60
1	AA	1192	G	C8-N9-C4	5.51	108.60	106.40
1	BA	2894	G	C2-N3-C4	-5.51	109.14	111.90
1	AA	450	G	C8-N9-C4	-5.51	104.20	106.40
1	AA	2828	C	N3-C4-C5	5.51	124.10	121.90
1	AA	1493	C	N3-C2-O2	-5.51	118.05	121.90
35	CE	12	LEU	CA-CB-CG	5.51	127.96	115.30
1	AA	54	G	C8-N9-C4	5.50	108.60	106.40
1	BA	1204	A	N1-C6-N6	5.50	121.90	118.60
1	BA	124	G	N9-C4-C5	-5.50	103.20	105.40
1	AA	2514	U	N3-C2-O2	5.50	126.05	122.20
1	AA	207	A	N1-C6-N6	5.50	121.90	118.60
1	AA	2453	A	C8-N9-C4	5.50	108.00	105.80
1	BA	60	G	C5-C6-O6	-5.50	125.30	128.60
1	AA	399	G	N9-C4-C5	-5.49	103.20	105.40
3	AD	274	ARG	C-N-CA	5.49	135.43	121.70
1	BA	271(H)	G	N9-C4-C5	-5.49	103.20	105.40
1	BA	335	C	C6-N1-C2	-5.49	118.10	120.30
1	BA	827	U	N3-C2-O2	5.49	126.05	122.20
1	BA	2523	G	N1-C6-O6	5.49	123.20	119.90
31	CA	1102	A	N3-C4-C5	-5.49	122.95	126.80
1	BA	975	C	N3-C4-N4	-5.49	114.16	118.00
1	AA	2523	G	C8-N9-C4	5.49	108.60	106.40
1	AA	845	G	C6-C5-N7	-5.49	127.11	130.40
1	BA	2022	U	N3-C4-O4	5.49	123.24	119.40
1	AA	1602	U	C4-C5-C6	5.49	122.99	119.70
1	BA	686	G	C4-C5-N7	5.49	113.00	110.80
31	CA	1123	A	N1-C6-N6	-5.49	115.31	118.60
1	AA	1204	A	N9-C4-C5	-5.48	103.61	105.80
1	BA	2490	G	C5-C6-N1	-5.48	108.76	111.50
1	AA	2200	C	C6-N1-C1'	-5.48	114.22	120.80
31	DA	736	C	C2-N1-C1'	5.48	124.83	118.80
1	BA	1033	U	C6-N1-C2	5.48	124.29	121.00
1	AA	2439	A	N9-C4-C5	-5.48	103.61	105.80
1	BA	1680	U	C6-N1-C2	-5.48	117.71	121.00
1	AA	465	G	C8-N9-C4	-5.47	104.21	106.40
1	BA	1622	G	N3-C2-N2	-5.47	116.07	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	2032	G	N7-C8-N9	5.47	115.84	113.10
31	CA	749	C	C5-C6-N1	5.47	123.73	121.00
31	CA	1122	U	C5-C6-N1	5.47	125.43	122.70
1	AA	652(F)	G	N7-C8-N9	5.47	115.83	113.10
1	BA	446	G	N1-C6-O6	5.47	123.18	119.90
31	CA	1492	A	C2'-C3'-O3'	5.47	122.45	113.70
1	AA	792	G	N9-C4-C5	-5.46	103.21	105.40
1	BA	1306	C	N3-C4-C5	5.46	124.09	121.90
1	AA	2187	G	C4-C5-N7	5.46	112.98	110.80
1	AA	2249	U	N3-C4-O4	5.46	123.22	119.40
31	CA	1012	U	N1-C2-O2	-5.46	118.98	122.80
1	AA	2560	C	N3-C4-C5	5.46	124.08	121.90
1	BA	2155	G	C5-C6-O6	-5.46	125.33	128.60
31	DA	1229	A	C6-N1-C2	5.46	121.88	118.60
1	AA	130	C	C6-N1-C1'	-5.46	114.25	120.80
1	AA	1210	A	C6-C5-N7	-5.46	128.48	132.30
1	AA	1784	A	C5-C6-N1	-5.46	114.97	117.70
1	BA	2482	G	N3-C4-C5	-5.46	125.87	128.60
1	BA	2504	U	C5-C6-N1	5.46	125.43	122.70
1	AA	1355	G	N7-C8-N9	5.45	115.83	113.10
31	CA	1030	C	C6-N1-C2	-5.45	118.12	120.30
1	BA	1343	G	C8-N9-C1'	-5.45	119.91	127.00
1	AA	71	A	N7-C8-N9	5.45	116.53	113.80
1	AA	1204	A	C6-C5-N7	-5.45	128.49	132.30
1	AA	1328	G	N9-C4-C5	-5.45	103.22	105.40
1	AA	1673	U	C6-N1-C2	5.45	124.27	121.00
1	AA	2440	C	C5-C4-N4	5.45	124.02	120.20
1	AA	1333	C	N1-C2-O2	5.45	122.17	118.90
1	AA	2448	A	N1-C6-N6	5.45	121.87	118.60
1	BA	1187	G	C8-N9-C4	-5.45	104.22	106.40
1	AA	949	C	C6-N1-C2	5.45	122.48	120.30
1	BA	1022	G	C8-N9-C1'	5.45	134.08	127.00
1	BA	1306	C	C5-C4-N4	-5.44	116.39	120.20
1	BA	1982	C	C5-C6-N1	5.44	123.72	121.00
31	CA	226	G	N7-C8-N9	-5.44	110.38	113.10
31	CA	738	C	C6-N1-C2	-5.44	118.12	120.30
1	BA	944	G	C8-N9-C1'	-5.44	119.93	127.00
1	BA	1721	G	C4-C5-N7	5.44	112.97	110.80
1	AA	1131	G	C8-N9-C1'	5.43	134.06	127.00
1	BA	56	A	N9-C4-C5	-5.43	103.63	105.80
1	BA	1313	U	C6-N1-C1'	-5.43	113.59	121.20
2	BB	92	C	C5-C6-N1	5.43	123.72	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	802	A	N1-C6-N6	5.43	121.86	118.60
1	AA	1644	C	C2-N1-C1'	5.43	124.78	118.80
1	BA	79	G	N3-C4-C5	5.43	131.32	128.60
31	CA	1030	C	C5-C6-N1	5.43	123.72	121.00
1	AA	1541	G	C4-C5-N7	-5.43	108.63	110.80
1	AA	1698	A	C5-C6-N1	-5.43	114.98	117.70
1	BA	450	G	C5-C6-O6	5.43	131.86	128.60
1	BA	1671	U	C6-N1-C2	-5.43	117.74	121.00
1	BA	2590	A	N3-C4-C5	5.43	130.60	126.80
1	AA	1190	G	C8-N9-C1'	5.43	134.06	127.00
1	BA	774	A	C8-N9-C4	-5.43	103.63	105.80
31	CA	181	G	C8-N9-C4	-5.43	104.23	106.40
31	DA	1311	G	N3-C2-N2	-5.43	116.10	119.90
1	AA	1382	G	C4-N9-C1'	-5.42	119.45	126.50
1	AA	2742	C	C6-N1-C2	5.42	122.47	120.30
1	BA	109	G	C4-C5-N7	-5.42	108.63	110.80
31	DA	76	C	N3-C4-C5	5.42	124.07	121.90
1	BA	1026	U	C2-N1-C1'	5.42	124.21	117.70
1	AA	2022	U	N1-C2-N3	-5.42	111.65	114.90
1	BA	1183	G	C2-N3-C4	-5.42	109.19	111.90
1	AA	1432	C	N3-C2-O2	5.42	125.69	121.90
1	AA	2022	U	C5-C4-O4	-5.42	122.65	125.90
1	BA	1914	C	P-O3'-C3'	-5.42	113.20	119.70
1	AA	460	A	C8-N9-C4	5.42	107.97	105.80
1	AA	141	A	N7-C8-N9	5.41	116.51	113.80
1	AA	945	A	C4-C5-N7	5.41	113.41	110.70
1	BA	865	C	N1-C2-N3	-5.41	115.41	119.20
31	DA	460	G	N7-C8-N9	5.41	115.81	113.10
1	BA	1531	C	C2-N1-C1'	5.41	124.75	118.80
1	BA	1671	U	C4-C5-C6	5.41	122.95	119.70
1	AA	676	A	C4-N9-C1'	5.41	136.04	126.30
1	BA	188	G	N9-C4-C5	-5.41	103.24	105.40
1	AA	1690	A	C8-N9-C4	5.41	107.96	105.80
31	CA	521	G	N1-C6-O6	-5.41	116.66	119.90
1	AA	988	A	N9-C4-C5	-5.41	103.64	105.80
1	AA	666	G	C6-C5-N7	-5.41	127.16	130.40
1	AA	1571	A	N1-C6-N6	-5.41	115.36	118.60
1	AA	1941	C	C2-N1-C1'	5.41	124.75	118.80
1	BA	2382	G	N9-C4-C5	-5.41	103.24	105.40
1	AA	117	G	N1-C6-O6	-5.40	116.66	119.90
1	BA	1678	G	N3-C2-N2	-5.40	116.12	119.90
1	BA	1426	G	C5-C6-O6	-5.40	125.36	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
31	CA	243	A	C8-N9-C4	-5.40	103.64	105.80
1	BA	92	A	N7-C8-N9	5.40	116.50	113.80
1	AA	546	C	C2-N1-C1'	5.40	124.74	118.80
1	AA	856	C	C5-C6-N1	5.39	123.70	121.00
1	BA	205	G	N3-C2-N2	5.39	123.68	119.90
1	AA	2658	C	N3-C2-O2	-5.39	118.12	121.90
1	BA	676	A	C8-N9-C4	-5.39	103.64	105.80
31	CA	1282	C	N3-C4-C5	5.39	124.06	121.90
31	CA	175	C	N1-C2-O2	5.39	122.14	118.90
31	CA	204	U	C2-N1-C1'	5.39	124.17	117.70
1	AA	450	G	N9-C4-C5	5.39	107.56	105.40
1	AA	2081	C	C6-N1-C2	5.39	122.46	120.30
1	AA	2512	C	N1-C2-O2	-5.39	115.67	118.90
1	BA	2271	G	N3-C4-C5	-5.39	125.91	128.60
1	BA	2581	G	N9-C4-C5	5.39	107.56	105.40
31	DA	1494	G	C6-N1-C2	-5.39	121.87	125.10
1	AA	1109	C	C6-N1-C2	-5.38	118.15	120.30
31	DA	895	G	N1-C6-O6	5.38	123.13	119.90
1	AA	825	C	C6-N1-C2	-5.38	118.15	120.30
1	BA	857	C	C6-N1-C2	-5.38	118.15	120.30
1	BA	1300	U	P-O3'-C3'	5.38	126.16	119.70
1	BA	530	G	C8-N9-C4	-5.38	104.25	106.40
1	BA	995	C	C6-N1-C1'	5.38	127.26	120.80
31	CA	221	C	C6-N1-C2	-5.38	118.15	120.30
1	BA	788	A	C8-N9-C4	5.38	107.95	105.80
31	CA	1323	G	N3-C4-N9	5.38	129.23	126.00
1	AA	2010	G	N1-C6-O6	5.38	123.13	119.90
31	DA	1030(B)	C	C5-C6-N1	5.38	123.69	121.00
31	CA	1149	C	N3-C4-C5	-5.38	119.75	121.90
1	BA	196	A	N9-C4-C5	-5.37	103.65	105.80
1	BA	528	A	C4-C5-N7	5.37	113.39	110.70
1	BA	1044	G	N3-C4-C5	-5.37	125.91	128.60
1	BA	2275	C	N3-C4-C5	5.37	124.05	121.90
1	BA	2585	U	C2-N1-C1'	5.37	124.15	117.70
1	AA	2058	A	C8-N9-C4	-5.37	103.65	105.80
1	AA	1840	G	C4-N9-C1'	5.37	133.48	126.50
1	BA	1925	C	C2-N1-C1'	-5.37	112.89	118.80
31	DA	1023	G	N3-C4-C5	-5.37	125.92	128.60
1	AA	752	A	C8-N9-C4	-5.37	103.65	105.80
3	AD	239	ARG	N-CA-C	-5.37	96.51	111.00
1	BA	2249	U	C5-C6-N1	5.37	125.38	122.70
31	CA	435	C	C6-N1-C2	-5.37	118.15	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1660	C	C2-N1-C1'	-5.36	112.90	118.80
31	DA	992	U	C5-C6-N1	5.36	125.38	122.70
1	BA	784	A	N3-C4-N9	-5.36	123.11	127.40
31	CA	1123	A	C6-C5-N7	5.36	136.05	132.30
1	AA	1107	G	C8-N9-C1'	-5.36	120.03	127.00
1	BA	2335	A	N1-C6-N6	-5.36	115.39	118.60
31	DA	354	G	C6-C5-N7	-5.36	127.19	130.40
1	AA	2238	G	C8-N9-C4	-5.36	104.26	106.40
1	BA	1271	G	C5-C6-N1	-5.36	108.82	111.50
1	BA	1344	G	N1-C6-O6	5.36	123.11	119.90
1	BA	1671	U	N1-C2-O2	-5.36	119.05	122.80
1	BA	1674	G	C8-N9-C1'	-5.36	120.04	127.00
1	BA	1938	A	C6-C5-N7	-5.36	128.55	132.30
31	DA	519	C	C6-N1-C2	-5.36	118.16	120.30
10	AO	8	LEU	CA-CB-CG	5.35	127.61	115.30
1	BA	1783	A	N9-C4-C5	5.35	107.94	105.80
31	DA	697	U	C5-C6-N1	-5.35	120.02	122.70
1	AA	1721	G	C2-N3-C4	5.35	114.58	111.90
1	BA	456	C	N3-C4-C5	5.35	124.04	121.90
31	DA	117	G	N1-C6-O6	5.35	123.11	119.90
1	BA	2283	C	N3-C2-O2	5.35	125.64	121.90
31	CA	501	C	C6-N1-C2	-5.35	118.16	120.30
1	AA	601	C	C6-N1-C2	5.34	122.44	120.30
1	BA	792	G	N9-C4-C5	-5.34	103.26	105.40
1	AA	1339	G	N9-C4-C5	-5.34	103.26	105.40
1	BA	728	G	C8-N9-C4	5.34	108.54	106.40
1	BA	1779	U	C6-N1-C1'	5.34	128.68	121.20
31	DA	1123	A	C6-C5-N7	5.34	136.04	132.30
1	AA	2599	G	N7-C8-N9	-5.34	110.43	113.10
1	BA	2489	G	C8-N9-C4	5.34	108.54	106.40
1	AA	566	U	N3-C4-O4	-5.34	115.67	119.40
2	BB	7	G	C5-C6-O6	-5.33	125.40	128.60
1	AA	85	G	C8-N9-C4	5.33	108.53	106.40
1	AA	2693	A	C8-N9-C4	5.33	107.93	105.80
1	BA	1050	A	N7-C8-N9	5.33	116.47	113.80
31	DA	435	C	C5-C6-N1	5.33	123.67	121.00
1	BA	784	A	N1-C6-N6	-5.33	115.40	118.60
1	BA	2785	C	C5-C6-N1	5.33	123.66	121.00
1	AA	596	G	N9-C4-C5	5.33	107.53	105.40
1	AA	2013	A	N1-C6-N6	5.33	121.80	118.60
1	AA	1955	U	C6-N1-C2	5.32	124.19	121.00
1	BA	2449	U	C5-C4-O4	-5.32	122.71	125.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
31	DA	1495	U	C6-N1-C2	-5.32	117.81	121.00
1	AA	2484	G	C4-N9-C1'	5.32	133.41	126.50
1	BA	933	A	N1-C6-N6	5.32	121.79	118.60
31	CA	443	C	C2-N1-C1'	5.32	124.65	118.80
1	BA	1332	G	C8-N9-C1'	5.32	133.91	127.00
31	CA	1282	C	C2-N1-C1'	5.32	124.65	118.80
31	DA	1501	C	C6-N1-C2	5.32	122.43	120.30
1	AA	139(A)	G	C6-C5-N7	-5.31	127.21	130.40
1	AA	188	G	N9-C4-C5	-5.31	103.28	105.40
1	AA	831	G	C5-C6-N1	-5.31	108.84	111.50
1	AA	1616	A	N9-C4-C5	-5.31	103.67	105.80
1	BA	1193	G	N1-C6-O6	5.31	123.09	119.90
1	AA	645	C	C6-N1-C2	-5.31	118.18	120.30
1	BA	24	G	C4-C5-N7	-5.31	108.68	110.80
1	BA	1779	U	C2-N3-C4	-5.31	123.81	127.00
31	DA	1151	A	N1-C6-N6	-5.31	115.41	118.60
1	BA	1049	C	C3'-C2'-C1'	5.31	105.75	101.50
1	BA	1279	G	C8-N9-C4	-5.31	104.28	106.40
1	AA	1740	G	C8-N9-C4	5.30	108.52	106.40
1	AA	1022	G	N1-C2-N2	5.30	120.97	116.20
1	AA	2493	U	C5-C6-N1	-5.30	120.05	122.70
1	AA	1285	G	C4-C5-N7	-5.30	108.68	110.80
31	CA	460	G	C6-C5-N7	-5.30	127.22	130.40
1	BA	1820	U	C5-C4-O4	5.30	129.08	125.90
31	DA	1311	G	N9-C4-C5	5.30	107.52	105.40
1	AA	130	C	C2-N3-C4	-5.30	117.25	119.90
1	AA	114	U	C6-N1-C1'	-5.30	113.79	121.20
1	AA	332	A	C8-N9-C4	5.30	107.92	105.80
1	AA	1115	G	C8-N9-C4	5.30	108.52	106.40
1	BA	109	G	N1-C6-O6	-5.30	116.72	119.90
1	BA	184	C	C6-N1-C2	-5.30	118.18	120.30
31	CA	1493	A	N7-C8-N9	-5.29	111.15	113.80
31	DA	316	G	C5-C6-O6	-5.29	125.42	128.60
1	BA	723	G	C6-C5-N7	-5.29	127.22	130.40
31	CA	1029	C	C6-N1-C1'	5.29	127.15	120.80
1	AA	2447	G	C4-N9-C1'	-5.29	119.62	126.50
1	BA	1519	G	C8-N9-C4	-5.29	104.28	106.40
1	BA	2863	C	C6-N1-C2	5.29	122.42	120.30
1	BA	83	G	N3-C4-N9	-5.29	122.83	126.00
1	BA	1721	G	C5-N7-C8	-5.29	101.66	104.30
1	BA	2121	G	N9-C4-C5	5.29	107.52	105.40
31	CA	172	A	C8-N9-C4	-5.29	103.69	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
31	DA	1492	A	N9-C4-C5	5.29	107.92	105.80
1	AA	1575	C	C6-N1-C2	5.29	122.42	120.30
31	CA	171	A	C8-N9-C4	-5.29	103.69	105.80
31	CA	953	G	N3-C4-N9	5.29	129.17	126.00
2	BB	12	C	N3-C2-O2	-5.28	118.20	121.90
1	AA	575	A	N1-C6-N6	5.28	121.77	118.60
31	DA	1409	C	C6-N1-C2	-5.28	118.19	120.30
1	BA	1024	G	N3-C4-N9	5.28	129.17	126.00
1	BA	1721	G	N3-C2-N2	5.28	123.60	119.90
1	BA	2141	G	C6-N1-C2	5.28	128.27	125.10
31	CA	1129	C	N1-C2-O2	5.28	122.07	118.90
1	AA	12	U	C6-N1-C2	-5.28	117.83	121.00
1	BA	652(H)	C	C2-N3-C4	5.28	122.54	119.90
1	BA	945	A	C5-C6-N6	-5.28	119.48	123.70
1	BA	995	C	C2-N1-C1'	-5.28	112.99	118.80
1	BA	451	C	C6-N1-C2	5.28	122.41	120.30
1	BA	639	U	N3-C2-O2	-5.27	118.51	122.20
1	AA	2484	G	C6-C5-N7	-5.27	127.24	130.40
1	BA	1373	A	C8-N9-C4	5.27	107.91	105.80
31	CA	1037	C	C6-N1-C2	-5.27	118.19	120.30
49	DS	9	VAL	N-CA-C	5.27	125.23	111.00
1	AA	221	A	C3'-C2'-C1'	5.27	105.71	101.50
1	AA	2501	C	C6-N1-C2	5.27	122.41	120.30
1	BA	1187	G	N9-C4-C5	5.27	107.51	105.40
31	CA	460	G	C8-N9-C4	-5.27	104.29	106.40
31	DA	1017	G	C5-C6-O6	5.27	131.76	128.60
1	AA	34	C	C2-N1-C1'	5.26	124.59	118.80
1	AA	1339	G	C8-N9-C4	5.26	108.51	106.40
1	AA	465	G	C4-N9-C1'	5.26	133.34	126.50
1	AA	587	C	C6-N1-C2	5.26	122.41	120.30
1	AA	1355	G	C8-N9-C4	-5.26	104.30	106.40
1	AA	1609	A	C8-N9-C4	-5.26	103.70	105.80
1	BA	139(A)	G	C4-C5-N7	5.26	112.90	110.80
1	BA	2489	G	C4-C5-N7	5.26	112.90	110.80
3	BD	60	ARG	NE-CZ-NH1	-5.26	117.67	120.30
1	BA	2791	C	C5-C6-N1	5.26	123.63	121.00
31	CA	1027	C	N3-C4-C5	-5.26	119.80	121.90
1	BA	1963	U	N1-C2-O2	5.25	126.48	122.80
1	AA	1120	G	N3-C4-N9	-5.25	122.85	126.00
1	AA	1259	G	C8-N9-C4	5.25	108.50	106.40
1	BA	652(O)	C	N1-C2-O2	5.25	122.05	118.90
31	DA	1112	C	C6-N1-C2	-5.25	118.20	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	827	U	N1-C2-O2	-5.25	119.12	122.80
1	AA	2508	G	C8-N9-C4	5.25	108.50	106.40
2	BB	84	C	C2-N1-C1'	-5.25	113.03	118.80
31	CA	1124	G	C6-N1-C2	5.25	128.25	125.10
1	AA	950	G	N3-C4-C5	-5.25	125.97	128.60
1	AA	2680	C	N3-C4-C5	5.25	124.00	121.90
1	BA	335	C	C5-C6-N1	5.25	123.62	121.00
1	BA	750	A	C8-N9-C4	-5.25	103.70	105.80
31	CA	1402	C	C6-N1-C2	-5.25	118.20	120.30
1	BA	680	G	N3-C2-N2	-5.24	116.23	119.90
1	BA	777	A	N1-C2-N3	5.24	131.92	129.30
1	BA	1698	A	C5-N7-C8	-5.24	101.28	103.90
1	BA	2058	A	C8-N9-C4	-5.24	103.70	105.80
1	BA	2449	U	N3-C4-O4	5.24	123.07	119.40
1	AA	693	C	N3-C4-C5	5.24	124.00	121.90
1	BA	1340	U	C5-C6-N1	-5.24	120.08	122.70
1	BA	1779	U	C6-N1-C2	5.24	124.14	121.00
1	AA	351	G	N1-C6-O6	5.24	123.05	119.90
1	BA	2306	C	C5-C6-N1	5.24	123.62	121.00
31	DA	376	G	C8-N9-C4	5.24	108.50	106.40
1	AA	2866	U	C6-N1-C1'	-5.24	113.87	121.20
1	AA	1963	U	N1-C2-O2	5.24	126.47	122.80
1	AA	2439	A	N7-C8-N9	5.24	116.42	113.80
1	AA	453	C	C5-C6-N1	-5.23	118.38	121.00
31	DA	76	C	C6-N1-C1'	-5.23	114.52	120.80
1	AA	271(M)	G	C8-N9-C1'	-5.23	120.20	127.00
1	AA	679	C	N1-C2-O2	-5.23	115.76	118.90
1	AA	2607	G	C4-N9-C1'	5.23	133.30	126.50
1	BA	1116	C	C6-N1-C1'	-5.23	114.52	120.80
1	BA	1955	U	C2-N1-C1'	-5.23	111.42	117.70
31	DA	1491	G	C4-C5-N7	-5.23	108.71	110.80
1	AA	1142(A)	A	N7-C8-N9	5.23	116.42	113.80
1	BA	2464	C	C4-C5-C6	-5.23	114.78	117.40
1	AA	2061	G	N9-C4-C5	-5.23	103.31	105.40
1	BA	970	C	C6-N1-C2	5.23	122.39	120.30
1	BA	2036	C	C5-C6-N1	5.23	123.61	121.00
1	BA	2790	A	C2-N3-C4	5.23	113.21	110.60
2	BB	109	C	C6-N1-C2	5.23	122.39	120.30
31	CA	1228	C	C6-N1-C2	-5.23	118.21	120.30
31	CA	105	G	C4-N9-C1'	-5.22	119.71	126.50
1	AA	1223	G	C8-N9-C4	-5.22	104.31	106.40
1	BA	61	G	N3-C2-N2	-5.22	116.25	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	2375	G	N7-C8-N9	-5.22	110.49	113.10
31	DA	442	C	C5-C6-N1	5.22	123.61	121.00
31	DA	442	C	C6-N1-C2	-5.22	118.21	120.30
31	DA	1054	C	N1-C2-O2	-5.22	115.77	118.90
1	BA	2744	G	N9-C4-C5	-5.22	103.31	105.40
2	BB	13	A	N1-C6-N6	-5.22	115.47	118.60
2	BB	116	G	C8-N9-C1'	-5.22	120.22	127.00
1	AA	2191	G	C4-C5-N7	5.22	112.89	110.80
31	DA	257	G	C8-N9-C4	5.22	108.49	106.40
1	AA	856	C	C3'-C2'-C1'	-5.22	97.33	101.50
7	AH	71	LEU	CA-CB-CG	5.22	127.30	115.30
1	BA	2137	C	N3-C4-C5	-5.22	119.81	121.90
1	BA	2429	G	C4-C5-N7	-5.22	108.71	110.80
1	AA	139(A)	G	N3-C4-N9	5.21	129.13	126.00
1	AA	2085	C	C6-N1-C2	5.21	122.39	120.30
11	AP	26	GLY	C-N-CA	5.21	134.73	121.70
1	BA	271(M)	G	C6-N1-C2	-5.21	121.97	125.10
1	BA	475	U	C2-N1-C1'	5.21	123.96	117.70
1	AA	2512	C	N3-C2-O2	5.21	125.55	121.90
1	AA	2601	C	N1-C2-O2	-5.21	115.77	118.90
31	DA	927	G	C5-C6-O6	5.21	131.73	128.60
1	BA	1689	A	N1-C6-N6	-5.21	115.47	118.60
1	AA	2606	C	N3-C4-C5	5.21	123.98	121.90
1	BA	254	G	N3-C4-C5	-5.21	126.00	128.60
1	BA	2706	G	C8-N9-C4	5.21	108.48	106.40
31	CA	1091	U	C6-N1-C2	-5.21	117.88	121.00
31	DA	200	G	C8-N9-C4	5.21	108.48	106.40
31	DA	1149	C	N3-C4-C5	-5.21	119.82	121.90
1	AA	1784	A	C8-N9-C4	5.21	107.88	105.80
1	BA	652(T)	C	C5-C6-N1	5.21	123.60	121.00
1	AA	2759	G	N9-C4-C5	-5.21	103.32	105.40
1	BA	58	G	N1-C2-N3	5.20	127.02	123.90
31	CA	1026	G	N7-C8-N9	5.20	115.70	113.10
1	BA	2062	A	N1-C6-N6	5.20	121.72	118.60
40	DJ	90	LEU	C-N-CD	-5.20	109.16	120.60
1	AA	2818	G	C8-N9-C4	5.20	108.48	106.40
31	DA	616	G	C4-N9-C1'	5.20	133.26	126.50
1	BA	114	U	N3-C2-O2	-5.20	118.56	122.20
1	AA	928	G	C4-C5-N7	5.20	112.88	110.80
1	AA	1371	G	N1-C6-O6	5.20	123.02	119.90
1	AA	2382	G	N3-C2-N2	5.20	123.54	119.90
1	BA	1037	G	C8-N9-C4	5.19	108.48	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	81	G	N1-C6-O6	5.19	123.02	119.90
1	AA	1190	G	N1-C6-O6	-5.19	116.78	119.90
1	BA	649	G	N3-C2-N2	-5.19	116.27	119.90
3	BD	244	ARG	N-CA-C	-5.19	96.98	111.00
1	AA	792	G	C8-N9-C1'	-5.19	120.25	127.00
1	AA	2607	G	C8-N9-C1'	-5.19	120.25	127.00
1	BA	2382	G	C2-N3-C4	-5.19	109.30	111.90
31	CA	1351	U	C5-C6-N1	5.19	125.30	122.70
1	BA	474	G	C8-N9-C4	-5.19	104.33	106.40
1	BA	1231	G	N3-C4-N9	-5.19	122.89	126.00
1	AA	511	U	C2-N1-C1'	5.19	123.92	117.70
1	AA	2757	A	C8-N9-C4	-5.18	103.73	105.80
31	DA	1417	G	C5-C6-N1	-5.18	108.91	111.50
1	AA	2253	G	N1-C6-O6	5.18	123.01	119.90
1	BA	690	G	N7-C8-N9	-5.18	110.51	113.10
1	BA	1313	U	N1-C2-O2	5.18	126.43	122.80
31	CA	749	C	C6-N1-C2	-5.18	118.23	120.30
1	BA	2102	U	N1-C2-O2	5.18	126.42	122.80
1	AA	1039	G	N9-C4-C5	-5.18	103.33	105.40
1	AA	1109	C	N3-C4-N4	5.18	121.62	118.00
1	BA	2176	A	C8-N9-C4	-5.18	103.73	105.80
1	AA	34	C	N1-C2-O2	5.17	122.00	118.90
1	AA	2055	C	C6-N1-C2	5.17	122.37	120.30
1	BA	2032	G	C8-N9-C1'	5.17	133.73	127.00
1	BA	2894	G	C4-N9-C1'	5.17	133.23	126.50
31	CA	1091	U	N1-C2-O2	5.17	126.42	122.80
1	AA	1615	C	C6-N1-C2	5.17	122.37	120.30
1	AA	1653	G	C8-N9-C1'	-5.17	120.28	127.00
1	BA	956	G	N1-C6-O6	5.17	123.00	119.90
1	BA	2028	U	N1-C2-O2	-5.17	119.18	122.80
1	BA	1274	A	C8-N9-C4	5.17	107.87	105.80
1	BA	2106	G	C6-N1-C2	5.17	128.20	125.10
40	CJ	90	LEU	C-N-CD	-5.17	109.23	120.60
1	BA	1258	C	C6-N1-C2	5.17	122.37	120.30
1	AA	1982	C	C6-N1-C1'	-5.17	114.60	120.80
1	AA	2241	A	N7-C8-N9	-5.17	111.22	113.80
1	BA	1721	G	N7-C8-N9	5.17	115.68	113.10
1	BA	856	C	C3'-C2'-C1'	-5.17	97.37	101.50
1	AA	139(A)	G	C5-N7-C8	-5.16	101.72	104.30
1	AA	1863	G	C8-N9-C4	5.16	108.47	106.40
1	AA	2099	U	C6-N1-C2	-5.16	117.90	121.00
1	AA	2105	C	C5-C6-N1	5.16	123.58	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	116	C	C5-C6-N1	-5.16	118.42	121.00
1	BA	1143	A	N1-C6-N6	5.16	121.70	118.60
1	BA	2311	A	C8-N9-C4	-5.16	103.73	105.80
1	BA	2501	C	C2-N1-C1'	-5.16	113.12	118.80
31	DA	972	C	C5-C6-N1	5.16	123.58	121.00
1	AA	566	U	C2-N3-C4	-5.16	123.90	127.00
1	AA	1698	A	N7-C8-N9	5.16	116.38	113.80
1	BA	565	C	N1-C2-O2	5.16	122.00	118.90
31	CA	1028	C	C6-N1-C2	-5.16	118.23	120.30
1	AA	1938	A	N1-C6-N6	5.16	121.70	118.60
2	AB	116	G	N9-C4-C5	-5.16	103.34	105.40
1	BA	761	A	N7-C8-N9	5.16	116.38	113.80
1	AA	139(A)	G	N7-C8-N9	5.16	115.68	113.10
1	AA	652(T)	C	N1-C2-O2	5.16	121.99	118.90
31	DA	1378	C	C6-N1-C2	-5.16	118.24	120.30
31	DA	940	C	C5-C6-N1	5.15	123.58	121.00
1	AA	1259	G	N7-C8-N9	-5.15	110.52	113.10
1	BA	572	A	C8-N9-C4	-5.15	103.74	105.80
1	BA	649	G	N1-C6-O6	5.15	122.99	119.90
1	AA	1678	G	C6-C5-N7	-5.15	127.31	130.40
1	BA	885	C	N3-C2-O2	-5.15	118.30	121.90
31	DA	97	G	N1-C6-O6	5.15	122.99	119.90
31	DA	460	G	C5-C6-O6	-5.15	125.51	128.60
31	DA	1002	G	C8-N9-C4	-5.15	104.34	106.40
1	BA	528	A	N3-C4-N9	-5.15	123.28	127.40
1	BA	937	U	C2-N1-C1'	-5.15	111.52	117.70
1	AA	2473	U	C6-N1-C1'	-5.15	113.99	121.20
1	BA	2503	A	C5-C6-N6	-5.15	119.58	123.70
31	DA	1138	G	N3-C4-C5	-5.15	126.03	128.60
1	AA	933	A	C5-N7-C8	-5.14	101.33	103.90
27	A5	54	GLY	N-CA-C	-5.14	100.24	113.10
1	BA	1678	G	C8-N9-C4	-5.14	104.34	106.40
1	BA	2249	U	C5-C4-O4	5.14	128.99	125.90
1	BA	2544	G	N3-C4-N9	5.14	129.09	126.00
1	BA	2894	G	C5-C6-N1	-5.14	108.93	111.50
31	CA	1034	G	C5-C6-N1	-5.14	108.93	111.50
31	DA	917	G	C8-N9-C4	-5.14	104.34	106.40
1	AA	1368	G	N3-C2-N2	-5.14	116.30	119.90
1	BA	33	U	C5-C4-O4	5.14	128.99	125.90
1	AA	1382	G	N3-C4-N9	-5.14	122.92	126.00
1	BA	2503	A	C2-N3-C4	5.14	113.17	110.60
31	CA	1502	A	N7-C8-N9	5.14	116.37	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	1314	C	N1-C2-O2	5.14	121.98	118.90
1	AA	1022	G	C8-N9-C1'	5.13	133.67	127.00
1	AA	1115	G	N3-C4-C5	5.13	131.17	128.60
31	DA	1255	G	C5-C6-O6	-5.13	125.52	128.60
31	DA	1184	G	C6-C5-N7	5.13	133.48	130.40
1	BA	933	A	C6-C5-N7	-5.13	128.71	132.30
1	BA	1930	G	C4-N9-C1'	-5.13	119.83	126.50
31	DA	1184	G	N3-C2-N2	-5.13	116.31	119.90
1	AA	2367	G	N9-C4-C5	-5.13	103.35	105.40
1	AA	2768	C	C6-N1-C2	-5.13	118.25	120.30
1	AA	1206	G	N1-C6-O6	5.13	122.98	119.90
1	AA	2318	G	C8-N9-C1'	-5.13	120.33	127.00
31	CA	1017	G	N3-C4-N9	-5.13	122.92	126.00
31	CA	1524	C	N3-C2-O2	5.13	125.49	121.90
31	DA	784	C	C6-N1-C2	5.13	122.35	120.30
1	AA	1882	C	C6-N1-C1'	-5.13	114.65	120.80
1	AA	2320	A	N1-C6-N6	-5.13	115.52	118.60
1	AA	2610	C	P-O3'-C3'	5.12	125.85	119.70
1	BA	1800	C	C2-N1-C1'	-5.12	113.16	118.80
31	CA	76	C	N1-C2-O2	5.12	121.97	118.90
31	CA	1032	G	N9-C4-C5	5.12	107.45	105.40
31	DA	1493	A	C4-N9-C1'	5.12	135.52	126.30
1	AA	190	A	N1-C6-N6	5.12	121.67	118.60
1	AA	389	G	N1-C6-O6	5.12	122.97	119.90
1	AA	1840	G	C8-N9-C1'	-5.12	120.34	127.00
1	AA	1992	G	N3-C4-C5	-5.12	126.04	128.60
31	DA	1184	G	N1-C2-N2	5.12	120.81	116.20
1	AA	62	C	C2-N1-C1'	-5.12	113.17	118.80
1	AA	885	C	N3-C2-O2	-5.12	118.31	121.90
1	BA	512	G	C8-N9-C1'	5.12	133.66	127.00
1	BA	1902	C	N1-C2-O2	5.12	121.97	118.90
1	BA	271(O)	C	N1-C2-O2	5.12	121.97	118.90
1	BA	1488	G	C4-N9-C1'	5.12	133.16	126.50
1	BA	949	C	N1-C2-O2	-5.12	115.83	118.90
1	BA	1656	C	C6-N1-C2	-5.12	118.25	120.30
1	AA	130	C	N3-C2-O2	-5.12	118.32	121.90
1	BA	2162	G	N3-C4-C5	-5.12	126.04	128.60
1	BA	2391	G	C4-N9-C1'	-5.12	119.85	126.50
1	BA	2689	U	N3-C2-O2	-5.12	118.62	122.20
34	CD	12	CYS	N-CA-C	-5.12	97.19	111.00
1	BA	758	C	C6-N1-C2	-5.11	118.25	120.30
1	BA	1773	A	C8-N9-C4	-5.11	103.75	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	855	G	C8-N9-C4	-5.11	104.36	106.40
2	BB	106	G	N1-C6-O6	5.11	122.97	119.90
2	BB	116	G	C4-N9-C1'	5.11	133.15	126.50
26	B4	42	PHE	C-N-CA	5.11	134.48	121.70
1	AA	71	A	C4-C5-N7	5.11	113.25	110.70
1	AA	190	A	N9-C4-C5	-5.11	103.76	105.80
1	BA	102	G	C4-C5-N7	-5.11	108.76	110.80
2	BB	92	C	C6-N1-C2	-5.11	118.26	120.30
31	CA	1123	A	C4-N9-C1'	-5.11	117.11	126.30
31	DA	299	G	C4-C5-N7	-5.11	108.76	110.80
1	AA	568	U	N3-C2-O2	5.11	125.77	122.20
1	AA	697	C	C2-N1-C1'	5.11	124.42	118.80
1	BA	1265	A	N7-C8-N9	5.11	116.35	113.80
31	DA	721	G	N1-C6-O6	5.11	122.96	119.90
31	DA	1229	A	C5-C6-N1	-5.11	115.15	117.70
1	BA	2685	G	N1-C2-N2	5.10	120.79	116.20
31	DA	1054	C	C2-N1-C1'	-5.10	113.19	118.80
1	AA	732	C	C6-N1-C2	5.10	122.34	120.30
31	CA	433	C	C6-N1-C2	-5.10	118.26	120.30
31	DA	1493	A	O4'-C1'-N9	5.10	112.28	108.20
1	BA	847	U	C5-C6-N1	-5.10	120.15	122.70
31	DA	1129	C	N1-C2-O2	5.10	121.96	118.90
31	DA	1255	G	N3-C4-N9	5.10	129.06	126.00
1	AA	2439	A	C2-N3-C4	-5.10	108.05	110.60
1	BA	143	G	C8-N9-C1'	5.10	133.63	127.00
1	BA	1779	U	N1-C2-O2	-5.10	119.23	122.80
1	BA	1992	G	P-O3'-C3'	5.10	125.82	119.70
1	BA	2482	G	C8-N9-C1'	-5.10	120.37	127.00
1	AA	210	C	C5-C6-N1	-5.10	118.45	121.00
1	AA	2419	U	N3-C4-O4	5.10	122.97	119.40
1	AA	2526	G	N3-C4-N9	-5.10	122.94	126.00
1	BA	2193	G	N1-C6-O6	5.10	122.96	119.90
1	BA	2430	A	C5-N7-C8	-5.10	101.35	103.90
2	BB	53	A	C6-N1-C2	-5.10	115.54	118.60
31	DA	1272	G	C6-C5-N7	-5.10	127.34	130.40
1	AA	60	G	C8-N9-C4	-5.10	104.36	106.40
1	BA	179	G	C6-C5-N7	-5.10	127.34	130.40
1	BA	574	C	C2-N1-C1'	-5.10	113.19	118.80
1	BA	2104	G	C4-N9-C1'	5.10	133.12	126.50
1	BA	2781	A	N1-C6-N6	-5.10	115.54	118.60
1	AA	211	A	C8-N9-C4	5.09	107.84	105.80
1	AA	2318	G	N3-C4-C5	-5.09	126.05	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	BB	37	C	N1-C2-O2	-5.09	115.84	118.90
31	CA	1292	U	C5-C6-N1	-5.09	120.15	122.70
31	DA	91	C	C6-N1-C2	-5.09	118.26	120.30
1	AA	2550	G	C8-N9-C1'	-5.09	120.38	127.00
1	AA	1399	C	N3-C4-C5	5.09	123.94	121.90
1	AA	1662	C	C6-N1-C2	-5.09	118.26	120.30
1	BA	657	U	C5-C4-O4	5.09	128.96	125.90
1	BA	2032	G	C4-C5-N7	5.09	112.84	110.80
1	BA	1578	U	N3-C2-O2	-5.09	118.64	122.20
1	BA	683	C	N1-C2-O2	5.09	121.95	118.90
1	BA	2802	G	C3'-C2'-C1'	-5.09	97.43	101.50
31	CA	413	G	C4-C5-N7	-5.09	108.77	110.80
1	AA	570	G	N3-C4-C5	-5.08	126.06	128.60
1	AA	577	G	C8-N9-C4	5.08	108.43	106.40
1	AA	631	A	C8-N9-C4	5.08	107.83	105.80
1	BA	546	C	C6-N1-C1'	-5.08	114.70	120.80
1	BA	978	G	N7-C8-N9	-5.08	110.56	113.10
31	DA	1249	C	C6-N1-C2	-5.08	118.27	120.30
1	AA	1350	C	N1-C2-O2	-5.08	115.85	118.90
1	BA	2299	G	C5-C6-N1	-5.08	108.96	111.50
31	CA	105	G	C8-N9-C1'	5.08	133.61	127.00
1	AA	560	C	C6-N1-C2	5.08	122.33	120.30
1	AA	2253	G	C6-C5-N7	-5.08	127.35	130.40
1	BA	2131	G	N7-C8-N9	5.08	115.64	113.10
31	CA	284	G	N1-C6-O6	5.08	122.95	119.90
31	CA	1030(B)	C	C6-N1-C1'	-5.08	114.71	120.80
1	BA	271(J)	C	N3-C2-O2	-5.08	118.35	121.90
1	BA	2452	C	C6-N1-C2	5.08	122.33	120.30
1	BA	330	A	C4-C5-N7	5.08	113.24	110.70
16	BU	74	LEU	CA-CB-CG	5.08	126.97	115.30
31	CA	117	G	N1-C6-O6	5.08	122.95	119.90
31	DA	993	G	C4-N9-C1'	5.08	133.10	126.50
1	AA	453	C	C6-N1-C2	5.07	122.33	120.30
1	AA	1930	G	C4-N9-C1'	-5.07	119.90	126.50
1	BA	70	G	C4-C5-N7	5.07	112.83	110.80
1	BA	271(Y)	U	C5-C6-N1	-5.07	120.16	122.70
1	BA	531	C	C6-N1-C2	5.07	122.33	120.30
1	BA	892	G	C5-C6-O6	-5.07	125.56	128.60
1	AA	1743	C	C6-N1-C2	5.07	122.33	120.30
1	AA	2822	G	N9-C4-C5	-5.07	103.37	105.40
1	BA	2687	U	C5-C4-O4	5.07	128.94	125.90
2	AB	30	C	N3-C4-C5	-5.07	119.87	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	139(A)	G	C5-C6-N1	5.07	114.03	111.50
1	BA	1047	G	N3-C2-N2	5.07	123.45	119.90
1	AA	1541	G	N9-C4-C5	5.06	107.43	105.40
2	AB	20	C	C2-N1-C1'	5.06	124.37	118.80
1	BA	1488	G	N7-C8-N9	5.06	115.63	113.10
1	BA	1899	G	N3-C2-N2	-5.06	116.36	119.90
1	BA	2708	G	N1-C6-O6	5.06	122.94	119.90
1	AA	2677	G	C8-N9-C4	5.06	108.42	106.40
31	CA	790	A	C2-N3-C4	-5.06	108.07	110.60
1	AA	2559	C	C6-N1-C2	5.06	122.32	120.30
1	BA	337	C	C6-N1-C2	5.06	122.32	120.30
1	BA	1636	C	C6-N1-C2	5.06	122.32	120.30
1	BA	1743	C	N3-C4-C5	-5.06	119.88	121.90
1	BA	2849	U	N3-C2-O2	5.06	125.74	122.20
1	BA	2874	C	C2-N1-C1'	5.06	124.36	118.80
1	BA	2318	G	C8-N9-C1'	-5.06	120.43	127.00
31	CA	19	C	C6-N1-C2	-5.06	118.28	120.30
1	AA	1328	G	C8-N9-C1'	-5.05	120.43	127.00
1	AA	1493	C	N1-C2-O2	5.05	121.93	118.90
1	BA	2286	A	N7-C8-N9	5.05	116.33	113.80
31	DA	1417	G	C4-N9-C1'	5.05	133.07	126.50
1	BA	528	A	C5-C6-N1	-5.05	115.17	117.70
1	BA	1477	A	C8-N9-C4	5.05	107.82	105.80
1	BA	1687	G	C8-N9-C4	-5.05	104.38	106.40
1	BA	2086	U	C5-C4-O4	5.05	128.93	125.90
1	AA	146	G	C8-N9-C4	5.05	108.42	106.40
1	AA	1359	A	C2-N3-C4	5.05	113.12	110.60
1	AA	1799	G	P-O3'-C3'	5.05	125.76	119.70
1	BA	741	G	N1-C6-O6	5.05	122.93	119.90
31	CA	324	G	N1-C6-O6	5.05	122.93	119.90
31	CA	596	C	N1-C2-O2	5.05	121.93	118.90
1	AA	113	G	N3-C4-N9	-5.05	122.97	126.00
1	AA	614	U	N3-C2-O2	-5.05	118.67	122.20
1	AA	1131	G	N7-C8-N9	-5.05	110.58	113.10
1	AA	2439	A	C5-C6-N6	-5.05	119.66	123.70
1	AA	2447	G	N3-C2-N2	-5.05	116.37	119.90
1	BA	1983	C	C2-N1-C1'	-5.05	113.25	118.80
1	BA	2894	G	N9-C4-C5	-5.05	103.38	105.40
31	DA	117	G	C4-C5-N7	5.05	112.82	110.80
31	DA	345	C	C6-N1-C1'	5.05	126.86	120.80
31	DA	839	U	C6-N1-C2	-5.04	117.97	121.00
1	BA	847	U	C6-N1-C1'	5.04	128.26	121.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1913	A	C2-N3-C4	5.04	113.12	110.60
1	AA	1176	G	C4-N9-C1'	-5.04	119.95	126.50
1	AA	1820	U	C2-N1-C1'	-5.04	111.66	117.70
2	AB	117	G	N1-C6-O6	5.04	122.92	119.90
31	DA	345	C	C5-C6-N1	-5.04	118.48	121.00
31	CA	93	G	N3-C4-N9	5.04	129.02	126.00
31	DA	3	G	N3-C4-N9	5.04	129.02	126.00
1	AA	2499	C	C6-N1-C2	5.04	122.31	120.30
1	AA	2501	C	C2-N1-C1'	-5.04	113.26	118.80
1	BA	652(G)	G	N3-C4-C5	-5.04	126.08	128.60
1	BA	1231	G	N1-C6-O6	5.04	122.92	119.90
1	AA	2585	U	N1-C2-O2	5.03	126.32	122.80
1	BA	2625	G	N1-C6-O6	-5.03	116.88	119.90
43	DM	65	LYS	C-N-CA	5.03	134.28	121.70
31	DA	972	C	C6-N1-C2	-5.03	118.29	120.30
31	DA	1255	G	C2-N3-C4	5.03	114.42	111.90
1	AA	1602	U	C6-N1-C2	-5.03	117.98	121.00
1	BA	204	A	C2-N3-C4	5.03	113.11	110.60
1	BA	2307	G	N7-C8-N9	5.03	115.61	113.10
1	AA	1001	A	C8-N9-C4	5.03	107.81	105.80
31	DA	560	U	C3'-C2'-C1'	5.03	105.52	101.50
31	DA	1126	U	C6-N1-C1'	-5.03	114.16	121.20
1	BA	2088	G	N3-C4-N9	-5.03	122.98	126.00
1	BA	2136	C	C5-C6-N1	5.03	123.51	121.00
2	BB	16	G	N1-C6-O6	5.03	122.92	119.90
31	CA	1029	C	N3-C4-N4	-5.03	114.48	118.00
31	DA	1017	G	C6-N1-C2	5.03	128.12	125.10
1	AA	749	C	N1-C2-O2	5.02	121.91	118.90
1	AA	2155	G	N1-C6-O6	5.02	122.91	119.90
31	DA	1033	G	N7-C8-N9	5.02	115.61	113.10
1	BA	220	G	N3-C4-C5	5.02	131.11	128.60
31	CA	353	A	C5-N7-C8	-5.02	101.39	103.90
1	BA	1022	G	N3-C4-C5	5.02	131.11	128.60
31	CA	353	A	C4-C5-N7	5.02	113.21	110.70
31	DA	397	A	C8-N9-C4	-5.02	103.79	105.80
1	AA	1698	A	N3-C4-C5	5.02	130.31	126.80
1	BA	528	A	C6-C5-N7	-5.02	128.79	132.30
1	BA	566	U	N3-C2-O2	5.02	125.71	122.20
1	BA	1416	G	N1-C6-O6	-5.02	116.89	119.90
1	BA	1187	G	N1-C2-N3	5.02	126.91	123.90
1	BA	1231	G	C2-N3-C4	-5.02	109.39	111.90
31	DA	1378	C	C2-N1-C1'	5.02	124.32	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	71	A	C6-C5-N7	-5.01	128.79	132.30
14	BS	82	ILE	C-N-CA	5.01	134.24	121.70
1	BA	525	U	C5-C6-N1	5.01	125.21	122.70
31	CA	1056	U	C2-N3-C4	5.01	130.01	127.00
1	AA	1882	C	C5-C6-N1	5.01	123.51	121.00
1	AA	1991	U	C5-C6-N1	-5.01	120.19	122.70
1	AA	2386	C	C5-C6-N1	-5.01	118.49	121.00
1	BA	2226	C	C6-N1-C2	5.01	122.31	120.30
1	AA	2056	G	C5-N7-C8	-5.01	101.80	104.30
1	AA	2439	A	C4-N9-C1'	5.01	135.32	126.30
1	AA	36	G	N1-C6-O6	5.01	122.91	119.90
1	AA	2061	G	N7-C8-N9	-5.01	110.60	113.10
31	CA	998	G	C4-C5-N7	-5.01	108.80	110.80
1	BA	748	G	C4-N9-C1'	-5.01	119.99	126.50
1	BA	1915	U	C5'-C4'-C3'	5.01	124.01	116.00
31	DA	1397	C	N3-C2-O2	-5.01	118.40	121.90
1	BA	2283	C	N1-C2-O2	-5.00	115.90	118.90
1	BA	2589	A	N3-C4-C5	5.00	130.30	126.80
1	AA	189	G	C8-N9-C4	5.00	108.40	106.40
1	AA	317	G	N1-C6-O6	5.00	122.90	119.90
1	AA	2685	G	C5-C6-O6	5.00	131.60	128.60
1	BA	2475	C	C6-N1-C2	5.00	122.30	120.30
13	BR	22	ARG	NE-CZ-NH1	-5.00	117.80	120.30
35	DE	12	LEU	CA-CB-CG	5.00	126.81	115.30
1	BA	1654	A	N1-C6-N6	-5.00	115.60	118.60
1	BA	2035	G	N3-C4-N9	-5.00	123.00	126.00
1	BA	2181	G	C4-N9-C1'	-5.00	120.00	126.50
1	BA	2287	A	N3-C4-C5	5.00	130.30	126.80
31	DA	97	G	C5-C6-O6	-5.00	125.60	128.60
31	DA	1123	A	C5-C6-N6	5.00	127.70	123.70
31	DA	1311	G	N3-C4-N9	-5.00	123.00	126.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
26	A4	42	PHE	Peptide
34	CD	11	LEU	Peptide
42	CL	26	ALA	Peptide
34	DD	11	LEU	Peptide
42	DL	26	ALA	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	AA	60900	0	30712	1406	0
1	BA	60900	0	30712	1060	0
2	AB	2574	0	1306	87	0
2	BB	2574	0	1306	35	0
3	AD	2136	0	2218	77	0
3	BD	2136	0	2218	81	0
4	AE	1555	0	1607	81	0
4	BE	1555	0	1607	54	0
5	AF	1576	0	1616	57	0
5	BF	1576	0	1616	59	0
6	AG	1368	0	1324	70	0
6	BG	1368	0	1324	61	0
7	AH	1317	0	1376	66	0
7	BH	1317	0	1376	33	0
8	AI	1046	0	1067	55	2
8	BI	1046	0	1067	47	0
9	AN	1112	0	1180	60	0
9	BN	1112	0	1180	34	0
10	AO	923	0	981	28	0
10	BO	923	0	981	23	0
11	AP	1119	0	1186	38	0
11	BP	1119	0	1186	40	0
12	AQ	1122	0	1179	43	0
12	BQ	1122	0	1179	47	0
13	AR	968	0	1033	43	0
13	BR	968	0	1033	33	0
14	AS	865	0	905	62	0
14	BS	865	0	905	54	0
15	AT	1063	0	1103	48	0
15	BT	1063	0	1103	37	0
16	AU	959	0	1019	31	0
16	BU	959	0	1019	20	0
17	AV	771	0	830	22	0
17	BV	771	0	830	15	0
18	AW	881	0	935	25	0
18	BW	881	0	935	24	0
19	AX	742	0	799	22	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
19	BX	742	0	799	23	0
20	AY	785	0	832	36	0
20	BY	785	0	832	30	0
21	AZ	1522	0	1511	56	0
21	BZ	1522	0	1511	47	0
22	A0	594	0	604	27	0
22	B0	594	0	604	21	0
23	A1	745	0	804	26	0
23	B1	745	0	804	26	0
24	A2	588	0	643	31	0
24	B2	588	0	643	18	0
25	A3	458	0	503	22	0
25	B3	458	0	503	6	0
26	A4	349	0	340	22	0
26	B4	349	0	340	16	0
27	A5	455	0	476	25	0
27	B5	455	0	476	18	0
28	A6	449	0	464	19	0
28	B6	449	0	466	17	0
29	A7	418	0	467	18	0
29	B7	418	0	467	14	0
30	A8	509	0	565	23	0
30	B8	509	0	565	20	0
31	CA	32208	0	16256	921	2
31	DA	32208	0	16254	923	0
32	CB	1777	0	1747	100	0
32	DB	1777	0	1747	95	0
33	CC	1450	0	1314	45	0
33	DC	1450	0	1314	58	0
34	CD	1520	0	1407	73	0
34	DD	1520	0	1406	85	0
35	CE	1105	0	1130	48	0
35	DE	1105	0	1130	54	0
36	CF	781	0	741	25	0
36	DF	781	0	741	29	0
37	CG	1167	0	1108	39	0
37	DG	1167	0	1108	46	0
38	CH	1045	0	1033	45	0
38	DH	1045	0	1033	53	0
39	CI	852	0	742	43	0
39	DI	852	0	742	52	0
40	CJ	659	0	552	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
40	DJ	659	0	552	37	0
41	CK	828	0	822	24	0
41	DK	828	0	822	32	0
42	CL	909	0	927	43	0
42	DL	909	0	927	38	0
43	CM	801	0	743	33	0
43	DM	801	0	743	37	0
44	CN	478	0	498	33	0
44	DN	478	0	497	30	0
45	CO	724	0	749	32	0
45	DO	724	0	749	27	0
46	CP	651	0	638	33	0
46	DP	651	0	638	28	0
47	CQ	823	0	891	33	0
47	DQ	823	0	891	22	0
48	CR	514	0	530	19	0
48	DR	514	0	530	17	0
49	CS	544	0	457	21	0
49	DS	544	0	457	26	0
50	CT	708	0	764	37	0
50	DT	708	0	764	26	0
51	CU	199	0	208	8	0
51	DU	199	0	208	6	0
52	AA	44	0	20	31	0
52	BA	44	0	20	23	0
53	AA	2	0	0	0	0
53	BA	2	0	0	0	0
All	All	279316	0	185722	7170	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 16.

All (7170) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:AA:3001:T8B:C13	31:DA:1492:A:H5''	1.47	1.41
34:DD:26:CYS:SG	34:DD:26:CYS:CB	2.09	1.41
34:DD:12:CYS:SG	34:DD:12:CYS:CB	2.16	1.34
52:AA:3001:T8B:H13	52:AA:3001:T8B:C22	1.58	1.33
52:BA:3001:T8B:H13	52:BA:3001:T8B:C22	1.58	1.32
1:BA:885:C:H42	1:BA:890:A:N6	1.28	1.32

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:BA:3001:T8B:H17	52:BA:3001:T8B:C19	1.66	1.25
52:AA:3001:T8B:H17	52:AA:3001:T8B:C19	1.66	1.24
1:BA:1913:A:O2'	52:BA:3001:T8B:O10	1.59	1.16
1:AA:885:C:H42	1:AA:890:A:N6	1.45	1.14
52:AA:3001:T8B:C13	31:DA:1492:A:C5'	2.26	1.12
52:AA:3001:T8B:H10	31:DA:1492:A:H5''	1.26	1.08
52:AA:3001:T8B:H8	31:DA:1492:A:H5''	1.23	1.07
1:BA:885:C:N4	1:BA:890:A:N6	2.02	1.06
1:BA:1913:A:O2'	52:BA:3001:T8B:C24	2.04	1.05
1:AA:2100:G:H1	1:AA:2189:U:H3	1.12	0.98
1:AA:1403:C:H5''	1:AA:1471:A:H1'	1.46	0.97
43:DM:65:LYS:HA	43:DM:66:LEU:HB2	1.44	0.97
31:DA:1028:C:N4	31:DA:1033:G:H1	1.62	0.97
31:DA:1502:A:H2	31:DA:1505:G:H1	1.12	0.96
1:BA:1049:C:HO2'	1:BA:1050:A:H8	1.06	0.96
1:BA:1332:G:N2	1:BA:1609:A:O2'	1.99	0.96
1:AA:1689:A:H62	1:AA:1698:A:H2	1.12	0.95
1:AA:631:A:OP1	11:AP:65:ARG:NH1	2.00	0.95
1:BA:1019:U:HO2'	1:BA:1021:A:H2	1.02	0.95
34:CD:189:PRO:HB3	34:CD:194:LEU:HD11	1.50	0.94
1:AA:1332:G:N2	1:AA:1609:A:O2'	1.99	0.94
1:BA:1913:A:O2'	52:BA:3001:T8B:C23	2.16	0.93
31:CA:1502:A:H2	31:CA:1505:G:H1	1.03	0.93
6:BG:105:LYS:NZ	26:B4:25:TYR:O	2.02	0.93
31:CA:939:G:H5''	37:CG:102:ARG:HH22	1.30	0.93
1:BA:83:G:N2	1:BA:102:G:O2'	2.01	0.93
4:AE:47:VAL:HG21	4:AE:86:PRO:HD2	1.50	0.93
1:AA:1915:U:C2	52:AA:3001:T8B:C28	2.52	0.93
52:BA:3001:T8B:C22	52:BA:3001:T8B:C19	2.37	0.92
52:AA:3001:T8B:H8	31:DA:1492:A:C5'	1.95	0.92
8:AI:101:LEU:HD22	8:AI:107:VAL:HG11	1.50	0.92
34:DD:32:ALA:O	34:DD:36:ARG:N	2.02	0.92
1:BA:1913:A:C2'	52:BA:3001:T8B:O10	2.18	0.91
31:DA:448:A:OP2	31:DA:485:G:N2	2.02	0.91
1:BA:1310:G:OP2	29:B7:9:ARG:NH1	2.02	0.91
52:AA:3001:T8B:C22	52:AA:3001:T8B:C19	2.37	0.91
9:AN:47:ALA:HB2	9:AN:112:LEU:HD11	1.50	0.91
39:DI:16:ARG:HB2	39:DI:64:THR:HG23	1.53	0.90
1:AA:885:C:N4	1:AA:890:A:N6	2.19	0.90
31:CA:165:C:H2'	31:CA:166:G:H8	1.34	0.90
1:BA:1913:A:HO2'	52:BA:3001:T8B:C24	1.82	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DD:31:CYS:O	34:DD:33:MET:N	2.05	0.90
1:BA:686:G:H5''	29:B7:11:LYS:HE2	1.53	0.90
31:DA:1321:C:H3'	31:DA:1322:C:H5''	1.54	0.89
31:DA:735:C:H2'	31:DA:736:C:H6	1.38	0.89
31:DA:677:U:H3	31:DA:713:G:H22	1.14	0.89
1:AA:819:A:OP2	1:AA:1187:G:N2	2.06	0.88
4:AE:143:ASN:HD22	4:AE:147:PRO:HD2	1.35	0.88
32:CB:88:ALA:HB2	32:CB:219:VAL:HG13	1.55	0.88
8:BI:1:MET:N	8:BI:21:VAL:O	2.05	0.88
1:AA:676:A:H8	1:AA:2069:G:H21	1.16	0.88
1:BA:2207:G:O2'	1:BA:2208:A:OP1	1.89	0.88
31:CA:1293:G:HO2'	31:CA:1294:G:H8	1.15	0.88
1:BA:1332:G:H5''	1:BA:1332:G:C8	2.09	0.88
31:CA:78:G:H1	31:CA:91:C:H42	1.16	0.88
31:CA:1203:C:H2'	31:CA:1204:A:H8	1.39	0.88
32:CB:115:LEU:HD13	32:CB:145:LEU:HB3	1.54	0.88
4:AE:128:SER:OG	4:AE:129:HIS:N	2.02	0.87
1:BA:102:G:O2'	1:BA:103:A:OP2	1.90	0.87
1:BA:819:A:OP2	1:BA:1187:G:N2	2.07	0.87
31:CA:35:G:O2'	42:CL:118:SER:O	1.91	0.87
31:DA:673:G:H2'	31:DA:674:G:C8	2.10	0.87
45:DO:25:THR:HG21	45:DO:70:LEU:HB2	1.57	0.87
23:A1:3:LYS:HB2	23:A1:61:ARG:HH12	1.39	0.87
1:BA:1530:C:O2'	1:BA:1531:C:O5'	1.92	0.87
20:BY:76:CYS:SG	20:BY:99:CYS:HB2	2.14	0.87
34:DD:26:CYS:CB	34:DD:31:CYS:SG	2.62	0.87
41:CK:79:SER:HA	41:CK:104:GLN:HB2	1.54	0.87
1:BA:1913:A:O2'	52:BA:3001:T8B:C21	2.23	0.87
3:BD:69:ARG:NH2	3:BD:128:GLY:O	2.07	0.87
15:BT:16:ARG:NH2	15:BT:83:ILE:O	2.07	0.87
44:DN:23:ARG:HG3	44:DN:24:CYS:H	1.39	0.86
3:AD:133:LEU:HA	3:AD:136:ILE:HD12	1.55	0.86
9:AN:24:GLY:HA2	9:AN:27:ALA:HB3	1.56	0.86
1:AA:2207:G:O2'	1:AA:2208:A:OP1	1.91	0.86
1:AA:1022:G:H22	1:AA:1142(A):A:H2	1.22	0.86
31:DA:78:G:H1	31:DA:91:C:H42	1.20	0.86
6:AG:41:GLN:HB3	6:AG:43:LEU:HD13	1.58	0.85
1:BA:676:A:H8	1:BA:2069:G:H21	1.22	0.85
1:BA:885:C:N4	1:BA:890:A:H61	1.69	0.85
1:AA:2206:G:H5'	1:AA:2207:G:N7	1.91	0.85
10:AO:35:VAL:HG11	10:AO:103:ALA:HB3	1.58	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AP:138:LEU:HD23	11:AP:145:PRO:HG3	1.57	0.85
31:CA:656:C:O2'	45:CO:28:GLN:NE2	2.08	0.85
31:DA:1071:C:OP1	35:DE:27:ARG:NH2	2.09	0.85
23:B1:64:ALA:HA	23:B1:67:ILE:HG13	1.56	0.85
21:BZ:119:GLU:OE2	21:BZ:122:ARG:NH1	2.10	0.85
1:AA:96:G:OP1	24:A2:46:GLN:NE2	2.10	0.85
1:AA:1530:C:O2'	1:AA:1531:C:O5'	1.94	0.85
21:AZ:33:LEU:HD23	21:AZ:90:VAL:HG21	1.58	0.85
52:AA:3001:T8B:C13	31:DA:1491:G:H3'	2.07	0.84
1:BA:1332:G:H5''	1:BA:1332:G:H8	1.40	0.84
3:AD:69:ARG:NH2	3:AD:128:GLY:O	2.11	0.84
1:BA:1364:G:OP2	23:B1:3:LYS:HG2	1.77	0.84
22:A0:10:THR:HG22	22:A0:12:ASN:H	1.42	0.84
15:BT:54:ARG:HA	15:BT:59:THR:HB	1.60	0.84
21:AZ:160:GLY:HA2	21:AZ:161:VAL:HB	1.59	0.84
31:CA:165:C:H2'	31:CA:166:G:C8	2.10	0.84
1:BA:139(A):G:N2	19:BX:44:GLU:OE1	2.11	0.84
6:BG:56:ALA:HB2	6:BG:153:ARG:HE	1.43	0.84
8:AI:77:LEU:HD11	8:AI:101:LEU:HG	1.59	0.84
31:DA:73:G:H1	31:DA:96:U:H3	1.26	0.84
4:BE:47:VAL:HG21	4:BE:86:PRO:HD2	1.58	0.83
18:BW:14:PRO:HG2	18:BW:78:GLU:HG2	1.58	0.83
31:CA:1177:G:N2	31:CA:1181:G:O6	2.12	0.83
31:DA:975:A:H4'	31:DA:976:G:H5''	1.61	0.83
33:DC:100:ALA:O	33:DC:102:ASN:ND2	2.12	0.83
1:AA:994:C:OP1	16:AU:53:ARG:NH2	2.11	0.82
1:BA:2306:C:H5''	1:BA:2307:G:H2'	1.59	0.82
35:DE:81:GLU:HG2	35:DE:90:VAL:HG13	1.62	0.82
41:DK:48:ILE:O	41:DK:50:TYR:N	2.12	0.82
52:AA:3001:T8B:H8	31:DA:1491:G:H3'	1.58	0.82
12:AQ:51:ARG:NH2	21:AZ:186:GLU:OE1	2.12	0.82
1:BA:1176:G:H1'	1:BA:1177:A:OP1	1.79	0.82
14:BS:96:GLY:H	14:BS:99:LYS:H	1.27	0.82
1:AA:2371:G:O2'	28:A6:46:HIS:ND1	2.12	0.82
27:B5:46:CYS:HG	27:B5:49:CYS:HG	0.83	0.81
38:DH:86:ILE:HG21	38:DH:133:LEU:HD13	1.62	0.81
1:AA:2683:C:OP1	15:AT:53:ARG:NH2	2.13	0.81
1:AA:1039:G:O6	1:AA:1116:C:N4	2.13	0.81
4:BE:105:THR:OG1	4:BE:199:ARG:NH2	2.14	0.81
43:CM:60:VAL:HG12	43:CM:66:LEU:HD11	1.63	0.81
3:AD:275:LYS:HG3	3:AD:276:LYS:H	1.43	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:CE:102:ALA:HB1	35:CE:106:PRO:HG2	1.62	0.81
31:CA:642:A:N3	38:CH:113:SER:OG	2.12	0.81
31:CA:664:G:H22	31:CA:741:G:H1	1.25	0.81
31:CA:1279:A:O2'	31:CA:1281:U:OP2	1.99	0.81
12:AQ:31:ASP:HA	12:AQ:134:ARG:HG3	1.62	0.80
1:BA:631:A:OP1	11:BP:65:ARG:NH1	2.14	0.80
31:DA:1118:C:H1'	31:DA:1179:A:C4	2.16	0.80
1:AA:71:A:C2	19:AX:31:HIS:HE1	2.00	0.80
32:DB:187:LEU:HA	32:DB:201:ILE:HB	1.62	0.80
52:AA:3001:T8B:H18	52:AA:3001:T8B:O12	1.80	0.80
52:BA:3001:T8B:H18	52:BA:3001:T8B:O12	1.80	0.80
1:BA:2190:G:H2'	1:BA:2191:G:H5''	1.63	0.80
18:BW:78:GLU:OE1	18:BW:99:ARG:NH1	2.14	0.80
31:CA:662:G:H2'	31:CA:663:A:C8	2.17	0.80
32:CB:25:ASN:O	32:CB:27:LYS:N	2.15	0.80
32:CB:69:LEU:HB3	32:CB:162:ILE:HG22	1.62	0.80
16:AU:92:ARG:HA	16:AU:95:LEU:HB2	1.64	0.80
21:AZ:5:LEU:HD23	21:AZ:47:VAL:HG21	1.61	0.80
32:CB:116:GLU:HA	32:CB:119:GLU:HB2	1.63	0.80
4:BE:128:SER:OG	4:BE:129:HIS:N	2.12	0.80
47:CQ:66:SER:O	47:CQ:70:ARG:NH1	2.15	0.80
36:DF:11:ASN:HB3	36:DF:14:LEU:HG	1.64	0.80
13:BR:103:ARG:NH1	13:BR:108:GLY:O	2.14	0.79
34:DD:24:GLU:O	34:DD:27:TYR:HB2	1.81	0.79
1:AA:1310:G:OP2	29:A7:9:ARG:NH1	2.15	0.79
1:AA:1815:A:OP2	3:AD:54:ARG:NH2	2.15	0.79
1:BA:2306:C:H3'	1:BA:2307:G:H8	1.48	0.79
39:CI:4:TYR:HB2	39:CI:19:LEU:HB2	1.64	0.79
3:AD:108:PRO:HB3	3:AD:143:HIS:HE1	1.48	0.79
32:DB:84:GLU:HB3	32:DB:219:VAL:HG21	1.62	0.79
41:DK:22:HIS:HB3	41:DK:29:ILE:HB	1.63	0.79
1:AA:1315:C:H42	1:AA:1337:G:H1	1.31	0.79
3:AD:108:PRO:HB3	3:AD:143:HIS:CE1	2.18	0.79
1:BA:1913:A:N6	31:CA:1493:A:H5'	1.98	0.79
39:CI:83:ARG:HA	39:CI:86:VAL:HG22	1.64	0.79
1:AA:139(A):G:N2	19:AX:44:GLU:OE1	2.15	0.79
41:DK:79:SER:HA	41:DK:104:GLN:HB2	1.64	0.79
1:BA:1914:C:O2'	1:BA:1915:U:H5'	1.82	0.79
48:CR:32:ARG:HA	48:CR:69:THR:HG21	1.65	0.79
31:CA:1128:C:O2'	31:CA:1130:A:N7	2.16	0.78
41:DK:99:GLN:HG2	41:DK:105:VAL:HG21	1.64	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:1914:C:O2'	1:BA:1915:U:P	2.40	0.78
21:BZ:160:GLY:HA2	21:BZ:161:VAL:HB	1.65	0.78
27:B5:16:ARG:NH1	27:B5:17:ASP:OD1	2.16	0.78
1:BA:2306:C:H3'	1:BA:2307:G:C8	2.18	0.78
1:BA:1379:A:H4'	1:BA:1380:G:OP2	1.82	0.78
1:BA:2683:C:O2	10:BO:70:LYS:NZ	2.15	0.78
2:AB:49:C:H2'	2:AB:50:G:H8	1.48	0.78
1:BA:301:G:OP2	20:BY:84:ARG:NH2	2.16	0.78
15:BT:118:ARG:HH11	15:BT:118:ARG:HG3	1.47	0.78
1:AA:903:C:H2'	1:AA:904:C:H6	1.49	0.78
1:AA:2306:C:H5''	1:AA:2307:G:H2'	1.66	0.78
31:DA:735:C:H2'	31:DA:736:C:C6	2.19	0.78
1:AA:571:A:H5'	1:AA:2030:A:H62	1.48	0.78
15:BT:24:PRO:HA	15:BT:49:VAL:HG22	1.66	0.78
1:AA:1332:G:N2	1:AA:1609:A:HO2'	1.82	0.78
31:CA:922:G:H4'	35:CE:20:GLN:HA	1.66	0.77
1:BA:252:G:OP2	11:BP:50:ARG:NH1	2.17	0.77
46:CP:51:VAL:HG12	46:CP:53:VAL:H	1.48	0.77
31:DA:1228:C:H2'	31:DA:1229:A:H8	1.47	0.77
52:AA:3001:T8B:H10	31:DA:1492:A:C5'	2.05	0.77
9:AN:20:GLY:HA2	9:AN:61:ARG:HD3	1.67	0.77
1:BA:1047:G:H2'	1:BA:1110:G:H22	1.47	0.77
1:BA:1204:A:H2	1:BA:1241:A:H62	1.33	0.77
31:CA:986:A:H1'	49:CS:55:LYS:HA	1.65	0.77
31:DA:344:A:H5''	31:DA:345:C:H5	1.49	0.77
31:DA:564:C:O2'	38:DH:91:ARG:NH2	2.16	0.77
31:CA:272:C:H2'	31:CA:273:A:H8	1.49	0.77
39:CI:71:SER:HA	39:CI:74:ILE:HD12	1.65	0.77
31:DA:922:G:H4'	35:DE:20:GLN:HA	1.65	0.77
35:DE:75:THR:OG1	35:DE:76:ILE:N	2.17	0.77
5:BF:185:ASP:HA	5:BF:188:ARG:HD3	1.66	0.77
7:BH:40:GLU:OE2	7:BH:60:ARG:NH1	2.17	0.77
14:BS:25:ARG:NH1	14:BS:42:ASP:OD2	2.18	0.77
1:AA:686:G:H5''	29:A7:11:LYS:HE2	1.65	0.77
1:AA:1816:G:O6	3:AD:35:LYS:NZ	2.13	0.77
41:CK:48:ILE:O	41:CK:50:TYR:N	2.18	0.77
1:AA:2839:G:H5'	13:AR:46:GLY:HA2	1.65	0.77
1:BA:1798:U:H5'	3:BD:259:THR:HG22	1.65	0.77
10:BO:35:VAL:HG11	10:BO:103:ALA:HB3	1.67	0.77
18:BW:18:ARG:HG3	18:BW:76:VAL:HB	1.65	0.77
1:BA:1341:U:OP2	1:BA:1394:U:O2'	2.03	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:BN:89:LYS:O	9:BN:93:THR:OG1	2.04	0.76
1:AA:1412:A:H2'	1:AA:1413:G:H8	1.51	0.76
1:BA:2820:A:OP2	13:BR:2:ARG:NH2	2.19	0.76
2:AB:28:C:H2'	2:AB:29:A:C8	2.21	0.76
5:AF:167:ALA:HB1	5:AF:173:VAL:HG11	1.68	0.76
35:DE:122:GLU:O	35:DE:126:ARG:NH1	2.18	0.76
43:DM:6:GLY:HA3	43:DM:67:GLU:HB2	1.67	0.76
15:AT:54:ARG:HA	15:AT:59:THR:HB	1.65	0.76
1:BA:2562:U:H1'	10:BO:23:ARG:HH11	1.51	0.76
1:AA:301:G:OP2	20:AY:84:ARG:NH2	2.18	0.76
1:AA:1174:A:H5'	1:AA:1177:A:N6	1.99	0.76
1:AA:2329:G:H21	22:A0:41:ARG:HG3	1.50	0.76
1:AA:1176:G:H1'	1:AA:1177:A:OP1	1.85	0.76
1:BA:2100:G:H1	1:BA:2189:U:H3	1.32	0.76
7:BH:70:THR:O	7:BH:72:ILE:N	2.18	0.76
36:DF:61:LEU:HB3	36:DF:63:TYR:HE2	1.51	0.76
3:AD:148:GLU:HB2	3:AD:151:LYS:HD2	1.68	0.76
1:BA:1045:A:H1'	1:BA:1047:G:C2	2.21	0.76
21:BZ:151:HIS:O	21:BZ:153:SER:N	2.19	0.76
1:BA:2134:A:N3	1:BA:2159:G:O2'	2.18	0.76
1:BA:2287:A:H62	1:BA:2344:U:H3	1.31	0.76
31:CA:310:G:OP2	46:CP:27:LYS:NZ	2.19	0.75
1:AA:102:G:O2'	1:AA:103:A:OP2	2.04	0.75
1:BA:83:G:H22	1:BA:102:G:HO2'	1.32	0.75
35:CE:50:GLU:HB2	35:CE:53:LEU:HD13	1.68	0.75
43:CM:65:LYS:HA	43:CM:66:LEU:HB2	1.68	0.75
31:DA:1147:C:HO2'	39:DI:5:TYR:HH	1.33	0.75
31:CA:433:C:H2'	31:CA:434:U:H6	1.52	0.75
26:A4:7:PRO:HB2	26:A4:27:THR:HG21	1.68	0.75
1:BA:1531:C:H42	1:BA:1538:G:H1	1.34	0.75
1:AA:2329:G:N2	22:A0:41:ARG:HG3	2.02	0.75
1:BA:1913:A:H2'	52:BA:3001:T8B:O2	1.87	0.75
21:AZ:69:THR:HG22	21:AZ:90:VAL:HA	1.67	0.75
31:DA:171:A:H2'	31:DA:172:A:C8	2.22	0.75
31:DA:974:A:OP2	44:DN:29:ARG:NH2	2.19	0.75
1:BA:9:U:H3	1:BA:2629:A:H2	1.35	0.75
31:CA:559:A:OP1	35:CE:126:ARG:NH2	2.20	0.75
31:CA:991:U:O2'	31:CA:992:U:OP2	2.05	0.75
33:CC:177:THR:HB	33:CC:180:ALA:HB2	1.66	0.75
31:DA:1013:G:N2	31:DA:1016:A:OP2	2.19	0.75
1:AA:1178:C:H2'	1:AA:1179:C:H6	1.52	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AE:105:THR:OG1	4:AE:199:ARG:NH2	2.19	0.75
13:BR:20:LEU:HD21	13:BR:40:LYS:HD3	1.68	0.75
47:CQ:67:LYS:HA	47:CQ:70:ARG:HH12	1.52	0.75
1:AA:1980:G:O2'	1:AA:1982:C:OP2	2.05	0.75
1:BA:71:A:C2	19:BX:31:HIS:HE1	2.05	0.75
1:BA:1049:C:O2'	1:BA:1050:A:H8	1.69	0.75
31:DA:737:A:H2'	31:DA:738:C:C6	2.21	0.75
31:CA:975:A:H4'	31:CA:976:G:H5''	1.68	0.74
1:BA:779:U:OP1	3:BD:49:ILE:HG13	1.86	0.74
34:DD:72:GLU:OE1	34:DD:207:TYR:OH	2.05	0.74
46:DP:74:LEU:HB3	46:DP:79:VAL:HG11	1.68	0.74
12:AQ:39:PRO:HD3	12:AQ:99:PRO:HG3	1.69	0.74
1:BA:1858:G:O2'	1:BA:1884:A:N6	2.20	0.74
1:BA:2131:G:OP1	1:BA:2132:U:H3'	1.86	0.74
31:CA:932:C:H5'	37:CG:4:ARG:HG2	1.68	0.74
31:CA:1001:A:H61	31:CA:1040:U:H3	1.34	0.74
3:BD:8:PRO:HB3	3:BD:14:ARG:HB2	1.69	0.74
31:CA:193:C:H2'	31:CA:194:C:H6	1.51	0.74
31:DA:460:G:O6	31:DA:470:C:H5''	1.88	0.74
31:DA:1004:A:H8	31:DA:1025:U:H3	1.34	0.74
10:AO:2:ILE:HD12	10:AO:6:THR:HG21	1.70	0.74
7:BH:33:LEU:HD21	7:BH:136:ILE:HG13	1.67	0.74
31:CA:1063:C:OP2	31:CA:1064:G:O2'	2.03	0.74
1:AA:994:C:O2'	1:AA:996:A:OP1	2.03	0.74
1:BA:102:G:OP1	24:B2:7:ARG:NH2	2.20	0.74
8:BI:78:THR:O	8:BI:104:GLN:NE2	2.21	0.74
1:AA:468:G:N7	29:A7:39:ARG:NH2	2.36	0.74
8:AI:92:VAL:HG13	8:AI:120:ILE:HB	1.69	0.74
27:A5:41:PRO:O	27:A5:44:THR:OG1	2.04	0.74
31:DA:1030:C:N3	31:DA:1031:G:N2	2.36	0.74
1:AA:885:C:N4	1:AA:890:A:H61	1.86	0.74
1:AA:1593:G:H2'	1:AA:1594:G:C8	2.23	0.74
3:AD:71:ASP:HB3	3:AD:103:ARG:HH22	1.53	0.74
41:CK:29:ILE:HG23	41:CK:44:SER:HB3	1.68	0.74
42:CL:27:LEU:O	42:CL:29:GLY:N	2.21	0.74
1:AA:143:G:H2'	1:AA:143(A):C:H6	1.52	0.73
1:AA:2318:G:O2'	1:AA:2319:G:OP1	2.04	0.73
31:DA:826:C:O2	31:DA:874:G:N2	2.18	0.73
31:DA:1128:C:O2'	31:DA:1130:A:N7	2.19	0.73
33:DC:33:LEU:O	33:DC:37:GLN:NE2	2.20	0.73
1:AA:587:C:OP2	11:AP:21:ARG:NH2	2.21	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2327:A:H2'	1:AA:2328:A:C8	2.22	0.73
31:DA:985:C:H2'	31:DA:986:A:H8	1.49	0.73
6:BG:76:SER:HA	6:BG:83:ARG:HA	1.71	0.73
31:DA:982:U:H5''	44:DN:6:LEU:HD21	1.69	0.73
2:AB:49:C:H2'	2:AB:50:G:C8	2.24	0.73
32:CB:201:ILE:HG21	32:CB:214:ILE:HG21	1.70	0.73
1:BA:1174:A:H4'	1:BA:1175:U:OP1	1.87	0.73
1:BA:1779:U:H5	1:BA:1784:A:N7	1.87	0.73
1:BA:2130:U:HO2'	1:BA:2133:G:HO2'	1.31	0.73
31:DA:56:U:H2'	31:DA:57:G:C8	2.23	0.73
32:DB:88:ALA:HB2	32:DB:219:VAL:HG13	1.68	0.73
1:BA:2327:A:H2'	1:BA:2328:A:C8	2.23	0.73
21:BZ:110:GLY:HA3	21:BZ:174:VAL:HG11	1.70	0.73
31:CA:1203:C:H2'	31:CA:1204:A:C8	2.22	0.73
1:AA:1779:U:H5	1:AA:1784:A:N7	1.86	0.73
2:AB:2:C:H2'	2:AB:3:C:H6	1.52	0.73
1:BA:1021:A:H2'	1:BA:1023:U:H5'	1.69	0.73
31:CA:73:G:H1	31:CA:96:U:H3	1.37	0.73
31:CA:545:C:H5''	34:CD:72:GLU:HG2	1.70	0.73
31:CA:939:G:H5''	37:CG:102:ARG:NH2	2.01	0.73
32:DB:116:GLU:HA	32:DB:119:GLU:HB2	1.68	0.73
2:AB:2:C:H2'	2:AB:3:C:C6	2.24	0.73
31:CA:78:G:H1	31:CA:91:C:N4	1.86	0.73
31:CA:343:U:O2'	31:CA:346:G:O6	2.02	0.73
31:CA:1353:G:OP1	51:CU:10:ARG:NH1	2.22	0.73
31:DA:390:C:O3'	46:DP:28:ARG:NH2	2.20	0.73
46:DP:53:VAL:HG13	46:DP:79:VAL:HG23	1.71	0.73
31:CA:1264:C:H2'	31:CA:1265:G:H8	1.54	0.73
21:BZ:19:ARG:NH1	21:BZ:84:GLU:O	2.22	0.73
31:CA:1030:C:N3	31:CA:1031:G:N2	2.37	0.73
32:CB:178:ARG:HH21	38:CH:74:PRO:HB3	1.53	0.73
32:DB:77:ALA:HB2	32:DB:211:ILE:HD13	1.71	0.73
1:BA:2777:G:H5''	1:BA:2778:A:H5'	1.70	0.72
31:DA:1030:C:N4	31:DA:1031:G:N1	2.34	0.72
11:AP:101:VAL:O	11:AP:103:ALA:N	2.22	0.72
1:BA:1420:U:O2'	1:BA:1421:G:OP1	2.07	0.72
32:DB:12:GLU:O	32:DB:16:HIS:ND1	2.22	0.72
1:BA:1913:A:C3'	52:BA:3001:T8B:O10	2.38	0.72
34:CD:32:ALA:O	34:CD:36:ARG:N	2.22	0.72
31:DA:1179:A:H4'	39:DI:103:THR:HA	1.71	0.72
1:AA:1049:C:HO2'	1:AA:1050:A:H8	1.35	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1801:G:OP2	3:AD:154:LYS:NZ	2.18	0.72
1:AA:2287:A:H62	1:AA:2344:U:H3	1.37	0.72
17:BV:60:GLU:HB2	17:BV:97:LYS:HE2	1.71	0.72
31:DA:426:G:OP1	34:DD:38:TYR:OH	2.04	0.72
31:DA:501:C:H2'	31:DA:502:G:C8	2.24	0.72
31:DA:656:C:O2'	45:DO:28:GLN:NE2	2.21	0.72
41:DK:110:ASP:HB3	48:DR:85:LEU:HB3	1.72	0.72
5:AF:185:ASP:HA	5:AF:188:ARG:HD3	1.70	0.72
6:BG:44:GLY:HA2	6:BG:88:ILE:HG22	1.72	0.72
40:DJ:35:SER:HB3	40:DJ:73:ASP:H	1.55	0.72
1:AA:607:U:OP1	5:AF:102:PRO:HA	1.88	0.72
1:AA:2562:U:H1'	10:AO:23:ARG:HH11	1.54	0.72
32:CB:84:GLU:HB3	32:CB:219:VAL:HG21	1.69	0.72
31:DA:973:G:H3'	31:DA:974:A:H5''	1.70	0.72
1:AA:958:U:O2	2:AB:90:A:O2'	2.05	0.72
1:AA:1204:A:H2	1:AA:1241:A:H62	1.35	0.72
1:AA:2130:U:O2'	1:AA:2133:G:O2'	2.08	0.72
1:AA:2316:C:H2'	1:AA:2317:C:C6	2.25	0.72
4:AE:111:ARG:HG3	4:AE:160:TYR:CD1	2.25	0.72
7:AH:118:PRO:HG2	7:AH:121:ILE:HG13	1.72	0.72
1:BA:1914:C:O2'	1:BA:1915:U:C5'	2.38	0.72
1:BA:1915:U:O4	1:BA:1916:A:C2	2.43	0.72
20:BY:23:ARG:HG2	20:BY:42:VAL:HG22	1.71	0.72
31:DA:473:G:H2'	31:DA:474:G:H8	1.55	0.72
32:DB:167:PRO:HG3	32:DB:188:ALA:HB2	1.71	0.72
6:AG:25:TYR:HB3	6:AG:30:GLU:HB2	1.71	0.72
11:BP:52:GLU:OE1	11:BP:55:ARG:NH1	2.22	0.72
31:CA:976:G:H5'	31:CA:1358:U:O2'	1.90	0.72
31:DA:1028:C:C2	31:DA:1033:G:N2	2.58	0.72
1:AA:1915:U:N3	52:AA:3001:T8B:C27	2.52	0.72
1:BA:1980:G:O2'	1:BA:1982:C:OP2	2.06	0.72
31:CA:344:A:H4'	31:CA:345:C:OP2	1.90	0.72
1:AA:2147:G:H2'	1:AA:2148:G:O4'	1.90	0.72
33:CC:19:GLU:O	33:CC:40:ARG:NH2	2.23	0.72
31:DA:1030:C:H42	31:DA:1031:G:H1	1.37	0.72
25:B3:8:LEU:HD13	25:B3:31:LEU:HD23	1.71	0.71
31:CA:1224:G:N2	43:CM:100:GLY:O	2.22	0.71
1:AA:1359:A:H61	1:AA:1372:U:H3	1.38	0.71
4:AE:143:ASN:HD22	4:AE:147:PRO:CD	2.03	0.71
12:AQ:32:TYR:CE2	12:AQ:133:ARG:HG3	2.26	0.71
44:DN:40:CYS:O	44:DN:42:ILE:N	2.22	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2592:G:O6	1:AA:2601:C:N4	2.20	0.71
8:AI:105:HIS:CD2	8:AI:105:HIS:H	2.05	0.71
1:BA:1210:A:H5'	1:BA:1210:A:H8	1.55	0.71
1:BA:1247:A:OP1	5:BF:95:ARG:NH2	2.22	0.71
1:BA:1359:A:H61	1:BA:1372:U:H3	1.36	0.71
1:BA:2296:U:OP2	14:BS:9:ARG:NH2	2.22	0.71
41:CK:82:VAL:HB	41:CK:108:ILE:HG12	1.71	0.71
44:CN:24:CYS:SG	44:CN:25:VAL:N	2.63	0.71
38:DH:12:ARG:HD2	38:DH:26:VAL:HG12	1.71	0.71
1:AA:252:G:OP2	11:AP:50:ARG:NH1	2.23	0.71
1:AA:882:G:H1	1:AA:894:C:H42	1.36	0.71
1:BA:2287:A:N6	1:BA:2344:U:H3	1.89	0.71
31:DA:344:A:H5''	31:DA:345:C:C5	2.25	0.71
31:DA:1073:U:H2'	31:DA:1074:G:H8	1.54	0.71
31:DA:1329:A:H5''	43:DM:26:GLY:H	1.55	0.71
1:AA:885:C:H42	1:AA:890:A:H61	1.39	0.71
14:AS:96:GLY:H	14:AS:99:LYS:H	1.35	0.71
28:B6:30:THR:O	28:B6:30:THR:OG1	2.07	0.71
31:CA:1502:A:H2	31:CA:1505:G:N1	1.84	0.71
42:CL:89:ARG:HA	42:CL:97:ARG:HA	1.73	0.71
31:DA:1032:G:H2'	31:DA:1033:G:C8	2.25	0.71
37:DG:113:GLU:OE2	37:DG:122:HIS:ND1	2.21	0.71
12:AQ:138:ASP:OD2	21:AZ:81:ARG:NH1	2.23	0.71
1:BA:2115:G:O2'	1:BA:2166:G:N2	2.24	0.71
33:DC:177:THR:HB	33:DC:180:ALA:HB2	1.73	0.71
1:BA:1109:C:H5	1:BA:1110:G:C6	2.09	0.71
31:CA:1264:C:H2'	31:CA:1265:G:C8	2.26	0.71
32:CB:184:VAL:N	32:CB:198:ASP:OD2	2.23	0.71
43:DM:10:PRO:HG2	43:DM:18:ALA:HB1	1.72	0.71
1:AA:579:G:H2'	1:AA:580:C:C6	2.26	0.71
1:AA:676:A:H2	1:AA:802:A:H61	1.38	0.71
31:CA:1030:C:N4	31:CA:1031:G:N1	2.39	0.71
35:CE:139:LEU:HA	35:CE:142:LEU:HD12	1.72	0.71
31:DA:953:G:C2	31:DA:954:G:H1'	2.26	0.71
35:DE:88:LYS:HB3	35:DE:123:LEU:HB2	1.71	0.71
1:AA:2777:G:H5''	1:AA:2778:A:H5'	1.73	0.71
27:A5:40:LYS:HG2	27:A5:46:CYS:HB2	1.72	0.71
43:DM:19:LEU:HD21	43:DM:56:LEU:HD21	1.73	0.71
1:AA:2206:G:H3'	1:AA:2207:G:C8	2.26	0.71
1:BA:956:G:OP2	12:BQ:14:ARG:NH2	2.24	0.71
31:CA:503:C:OP2	42:CL:116:SER:HB3	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:1030(C):G:H2'	31:DA:1030(D):A:H8	1.56	0.71
1:BA:2115:G:H21	1:BA:2171:A:H61	1.37	0.70
48:CR:53:ARG:HH21	48:CR:60:ALA:N	1.89	0.70
31:DA:664:G:H22	31:DA:741:G:H1	1.38	0.70
31:DA:1237:C:H3'	31:DA:1336:C:H41	1.55	0.70
44:DN:7:ILE:HA	44:DN:23:ARG:HE	1.55	0.70
30:B8:6:THR:HG22	30:B8:63:PRO:HD2	1.73	0.70
31:CA:627:G:H2'	31:CA:628:G:H8	1.56	0.70
31:CA:954:G:H21	31:CA:1227:A:H62	1.38	0.70
1:BA:833:U:O2	11:BP:55:ARG:NH2	2.23	0.70
20:AY:28:LYS:HG3	20:AY:40:GLU:HG2	1.73	0.70
1:BA:1140:C:O3'	9:BN:25:ARG:NH1	2.24	0.70
35:CE:75:THR:OG1	35:CE:76:ILE:N	2.24	0.70
45:CO:39:LEU:O	45:CO:42:HIS:N	2.24	0.70
31:DA:946:A:H2'	31:DA:947:G:C8	2.26	0.70
29:A7:35:ARG:HG3	29:A7:42:LEU:HD11	1.74	0.70
1:BA:1174:A:H1'	1:BA:1175:U:H5''	1.72	0.70
1:BA:1427:A:H4'	1:BA:1428:C:O5'	1.91	0.70
37:CG:145:ALA:O	37:CG:147:ALA:N	2.24	0.70
49:CS:31:ILE:HG23	49:CS:49:ILE:HA	1.73	0.70
31:DA:78:G:H1	31:DA:91:C:N4	1.88	0.70
32:DB:134:GLU:O	32:DB:137:ARG:NH2	2.24	0.70
34:DD:53:ASP:O	34:DD:57:ARG:NH1	2.23	0.70
38:DH:10:LEU:HD22	38:DH:83:ILE:HD11	1.74	0.70
1:AA:1336:A:H2'	1:AA:1337:G:C8	2.27	0.70
1:AA:2051:A:OP1	4:AE:137:HIS:ND1	2.24	0.70
35:DE:126:ARG:HA	35:DE:131:ILE:HD11	1.73	0.70
8:AI:96:ASP:O	8:AI:100:ALA:N	2.22	0.70
12:BQ:34:LEU:HD11	12:BQ:129:THR:HB	1.74	0.70
31:CA:946:A:H2'	31:CA:947:G:C8	2.26	0.70
34:CD:13:ARG:O	34:CD:15:GLU:N	2.24	0.70
34:CD:41:GLY:O	34:CD:43:HIS:N	2.24	0.70
1:AA:2166:G:N2	1:AA:2172:U:O4	2.24	0.70
1:BA:1817:G:OP1	3:BD:88:ARG:NH2	2.24	0.70
3:BD:275:LYS:HG3	3:BD:276:LYS:H	1.54	0.70
37:CG:111:ARG:CZ	37:CG:122:HIS:HB3	2.22	0.70
47:CQ:58:GLU:HG3	47:CQ:77:VAL:HG21	1.74	0.70
31:DA:97:G:O2'	31:DA:98:G:H5''	1.90	0.70
32:DB:25:ASN:O	32:DB:27:LYS:N	2.24	0.70
37:DG:46:ALA:HB2	37:DG:117:ALA:HB1	1.73	0.70
1:AA:335:C:H4'	20:AY:73:ARG:HD3	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1210:A:H8	1:AA:1210:A:H5'	1.56	0.70
26:A4:42:PHE:HB3	26:A4:43:TYR:HB2	1.73	0.70
44:CN:21:TYR:OH	44:CN:23:ARG:NH2	2.25	0.70
31:DA:600:C:H2'	31:DA:601:C:C6	2.27	0.70
1:AA:1340:U:OP1	19:AX:16:LYS:NZ	2.24	0.70
35:DE:68:GLU:HG2	35:DE:70:PRO:HG3	1.74	0.70
45:DO:70:LEU:HD11	45:DO:77:ARG:HB3	1.74	0.70
1:AA:2659:G:P	7:AH:158:HIS:HE2	2.15	0.69
31:CA:235:C:H5'	47:CQ:70:ARG:HG2	1.74	0.69
1:AA:796:C:H2'	1:AA:797:C:C6	2.28	0.69
11:BP:38:GLN:HA	11:BP:41:ARG:HG2	1.73	0.69
1:AA:143:G:H4'	19:AX:35:THR:HG21	1.73	0.69
4:AE:101:ARG:NH1	4:AE:169:ASN:O	2.22	0.69
5:BF:117:ARG:HH21	5:BF:187:VAL:HA	1.57	0.69
5:AF:53:THR:HG23	5:AF:55:GLY:H	1.58	0.69
1:BA:2285:C:OP2	28:B6:6:ARG:NH1	2.26	0.69
31:DA:445:G:H2'	31:DA:446:G:H8	1.57	0.69
31:CA:627:G:H2'	31:CA:628:G:C8	2.27	0.69
34:CD:31:CYS:O	34:CD:33:MET:N	2.25	0.69
1:AA:1412:A:H2'	1:AA:1413:G:C8	2.27	0.69
50:CT:30:LYS:HA	50:CT:33:ILE:HD12	1.75	0.69
1:AA:1935:G:H1'	1:AA:1964:G:N2	2.07	0.69
1:BA:885:C:H3'	1:BA:886:C:H5''	1.73	0.69
2:BB:49:C:OP1	14:BS:97:ARG:N	2.24	0.69
19:BX:53:LYS:HB3	19:BX:82:GLN:HB3	1.74	0.69
31:CA:262:A:H2'	31:CA:263:A:C8	2.27	0.69
31:CA:373:A:H2'	31:CA:374:A:H8	1.57	0.69
31:CA:1086:U:H2'	31:CA:1087:G:H8	1.55	0.69
31:DA:1028:C:N4	31:DA:1033:G:N1	2.39	0.69
34:DD:189:PRO:HB3	34:DD:194:LEU:HD11	1.75	0.69
1:AA:1364:G:OP2	23:A1:3:LYS:HG2	1.93	0.69
1:AA:1823:G:OP1	3:AD:54:ARG:NH1	2.25	0.69
1:BA:2894:G:H8	1:BA:2894:G:O5'	1.74	0.69
20:BY:76:CYS:SG	20:BY:99:CYS:CB	2.81	0.69
38:CH:124:ALA:O	38:CH:128:GLY:N	2.26	0.69
31:DA:56:U:H2'	31:DA:57:G:H8	1.57	0.69
31:DA:1053:G:N7	31:DA:1200:C:H5'	2.08	0.69
31:DA:1346:A:N1	31:DA:1374:A:H5''	2.07	0.69
1:AA:271(F):C:H2'	1:AA:271(G):C:H6	1.57	0.69
1:AA:1049:C:O2'	1:AA:1050:A:H8	1.76	0.69
1:AA:1109:C:H5	1:AA:1110:G:C6	2.11	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1430:C:H2'	1:AA:1431:U:C6	2.27	0.69
6:AG:15:VAL:HG13	6:AG:175:LEU:HB3	1.74	0.69
1:BA:1913:A:O2'	52:BA:3001:T8B:O9	2.11	0.69
1:BA:2445:G:OP1	5:BF:74:ARG:NH2	2.26	0.69
1:BA:2815:C:H5'	27:B5:29:THR:HG21	1.73	0.69
31:DA:833:U:H2'	31:DA:834:C:H6	1.56	0.69
1:AA:641:C:O2'	1:AA:2350:C:OP1	2.06	0.69
1:AA:947:G:H2'	1:AA:948:G:C8	2.28	0.69
1:AA:2303:G:N2	1:AA:2313:C:O2	2.26	0.69
1:BA:517:C:OP1	27:B5:16:ARG:NH2	2.26	0.69
22:B0:27:GLU:HG3	22:B0:68:GLU:HA	1.74	0.69
31:DA:1030:C:C4	31:DA:1030(A):G:H1'	2.28	0.69
1:AA:794:G:H2'	1:AA:795:C:C6	2.28	0.68
1:AA:2445:G:OP1	5:AF:74:ARG:NH2	2.26	0.68
1:BA:651:G:OP1	30:B8:19:SER:OG	2.09	0.68
1:BA:652(G):G:H2'	1:BA:652(H):C:C6	2.28	0.68
35:CE:100:VAL:O	35:CE:107:ARG:NH2	2.24	0.68
35:CE:110:LEU:HD13	35:CE:118:ILE:HG21	1.74	0.68
1:AA:1786:A:H1'	1:AA:1938:A:N6	2.07	0.68
13:AR:33:ARG:NH2	27:A5:57:VAL:O	2.22	0.68
12:BQ:31:ASP:OD1	12:BQ:134:ARG:NH1	2.26	0.68
31:DA:878:G:H5'	38:DH:89:PRO:HG2	1.75	0.68
1:AA:271(K):U:H4'	1:AA:271(L):U:OP2	1.92	0.68
1:AA:517:C:OP1	27:A5:16:ARG:NH2	2.26	0.68
8:AI:1:MET:N	8:AI:1:MET:SD	2.67	0.68
1:BA:90:U:O2'	1:BA:92:A:O4'	2.10	0.68
1:BA:2104:G:N7	1:BA:2186:G:N2	2.42	0.68
31:CA:1291:G:H4'	39:CI:39:GLY:HA3	1.74	0.68
3:AD:8:PRO:HB3	3:AD:14:ARG:HB2	1.75	0.68
30:A8:61:LEU:O	30:A8:63:PRO:HD3	1.93	0.68
1:BA:1518:U:H2'	1:BA:1519:G:O4'	1.93	0.68
41:CK:34:ASP:HB3	41:CK:40:ILE:HD11	1.75	0.68
31:DA:750:G:N3	45:DO:23:GLY:HA3	2.07	0.68
31:DA:1028:C:N3	31:DA:1033:G:N2	2.40	0.68
31:DA:1279:A:O2'	31:DA:1281:U:OP2	2.11	0.68
1:AA:143:G:H2'	1:AA:143(A):C:C6	2.28	0.68
1:AA:330:A:H2	1:AA:1210:A:H2'	1.58	0.68
1:BA:1529:G:C6	1:BA:1530:C:N4	2.60	0.68
1:BA:1915:U:C5	1:BA:1916:A:C5	2.82	0.68
8:BI:104:GLN:HB2	8:BI:105:HIS:HD2	1.59	0.68
31:CA:269:C:H2'	31:CA:270:A:C8	2.29	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:343:U:O2'	31:DA:346:G:O6	2.06	0.68
1:BA:1225:G:OP1	17:BV:69:LYS:NZ	2.22	0.68
32:CB:195:ASP:O	38:CH:68:ARG:NH2	2.26	0.68
41:CK:99:GLN:HG2	41:CK:105:VAL:HG21	1.76	0.68
1:BA:1530:C:N4	1:BA:1539:G:H1	1.92	0.68
31:CA:674:G:H2'	31:CA:675:A:H8	1.59	0.68
31:CA:1100:C:O2'	31:CA:1102:A:OP1	2.08	0.68
32:DB:201:ILE:HG21	32:DB:214:ILE:HG21	1.75	0.68
38:DH:73:ASP:OD2	38:DH:75:ARG:HD3	1.93	0.68
1:BA:249:C:O2	30:B8:12:LYS:NZ	2.20	0.68
4:BE:181:LEU:HD21	15:BT:6:LEU:HD12	1.75	0.68
40:CJ:61:GLU:OE2	44:CN:49:HIS:NE2	2.27	0.68
31:DA:176:C:OP1	50:DT:29:LYS:NZ	2.17	0.68
31:DA:1347:G:N2	31:DA:1373:G:H2'	2.09	0.68
35:DE:78:HIS:HE1	35:DE:142:LEU:HA	1.59	0.68
2:AB:55:U:O3'	6:AG:27:ASN:ND2	2.25	0.68
3:AD:206:LEU:HD22	3:AD:211:ARG:HG2	1.76	0.68
8:AI:124:GLY:N	8:AI:144:VAL:HG13	2.09	0.68
18:AW:18:ARG:HG3	18:AW:76:VAL:HB	1.76	0.68
31:CA:600:C:H2'	31:CA:601:C:C6	2.29	0.68
31:CA:1030(D):A:H62	31:CA:1031:G:H21	1.40	0.68
38:CH:4:ASP:OD1	38:CH:85:ARG:NH1	2.23	0.68
1:AA:2104:G:N7	1:AA:2186:G:N2	2.41	0.67
7:AH:153:LYS:HG3	7:AH:154:PRO:HD2	1.76	0.67
1:BA:2164:C:H3'	1:BA:2165:G:H8	1.58	0.67
1:BA:2394:C:OP2	30:B8:30:ARG:NH1	2.26	0.67
2:BB:66:A:H61	2:BB:108:U:H2'	1.59	0.67
31:CA:1259:C:O2'	31:CA:1283:G:N2	2.27	0.67
1:AA:729:G:O5'	3:AD:208:LYS:NZ	2.26	0.67
4:AE:56:PRO:HG3	4:AE:74:PRO:HG2	1.75	0.67
1:BA:2318:G:O2'	1:BA:2319:G:OP1	2.10	0.67
31:DA:96:U:O2'	31:DA:97:G:OP2	2.12	0.67
1:AA:910:A:H62	12:AQ:12:GLN:HA	1.59	0.67
1:AA:1047:G:H2'	1:AA:1110:G:H22	1.59	0.67
31:CA:991:U:HO2'	31:CA:992:U:P	2.18	0.67
32:CB:194:PRO:O	32:CB:196:LEU:N	2.27	0.67
39:CI:42:ARG:O	39:CI:44:VAL:N	2.25	0.67
1:AA:272(B):G:H2'	1:AA:272(C):G:H8	1.59	0.67
1:AA:1366:A:OP1	23:A1:3:LYS:NZ	2.28	0.67
1:BA:652(B):A:H61	1:BA:655:A:H2	1.41	0.67
1:BA:1915:U:C4	1:BA:1916:A:C4	2.82	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BG:43:LEU:HD11	6:BG:153:ARG:HG2	1.76	0.67
8:BI:61:ARG:HB3	8:BI:133:HIS:HD2	1.58	0.67
31:CA:518:C:O2'	31:CA:530:G:N2	2.27	0.67
31:DA:400:C:H5''	34:DD:73:ARG:HH22	1.59	0.67
31:DA:1032:G:H2'	31:DA:1033:G:H8	1.60	0.67
1:AA:668:G:H5'	1:AA:669:G:OP2	1.93	0.67
28:A6:30:THR:OG1	28:A6:30:THR:O	2.11	0.67
12:BQ:43:THR:HG22	12:BQ:94:VAL:HG12	1.77	0.67
32:CB:87:ARG:HE	32:CB:233:SER:HB2	1.59	0.67
49:DS:11:VAL:HG23	49:DS:38:SER:HB2	1.76	0.67
1:AA:947:G:H2'	1:AA:948:G:H8	1.60	0.67
1:AA:2786:U:O2'	4:AE:62:PRO:O	2.10	0.67
8:AI:97:ILE:HA	8:AI:100:ALA:HB3	1.77	0.67
22:A0:27:GLU:HG3	22:A0:68:GLU:HA	1.74	0.67
1:BA:607:U:OP1	5:BF:102:PRO:HA	1.95	0.67
31:CA:1320:C:H5'	49:CS:70:LYS:HD3	1.77	0.67
31:DA:942:G:H21	39:DI:124:GLN:HE22	1.40	0.67
1:AA:885:C:H3'	1:AA:886:C:H5''	1.76	0.67
1:AA:2136:C:N4	1:AA:2155:G:H1	1.93	0.67
5:AF:24:LEU:HD21	5:AF:199:TRP:HH2	1.60	0.67
31:CA:243:A:H4'	31:CA:244:U:O5'	1.94	0.67
39:CI:26:VAL:HG22	39:CI:61:ALA:HB3	1.76	0.67
31:DA:801:U:H2'	31:DA:802:A:H8	1.59	0.67
41:DK:48:ILE:HD11	41:DK:64:ALA:HA	1.77	0.67
1:AA:1174:A:H1'	1:AA:1175:U:H5''	1.76	0.67
33:CC:152:ILE:HB	33:CC:199:LYS:HB2	1.75	0.67
34:CD:18:LYS:NZ	34:CD:31:CYS:HB3	2.10	0.67
1:AA:2104:G:N2	1:AA:2105:C:C2	2.62	0.67
1:AA:2349:G:H5'	1:AA:2350:C:OP2	1.95	0.67
1:BA:2820:A:OP1	13:BR:4:LEU:HD23	1.94	0.67
31:CA:1391:U:H2'	31:CA:1392:G:C8	2.30	0.67
1:AA:652(G):G:H2'	1:AA:652(H):C:C6	2.30	0.67
1:BA:1429:G:H2'	1:BA:1430:C:C6	2.30	0.67
1:BA:1639:U:H2'	1:BA:1640:C:H5''	1.75	0.67
27:B5:36:CYS:O	27:B5:37:LYS:HD3	1.94	0.67
45:CO:25:THR:HG21	45:CO:70:LEU:HB2	1.77	0.67
31:DA:17:U:O2'	31:DA:1079:G:N3	2.28	0.67
1:AA:2133:G:O2'	1:AA:2158:A:N1	2.28	0.66
1:BA:1505:C:H2'	1:BA:1506:C:H6	1.59	0.66
31:CA:1372:U:OP1	39:CI:72:GLY:N	2.28	0.66
32:DB:91:PRO:HG2	32:DB:155:LEU:HD23	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:DI:3:GLN:HG2	39:DI:20:ARG:HG2	1.77	0.66
1:AA:1531:C:H42	1:AA:1538:G:H1	1.42	0.66
2:AB:48:A:H4'	14:AS:95:HIS:HD2	1.61	0.66
1:BA:993:G:OP1	16:BU:50:ARG:NH2	2.29	0.66
1:BA:2108:C:H42	1:BA:2181:G:H1	1.41	0.66
9:BN:56:ASN:HA	9:BN:125:GLY:H	1.60	0.66
31:CA:1031:G:H2'	31:CA:1032:G:C8	2.30	0.66
34:CD:108:LEU:HB3	34:CD:110:PHE:HE1	1.60	0.66
31:DA:625:G:H2'	31:DA:626:U:C6	2.30	0.66
1:AA:9:U:H3	1:AA:2629:A:H2	1.42	0.66
1:AA:1753:G:H5''	15:AT:95:ARG:HE	1.60	0.66
5:AF:22:ALA:HB1	5:AF:24:LEU:HD22	1.77	0.66
8:AI:5:LEU:HD11	8:AI:19:VAL:HG22	1.77	0.66
1:BA:1913:A:N6	31:CA:1493:A:C5'	2.58	0.66
14:BS:14:VAL:O	14:BS:18:ILE:HG12	1.95	0.66
31:CA:308:C:H2'	31:CA:309:G:H8	1.59	0.66
32:CB:135:GLN:HA	32:CB:138:LEU:HD12	1.77	0.66
40:CJ:54:PHE:HD2	40:CJ:55:LYS:HD2	1.58	0.66
31:DA:664:G:P	48:DR:64:ARG:HH21	2.19	0.66
9:AN:128:HIS:O	9:AN:131:GLN:NE2	2.28	0.66
20:AY:99:CYS:CB	20:AY:102:CYS:HG	2.07	0.66
31:CA:991:U:H4'	31:CA:992:U:O5'	1.96	0.66
31:DA:1190:G:OP1	33:DC:5:ILE:N	2.28	0.66
1:AA:1754:C:OP1	15:AT:96:ARG:NH1	2.28	0.66
7:BH:137:ASP:HB3	7:BH:140:LYS:HB3	1.78	0.66
31:CA:222:U:H2'	31:CA:223:U:C6	2.31	0.66
33:CC:43:LEU:HD22	33:CC:47:LEU:HD12	1.78	0.66
32:DB:20:GLU:O	32:DB:40:HIS:N	2.27	0.66
1:AA:2755:C:HO2'	1:AA:2756:U:H6	1.44	0.66
1:AA:2823:A:OP1	4:AE:159:HIS:NE2	2.22	0.66
33:CC:15:THR:HG21	33:CC:181:ASN:HA	1.76	0.66
31:DA:538:G:OP2	42:DL:115:LYS:HB2	1.96	0.66
35:DE:50:GLU:HB2	35:DE:53:LEU:HD13	1.77	0.66
1:AA:2693:A:H2'	1:AA:2694:G:H8	1.60	0.66
1:AA:2884:U:H1'	27:A5:53:ALA:HB2	1.76	0.66
3:AD:242:ARG:N	3:AD:242:ARG:HD3	2.11	0.66
31:CA:413:G:N2	31:CA:428:G:H1'	2.11	0.66
31:CA:965:A:H5'	31:CA:969:A:O4'	1.96	0.66
32:CB:21:ARG:HB3	32:CB:39:ILE:HG12	1.77	0.66
31:DA:445:G:H2'	31:DA:446:G:C8	2.30	0.66
38:DH:29:SER:HB3	38:DH:32:LYS:HD2	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:AW:88:ARG:NH1	18:AW:94:ASP:OD1	2.28	0.66
5:BF:185:ASP:OD1	5:BF:188:ARG:NH1	2.27	0.66
31:CA:618:C:N3	31:CA:622:A:N6	2.44	0.66
31:CA:1228:C:H5'	43:CM:114:ARG:HB2	1.77	0.66
39:CI:46:ALA:O	39:CI:49:PRO:HD2	1.96	0.66
31:DA:1237:C:O2'	31:DA:1300:G:N2	2.25	0.66
1:AA:83:G:H22	1:AA:102:G:HO2'	1.43	0.66
9:AN:30:ILE:HG23	9:AN:52:VAL:HG11	1.78	0.66
31:CA:1025:U:O2	31:CA:1036:G:O6	2.14	0.66
34:CD:9:CYS:HA	34:CD:12:CYS:HB2	1.78	0.66
37:DG:150:ALA:HA	41:DK:59:TYR:HB3	1.78	0.66
1:AA:1113:U:H2'	1:AA:1114:G:C8	2.31	0.66
1:AA:1143:A:OP1	9:AN:25:ARG:NH2	2.29	0.66
32:CB:163:PHE:CD2	32:CB:185:ILE:HG13	2.30	0.66
31:DA:243:A:H4'	31:DA:244:U:O5'	1.94	0.66
31:DA:1286:A:H2'	31:DA:1287:A:H4'	1.77	0.66
32:DB:81:VAL:HG22	32:DB:215:LEU:HD11	1.77	0.66
39:DI:9:ARG:HB2	39:DI:104:ARG:HE	1.59	0.66
1:AA:2377:A:H2'	1:AA:2378:A:C8	2.31	0.65
26:A4:13:ARG:N	26:A4:29:PRO:O	2.25	0.65
1:BA:641:C:O2'	1:BA:2350:C:OP1	2.10	0.65
1:BA:673:C:H5''	5:BF:81:PRO:HD2	1.78	0.65
1:BA:2113:U:H2'	1:BA:2114:A:O4'	1.96	0.65
31:DA:308:C:H2'	31:DA:309:G:H8	1.60	0.65
31:DA:983:A:N1	31:DA:1222:G:N2	2.43	0.65
1:AA:903:C:H2'	1:AA:904:C:C6	2.30	0.65
1:AA:1300:U:H4'	1:AA:1301:A:H5''	1.78	0.65
1:AA:1359:A:N1	1:AA:1372:U:O4	2.30	0.65
1:AA:2296:U:OP2	14:AS:9:ARG:NH2	2.28	0.65
6:AG:108:ASN:HA	26:A4:37:SER:HB3	1.78	0.65
7:AH:137:ASP:HB3	7:AH:140:LYS:HB3	1.78	0.65
15:AT:27:THR:HB	15:AT:89:VAL:HG22	1.78	0.65
21:AZ:156:LYS:HD2	21:AZ:158:PRO:HD3	1.78	0.65
1:BA:141:A:H8	1:BA:1408:C:HO2'	1.44	0.65
1:BA:2349:G:H5'	1:BA:2350:C:OP2	1.96	0.65
31:CA:542:G:H5'	34:CD:41:GLY:HA3	1.78	0.65
31:CA:580:U:H2'	31:CA:581:G:O4'	1.96	0.65
31:CA:673:G:H2'	31:CA:674:G:C8	2.31	0.65
31:CA:1300:G:O2'	31:CA:1301:U:O5'	2.13	0.65
48:CR:56:THR:HB	48:CR:58:LEU:HD13	1.78	0.65
1:AA:384:U:H2'	1:AA:385:C:H6	1.61	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1462:C:H4'	1:AA:2703:C:H5'	1.78	0.65
31:CA:1435:G:H2'	31:CA:1436:U:C6	2.31	0.65
32:DB:186:ALA:O	32:DB:201:ILE:N	2.27	0.65
1:AA:919:G:N2	1:AA:2269:A:OP2	2.30	0.65
1:AA:1427:A:H4'	1:AA:1428:C:O5'	1.94	0.65
5:AF:110:LEU:HD21	5:AF:181:LEU:HG	1.77	0.65
13:AR:56:LYS:NZ	13:AR:90:ARG:O	2.29	0.65
25:A3:10:LYS:HB3	25:A3:53:LEU:HA	1.78	0.65
25:A3:22:ALA:O	25:A3:26:LEU:HG	1.96	0.65
20:BY:99:CYS:CB	20:BY:102:CYS:HG	2.09	0.65
31:DA:158:G:N2	31:DA:163:C:O2	2.29	0.65
1:BA:1798:U:C5'	3:BD:259:THR:HG22	2.25	0.65
21:BZ:72:ARG:NH2	21:BZ:97:GLU:O	2.29	0.65
34:CD:65:ARG:HG2	34:CD:75:PHE:CD1	2.32	0.65
31:DA:1411:C:H2'	31:DA:1412:C:H6	1.62	0.65
32:DB:24:TRP:HZ3	32:DB:29:ALA:HB2	1.59	0.65
1:AA:637:A:H8	11:AP:117:GLU:HG3	1.62	0.65
1:AA:1315:C:N4	1:AA:1337:G:H1	1.95	0.65
1:AA:2174:C:H2'	1:AA:2175:C:H6	1.62	0.65
2:AB:44:G:N3	2:AB:47:C:N4	2.43	0.65
17:AV:35:LEU:HB2	17:AV:57:VAL:HG23	1.78	0.65
1:BA:2114:A:H3'	1:BA:2115:G:C8	2.32	0.65
31:CA:192:U:H4'	50:CT:57:ARG:HD2	1.79	0.65
31:CA:1356:G:H2'	31:CA:1357:A:C8	2.32	0.65
34:CD:129:ASN:HD21	34:CD:144:ASP:HA	1.61	0.65
31:DA:189(F):U:O2	47:DQ:63:ARG:NH2	2.30	0.65
33:DC:77:ILE:HA	33:DC:84:ILE:H	1.62	0.65
39:DI:40:LEU:HB2	39:DI:43:ALA:HB2	1.79	0.65
1:AA:1019:U:H3	1:AA:1142(A):A:H62	1.43	0.65
1:AA:2297:C:O2	1:AA:2321:G:N2	2.15	0.65
1:BA:642:G:H21	1:BA:646:A:H2	1.42	0.65
31:CA:1062:U:H2'	31:CA:1063:C:C6	2.32	0.65
31:CA:1318:A:H2'	31:CA:1319:A:H5''	1.78	0.65
37:CG:89:MET:HG2	37:CG:155:ARG:HG3	1.78	0.65
31:DA:342:C:H2'	31:DA:343:U:O4'	1.97	0.65
31:DA:1178:G:N2	31:DA:1181:G:OP2	2.23	0.65
1:AA:299:A:H5''	20:AY:86:ARG:HH21	1.61	0.65
3:AD:125:ILE:HB	36:DF:81:ILE:HD11	1.79	0.65
1:BA:384:U:H2'	1:BA:385:C:H6	1.60	0.65
1:BA:1027:A:C2	1:BA:2488:A:H5'	2.32	0.65
1:BA:1448:G:H4'	1:BA:1542:A:OP1	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:BP:59:LEU:HD11	30:B8:10:ALA:HB2	1.79	0.65
13:BR:44:LEU:HD22	13:BR:48:VAL:HG23	1.79	0.65
31:CA:15:G:H4'	35:CE:24:ARG:HH12	1.62	0.65
31:DA:160:A:O5'	31:DA:160:A:H8	1.80	0.65
32:DB:155:LEU:HD11	32:DB:159:PRO:HD3	1.77	0.65
35:DE:147:ASP:OD2	35:DE:147:ASP:N	2.25	0.65
5:AF:184:TYR:CE2	5:AF:188:ARG:HD2	2.32	0.65
3:BD:108:PRO:HB3	3:BD:143:HIS:HE1	1.62	0.65
31:CA:959:A:HO2'	31:CA:984:C:HO2'	1.45	0.65
34:CD:18:LYS:HZ2	34:CD:31:CYS:HB3	1.62	0.65
35:CE:147:ASP:OD2	35:CE:147:ASP:N	2.26	0.65
32:DB:17:PHE:H	32:DB:17:PHE:HD2	1.43	0.65
1:AA:1766:U:H2'	1:AA:1767:C:H6	1.62	0.65
32:CB:16:HIS:HB3	32:CB:210:SER:HA	1.78	0.65
31:DA:1106:G:H5''	33:DC:172:ARG:HG2	1.79	0.65
1:AA:247:G:H4'	1:AA:386:G:C5	2.32	0.64
3:AD:16:MET:HG3	3:AD:206:LEU:O	1.96	0.64
1:BA:389:G:N1	11:BP:70:GLN:HG3	2.12	0.64
1:BA:1495:A:H2'	1:BA:1496:A:C8	2.32	0.64
1:BA:1689:A:H62	1:BA:1698:A:H2	1.42	0.64
31:CA:93:G:O2'	31:CA:96:U:OP2	2.15	0.64
31:CA:344:A:H5''	31:CA:345:C:H5	1.62	0.64
31:CA:933:G:O6	37:CG:3:ARG:NH2	2.31	0.64
47:CQ:45:HIS:HB2	47:CQ:65:ILE:HD13	1.79	0.64
31:DA:1352:C:H2'	31:DA:1353:G:C8	2.32	0.64
31:DA:1366:C:O2'	40:DJ:60:ARG:NH1	2.28	0.64
33:DC:5:ILE:HD11	44:DN:49:HIS:CE1	2.32	0.64
1:AA:2287:A:N6	1:AA:2344:U:H3	1.95	0.64
15:AT:64:ARG:HB2	15:AT:73:GLU:HG2	1.80	0.64
20:AY:76:CYS:HB2	20:AY:106:LEU:HD21	1.77	0.64
1:BA:336:C:O2'	20:BY:35:TYR:OH	2.15	0.64
1:BA:535:C:O3'	16:BU:53:ARG:NH1	2.29	0.64
1:BA:2079:U:OP1	23:B1:21:ARG:NH2	2.30	0.64
8:BI:5:LEU:HD11	8:BI:19:VAL:HG22	1.78	0.64
31:CA:584:G:H5'	47:CQ:91:ARG:HH22	1.62	0.64
1:AA:95:G:H4'	24:A2:46:GLN:H	1.61	0.64
1:AA:2158:A:H4'	1:AA:2159:G:H5'	1.79	0.64
21:AZ:5:LEU:HD21	21:AZ:44:PHE:HA	1.79	0.64
1:BA:857:C:H4'	22:B0:23:VAL:HG21	1.77	0.64
31:CA:100:C:H2'	31:CA:101:A:C8	2.31	0.64
32:CB:87:ARG:HH21	32:CB:233:SER:HB2	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:373:A:H2'	31:DA:374:A:H8	1.63	0.64
31:DA:674:G:H2'	31:DA:675:A:H8	1.63	0.64
1:AA:794:G:H2'	1:AA:795:C:H6	1.61	0.64
1:AA:2074:U:H2'	1:AA:2075:U:C6	2.32	0.64
15:AT:24:PRO:HA	15:AT:49:VAL:HG22	1.77	0.64
1:BA:1903:G:OP1	3:BD:241:PRO:HB2	1.97	0.64
31:CA:1306:A:H1'	31:CA:1332:A:C2	2.33	0.64
37:CG:150:ALA:HA	41:CK:59:TYR:HB3	1.78	0.64
31:DA:1356:G:H2'	31:DA:1357:A:C8	2.32	0.64
49:DS:69:HIS:HD1	49:DS:74:PHE:HZ	1.43	0.64
1:AA:2022:U:O2'	1:AA:2617:C:H5'	1.98	0.64
1:AA:2820:A:OP2	13:AR:2:ARG:NH2	2.27	0.64
5:AF:53:THR:HG22	5:AF:56:GLU:HG3	1.80	0.64
21:AZ:110:GLY:HA3	21:AZ:174:VAL:HG11	1.78	0.64
1:BA:223:A:O2'	1:BA:420:C:O2	2.15	0.64
1:BA:271(K):U:O2'	1:BA:271(M):G:N2	2.31	0.64
31:CA:625:G:H2'	31:CA:626:U:C6	2.32	0.64
31:CA:688:G:H2'	31:CA:689:C:H6	1.61	0.64
31:DA:165:C:H2'	31:DA:166:G:C8	2.33	0.64
34:DD:41:GLY:O	34:DD:43:HIS:N	2.31	0.64
1:AA:987:G:O2'	1:AA:1000:A:N3	2.28	0.64
7:BH:90:LYS:NZ	7:BH:159:GLU:OE1	2.29	0.64
20:BY:99:CYS:HB3	20:BY:102:CYS:SG	2.38	0.64
31:CA:166:G:H2'	31:CA:167:G:C8	2.32	0.64
31:CA:1293:G:O2'	31:CA:1294:G:H8	1.81	0.64
32:CB:187:LEU:HA	32:CB:201:ILE:HB	1.79	0.64
31:DA:964:A:N3	31:DA:969:A:O2'	2.25	0.64
1:AA:96:G:P	24:A2:46:GLN:HE21	2.20	0.64
10:AO:115:VAL:HG13	10:AO:121:VAL:HG21	1.80	0.64
1:BA:604:G:OP2	11:BP:90:ARG:NH1	2.31	0.64
1:BA:1187:G:H5''	17:BV:81:TYR:CE2	2.33	0.64
1:BA:2567:G:H2'	1:BA:2568:C:C6	2.33	0.64
1:AA:2126:A:H4'	1:AA:2127:G:O5'	1.97	0.64
6:AG:11:TYR:CZ	6:AG:16:ARG:HD3	2.32	0.64
7:AH:24:VAL:HG22	7:AH:35:VAL:HB	1.80	0.64
22:A0:31:VAL:HG11	22:A0:37:LEU:HD21	1.80	0.64
1:BA:2222:G:O2'	3:BD:148:GLU:HG2	1.97	0.64
26:B4:7:PRO:HB2	26:B4:27:THR:HG21	1.79	0.64
37:CG:41:ARG:O	37:CG:45:ASP:N	2.30	0.64
31:DA:105:G:H2'	31:DA:106:C:C6	2.33	0.64
33:DC:142:MET:HA	33:DC:146:ALA:HB3	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:652(Q):G:H2'	1:AA:652(R):C:H5'	1.80	0.64
1:AA:2158:A:H1'	1:AA:2159:G:C8	2.33	0.64
13:AR:101:ALA:HA	27:A5:44:THR:HG21	1.78	0.64
25:A3:6:VAL:HG13	25:A3:56:VAL:HG22	1.78	0.64
1:BA:1403:C:H5''	1:BA:1471:A:H1'	1.79	0.64
1:BA:2126:A:H4'	1:BA:2127:G:O5'	1.98	0.64
19:BX:31:HIS:CD2	19:BX:33:LYS:H	2.16	0.64
31:DA:1130:A:O5'	39:DI:20:ARG:NH2	2.29	0.64
34:DD:60:GLU:OE1	34:DD:199:ASN:N	2.28	0.64
1:AA:1187:G:H5''	17:AV:81:TYR:CE2	2.32	0.64
1:AA:1300:U:H4'	1:AA:1301:A:C5'	2.28	0.64
1:AA:2693:A:H2'	1:AA:2694:G:C8	2.33	0.64
15:AT:9:LEU:O	15:AT:12:SER:OG	2.10	0.64
1:BA:1266:G:O5'	18:BW:15:ARG:NH2	2.31	0.64
31:DA:1118:C:OP1	39:DI:9:ARG:NH1	2.31	0.64
1:BA:2099:U:H3	1:BA:2190:G:H1	1.45	0.63
4:BE:143:ASN:HB2	4:BE:147:PRO:HD2	1.78	0.63
10:BO:115:VAL:HG13	10:BO:121:VAL:HG21	1.79	0.63
31:CA:590:C:H2'	31:CA:591:U:C6	2.33	0.63
31:CA:1458:G:OP1	50:CT:35:THR:OG1	2.07	0.63
32:CB:18:GLY:HA2	32:CB:42:ILE:HG13	1.81	0.63
33:DC:11:ARG:HH21	33:DC:180:ALA:HB3	1.62	0.63
1:AA:644:A:H4'	1:AA:645:C:C5	2.33	0.63
1:AA:805:G:N2	1:AA:829:A:OP1	2.30	0.63
31:DA:954:G:H21	31:DA:1227:A:H62	1.46	0.63
32:DB:87:ARG:NH1	32:DB:220:ASP:OD1	2.31	0.63
34:DD:14:ARG:HB2	34:DD:40:PRO:HD2	1.80	0.63
1:AA:1332:G:H21	1:AA:1610:A:H8	1.45	0.63
1:AA:1430:C:H2'	1:AA:1431:U:H6	1.62	0.63
1:AA:2307:G:H8	1:AA:2307:G:OP1	1.80	0.63
31:CA:233:C:H2'	31:CA:234:C:H6	1.63	0.63
31:CA:750:G:N3	45:CO:23:GLY:HA3	2.13	0.63
31:CA:1379:G:N2	31:CA:1381:U:O4	2.30	0.63
32:CB:219:VAL:HA	32:CB:222:ILE:HD12	1.80	0.63
34:CD:108:LEU:HB3	34:CD:110:PHE:CE1	2.33	0.63
31:DA:165:C:H2'	31:DA:166:G:H8	1.63	0.63
31:DA:559:A:OP1	35:DE:126:ARG:NH2	2.31	0.63
1:AA:2567:G:H2'	1:AA:2568:C:C6	2.33	0.63
5:AF:137:LYS:HA	5:AF:140:LEU:HD23	1.81	0.63
28:A6:18:ARG:HG3	28:A6:42:TRP:CD1	2.32	0.63
8:BI:92:VAL:HG13	8:BI:120:ILE:HB	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:193:C:H2'	31:CA:194:C:C6	2.32	0.63
39:CI:5:TYR:O	39:CI:87:GLN:NE2	2.31	0.63
1:AA:107:C:H2'	1:AA:108:U:H6	1.63	0.63
1:AA:1688:U:H1'	1:AA:1701:A:C6	2.33	0.63
1:AA:2125:G:H22	1:AA:2172:U:H3'	1.63	0.63
1:BA:443:A:H1'	1:BA:1201:C:O4'	1.99	0.63
6:BG:41:GLN:HE22	6:BG:153:ARG:HB3	1.63	0.63
31:DA:1502:A:H2	31:DA:1505:G:N1	1.91	0.63
35:DE:126:ARG:HG2	35:DE:126:ARG:HH11	1.63	0.63
43:DM:60:VAL:HA	43:DM:64:TRP:HZ3	1.63	0.63
1:AA:322:A:OP2	5:AF:169:ASN:HB2	1.98	0.63
1:AA:1021:A:H62	1:AA:1141:U:H3	1.46	0.63
1:BA:639:U:H2'	1:BA:640:C:C6	2.33	0.63
1:BA:676:A:H2	1:BA:802:A:H61	1.47	0.63
1:BA:2166:G:N2	1:BA:2172:U:O4	2.31	0.63
33:CC:150:LYS:HB3	33:CC:173:VAL:HG21	1.81	0.63
49:CS:50:ALA:HB1	49:CS:57:HIS:HB3	1.79	0.63
31:DA:985:C:H2'	31:DA:986:A:C8	2.32	0.63
31:DA:1025:U:C2	31:DA:1036:G:O6	2.51	0.63
46:DP:72:ARG:HH21	46:DP:73:LEU:HD21	1.64	0.63
1:AA:879:G:H2'	1:AA:880:G:O4'	1.97	0.63
28:A6:16:CYS:HB3	28:A6:43:CYS:SG	2.39	0.63
8:BI:93:THR:HG22	8:BI:119:PRO:HB3	1.81	0.63
9:BN:20:GLY:HA2	9:BN:61:ARG:HD3	1.79	0.63
12:BQ:135:ASP:OD2	21:BZ:49:ARG:NH2	2.31	0.63
31:CA:677:U:H3	31:CA:713:G:H22	1.46	0.63
31:CA:1251:A:H2'	31:CA:1252:A:C8	2.34	0.63
31:DA:430:A:OP1	34:DD:9:CYS:HB2	1.98	0.63
32:DB:47:THR:HA	32:DB:202:PRO:HG2	1.81	0.63
1:AA:2198:A:O2'	1:AA:2224:G:N2	2.32	0.63
30:A8:39:LYS:HA	30:A8:42:ARG:NH1	2.13	0.63
1:BA:1503:U:H2'	1:BA:1504:C:C6	2.34	0.63
8:BI:12:LEU:HD22	8:BI:19:VAL:HG21	1.81	0.63
13:BR:55:ALA:HB2	13:BR:79:LEU:HD13	1.80	0.63
31:CA:735:C:H2'	31:CA:736:C:H6	1.64	0.63
32:CB:20:GLU:O	32:CB:40:HIS:N	2.31	0.63
39:CI:24:GLY:HA2	39:CI:60:ASP:HA	1.78	0.63
31:DA:1435:G:H2'	31:DA:1436:U:C6	2.33	0.63
1:AA:83:G:N2	1:AA:102:G:O2'	2.31	0.63
1:AA:1047:G:H2'	1:AA:1110:G:H1	1.64	0.63
1:AA:2611:U:C4	27:A5:3:LYS:HG2	2.34	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:28:C:H2'	2:AB:29:A:H8	1.63	0.63
2:BB:31:C:O2'	2:BB:53:A:N6	2.31	0.63
31:CA:132:C:H2'	31:CA:133:U:H6	1.64	0.63
31:CA:176:C:H2'	31:CA:177:C:C6	2.33	0.63
31:DA:737:A:H1'	36:DF:73:ASN:HD21	1.63	0.63
33:DC:155:GLY:O	33:DC:157:ILE:N	2.32	0.63
1:AA:1012:U:H5	9:AN:28:THR:HG21	1.64	0.62
1:AA:1833:U:O2'	1:AA:1969:A:N1	2.29	0.62
1:BA:1426:G:O6	3:BD:31:LYS:NZ	2.32	0.62
32:DB:71:VAL:HA	32:DB:93:VAL:HG23	1.81	0.62
39:DI:43:ALA:C	39:DI:45:ALA:HB2	2.19	0.62
1:AA:375:C:H2'	1:AA:376:C:C6	2.34	0.62
1:AA:1174:A:H4'	1:AA:1175:U:OP1	1.98	0.62
1:AA:1332:G:N2	1:AA:1610:A:C8	2.67	0.62
1:AA:1537:G:H2'	1:AA:1538:G:H8	1.64	0.62
1:BA:271(M):G:O2'	1:BA:271(N):U:O5'	2.16	0.62
1:BA:878:A:H2'	1:BA:879:G:H5'	1.81	0.62
52:BA:3001:T8B:H13	52:BA:3001:T8B:H17	0.73	0.62
14:BS:26:LEU:HD13	14:BS:87:PHE:HD1	1.64	0.62
31:CA:1128:C:H1'	31:CA:1146:A:H61	1.64	0.62
35:CE:78:HIS:HE1	35:CE:142:LEU:HA	1.64	0.62
31:DA:663:A:O3'	48:DR:64:ARG:NH2	2.31	0.62
1:AA:95:G:O2'	24:A2:46:GLN:HA	2.00	0.62
1:AA:392:C:H5''	1:AA:409:C:H5''	1.81	0.62
1:AA:1266:G:O5'	18:AW:15:ARG:NH2	2.32	0.62
1:AA:1420:U:O2'	1:AA:1421:G:OP1	2.17	0.62
1:AA:2854:G:H2'	1:AA:2855:C:C6	2.33	0.62
8:AI:95:LYS:O	8:AI:98:ALA:HB3	1.99	0.62
11:AP:121:LYS:O	11:AP:123:LEU:N	2.32	0.62
14:AS:34:HIS:ND1	14:AS:53:SER:OG	2.32	0.62
12:BQ:21:THR:HG21	12:BQ:101:ARG:HB2	1.79	0.62
31:CA:67:C:H2'	31:CA:68:G:C8	2.35	0.62
31:DA:1255:G:O2'	31:DA:1259:C:O4'	2.13	0.62
1:AA:576:U:H2'	1:AA:577:G:C8	2.34	0.62
1:AA:2305:A:H2	6:AG:152:LEU:HD11	1.64	0.62
1:BA:242:G:O2'	1:BA:254:G:O6	2.13	0.62
31:CA:272:C:H2'	31:CA:273:A:C8	2.34	0.62
34:CD:32:ALA:O	34:CD:35:ARG:N	2.33	0.62
31:DA:748:C:H4'	31:DA:749:C:O5'	1.99	0.62
32:DB:52:GLU:O	32:DB:56:ARG:HG2	1.99	0.62
33:DC:40:ARG:NH2	33:DC:55:VAL:O	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:186:G:H2'	1:AA:187:G:H8	1.65	0.62
1:AA:1530:C:N4	1:AA:1539:G:H1	1.98	0.62
14:AS:34:HIS:O	14:AS:97:ARG:NH2	2.32	0.62
1:BA:2125:G:H22	1:BA:2172:U:H3'	1.64	0.62
20:BY:79:CYS:HB2	20:BY:81:LYS:H	1.63	0.62
31:CA:187:C:H2'	31:CA:188:C:C6	2.35	0.62
31:DA:1201:A:H4'	31:DA:1202:G:O5'	2.00	0.62
34:DD:134:ASP:O	34:DD:136:PRO:HD3	1.99	0.62
1:AA:639:U:H2'	1:AA:640:C:C6	2.33	0.62
1:AA:1138:G:N3	9:AN:106:MET:HE2	2.15	0.62
1:AA:2233:U:H2'	1:AA:2234:G:C8	2.34	0.62
1:BA:644:A:H4'	1:BA:645:C:C5	2.34	0.62
1:BA:1434:A:H61	1:BA:1558:A:N6	1.97	0.62
1:BA:2104:G:N2	1:BA:2105:C:C2	2.67	0.62
18:BW:10:VAL:HG12	18:BW:12:ILE:HG22	1.82	0.62
31:CA:688:G:H2'	31:CA:689:C:C6	2.34	0.62
31:DA:801:U:H2'	31:DA:802:A:C8	2.33	0.62
34:DD:128:VAL:HG12	34:DD:129:ASN:HD22	1.64	0.62
35:DE:143:ARG:NH1	38:DH:77:GLU:OE1	2.32	0.62
38:DH:96:GLY:N	38:DH:99:GLU:OE2	2.27	0.62
1:AA:994:C:OP2	16:AU:54:LYS:NZ	2.24	0.62
7:AH:30:LYS:HG3	7:AH:80:SER:O	1.98	0.62
13:AR:23:ASN:OD1	13:AR:23:ASN:N	2.32	0.62
1:BA:855:G:O2'	22:B0:27:GLU:OE2	2.12	0.62
31:CA:1126:U:H1'	31:CA:1280:A:C5	2.35	0.62
31:CA:1215:G:H2'	31:CA:1216:G:H8	1.64	0.62
32:CB:100:GLY:N	32:CB:176:GLU:OE2	2.30	0.62
39:CI:44:VAL:N	39:CI:45:ALA:HA	2.15	0.62
46:CP:29:ASP:OD2	46:CP:29:ASP:N	2.31	0.62
31:DA:458:C:H2'	31:DA:460:G:H8	1.65	0.62
8:BI:61:ARG:HB3	8:BI:133:HIS:CD2	2.33	0.62
10:BO:120:GLU:HG2	10:BO:122:LEU:HG	1.81	0.62
14:BS:96:GLY:N	14:BS:99:LYS:HB3	2.14	0.62
31:DA:618:C:N3	31:DA:622:A:N6	2.47	0.62
31:DA:1071:C:H5''	35:DE:49:PRO:HG3	1.81	0.62
31:DA:1289:A:H3'	31:DA:1290:G:H8	1.65	0.62
41:DK:62:GLN:NE2	41:DK:93:GLN:OE1	2.32	0.62
1:AA:1529:G:C6	1:AA:1530:C:N4	2.68	0.62
1:AA:2735:G:H2'	1:AA:2736:G:H8	1.62	0.62
1:BA:571:A:H5'	1:BA:2030:A:H62	1.64	0.62
1:BA:1568:G:H5''	3:BD:61:LEU:HD22	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:BR:28:LEU:HD12	13:BR:48:VAL:HG21	1.82	0.62
40:CJ:61:GLU:OE1	44:CN:58:LYS:NZ	2.28	0.62
31:DA:473:G:H2'	31:DA:474:G:C8	2.34	0.62
1:AA:747:U:O2	1:AA:2014:A:H1'	2.00	0.62
1:AA:1530:C:H1'	1:AA:1531:C:OP1	1.99	0.62
1:AA:2462:U:H2'	1:AA:2463:C:O4'	1.99	0.62
5:AF:103:LYS:HA	5:AF:106:ARG:HG3	1.82	0.62
1:BA:265:A:H1'	1:BA:266:G:O4'	1.99	0.62
1:BA:1227:G:OP1	16:BU:13:LYS:NZ	2.32	0.62
1:BA:1839:G:H8	1:BA:1839:G:H5'	1.64	0.62
1:BA:2314:C:H2'	1:BA:2315:G:C8	2.35	0.62
1:BA:2483:C:N3	12:BQ:124:LYS:NZ	2.47	0.62
14:BS:96:GLY:HA2	14:BS:100:ALA:H	1.65	0.62
31:CA:475:G:O2'	31:CA:476:G:H5'	2.00	0.62
31:CA:662:G:H2'	31:CA:663:A:H8	1.62	0.62
31:CA:877:C:H5''	38:CH:88:LYS:HD3	1.80	0.62
42:CL:27:LEU:C	42:CL:29:GLY:H	2.04	0.62
32:DB:74:LYS:O	32:DB:76:GLN:N	2.32	0.62
32:DB:135:GLN:HA	32:DB:138:LEU:HD12	1.82	0.62
33:DC:48:TYR:O	33:DC:50:ALA:N	2.33	0.62
1:AA:271(K):U:O2'	1:AA:271(M):G:N2	2.32	0.61
1:AA:1283:G:N2	1:AA:1285:G:H3'	2.15	0.61
1:AA:2115:G:O2'	1:AA:2166:G:N2	2.33	0.61
1:AA:2122:U:H2'	1:AA:2123:G:H8	1.65	0.61
21:AZ:160:GLY:HA2	21:AZ:161:VAL:CB	2.29	0.61
1:BA:652(Q):G:H2'	1:BA:652(R):C:H5'	1.82	0.61
1:BA:2022:U:O2'	1:BA:2617:C:H5'	1.99	0.61
3:BD:145:VAL:HG12	3:BD:146:GLU:O	2.00	0.61
14:BS:58:LEU:HB2	14:BS:59:LYS:HB2	1.82	0.61
26:B4:42:PHE:HB3	26:B4:43:TYR:HB2	1.81	0.61
31:CA:427:U:OP1	34:CD:13:ARG:NH2	2.32	0.61
32:DB:12:GLU:C	32:DB:16:HIS:HD1	2.04	0.61
1:AA:855:G:O2'	22:A0:27:GLU:OE2	2.15	0.61
1:AA:1470:G:N2	1:AA:1520:G:OP2	2.33	0.61
1:BA:184:C:H2'	1:BA:185:U:C6	2.35	0.61
1:BA:754:C:H2'	1:BA:755:C:H6	1.64	0.61
21:BZ:69:THR:HG22	21:BZ:90:VAL:HA	1.81	0.61
31:CA:159:G:H2'	31:CA:161:A:OP2	2.00	0.61
31:CA:537:G:H2'	31:CA:538:G:C8	2.35	0.61
31:CA:955:U:H1'	31:CA:1227:A:H61	1.65	0.61
33:CC:58:GLU:HB2	33:CC:65:ALA:HB3	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:CM:23:TYR:HB3	43:CM:67:GLU:HG2	1.82	0.61
5:AF:7:TYR:HB2	5:AF:22:ALA:HB3	1.82	0.61
1:BA:84:A:N6	1:BA:102:G:H1'	2.15	0.61
3:BD:77:ALA:HB2	3:BD:97:TYR:CD2	2.35	0.61
31:CA:1155:G:H2'	31:CA:1156:G:C8	2.35	0.61
32:DB:163:PHE:HA	32:DB:185:ILE:HG12	1.83	0.61
35:DE:145:LYS:O	35:DE:149:GLU:HG2	1.99	0.61
1:AA:1022:G:N7	9:AN:66:LYS:HE2	2.15	0.61
1:AA:1506:C:H2'	1:AA:1507:A:H5'	1.82	0.61
1:AA:2563:U:H4'	10:AO:28:SER:HA	1.81	0.61
1:AA:2887:U:H2'	1:AA:2888:C:C6	2.36	0.61
2:AB:75:G:H5''	2:AB:75:G:H8	1.64	0.61
8:AI:116:LEU:HD13	8:AI:119:PRO:HA	1.81	0.61
23:A1:10:LYS:NZ	23:A1:65:SER:OG	2.34	0.61
1:BA:528:A:C2	1:BA:2043:C:H4'	2.34	0.61
3:BD:35:LYS:HG2	3:BD:36:PRO:HD2	1.81	0.61
5:BF:33:LEU:HD13	5:BF:112:MET:HE2	1.82	0.61
34:CD:22:LYS:HB2	34:CD:26:CYS:HB2	1.82	0.61
51:CU:12:LYS:O	51:CU:16:GLY:N	2.34	0.61
31:DA:262:A:H2'	31:DA:263:A:C8	2.35	0.61
31:DA:624:C:H2'	31:DA:625:G:H8	1.65	0.61
31:DA:1307:U:OP1	43:DM:101:GLN:NE2	2.34	0.61
1:AA:2294:C:OP2	14:AS:13:ARG:NH1	2.34	0.61
2:AB:24:G:H4'	2:AB:25:A:C8	2.36	0.61
1:BA:1021:A:H3'	1:BA:1021:A:C8	2.36	0.61
1:BA:1300:U:H4'	1:BA:1301:A:C5'	2.31	0.61
10:BO:8:LEU:HB2	10:BO:19:ILE:HG13	1.82	0.61
31:CA:600:C:OP1	38:CH:97:VAL:N	2.27	0.61
31:DA:246:A:N1	31:DA:278:G:O2'	2.27	0.61
31:DA:256:U:H2'	31:DA:257:G:C8	2.36	0.61
31:DA:353:A:H8	31:DA:353:A:H5'	1.64	0.61
34:DD:100:ARG:NH1	34:DD:137:SER:HB3	2.15	0.61
44:DN:29:ARG:HD3	44:DN:40:CYS:HB2	1.83	0.61
1:AA:2164:C:H3'	1:AA:2165:G:H8	1.64	0.61
1:AA:2420:C:P	30:A8:33:ASN:H	2.23	0.61
13:AR:100:LEU:HD11	13:AR:113:LEU:HD23	1.81	0.61
1:BA:2789:C:O2'	1:BA:2790:A:O2'	2.18	0.61
10:BO:2:ILE:HD12	10:BO:6:THR:HG21	1.81	0.61
28:B6:26:ASN:HB3	28:B6:29:ASN:HB2	1.82	0.61
31:CA:620:C:H2'	31:CA:621:A:O4'	2.00	0.61
31:CA:1355:G:H2'	31:CA:1356:G:H8	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:263:A:OP1	50:DT:79:ARG:NH1	2.32	0.61
32:DB:134:GLU:HA	32:DB:137:ARG:HE	1.65	0.61
32:DB:219:VAL:HA	32:DB:222:ILE:HD12	1.81	0.61
1:AA:652(G):G:H2'	1:AA:652(H):C:H6	1.66	0.61
1:AA:1429:G:H2'	1:AA:1430:C:C6	2.35	0.61
1:AA:2108:C:H42	1:AA:2181:G:H1	1.46	0.61
15:AT:56:GLY:O	15:AT:59:THR:HG23	2.00	0.61
31:CA:1141:C:H2'	31:CA:1142:G:O4'	2.01	0.61
32:CB:185:ILE:HG22	32:CB:199:TYR:HB2	1.82	0.61
31:DA:224:C:H2'	31:DA:225:C:C6	2.36	0.61
35:DE:36:ASP:OD2	35:DE:38:GLN:N	2.32	0.61
1:AA:84:A:N6	1:AA:102:G:H1'	2.16	0.61
1:AA:1495:A:H2'	1:AA:1496:A:C8	2.35	0.61
4:AE:38:THR:O	4:AE:42:ASP:N	2.31	0.61
7:AH:56:SER:OG	7:AH:61:HIS:ND1	2.28	0.61
1:BA:882:G:H1	1:BA:894:C:H42	1.49	0.61
1:BA:1914:C:HO2'	1:BA:1915:U:P	2.24	0.61
33:CC:134:ILE:HG22	33:CC:168:ALA:HB3	1.82	0.61
31:DA:811:C:O2'	31:DA:901:A:N1	2.34	0.61
1:AA:1106:A:H4'	1:AA:1107:G:OP2	2.01	0.61
14:AS:26:LEU:HD22	14:AS:87:PHE:CE1	2.36	0.61
30:A8:33:ASN:HA	30:A8:36:LYS:HD2	1.83	0.61
1:BA:1048:A:OP2	1:BA:1109:C:N4	2.34	0.61
1:BA:2390:U:P	30:B8:35:GLN:HE22	2.23	0.61
52:BA:3001:T8B:C6	31:CA:1409:C:C2	2.83	0.61
5:BF:101:LEU:HD12	5:BF:102:PRO:HD2	1.82	0.61
31:CA:559:A:H4'	31:CA:560:U:H3'	1.81	0.61
31:CA:1073:U:H2'	31:CA:1074:G:H8	1.65	0.61
31:DA:235:C:H5'	47:DQ:70:ARG:HG2	1.81	0.61
31:DA:1210:C:H3'	31:DA:1211:U:H5''	1.80	0.61
14:AS:63:THR:O	14:AS:66:ALA:HB3	2.01	0.61
20:AY:9:LYS:NZ	20:AY:28:LYS:O	2.31	0.61
26:A4:9:LEU:HD23	26:A4:27:THR:HG23	1.83	0.61
14:BS:15:ARG:O	14:BS:19:LYS:HG2	2.01	0.61
32:CB:102:LEU:HB3	32:CB:180:LEU:HD12	1.82	0.61
34:CD:170:VAL:HG22	34:CD:171:GLY:H	1.66	0.61
47:CQ:31:LEU:HD23	47:CQ:32:TYR:CZ	2.36	0.61
31:DA:171:A:H2'	31:DA:172:A:H8	1.66	0.61
32:DB:115:LEU:HD13	32:DB:145:LEU:HB3	1.83	0.61
1:AA:30:G:H2'	1:AA:31:C:C6	2.36	0.60
1:AA:854:G:H2'	1:AA:855:G:C8	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AI:106:GLY:HA2	8:AI:107:VAL:HB	1.83	0.60
1:BA:1359:A:N1	1:BA:1372:U:O4	2.34	0.60
1:BA:1665:A:H4'	10:BO:67:LYS:HB2	1.81	0.60
31:CA:828:A:H2'	31:CA:829:G:O4'	2.01	0.60
34:CD:14:ARG:HB2	34:CD:40:PRO:HD2	1.81	0.60
1:AA:1796:U:H2'	1:AA:1797:C:C6	2.36	0.60
1:AA:2079:U:OP1	23:A1:21:ARG:NH2	2.35	0.60
1:AA:2357:U:OP1	22:A0:20:ARG:NH1	2.34	0.60
1:AA:2820:A:OP1	13:AR:4:LEU:HD23	2.02	0.60
4:AE:201:THR:OG1	4:AE:202:LYS:N	2.34	0.60
7:AH:105:LEU:HD21	7:AH:148:ILE:HG23	1.83	0.60
21:AZ:54:HIS:ND1	21:AZ:101:PRO:HG3	2.16	0.60
30:A8:8:LYS:HB3	30:A8:12:LYS:HE3	1.83	0.60
1:BA:1045:A:N3	1:BA:1045:A:H2'	2.17	0.60
1:BA:1915:U:C4	1:BA:1916:A:C5	2.88	0.60
31:DA:166:G:H2'	31:DA:167:G:H8	1.65	0.60
49:DS:52:TYR:HB2	49:DS:57:HIS:CE1	2.35	0.60
1:AA:1046:A:O2'	1:AA:1047:G:OP2	2.19	0.60
1:AA:2324:C:H5''	1:AA:2325:G:H5'	1.82	0.60
6:AG:16:ARG:NE	6:AG:31:VAL:HG21	2.16	0.60
15:AT:15:VAL:HG13	15:AT:79:HIS:CE1	2.37	0.60
27:A5:16:ARG:NH1	27:A5:17:ASP:OD1	2.34	0.60
1:BA:848:G:H2'	1:BA:849:A:C8	2.36	0.60
1:BA:2408:U:H2'	1:BA:2409:G:C8	2.36	0.60
31:CA:1086:U:H3	31:CA:1099:G:H22	1.49	0.60
31:DA:976:G:H5'	31:DA:1358:U:O2'	2.01	0.60
32:DB:102:LEU:HD23	32:DB:182:ILE:HD12	1.83	0.60
33:DC:7:PRO:HG3	33:DC:201:TYR:CE2	2.36	0.60
1:AA:848:G:H2'	1:AA:849:A:C8	2.36	0.60
1:AA:2131:G:OP1	1:AA:2132:U:H3'	2.02	0.60
14:AS:52:SER:HB2	14:AS:55:ALA:H	1.65	0.60
17:AV:60:GLU:HB2	17:AV:97:LYS:HE2	1.83	0.60
23:B1:19:GLN:O	23:B1:35:THR:HG22	2.01	0.60
28:B6:11:LEU:HB3	28:B6:49:HIS:HB3	1.83	0.60
32:CB:178:ARG:HH22	38:CH:68:ARG:HH12	1.47	0.60
38:CH:29:SER:HB3	38:CH:32:LYS:HD2	1.83	0.60
31:DA:1259:C:N4	31:DA:1260:C:O2	2.35	0.60
31:DA:1411:C:H2'	31:DA:1412:C:C6	2.36	0.60
5:AF:20:LEU:HD22	5:AF:21:ALA:HB3	1.83	0.60
6:AG:138:GLN:HE22	6:AG:153:ARG:HB2	1.66	0.60
11:AP:138:LEU:HG	11:AP:143:GLY:HA3	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:1040:C:H2'	1:BA:1041:C:O4'	2.02	0.60
1:BA:2319:G:N1	14:BS:3:ARG:HA	2.16	0.60
31:CA:1118:C:H1'	31:CA:1179:A:C4	2.36	0.60
32:DB:16:HIS:HB3	32:DB:210:SER:HA	1.82	0.60
41:DK:34:ASP:HB3	41:DK:40:ILE:HD11	1.84	0.60
1:AA:185:U:H4'	1:AA:218:A:H4'	1.84	0.60
1:AA:854:G:H2'	1:AA:855:G:H8	1.66	0.60
5:AF:183:VAL:O	5:AF:187:VAL:HG23	2.01	0.60
7:AH:159:GLU:HG3	7:AH:169:VAL:HG21	1.83	0.60
1:BA:1300:U:H4'	1:BA:1301:A:H5''	1.82	0.60
1:BA:1779:U:C5	1:BA:1784:A:N7	2.70	0.60
1:BA:2136:C:N4	1:BA:2155:G:H1	1.99	0.60
16:BU:92:ARG:HA	16:BU:95:LEU:HB2	1.82	0.60
28:B6:18:ARG:HG3	28:B6:42:TRP:CD1	2.36	0.60
31:CA:400:C:H5''	34:CD:73:ARG:HH22	1.65	0.60
31:CA:460:G:O6	31:CA:470:C:H5''	2.01	0.60
31:CA:1030:C:N4	31:CA:1031:G:C6	2.70	0.60
31:CA:1327:C:H2'	31:CA:1328:C:C6	2.36	0.60
32:CB:77:ALA:HB2	32:CB:211:ILE:HD13	1.84	0.60
36:CF:10:LEU:HD23	36:CF:61:LEU:HD13	1.84	0.60
31:DA:1097:C:H1'	31:DA:1169:A:C2	2.36	0.60
1:AA:864:G:C6	1:AA:865:C:N4	2.69	0.60
1:AA:958:U:OP2	12:AQ:14:ARG:NH1	2.35	0.60
1:AA:1790:C:H5''	1:AA:1791:A:OP1	2.00	0.60
2:AB:32:C:C2	2:AB:51:G:N2	2.70	0.60
14:AS:102:ALA:HA	14:AS:105:ALA:HB3	1.84	0.60
1:BA:576:U:H2'	1:BA:577:G:C8	2.37	0.60
1:BA:2133:G:O2'	1:BA:2158:A:N1	2.35	0.60
31:CA:537:G:H2'	31:CA:538:G:H8	1.66	0.60
31:DA:35:G:O2'	42:DL:118:SER:O	2.19	0.60
31:DA:1228:C:H2'	31:DA:1229:A:C8	2.34	0.60
33:DC:141:VAL:HG11	33:DC:202:ILE:HD13	1.84	0.60
1:AA:531:C:H4'	1:AA:532:A:H5''	1.84	0.60
1:AA:729:G:C5	3:AD:208:LYS:HB2	2.37	0.60
1:AA:729:G:OP1	3:AD:10:THR:OG1	2.12	0.60
2:AB:94:C:H2'	2:AB:95:C:H6	1.67	0.60
1:BA:2137:C:O2	1:BA:2137:C:H2'	2.00	0.60
6:BG:27:ASN:HB3	6:BG:30:GLU:HG3	1.84	0.60
21:BZ:160:GLY:HA2	21:BZ:161:VAL:CB	2.31	0.60
27:B5:16:ARG:HH11	27:B5:16:ARG:HG2	1.66	0.60
31:CA:186:C:H2'	31:CA:187:C:H6	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:685:G:O2'	31:CA:686:U:H5'	2.02	0.60
31:CA:1207:G:H2'	31:CA:1208:C:C6	2.36	0.60
31:CA:1284:C:H3'	31:CA:1285:A:H8	1.67	0.60
31:DA:67:C:H2'	31:DA:68:G:C8	2.36	0.60
31:DA:437:U:H5''	34:DD:155:LEU:HD11	1.83	0.60
31:DA:1367:C:H4'	40:DJ:48:THR:HG21	1.83	0.60
35:DE:36:ASP:OD2	35:DE:37:ARG:N	2.35	0.60
6:AG:66:GLN:HG2	26:A4:1:MET:HE3	1.81	0.60
11:AP:88:LEU:HD11	11:AP:114:ILE:HD12	1.82	0.60
6:BG:20:ILE:O	6:BG:24:GLY:N	2.29	0.60
8:BI:83:ALA:HA	8:BI:89:TYR:HD1	1.67	0.60
31:CA:433:C:H2'	31:CA:434:U:C6	2.35	0.60
35:CE:53:LEU:H	35:CE:53:LEU:HD12	1.66	0.60
31:DA:552:U:O3'	42:DL:87:GLY:HA3	2.01	0.60
31:DA:1176:A:H2'	31:DA:1177:G:C8	2.37	0.60
31:DA:1262:C:H2'	31:DA:1263:C:C6	2.36	0.60
35:DE:100:VAL:O	35:DE:107:ARG:NH2	2.35	0.60
40:DJ:9:ARG:HB2	40:DJ:95:GLU:HB3	1.84	0.60
44:DN:32:SER:HB3	44:DN:41:ARG:HB3	1.83	0.60
1:AA:300:A:P	20:AY:86:ARG:HH22	2.24	0.60
3:AD:142:VAL:HG23	3:AD:193:VAL:HA	1.84	0.60
16:AU:76:TYR:CZ	16:AU:80:ILE:HG13	2.37	0.60
1:BA:910:A:H62	12:BQ:12:GLN:HA	1.65	0.60
34:CD:60:GLU:HG2	34:CD:202:LEU:HB2	1.83	0.60
34:DD:9:CYS:HA	34:DD:12:CYS:H	1.67	0.60
1:AA:656:G:H2'	1:AA:657:U:O4'	2.01	0.59
1:AA:2285:C:OP2	28:A6:6:ARG:NH1	2.34	0.59
2:AB:66:A:H61	2:AB:109:C:H5''	1.66	0.59
15:AT:95:ARG:HG2	15:AT:95:ARG:HH11	1.66	0.59
1:BA:2712:U:H2'	1:BA:2714:G:H5''	1.84	0.59
9:BN:56:ASN:H	9:BN:125:GLY:HA3	1.67	0.59
31:CA:353:A:H5'	31:CA:353:A:H8	1.67	0.59
31:CA:713:G:H2'	31:CA:714:G:C8	2.36	0.59
31:CA:757:U:H2'	31:CA:758:G:O4'	2.01	0.59
31:CA:1324:A:H2'	31:CA:1325:C:C6	2.37	0.59
31:DA:154:C:H2'	31:DA:155:C:H6	1.66	0.59
31:DA:746:A:H2'	31:DA:747:C:H6	1.67	0.59
31:DA:1305:G:N2	31:DA:1331:G:HO2'	1.99	0.59
2:AB:43:C:H5''	26:A4:1:MET:HG3	1.83	0.59
8:AI:123:LEU:HB3	8:AI:144:VAL:HG22	1.83	0.59
1:BA:221:A:N1	1:BA:265:A:O2'	2.34	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:1492:A:OP1	42:CL:47:LYS:N	2.27	0.59
31:DA:1286:A:N6	31:DA:1354:C:O3'	2.35	0.59
31:DA:1346:A:H61	31:DA:1374:A:H3'	1.67	0.59
42:DL:34:ARG:O	42:DL:61:THR:HG23	2.03	0.59
46:DP:28:ARG:NH1	46:DP:29:ASP:OD2	2.35	0.59
1:AA:2615:U:H2'	1:AA:2616:C:H6	1.67	0.59
1:AA:2802:G:H2'	1:AA:2803:C:C6	2.38	0.59
7:AH:11:VAL:HG21	7:AH:50:VAL:HG23	1.83	0.59
14:AS:27:SER:HA	14:AS:88:ASP:HB3	1.85	0.59
20:AY:79:CYS:HB2	20:AY:81:LYS:H	1.68	0.59
1:BA:910:A:C5	12:BQ:13:GLN:HG3	2.37	0.59
1:BA:2683:C:OP1	15:BT:53:ARG:NH2	2.35	0.59
6:BG:126:ASP:HB3	6:BG:130:ASN:H	1.67	0.59
31:CA:116:A:H61	31:CA:313:A:H1'	1.68	0.59
31:CA:171:A:H2'	31:CA:172:A:C8	2.37	0.59
31:CA:503:C:OP1	42:CL:119:LYS:NZ	2.26	0.59
31:CA:1030:C:C4	31:CA:1030(A):G:H1'	2.37	0.59
34:CD:110:PHE:H	34:CD:110:PHE:HD1	1.50	0.59
31:DA:166:G:H2'	31:DA:167:G:C8	2.38	0.59
31:DA:501:C:H2'	31:DA:502:G:H8	1.65	0.59
31:DA:625:G:H2'	31:DA:626:U:H6	1.65	0.59
1:AA:1379:A:H4'	1:AA:1380:G:OP2	2.03	0.59
3:AD:274:ARG:HA	3:AD:275:LYS:HB2	1.85	0.59
1:BA:1047:G:H2'	1:BA:1110:G:H1	1.67	0.59
1:BA:1331:A:HO2'	1:BA:1332:G:H8	1.50	0.59
1:BA:1420:U:HO2'	1:BA:1421:G:P	2.25	0.59
1:BA:2156:G:O6	1:BA:2157:G:N2	2.33	0.59
1:BA:2408:U:H2'	1:BA:2409:G:H8	1.68	0.59
1:BA:2420:C:H5'	28:B6:54:ILE:HD11	1.83	0.59
31:CA:953:G:H5''	31:CA:965:A:H61	1.67	0.59
31:CA:1256:A:H5''	31:CA:1257:U:OP1	2.01	0.59
34:CD:149:ALA:O	34:CD:152:SER:N	2.32	0.59
31:DA:382:A:H2'	31:DA:383:A:H8	1.66	0.59
37:DG:43:PHE:O	37:DG:47:CYS:N	2.30	0.59
37:DG:62:PHE:HA	37:DG:124:LEU:HD22	1.83	0.59
39:DI:28:VAL:HG22	39:DI:63:ILE:HB	1.84	0.59
1:AA:2723:C:H4'	13:AR:1:MET:HG3	1.83	0.59
19:AX:12:VAL:HG22	19:AX:29:TRP:CE2	2.37	0.59
1:BA:784:A:C5	3:BD:229:VAL:HG21	2.38	0.59
1:BA:1178:C:H2'	1:BA:1179:C:C6	2.38	0.59
52:BA:3001:T8B:C10	31:CA:1491:G:C6	2.86	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BF:197:ASP:OD2	5:BF:197:ASP:N	2.35	0.59
21:BZ:44:PHE:CZ	21:BZ:86:VAL:HG11	2.37	0.59
31:CA:105:G:H2'	31:CA:106:C:C6	2.37	0.59
31:DA:1003:G:H21	31:DA:1038:C:N4	2.01	0.59
36:DF:3:ARG:HB3	36:DF:93:SER:HB2	1.85	0.59
43:DM:24:GLY:O	43:DM:25:ILE:HG13	2.03	0.59
1:AA:296:C:H2'	1:AA:297:C:H6	1.68	0.59
1:BA:207:A:H2'	1:BA:208:C:O4'	2.03	0.59
15:BT:16:ARG:HH21	15:BT:83:ILE:HB	1.67	0.59
31:CA:59:A:H5''	31:CA:60:A:H5''	1.84	0.59
31:CA:93:G:H1'	31:CA:96:U:O5'	2.03	0.59
31:CA:417:C:H2'	31:CA:418:C:C6	2.37	0.59
31:CA:1187:G:H2'	31:CA:1188:A:C8	2.37	0.59
31:CA:1226:C:H4'	49:CS:80:TYR:CZ	2.38	0.59
31:DA:560:U:H4'	31:DA:561:U:O5'	2.02	0.59
31:DA:1063:C:H3'	31:DA:1064:G:H2'	1.85	0.59
1:AA:1448:G:H2'	1:AA:1449:A:C8	2.37	0.59
1:AA:1849:G:H2'	1:AA:1850:G:H8	1.67	0.59
31:CA:1279:A:O2'	31:CA:1282:C:N4	2.35	0.59
45:CO:56:LEU:O	45:CO:60:VAL:HG23	2.02	0.59
31:DA:93:G:H1'	31:DA:96:U:O5'	2.02	0.59
31:DA:187:C:H2'	31:DA:188:C:C6	2.38	0.59
31:DA:1360:A:C5	44:DN:18:VAL:HG12	2.38	0.59
1:AA:1155:A:O2'	1:AA:1156:A:H2'	2.03	0.59
19:AX:31:HIS:CD2	19:AX:33:LYS:H	2.20	0.59
12:BQ:1:MET:O	12:BQ:2:LEU:HB2	2.02	0.59
31:CA:142:G:H2'	31:CA:143:A:H8	1.68	0.59
31:CA:1009:G:O6	31:CA:1020:U:O2	2.21	0.59
31:CA:1126:U:H1'	31:CA:1280:A:C4	2.38	0.59
31:CA:1202:G:H2'	31:CA:1203:C:O4'	2.01	0.59
31:CA:1325:C:H2'	31:CA:1326:C:H6	1.68	0.59
31:CA:1330:U:O4	31:CA:1331:G:N1	2.36	0.59
31:CA:1392:G:H21	31:CA:1502:A:H8	1.50	0.59
48:DR:73:ALA:HB1	48:DR:78:LEU:HB2	1.84	0.59
1:AA:2615:U:H2'	1:AA:2616:C:C6	2.38	0.59
1:AA:2646:C:H2'	1:AA:2647:U:O4'	2.03	0.59
30:A8:39:LYS:HA	30:A8:42:ARG:HH12	1.67	0.59
1:BA:2473:U:O2	1:BA:2473:U:H2'	2.03	0.59
15:BT:118:ARG:HG3	15:BT:118:ARG:NH1	2.16	0.59
31:CA:993:G:O6	31:CA:1045:C:N4	2.31	0.59
31:DA:955:U:H2'	31:DA:956:U:H6	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:1318:A:H4'	49:DS:10:PHE:CE2	2.38	0.59
1:AA:2061:G:H5''	1:AA:2503:A:C2	2.37	0.59
4:AE:9:VAL:HG22	4:AE:25:VAL:HB	1.85	0.59
1:BA:857:C:OP2	22:B0:77:ARG:NH2	2.36	0.59
1:BA:1506:C:H2'	1:BA:1507:A:H5'	1.84	0.59
1:BA:2701:C:H2'	1:BA:2702:U:H2'	1.85	0.59
31:CA:308:C:H2'	31:CA:309:G:C8	2.37	0.59
31:CA:590:C:H2'	31:CA:591:U:H6	1.68	0.59
31:CA:1155:G:H2'	31:CA:1156:G:H8	1.66	0.59
31:CA:1276:G:H1'	31:CA:1282:C:O2'	2.03	0.59
38:CH:121:ASP:N	38:CH:121:ASP:OD1	2.34	0.59
31:DA:397:A:H3'	31:DA:397:A:N3	2.18	0.59
31:DA:971:G:P	31:DA:1231:G:H21	2.25	0.59
31:DA:1250:A:C2	31:DA:1370:G:H1'	2.38	0.59
35:DE:110:LEU:HD13	35:DE:118:ILE:HG21	1.85	0.59
1:AA:500:G:N1	1:AA:503:A:OP2	2.36	0.58
1:AA:2023:G:H5'	1:AA:2617:C:H4'	1.85	0.58
1:AA:2133:G:H1'	1:AA:2158:A:H61	1.68	0.58
7:AH:33:LEU:HD21	7:AH:136:ILE:HG13	1.85	0.58
22:A0:10:THR:HG22	22:A0:12:ASN:N	2.15	0.58
1:BA:1364:G:OP1	23:B1:2:SER:HA	2.03	0.58
8:BI:40:THR:O	8:BI:44:LEU:HB2	2.03	0.58
8:BI:77:LEU:HB3	8:BI:142:VAL:HG12	1.85	0.58
14:BS:52:SER:HB2	14:BS:55:ALA:H	1.68	0.58
14:BS:82:ILE:HA	14:BS:83:LYS:CB	2.33	0.58
45:CO:39:LEU:HD13	45:CO:56:LEU:HB2	1.83	0.58
1:AA:652(B):A:H61	1:AA:655:A:H2	1.51	0.58
1:AA:1671:U:HO2'	1:AA:1673:U:H5	1.49	0.58
12:AQ:8:LYS:HA	21:AZ:196:VAL:HB	1.84	0.58
14:AS:96:GLY:N	14:AS:99:LYS:H	2.01	0.58
19:AX:2:LYS:NZ	19:AX:38:GLU:OE2	2.31	0.58
1:BA:776:G:H4'	1:BA:777:A:O5'	2.03	0.58
1:BA:1239:G:H2'	1:BA:1240:U:O4'	2.04	0.58
2:BB:28:C:OP1	14:BS:36:TYR:OH	2.13	0.58
31:CA:1490:C:O2'	31:CA:1491:G:H5'	2.03	0.58
32:CB:17:PHE:H	32:CB:17:PHE:HD2	1.50	0.58
38:CH:86:ILE:HG13	38:CH:133:LEU:HD22	1.85	0.58
31:DA:971:G:OP2	31:DA:1231:G:N2	2.35	0.58
31:DA:1399:C:C2	31:DA:1502:A:N6	2.71	0.58
31:DA:1492:A:H4'	31:DA:1493:A:OP1	2.02	0.58
42:DL:84:LEU:HD23	42:DL:105:TYR:HE1	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:322:A:OP1	5:AF:168:ARG:NH1	2.35	0.58
10:AO:43:VAL:HG23	10:AO:56:ASP:O	2.03	0.58
20:AY:99:CYS:CB	20:AY:102:CYS:SG	2.91	0.58
21:AZ:119:GLU:OE2	21:AZ:122:ARG:NH1	2.36	0.58
1:BA:1339:G:H5''	19:BX:16:LYS:HD2	1.84	0.58
1:BA:1916:A:N6	52:BA:3001:T8B:O1	2.35	0.58
31:CA:583:A:N6	31:CA:758:G:O2'	2.36	0.58
31:CA:677:U:H2'	31:CA:678:U:H6	1.67	0.58
31:CA:1493:A:N3	31:CA:1493:A:H2'	2.17	0.58
45:CO:18:PHE:HB2	45:CO:19:PRO:HD2	1.85	0.58
50:CT:77:ALA:O	50:CT:81:LYS:HG3	2.04	0.58
31:DA:1113:C:H42	31:DA:1187:G:H1	1.48	0.58
31:DA:1250:A:H2	31:DA:1370:G:H1'	1.67	0.58
35:DE:102:ALA:HB1	35:DE:106:PRO:HG2	1.84	0.58
37:DG:46:ALA:HB1	37:DG:121:ALA:HB2	1.83	0.58
1:AA:1882:C:H5'	1:AA:1883:G:OP2	2.04	0.58
1:AA:2125:G:N2	1:AA:2172:U:H3'	2.17	0.58
1:AA:2471:C:N4	1:AA:2476:A:O2'	2.36	0.58
2:AB:55:U:H2'	2:AB:56:G:O4'	2.04	0.58
1:BA:213:A:H2'	1:BA:214:G:O4'	2.03	0.58
1:BA:1153:C:OP1	16:BU:92:ARG:NH1	2.36	0.58
10:BO:80:ASP:OD1	15:BT:64:ARG:NH2	2.36	0.58
31:CA:878:G:H5'	38:CH:89:PRO:HG2	1.85	0.58
31:CA:1109:C:H2'	31:CA:1110:A:O4'	2.04	0.58
31:CA:1273:G:H3'	31:CA:1274:G:C8	2.38	0.58
35:CE:36:ASP:OD2	35:CE:38:GLN:N	2.37	0.58
31:DA:20:U:H2'	31:DA:21:G:O4'	2.02	0.58
31:DA:1071:C:H2'	31:DA:1072:G:H8	1.67	0.58
1:AA:690:G:N2	1:AA:773:U:C2	2.72	0.58
1:BA:2299:G:C2	1:BA:2318:G:H8	2.21	0.58
33:CC:182:ILE:HA	33:CC:202:ILE:O	2.03	0.58
31:DA:102:G:H2'	31:DA:103:C:H6	1.68	0.58
31:DA:235:C:H2'	31:DA:236:G:H8	1.69	0.58
31:DA:1305:G:H1	31:DA:1331:G:HO2'	1.48	0.58
46:DP:55:ARG:O	46:DP:58:TYR:N	2.36	0.58
1:AA:1639:U:H2'	1:AA:1640:C:H5''	1.85	0.58
1:AA:2262:U:P	22:A0:19:LYS:HE2	2.44	0.58
20:AY:99:CYS:HB3	20:AY:102:CYS:SG	2.42	0.58
28:A6:9:LEU:HD21	28:A6:25:LYS:HB3	1.85	0.58
1:BA:754:C:H2'	1:BA:755:C:C6	2.38	0.58
1:BA:1372:U:H6	1:BA:1372:U:O5'	1.87	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:1882:C:H5'	1:BA:1883:G:OP2	2.04	0.58
31:CA:438:G:O2'	31:CA:494:U:O4	2.18	0.58
36:CF:22:GLU:OE2	36:CF:82:ARG:HG2	2.04	0.58
37:CG:42:ILE:HG22	37:CG:120:ILE:HG13	1.86	0.58
45:CO:70:LEU:HD11	45:CO:77:ARG:HB3	1.84	0.58
31:DA:828:A:H2'	31:DA:829:G:O4'	2.04	0.58
31:DA:1345:U:O2	31:DA:1375:A:N6	2.37	0.58
40:DJ:8:LEU:N	40:DJ:70:ARG:O	2.36	0.58
1:AA:272(B):G:H2'	1:AA:272(C):G:C8	2.39	0.58
1:AA:2887:U:H2'	1:AA:2888:C:H6	1.69	0.58
4:AE:101:ARG:NH2	4:AE:171:GLU:HB2	2.19	0.58
1:BA:90:U:HO2'	1:BA:92:A:H8	1.44	0.58
1:BA:185:U:H4'	1:BA:218:A:H4'	1.86	0.58
1:BA:831:G:O2'	11:BP:38:GLN:HG2	2.04	0.58
1:BA:1530:C:H42	1:BA:1539:G:H1	1.49	0.58
1:BA:1899:G:H2'	1:BA:1899:G:N3	2.18	0.58
1:BA:2357:U:OP1	22:B0:20:ARG:NH1	2.31	0.58
18:BW:86:LEU:HD22	18:BW:96:ILE:HD11	1.86	0.58
20:BY:86:ARG:NH1	20:BY:100:ALA:HB1	2.18	0.58
31:CA:630:G:H2'	31:CA:631:G:H8	1.67	0.58
31:DA:458:C:H2'	31:DA:460:G:C8	2.38	0.58
31:DA:1147:C:O2'	39:DI:5:TYR:OH	2.14	0.58
47:DQ:45:HIS:HB2	47:DQ:65:ILE:HD13	1.85	0.58
49:DS:45:VAL:HA	49:DS:62:ILE:HG22	1.86	0.58
1:AA:1405:U:H2'	1:AA:1406:U:C6	2.38	0.58
1:AA:1537:G:H2'	1:AA:1538:G:C8	2.38	0.58
1:AA:2393:A:H5'	11:AP:63:PRO:HB3	1.84	0.58
1:BA:443:A:H5''	1:BA:444:C:OP1	2.03	0.58
1:BA:675:A:OP1	5:BF:63:LYS:NZ	2.29	0.58
4:BE:71:GLY:HA2	4:BE:72:VAL:O	2.04	0.58
6:BG:114:ILE:HB	6:BG:117:PHE:HB2	1.84	0.58
50:CT:87:LYS:O	50:CT:91:LEU:HG	2.04	0.58
31:DA:198:G:O6	31:DA:219:C:N4	2.37	0.58
31:DA:359:U:H2'	31:DA:360:A:C8	2.39	0.58
31:DA:1217:C:H2'	31:DA:1218:C:O4'	2.03	0.58
32:DB:100:GLY:N	32:DB:176:GLU:OE2	2.31	0.58
38:DH:86:ILE:HG13	38:DH:133:LEU:HD22	1.85	0.58
1:AA:271(E):U:H2'	1:AA:271(F):C:C6	2.39	0.58
1:AA:1913:A:O2'	52:AA:3001:T8B:C21	2.52	0.58
7:AH:12:PRO:O	7:AH:15:VAL:HG13	2.03	0.58
14:AS:82:ILE:HA	14:AS:83:LYS:CB	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:1047:G:O2'	1:BA:1048:A:O5'	2.21	0.58
1:BA:1688:U:H1'	1:BA:1701:A:C6	2.39	0.58
1:BA:2139:C:H42	1:BA:2152:G:H1	1.51	0.58
4:BE:72:VAL:HA	4:BE:73:GLU:HB3	1.86	0.58
4:BE:110:GLY:HA2	4:BE:161:GLY:HA3	1.85	0.58
20:BY:99:CYS:CB	20:BY:102:CYS:SG	2.92	0.58
23:B1:85:LEU:HB3	23:B1:89:GLU:HG3	1.85	0.58
31:CA:359:U:H2'	31:CA:360:A:C8	2.38	0.58
31:CA:521:G:O6	31:CA:529:G:C2	2.57	0.58
43:CM:5:ALA:HA	43:CM:61:GLU:HG2	1.85	0.58
31:DA:272:C:H2'	31:DA:273:A:H8	1.68	0.58
31:DA:373:A:H61	31:DA:391:G:H1'	1.69	0.58
43:DM:69:GLU:O	43:DM:71:ARG:N	2.32	0.58
1:AA:1140:C:O3'	9:AN:25:ARG:NH1	2.37	0.58
1:AA:2783:G:H2'	1:AA:2784:C:C6	2.39	0.58
52:AA:3001:T8B:H13	52:AA:3001:T8B:H17	0.73	0.58
2:AB:21:G:H2'	2:AB:22:U:O4'	2.03	0.58
11:AP:80:TYR:HE1	11:AP:128:HIS:HB3	1.69	0.58
11:AP:83:VAL:HG13	11:AP:112:LEU:HD21	1.84	0.58
14:AS:67:ARG:HG2	14:AS:71:ARG:NH2	2.18	0.58
27:A5:47:PRO:O	27:A5:60:VAL:HG21	2.03	0.58
1:BA:2108:C:N4	1:BA:2181:G:H1	2.01	0.58
1:BA:2165:G:H2'	1:BA:2166:G:C8	2.39	0.58
1:BA:2272:U:H5''	1:BA:2273:A:OP1	2.03	0.58
1:BA:2292:C:OP1	14:BS:17:ARG:NH2	2.33	0.58
1:BA:2328:A:H2'	1:BA:2329:G:C8	2.39	0.58
31:CA:501:C:H1'	31:CA:549:C:H1'	1.85	0.58
31:CA:994:A:N1	31:CA:1047:G:H4'	2.19	0.58
39:CI:34:ASN:N	39:CI:34:ASN:OD1	2.36	0.58
31:DA:232:G:H1'	31:DA:262:A:N1	2.19	0.58
31:DA:1285:A:H1'	31:DA:1286:A:OP2	2.04	0.58
1:AA:1341:U:OP2	1:AA:1394:U:O2'	2.19	0.57
2:AB:53:A:H2'	2:AB:54:G:O4'	2.04	0.57
2:AB:105:A:OP1	21:AZ:72:ARG:NH1	2.37	0.57
8:AI:87:LYS:HG3	8:AI:121:LYS:O	2.04	0.57
18:AW:75:TYR:CZ	18:AW:104:THR:HG21	2.39	0.57
1:BA:141:A:H8	1:BA:1408:C:O2'	1.87	0.57
1:BA:728:G:H4'	3:BD:13:ARG:HD3	1.85	0.57
1:BA:1761:C:H2'	1:BA:1762:A:H5''	1.86	0.57
5:BF:22:ALA:HB1	5:BF:24:LEU:HD22	1.85	0.57
31:CA:375:U:C2	31:CA:376:G:C8	2.92	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:633:G:H2'	31:CA:634:C:C6	2.38	0.57
31:CA:696:A:N1	31:CA:797:C:O2'	2.29	0.57
33:CC:74:GLY:C	33:CC:76:VAL:H	2.08	0.57
31:DA:590:C:H2'	31:DA:591:U:C6	2.39	0.57
31:DA:1218:C:H2'	31:DA:1219:U:C6	2.39	0.57
1:AA:330:A:HO2'	1:AA:331:A:H8	1.52	0.57
1:AA:2036:C:H6	1:AA:2036:C:H5'	1.68	0.57
6:AG:105:LYS:NZ	26:A4:25:TYR:O	2.31	0.57
1:BA:2319:G:H22	14:BS:3:ARG:NE	2.02	0.57
1:BA:2405:G:O2'	1:BA:2411:A:N6	2.37	0.57
1:BA:2470:G:O6	1:BA:2476:A:O2'	2.18	0.57
13:BR:33:ARG:NH2	27:B5:57:VAL:O	2.34	0.57
31:CA:194:C:O3'	50:CT:68:LYS:HD2	2.04	0.57
31:CA:1005:A:N3	31:CA:1036:G:N2	2.50	0.57
32:CB:158:LEU:HG	32:CB:182:ILE:HD11	1.85	0.57
35:CE:78:HIS:CE1	35:CE:142:LEU:HD23	2.40	0.57
40:CJ:61:GLU:OE2	44:CN:45:ARG:NE	2.37	0.57
34:DD:121:VAL:HG22	34:DD:126:ILE:HG13	1.86	0.57
38:DH:51:VAL:HG21	38:DH:60:ARG:HG3	1.86	0.57
1:AA:1221(A):C:C2	1:AA:1229:G:C2	2.93	0.57
6:AG:50:ALA:C	6:AG:52:ILE:H	2.07	0.57
17:AV:40:LEU:HB2	17:AV:46:VAL:HG13	1.85	0.57
1:BA:141:A:C8	1:BA:1408:C:O2'	2.57	0.57
31:CA:691:G:H2'	31:CA:692:U:C6	2.39	0.57
31:CA:859:A:H2'	31:CA:860:A:O4'	2.04	0.57
31:CA:955:U:H1'	31:CA:1227:A:N6	2.19	0.57
31:DA:222:U:H2'	31:DA:223:U:C6	2.39	0.57
34:DD:26:CYS:HA	34:DD:31:CYS:HB2	1.86	0.57
40:DJ:8:LEU:HD12	40:DJ:20:ALA:HB2	1.86	0.57
1:AA:71:A:C2	19:AX:31:HIS:CE1	2.88	0.57
1:AA:459:U:H5''	29:A7:40:TRP:CD2	2.40	0.57
1:AA:1858:G:O2'	1:AA:1884:A:N6	2.37	0.57
6:AG:13:GLU:O	6:AG:14:GLU:HB2	2.03	0.57
7:AH:113:VAL:HG11	7:AH:151:ILE:HD13	1.85	0.57
13:AR:2:ARG:NH1	13:AR:5:LYS:O	2.37	0.57
23:A1:3:LYS:HB2	23:A1:61:ARG:NH1	2.14	0.57
1:BA:729:G:C5	3:BD:208:LYS:HB2	2.39	0.57
31:DA:1228:C:N4	43:DM:104:ARG:O	2.38	0.57
42:DL:24:VAL:HG13	42:DL:98:TYR:HE2	1.70	0.57
1:AA:263:C:H2'	1:AA:264:C:O4'	2.04	0.57
1:AA:2067:G:O2'	1:AA:2069:G:H5'	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2152:G:H2'	1:AA:2153:G:H8	1.70	0.57
6:AG:122:PRO:HG3	6:AG:180:PHE:HB3	1.86	0.57
28:A6:10:LEU:HD23	28:A6:22:ALA:HB2	1.87	0.57
1:BA:621:A:OP2	11:BP:108:LYS:NZ	2.37	0.57
1:BA:2127:G:N2	1:BA:2162:G:H1'	2.19	0.57
1:BA:2128:C:H42	1:BA:2160:G:H1	1.52	0.57
1:BA:2218:U:H4'	1:BA:2219:G:OP2	2.04	0.57
18:BW:40:ASN:O	18:BW:41:LYS:HG3	2.05	0.57
29:B7:47:ARG:HH11	29:B7:47:ARG:HG3	1.70	0.57
31:CA:969:A:OP1	40:CJ:55:LYS:NZ	2.37	0.57
32:CB:163:PHE:HA	32:CB:185:ILE:HG12	1.85	0.57
31:DA:186:C:H2'	31:DA:187:C:H6	1.69	0.57
32:DB:73:THR:HB	32:DB:95:GLN:O	2.05	0.57
34:DD:10:ARG:HG3	34:DD:11:LEU:HD23	1.86	0.57
35:DE:94:ALA:HB1	35:DE:98:THR:OG1	2.04	0.57
1:AA:1798:U:H5'	3:AD:259:THR:HG22	1.85	0.57
1:AA:2558:C:H2'	1:AA:2559:C:O4'	2.05	0.57
1:AA:2723:C:OP2	4:AE:109:LYS:NZ	2.35	0.57
9:AN:24:GLY:HA2	9:AN:27:ALA:CB	2.33	0.57
10:AO:102:VAL:HB	10:AO:106:LEU:HD12	1.87	0.57
15:AT:16:ARG:NH2	15:AT:83:ILE:O	2.38	0.57
21:AZ:24:LEU:HB2	21:AZ:41:LEU:HD23	1.87	0.57
1:BA:102:G:HO2'	1:BA:103:A:P	2.21	0.57
1:BA:652(Q):G:C2'	1:BA:652(R):C:H5'	2.35	0.57
1:BA:846:C:H4'	1:BA:847:U:O5'	2.04	0.57
1:BA:1021:A:H3'	1:BA:1021:A:H8	1.70	0.57
1:BA:2203:U:O2'	1:BA:2205:C:H5'	2.03	0.57
1:BA:2791:C:H2'	1:BA:2792:G:C8	2.39	0.57
19:BX:41:ASN:O	19:BX:45:THR:HG23	2.04	0.57
21:BZ:158:PRO:O	21:BZ:161:VAL:HG11	2.05	0.57
31:CA:719:C:O2'	48:CR:49:LYS:HB3	2.05	0.57
31:CA:757:U:O2'	31:CA:879:C:O2	2.21	0.57
31:CA:977:A:H2'	31:CA:978:A:H5''	1.86	0.57
31:CA:1131:G:H2'	31:CA:1132:C:C6	2.39	0.57
31:DA:6:G:C4	35:DE:119:LEU:HD11	2.39	0.57
31:DA:518:C:O2'	31:DA:530:G:N2	2.37	0.57
31:DA:1141:C:H2'	31:DA:1142:G:O4'	2.04	0.57
33:DC:7:PRO:O	33:DC:11:ARG:NH1	2.36	0.57
8:AI:27:ARG:HD2	23:A1:71:TYR:CE1	2.39	0.57
15:AT:16:ARG:HH21	15:AT:83:ILE:HB	1.70	0.57
18:AW:71:VAL:HA	18:AW:107:LEU:HD12	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:71:A:C2	19:BX:31:HIS:CE1	2.90	0.57
1:BA:652(B):A:O2'	1:BA:652(C):G:H5'	2.03	0.57
1:BA:1540:U:H2'	1:BA:1541:G:O4'	2.04	0.57
4:BE:116:VAL:HG13	4:BE:122:PHE:HB2	1.85	0.57
31:CA:41:G:H2'	31:CA:42:G:C8	2.40	0.57
31:CA:62:U:H2'	31:CA:63:C:C6	2.38	0.57
31:CA:448:A:OP2	31:CA:485:G:N2	2.21	0.57
31:CA:1036:G:H3'	31:CA:1037:C:C6	2.40	0.57
47:DQ:67:LYS:HA	47:DQ:70:ARG:HH12	1.69	0.57
1:AA:1488:G:H5'	1:AA:1489:U:OP2	2.04	0.57
1:AA:1518:U:H2'	1:AA:1519:G:O4'	2.03	0.57
1:AA:1889:A:H2'	1:AA:1890:A:C8	2.39	0.57
1:AA:2356:C:O3'	22:A0:20:ARG:HD3	2.04	0.57
1:AA:2849:U:OP2	15:AT:95:ARG:NH1	2.38	0.57
1:BA:272(G):C:N4	1:BA:272(H):C:H41	2.03	0.57
1:BA:652(D):C:H2'	1:BA:652(E):G:O4'	2.05	0.57
4:BE:174:ASP:OD2	4:BE:175:VAL:N	2.38	0.57
31:CA:954:G:N2	31:CA:1227:A:H62	2.03	0.57
37:CG:111:ARG:HB3	37:CG:113:GLU:HG2	1.87	0.57
44:CN:24:CYS:SG	44:CN:39:LEU:HA	2.45	0.57
31:DA:522:C:H41	42:DL:53:ARG:HH22	1.53	0.57
34:DD:158:ILE:O	34:DD:162:LEU:N	2.33	0.57
4:AE:112:GLY:O	4:AE:159:HIS:HA	2.04	0.57
1:BA:1815:A:P	3:BD:54:ARG:HH22	2.28	0.57
1:BA:2803:C:H2'	1:BA:2804:C:C6	2.39	0.57
31:DA:148:G:O2'	31:DA:149:A:H5'	2.04	0.57
31:DA:333:G:H4'	50:DT:16:HIS:CE1	2.40	0.57
31:DA:560:U:H5'	31:DA:566:G:N2	2.20	0.57
31:DA:590:C:H2'	31:DA:591:U:H6	1.70	0.57
31:DA:1010:G:N2	31:DA:1020:U:H1'	2.20	0.57
50:DT:64:ASP:OD1	50:DT:81:LYS:NZ	2.30	0.57
1:AA:29:U:H2'	1:AA:30:G:C8	2.40	0.57
1:AA:40:C:H2'	1:AA:41:C:C6	2.40	0.57
1:AA:1593:G:H2'	1:AA:1594:G:H8	1.66	0.57
1:AA:2151:G:H2'	1:AA:2152:G:O4'	2.05	0.57
1:AA:2299:G:N1	1:AA:2318:G:C8	2.73	0.57
2:AB:37:C:N4	2:AB:38:C:N3	2.51	0.57
6:AG:27:ASN:HB3	6:AG:30:GLU:HG3	1.86	0.57
24:A2:50:ILE:O	24:A2:51:ARG:HB3	2.04	0.57
29:A7:5:TRP:O	29:A7:6:GLN:NE2	2.38	0.57
30:A8:34:TRP:CE2	30:A8:35:GLN:HB3	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:2125:G:O2'	1:BA:2173:A:N6	2.38	0.57
3:BD:108:PRO:HB3	3:BD:143:HIS:CE1	2.39	0.57
6:BG:16:ARG:NE	6:BG:31:VAL:HG21	2.19	0.57
14:BS:27:SER:HA	14:BS:88:ASP:HB3	1.86	0.57
31:CA:975:A:N6	31:CA:1367:C:O4'	2.38	0.57
40:CJ:45:ARG:NH2	44:CN:36:PHE:HB2	2.20	0.57
49:CS:20:LEU:HG	49:CS:43:GLU:HG2	1.86	0.57
31:DA:93:G:O2'	31:DA:96:U:OP2	2.21	0.57
31:DA:1333:A:H3'	31:DA:1334:G:H8	1.70	0.57
43:DM:96:LEU:HD13	43:DM:97:PRO:HD2	1.86	0.57
1:AA:481:G:H4'	1:AA:482:A:H5'	1.87	0.56
1:AA:652(E):G:O6	1:AA:652(S):C:N4	2.38	0.56
1:AA:1540:U:H2'	1:AA:1541:G:O4'	2.05	0.56
1:AA:2114:A:H3'	1:AA:2115:G:C8	2.40	0.56
1:AA:2845:G:H2'	1:AA:2846:G:C8	2.40	0.56
1:BA:2144:U:O2	1:BA:2148:G:N1	2.38	0.56
1:BA:2463:C:C2'	1:BA:2464:C:H5'	2.35	0.56
13:BR:38:VAL:HG22	13:BR:112:ALA:HB2	1.87	0.56
18:BW:32:ALA:HB1	18:BW:51:LEU:HD11	1.87	0.56
32:CB:87:ARG:NE	32:CB:233:SER:HB2	2.19	0.56
32:CB:87:ARG:NH2	32:CB:233:SER:HB2	2.19	0.56
34:CD:12:CYS:O	34:CD:33:MET:HG2	2.05	0.56
31:DA:9:G:H2'	31:DA:10:A:C8	2.40	0.56
31:DA:568:G:N7	42:DL:5:PRO:HD3	2.19	0.56
31:DA:1183:A:H5''	31:DA:1184:G:OP2	2.05	0.56
42:DL:89:ARG:HA	42:DL:97:ARG:HA	1.87	0.56
46:DP:72:ARG:HE	46:DP:73:LEU:HD23	1.70	0.56
1:AA:1000:A:OP2	1:AA:1154:G:N1	2.20	0.56
1:AA:1915:U:O2	52:AA:3001:T8B:C28	2.54	0.56
1:AA:2584:U:H2'	1:AA:2585:U:H2'	1.87	0.56
1:AA:2590:A:H5''	3:AD:239:ARG:HE	1.69	0.56
3:AD:35:LYS:HG2	3:AD:36:PRO:HD2	1.86	0.56
3:AD:71:ASP:HB3	3:AD:103:ARG:NH2	2.18	0.56
1:BA:470:A:OP1	5:BF:59:TYR:HE2	1.88	0.56
1:BA:1021:A:H62	1:BA:1141:U:H3	1.51	0.56
1:BA:1210:A:H5'	1:BA:1210:A:C8	2.40	0.56
1:BA:1292:U:H2'	1:BA:1293:C:C6	2.40	0.56
1:BA:2820:A:O2'	1:BA:2821:A:OP1	2.23	0.56
10:BO:16:ALA:HB2	10:BO:52:VAL:HG21	1.87	0.56
31:CA:114:U:H2'	31:CA:115:G:C8	2.40	0.56
31:CA:835:U:OP1	48:CR:64:ARG:NH1	2.21	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:1008:C:H2'	31:DA:1009:G:O4'	2.05	0.56
31:DA:1016:A:H2'	31:DA:1017:G:O4'	2.05	0.56
1:AA:90:U:O2'	1:AA:92:A:O4'	2.23	0.56
1:AA:857:C:H4'	22:A0:23:VAL:HG21	1.87	0.56
1:AA:1175:U:O2'	1:AA:1176:G:H5'	2.04	0.56
1:AA:2134:A:H1'	1:AA:2159:G:H1'	1.87	0.56
1:AA:2305:A:C2	6:AG:152:LEU:HD11	2.40	0.56
1:AA:2494:G:C4	1:AA:2495:G:C8	2.93	0.56
1:AA:2751:G:H4'	7:AH:4:ILE:HD11	1.86	0.56
2:AB:79:C:H2'	2:AB:80:U:O4'	2.05	0.56
6:AG:103:LEU:O	6:AG:106:LEU:HB3	2.04	0.56
13:AR:20:LEU:HD21	13:AR:40:LYS:HD3	1.88	0.56
14:AS:11:LYS:HG3	14:AS:91:PRO:HD3	1.86	0.56
17:AV:76:LYS:HB2	17:AV:81:TYR:HB3	1.87	0.56
1:BA:330:A:H2	1:BA:1210:A:H2'	1.70	0.56
1:BA:384:U:H2'	1:BA:385:C:C6	2.40	0.56
1:BA:570:G:H2'	1:BA:2030:A:C5	2.40	0.56
1:BA:2096:U:H3	1:BA:2193:G:H1	1.53	0.56
9:BN:47:ALA:HB2	9:BN:112:LEU:HD11	1.87	0.56
31:CA:240:C:H2'	31:CA:241:C:H6	1.71	0.56
31:CA:519:C:H2'	31:CA:520:A:C8	2.41	0.56
31:CA:1309:G:H2'	31:CA:1310:G:O4'	2.05	0.56
31:CA:1504:G:P	31:CA:1504:G:H3'	2.45	0.56
32:CB:15:VAL:HG23	32:CB:209:ARG:HG2	1.88	0.56
33:CC:127:ARG:HH11	33:CC:127:ARG:HB3	1.70	0.56
34:CD:24:GLU:O	34:CD:27:TYR:HB2	2.05	0.56
36:CF:3:ARG:HB3	36:CF:93:SER:HB2	1.86	0.56
42:CL:83:VAL:HG13	42:CL:100:ILE:HG23	1.88	0.56
50:CT:12:ALA:O	50:CT:15:ARG:HB2	2.05	0.56
31:DA:475:G:O2'	31:DA:476:G:H5'	2.05	0.56
31:DA:628:G:H2'	31:DA:629:G:C8	2.41	0.56
31:DA:746:A:H2'	31:DA:747:C:C6	2.39	0.56
31:DA:826:C:H2'	31:DA:827:U:C6	2.40	0.56
31:DA:863:U:H2'	31:DA:865:A:OP2	2.06	0.56
31:DA:1001(A):G:H2'	31:DA:1002:G:O4'	2.06	0.56
31:DA:1111:A:H2'	31:DA:1112:C:H6	1.70	0.56
32:DB:55:PHE:HA	32:DB:58:ILE:HD12	1.87	0.56
1:AA:1365:A:H5''	23:A1:41:ARG:HH22	1.70	0.56
1:AA:1505:C:H2'	1:AA:1506:C:H6	1.70	0.56
14:AS:96:GLY:HA2	14:AS:97:ARG:C	2.26	0.56
25:A3:4:LEU:HD22	25:A3:56:VAL:HG11	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:65:C:H2'	1:BA:66:C:H6	1.70	0.56
1:BA:511:U:H5	1:BA:512:G:C5	2.23	0.56
23:B1:23:LYS:HB3	23:B1:29:GLY:HA3	1.88	0.56
31:CA:20:U:H2'	31:CA:21:G:O4'	2.05	0.56
31:DA:272:C:H2'	31:DA:273:A:C8	2.41	0.56
31:DA:512:U:H2'	31:DA:513:C:H6	1.69	0.56
31:DA:664:G:N2	31:DA:741:G:H1	2.04	0.56
31:DA:975:A:H5'	31:DA:975:A:H8	1.70	0.56
31:DA:1131:G:H2'	31:DA:1132:C:C6	2.41	0.56
1:AA:2316:C:H2'	1:AA:2317:C:H6	1.69	0.56
1:AA:2740:A:H2'	1:AA:2741:A:C8	2.41	0.56
24:A2:1:MET:N	24:A2:52:ASP:OD1	2.31	0.56
1:BA:2053:G:H5'	4:BE:144:ARG:O	2.05	0.56
1:BA:2147:G:H2'	1:BA:2148:G:O4'	2.05	0.56
1:BA:2349:G:H3'	1:BA:2350:C:H5''	1.88	0.56
1:BA:2355:C:H1'	22:B0:39:ARG:HH21	1.70	0.56
2:BB:48:A:H4'	14:BS:95:HIS:HD2	1.68	0.56
1:AA:203:C:H3'	1:AA:204:A:H5''	1.87	0.56
1:AA:861:A:N6	1:AA:916:G:O2'	2.37	0.56
1:AA:1915:U:C2	52:AA:3001:T8B:C27	2.88	0.56
12:AQ:62:GLY:HA2	21:AZ:116:VAL:HG21	1.87	0.56
1:BA:2803:C:H2'	1:BA:2804:C:H6	1.70	0.56
1:BA:2852:G:H2'	1:BA:2853:C:O4'	2.05	0.56
2:BB:31:C:H4'	6:BG:29:TRP:CH2	2.40	0.56
5:BF:24:LEU:HD21	5:BF:199:TRP:HH2	1.70	0.56
27:B5:46:CYS:CB	27:B5:49:CYS:HG	2.19	0.56
31:CA:715:A:H2'	31:CA:716:A:C8	2.40	0.56
37:CG:18:TYR:HD2	37:CG:59:LEU:HD12	1.71	0.56
31:DA:1412:C:H2'	31:DA:1413:A:C8	2.40	0.56
33:DC:175:LEU:HD12	33:DC:175:LEU:H	1.70	0.56
1:AA:1580:A:OP2	1:AA:1580:A:H8	1.88	0.56
1:AA:1638:C:H4'	1:AA:2710:C:O2	2.06	0.56
1:BA:455:C:N3	1:BA:473:G:H5'	2.20	0.56
1:BA:1453:U:OP1	13:BR:77:ARG:NH1	2.37	0.56
1:BA:1790:C:H5''	1:BA:1791:A:OP1	2.05	0.56
1:BA:2012:G:OP1	18:BW:11:ARG:NH2	2.39	0.56
1:BA:2036:C:H5'	1:BA:2036:C:H6	1.70	0.56
1:BA:2787:C:H1'	4:BE:62:PRO:HG3	1.87	0.56
31:CA:96:U:O2'	31:CA:97:G:OP2	2.19	0.56
31:CA:394:G:H2'	31:CA:395:C:H6	1.70	0.56
31:CA:434:U:H2'	31:CA:435:C:C6	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:458:C:H2'	31:CA:460:G:C8	2.41	0.56
34:DD:19:LEU:HD23	34:DD:21:LEU:HD11	1.88	0.56
35:DE:122:GLU:HG2	35:DE:131:ILE:HD12	1.87	0.56
49:DS:18:LYS:O	49:DS:22:LEU:N	2.37	0.56
1:AA:7:G:H2'	1:AA:8:A:C8	2.41	0.56
1:BA:458:G:C8	29:B7:37:LYS:HG2	2.40	0.56
1:BA:1175:U:O2'	1:BA:1176:G:H5'	2.06	0.56
1:BA:1971:A:C4	3:BD:241:PRO:HD3	2.40	0.56
1:BA:2106:G:N2	1:BA:2184:G:C4	2.74	0.56
9:BN:24:GLY:HA2	9:BN:27:ALA:HB3	1.88	0.56
31:CA:189(B):C:H2'	31:CA:189(C):C:C6	2.41	0.56
31:CA:458:C:H2'	31:CA:460:G:H8	1.71	0.56
35:CE:39:GLY:O	35:CE:69:VAL:HG13	2.06	0.56
31:DA:430:A:OP2	34:DD:8:VAL:HG12	2.06	0.56
31:DA:675:A:H2'	31:DA:676:A:H8	1.69	0.56
31:DA:792:A:H4'	31:DA:793:U:O5'	2.06	0.56
31:DA:1298:C:OP2	37:DG:114:ARG:NH2	2.39	0.56
31:DA:1348:U:H4'	39:DI:120:ARG:HD2	1.88	0.56
1:AA:652(Q):G:C2'	1:AA:652(R):C:H5'	2.36	0.56
1:AA:807:U:OP2	11:AP:41:ARG:NH2	2.38	0.56
1:AA:1339:G:H5''	19:AX:16:LYS:HD2	1.88	0.56
1:AA:1494:A:H2'	1:AA:1495:A:C8	2.41	0.56
1:AA:2108:C:N4	1:AA:2181:G:H1	2.04	0.56
1:AA:2784:C:H1'	4:AE:37:ARG:HH12	1.71	0.56
3:AD:145:VAL:HG11	3:AD:175:LEU:HD11	1.86	0.56
5:AF:197:ASP:OD2	5:AF:197:ASP:N	2.28	0.56
11:AP:59:LEU:HD11	30:A8:10:ALA:HB2	1.88	0.56
13:AR:24:GLN:HB3	13:AR:44:LEU:HD11	1.88	0.56
1:BA:286:C:H2'	1:BA:287:C:C6	2.41	0.56
1:BA:652(E):G:O6	1:BA:652(S):C:N4	2.39	0.56
4:BE:36:ARG:NH1	4:BE:85:ASN:OD1	2.39	0.56
31:CA:185:A:H2'	31:CA:186:C:C6	2.40	0.56
31:CA:1263:C:H2'	31:CA:1264:C:C6	2.41	0.56
31:CA:1346:A:H61	31:CA:1374:A:H3'	1.71	0.56
31:CA:1346:A:H5''	39:CI:120:ARG:NH1	2.20	0.56
31:DA:405:U:O4	34:DD:2:GLY:N	2.39	0.56
32:DB:104:ASN:O	32:DB:108:ILE:HG12	2.06	0.56
32:DB:194:PRO:O	32:DB:196:LEU:N	2.38	0.56
1:AA:555:U:O2'	1:AA:556:G:N7	2.36	0.56
4:AE:2:LYS:HB2	4:AE:95:ILE:HD12	1.87	0.56
20:AY:98:VAL:HG12	20:AY:105:ALA:HA	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:2125:G:N2	1:BA:2172:U:H3'	2.21	0.56
3:BD:46:GLN:HB2	3:BD:48:ARG:HG2	1.88	0.56
31:CA:1064:G:N2	31:CA:1191:A:OP2	2.39	0.56
31:CA:1070:U:H2'	31:CA:1071:C:C6	2.41	0.56
31:CA:1109:C:OP2	33:CC:176:HIS:ND1	2.39	0.56
31:DA:339:C:H2'	31:DA:340:U:C6	2.41	0.56
31:DA:624:C:H2'	31:DA:625:G:C8	2.40	0.56
31:DA:986:A:H1'	49:DS:54:GLY:O	2.05	0.56
31:DA:1130:A:P	39:DI:20:ARG:HH22	2.28	0.56
43:DM:19:LEU:HD12	43:DM:22:ILE:HD11	1.88	0.56
1:AA:1720:U:H2'	1:AA:1721:G:O4'	2.06	0.55
1:AA:2206:G:O2'	1:AA:2207:G:OP1	2.21	0.55
5:AF:61:GLY:HA2	5:AF:77:ASP:HB3	1.87	0.55
5:AF:143:ALA:HB1	5:AF:148:LEU:HB2	1.87	0.55
10:AO:77:ILE:HG13	15:AT:74:ARG:HG2	1.87	0.55
1:BA:577:G:OP1	1:BA:2502:G:O2'	2.20	0.55
1:BA:1047:G:H2'	1:BA:1110:G:N2	2.19	0.55
1:BA:2564:A:C2	1:BA:2647:U:H4'	2.40	0.55
8:BI:81:VAL:O	8:BI:146:ALA:HA	2.06	0.55
31:CA:41:G:H2'	31:CA:42:G:H8	1.70	0.55
31:CA:1347:G:O2'	31:CA:1348:U:OP2	2.24	0.55
38:CH:87:SER:HB2	38:CH:93:VAL:HB	1.88	0.55
31:DA:376:G:H2'	31:DA:377:G:H8	1.71	0.55
31:DA:920:U:H2'	31:DA:921:U:C6	2.41	0.55
31:DA:1241:G:H2'	31:DA:1242:C:C6	2.41	0.55
31:DA:1391:U:H2'	31:DA:1392:G:C8	2.41	0.55
32:DB:178:ARG:HH21	38:DH:74:PRO:HB3	1.70	0.55
35:DE:52:PRO:HD2	35:DE:53:LEU:HD12	1.89	0.55
1:AA:651:G:C6	1:AA:652:C:C4	2.95	0.55
1:AA:1178:C:H2'	1:AA:1179:C:C6	2.39	0.55
1:AA:1607:C:H4'	1:AA:1608:A:O5'	2.06	0.55
16:AU:106:PHE:HA	16:AU:109:LEU:HD12	1.88	0.55
1:BA:414:C:H2'	1:BA:415:A:C8	2.40	0.55
1:BA:458:G:O2'	29:B7:39:ARG:HD3	2.06	0.55
1:BA:1786:A:H1'	1:BA:1938:A:N6	2.21	0.55
5:BF:149:ASP:OD2	5:BF:149:ASP:N	2.32	0.55
6:BG:56:ALA:HA	6:BG:153:ARG:HH21	1.71	0.55
14:BS:59:LYS:HB3	14:BS:60:GLY:HA2	1.88	0.55
20:BY:76:CYS:HG	20:BY:99:CYS:HG	1.49	0.55
21:BZ:153:SER:OG	21:BZ:154:ASP:N	2.39	0.55
21:BZ:160:GLY:CA	21:BZ:161:VAL:HB	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:B1:86:SER:O	23:B1:89:GLU:HG2	2.06	0.55
31:CA:192:U:H2'	31:CA:193:C:H6	1.71	0.55
31:CA:1146:A:H2'	31:CA:1147:C:O4'	2.05	0.55
31:CA:1346:A:N1	31:CA:1374:A:H5''	2.21	0.55
46:CP:53:VAL:HG13	46:CP:79:VAL:HG23	1.87	0.55
31:DA:21:G:H2'	31:DA:22:G:C8	2.42	0.55
31:DA:586:C:O2'	31:DA:878:G:H4'	2.05	0.55
31:DA:833:U:H2'	31:DA:834:C:C6	2.39	0.55
31:DA:1062:U:H2'	31:DA:1063:C:C6	2.41	0.55
31:DA:1239:A:H4'	31:DA:1240:U:H5''	1.88	0.55
36:DF:61:LEU:HB3	36:DF:63:TYR:CE2	2.37	0.55
48:DR:73:ALA:HB3	48:DR:79:LEU:HD12	1.87	0.55
1:AA:1331:A:H2'	1:AA:1333:C:C5	2.40	0.55
1:AA:1332:G:H5''	1:AA:1332:G:H8	1.71	0.55
1:AA:2172:U:H4'	1:AA:2173:A:OP2	2.06	0.55
4:AE:107:THR:O	4:AE:190:GLY:HA3	2.06	0.55
21:AZ:151:HIS:C	21:AZ:153:SER:H	2.09	0.55
1:BA:2109:U:H2'	1:BA:2110:G:C8	2.41	0.55
1:BA:2163:C:OP2	1:BA:2164:C:H5	1.88	0.55
5:BF:103:LYS:HA	5:BF:106:ARG:HG3	1.88	0.55
31:CA:166:G:H2'	31:CA:167:G:H8	1.71	0.55
31:CA:767:A:H2'	31:CA:768:A:O4'	2.06	0.55
31:CA:1028:C:N4	31:CA:1033:G:H1	2.05	0.55
31:DA:954:G:N2	31:DA:1227:A:H62	2.04	0.55
31:DA:1329:A:H5''	43:DM:26:GLY:N	2.20	0.55
31:DA:1465:C:H2'	31:DA:1466:C:O4'	2.06	0.55
34:DD:100:ARG:NH2	34:DD:102:ASP:OD2	2.39	0.55
48:DR:32:ARG:HA	48:DR:69:THR:HG21	1.86	0.55
1:AA:674:G:H1'	5:AF:74:ARG:HD3	1.88	0.55
1:AA:775:G:N2	1:AA:794:G:H5'	2.22	0.55
1:AA:1410:G:H2'	1:AA:1411:C:H6	1.70	0.55
12:AQ:2:LEU:HB3	12:AQ:70:PRO:HG3	1.89	0.55
15:AT:55:ASN:H	15:AT:59:THR:HG22	1.71	0.55
20:AY:13:VAL:HG12	20:AY:74:PRO:HA	1.89	0.55
1:BA:1639:U:C2'	1:BA:1640:C:H5''	2.37	0.55
1:BA:1843:C:H5'	3:BD:253:GLN:NE2	2.21	0.55
1:BA:2282:G:H4'	1:BA:2283:C:O5'	2.06	0.55
1:BA:2469:A:H4'	12:BQ:56:ARG:HG2	1.86	0.55
30:B8:23:VAL:HG11	30:B8:47:LYS:HD3	1.87	0.55
31:CA:259:G:OP2	50:CT:83:ARG:HD3	2.06	0.55
31:CA:611:A:H2	31:CA:630:G:H22	1.52	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:864:A:H2'	31:CA:865:A:C8	2.42	0.55
32:CB:111:ARG:HH11	32:CB:111:ARG:HA	1.70	0.55
42:CL:11:VAL:HG11	47:CQ:36:ILE:HG21	1.87	0.55
42:CL:11:VAL:HG13	47:CQ:29:HIS:CD2	2.41	0.55
42:CL:32:PHE:HE1	42:CL:86:ARG:HG3	1.70	0.55
31:DA:522:C:H5''	42:DL:120:TYR:OH	2.07	0.55
31:DA:814:A:N7	31:DA:816:A:C4	2.74	0.55
31:DA:1512:U:H2'	31:DA:1513:A:H8	1.72	0.55
47:DQ:57:VAL:HA	47:DQ:77:VAL:HG23	1.88	0.55
1:AA:900:A:H2'	1:AA:901:A:O4'	2.06	0.55
1:AA:1153:C:H2'	1:AA:1154:G:O4'	2.07	0.55
1:AA:1332:G:H5''	1:AA:1332:G:C8	2.41	0.55
1:AA:1403:C:H5''	1:AA:1471:A:C1'	2.29	0.55
4:AE:115:GLY:O	4:AE:119:ARG:HB2	2.06	0.55
23:A1:73:LEU:HB3	23:A1:94:LEU:HD22	1.88	0.55
1:BA:1486:A:H2'	1:BA:1487:G:H8	1.70	0.55
1:BA:2115:G:H21	1:BA:2171:A:N6	2.05	0.55
6:BG:47:LYS:HB3	6:BG:82:LEU:HD11	1.87	0.55
7:BH:7:LEU:HG	7:BH:69:ARG:NH1	2.21	0.55
20:BY:35:TYR:CE2	20:BY:69:ALA:HB3	2.41	0.55
28:B6:6:ARG:NE	28:B6:24:GLU:OE1	2.25	0.55
31:CA:971:G:N2	31:CA:1363(A):A:OP2	2.31	0.55
31:DA:630:G:H2'	31:DA:631:G:H8	1.72	0.55
31:DA:1012:U:H2'	31:DA:1013:G:C8	2.41	0.55
34:DD:62:GLN:HB3	34:DD:66:ARG:HD2	1.87	0.55
1:AA:19:C:H2'	1:AA:20:C:H6	1.70	0.55
1:AA:1028:A:N6	1:AA:1125:G:H2'	2.21	0.55
1:AA:1500:G:H2'	1:AA:1501:C:C6	2.42	0.55
1:AA:1688:U:O2	1:AA:1700:A:H5'	2.07	0.55
1:AA:1792:G:H5'	3:AD:205:VAL:HG13	1.89	0.55
1:AA:1999:C:H4'	1:AA:2723:C:O2	2.05	0.55
14:AS:96:GLY:HA2	14:AS:100:ALA:H	1.72	0.55
1:BA:271(K):U:H4'	1:BA:271(L):U:OP2	2.07	0.55
1:BA:2839:G:H5'	13:BR:46:GLY:HA2	1.88	0.55
8:BI:94:ALA:HA	8:BI:97:ILE:HB	1.88	0.55
12:BQ:62:GLY:O	21:BZ:178:GLU:HG2	2.07	0.55
31:CA:342:C:H2'	31:CA:343:U:O4'	2.06	0.55
31:CA:1073:U:H2'	31:CA:1074:G:C8	2.40	0.55
31:CA:1121:U:H2'	31:CA:1122:U:C6	2.42	0.55
31:CA:1298:C:N4	37:CG:114:ARG:HD3	2.22	0.55
31:DA:32:A:H2'	31:DA:33:A:C8	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:1004:A:H8	31:DA:1025:U:N3	2.03	0.55
31:DA:1063:C:H2'	31:DA:1064:G:C8	2.42	0.55
31:DA:1238:A:C2	31:DA:1303:C:H4'	2.42	0.55
31:DA:1320:C:H1'	49:DS:73:GLU:N	2.20	0.55
33:DC:125:GLU:HG3	33:DC:189:ALA:HB1	1.88	0.55
37:DG:34:GLY:O	37:DG:36:LYS:N	2.38	0.55
43:DM:44:ARG:O	43:DM:46:LYS:N	2.40	0.55
48:DR:56:THR:HB	48:DR:58:LEU:HD13	1.88	0.55
1:AA:1312:U:H4'	1:AA:1313:U:O5'	2.07	0.55
1:AA:1378:A:OP1	29:A7:10:ARG:NH2	2.39	0.55
1:AA:1448:G:N2	1:AA:1463:C:O2	2.32	0.55
1:AA:1913:A:H2'	52:AA:3001:T8B:O2	2.07	0.55
1:AA:2171:A:H4'	1:AA:2172:U:OP1	2.07	0.55
1:AA:2384:G:OP2	22:A0:55:ARG:NH1	2.40	0.55
1:AA:2815:C:H5'	27:A5:29:THR:HG21	1.89	0.55
15:AT:13:ARG:HG3	15:AT:14:TYR:CE1	2.41	0.55
18:AW:86:LEU:HD22	18:AW:96:ILE:HD11	1.89	0.55
1:BA:729:G:OP2	3:BD:208:LYS:NZ	2.40	0.55
1:BA:1530:C:H1'	1:BA:1531:C:OP1	2.07	0.55
4:BE:29:GLY:HA3	4:BE:180:ASN:OD1	2.06	0.55
7:BH:11:VAL:HG21	7:BH:50:VAL:HG23	1.89	0.55
10:BO:68:GLU:OE2	10:BO:78:ARG:NH1	2.40	0.55
31:CA:1101:A:H4'	31:CA:1102:A:O5'	2.07	0.55
32:CB:102:LEU:HD23	32:CB:182:ILE:HD12	1.89	0.55
34:CD:173:TRP:CD1	34:CD:174:LEU:HG	2.41	0.55
44:CN:32:SER:HB3	44:CN:41:ARG:HB3	1.87	0.55
6:AG:56:ALA:HB2	6:AG:153:ARG:HE	1.71	0.55
7:AH:56:SER:HG	7:AH:61:HIS:HD1	1.52	0.55
12:BQ:138:ASP:OD2	21:BZ:81:ARG:NH1	2.35	0.55
14:BS:102:ALA:HA	14:BS:105:ALA:HB3	1.87	0.55
31:CA:115:G:H1'	31:CA:116:A:N7	2.21	0.55
31:CA:819:A:H4'	31:CA:820:U:OP2	2.06	0.55
31:CA:1124:G:H1	31:CA:1149:C:N4	2.05	0.55
31:CA:1229:A:OP2	43:CM:114:ARG:HD2	2.06	0.55
38:CH:86:ILE:HG21	38:CH:133:LEU:HD13	1.89	0.55
43:CM:102:ARG:HE	43:CM:105:THR:HG23	1.72	0.55
31:DA:1011:G:H1	31:DA:1018:C:N4	2.03	0.55
31:DA:1061:G:H5''	40:DJ:59:SER:HB2	1.88	0.55
36:DF:62:TRP:CH2	36:DF:64:GLN:HB2	2.42	0.55
37:DG:77:SER:HA	37:DG:86:GLN:HA	1.89	0.55
1:AA:359:A:H2'	1:AA:360:G:O4'	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:686:G:C8	29:A7:7:PRO:HA	2.42	0.55
1:AA:1406:U:H2'	1:AA:1407:C:C6	2.41	0.55
1:AA:1448:G:H4'	1:AA:1542:A:OP1	2.06	0.55
6:AG:61:ALA:HA	6:AG:66:GLN:O	2.07	0.55
1:BA:27:G:N2	1:BA:512:G:H1'	2.22	0.55
1:BA:1204:A:H2	1:BA:1241:A:N6	2.04	0.55
1:BA:1810:A:H2'	1:BA:1811:G:O4'	2.07	0.55
4:BE:28:ALA:HB3	4:BE:93:VAL:HG13	1.89	0.55
31:CA:115:G:H4'	31:CA:116:A:O5'	2.06	0.55
31:CA:344:A:H5''	31:CA:345:C:C5	2.42	0.55
31:CA:736:C:H2'	31:CA:737:A:C8	2.42	0.55
31:CA:1027:C:H2'	31:CA:1028:C:C5	2.41	0.55
31:CA:1292:U:O2'	31:CA:1293:G:H5'	2.06	0.55
31:CA:1414:U:H3	31:CA:1486:G:H1	1.55	0.55
36:CF:67:MET:SD	36:CF:75:LEU:HD12	2.46	0.55
43:CM:57:ARG:O	43:CM:61:GLU:HG3	2.07	0.55
31:DA:7:G:O2'	35:DE:120:THR:O	2.24	0.55
31:DA:1256:A:N6	31:DA:1278:U:O2	2.40	0.55
37:DG:126:ASP:HB3	37:DG:131:LYS:O	2.06	0.55
38:DH:6:ILE:O	38:DH:10:LEU:HG	2.07	0.55
39:DI:105:ASP:HB2	39:DI:107:ARG:HD3	1.89	0.55
1:AA:84:A:H61	1:AA:102:G:H1'	1.72	0.55
1:AA:90:U:O2'	1:AA:92:A:C8	2.56	0.55
1:AA:1653:G:N1	13:AR:11:ASN:OD1	2.38	0.55
4:AE:14:ILE:HG13	4:AE:21:VAL:HG13	1.89	0.55
26:A4:42:PHE:CB	26:A4:43:TYR:HB2	2.36	0.55
1:BA:34:C:H5''	1:BA:35:G:OP2	2.07	0.55
2:BB:19:G:H5'	2:BB:20:C:OP2	2.07	0.55
6:BG:56:ALA:HB2	6:BG:153:ARG:NE	2.17	0.55
15:BT:120:ARG:HA	15:BT:123:GLN:HG2	1.89	0.55
25:B3:39:ASP:OD2	25:B3:44:ARG:NH1	2.41	0.55
31:CA:652:U:O4	31:CA:752:G:O2'	2.21	0.55
31:CA:1347:G:H5''	39:CI:107:ARG:HB3	1.88	0.55
42:CL:53:ARG:HB3	42:CL:93:LEU:HD11	1.87	0.55
48:CR:47:THR:HG23	48:CR:49:LYS:HG3	1.89	0.55
31:DA:618:C:N4	31:DA:621:A:N7	2.54	0.55
31:DA:673:G:H2'	31:DA:674:G:H8	1.68	0.55
31:DA:1112:C:O2	33:DC:179:ARG:HG2	2.07	0.55
31:DA:1307:U:H2'	31:DA:1308:U:C6	2.42	0.55
32:DB:102:LEU:HB3	32:DB:180:LEU:HD12	1.88	0.55
38:DH:21:LYS:O	38:DH:63:LEU:HD23	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DH:121:ASP:N	38:DH:121:ASP:OD1	2.40	0.55
42:DL:90:VAL:O	42:DL:92:ASP:N	2.40	0.55
49:DS:32:LYS:HG2	49:DS:50:ALA:HB3	1.88	0.55
1:AA:9:U:N3	1:AA:2629:A:H2	2.05	0.54
1:AA:986:C:C2'	1:AA:987:G:H5'	2.37	0.54
4:AE:203:LYS:CB	4:AE:204:ALA:HA	2.37	0.54
7:AH:54:ARG:HD3	7:AH:65:HIS:ND1	2.22	0.54
9:AN:34:LEU:O	9:AN:49:GLY:HA3	2.07	0.54
13:AR:67:LEU:HD22	13:AR:76:VAL:HG21	1.88	0.54
1:BA:2299:G:N1	1:BA:2318:G:C8	2.75	0.54
2:BB:75:G:H5''	2:BB:75:G:H8	1.71	0.54
12:BQ:35:VAL:HG13	12:BQ:130:LYS:HB3	1.89	0.54
22:B0:55:ARG:NH1	22:B0:55:ARG:HB2	2.21	0.54
31:CA:539:A:H2'	31:CA:540:G:C8	2.43	0.54
31:CA:955:U:H2'	31:CA:956:U:C6	2.42	0.54
31:CA:1090:U:H2'	31:CA:1091:U:H6	1.72	0.54
32:CB:87:ARG:HH11	32:CB:219:VAL:HG12	1.72	0.54
35:CE:7:GLU:O	35:CE:8:GLU:HB3	2.07	0.54
38:CH:73:ASP:OD2	38:CH:75:ARG:HD3	2.08	0.54
31:DA:607:A:H2'	31:DA:608:A:O4'	2.07	0.54
31:DA:634:C:H2'	31:DA:635:G:H8	1.72	0.54
31:DA:642:A:N3	38:DH:113:SER:OG	2.37	0.54
31:DA:1090:U:H2'	31:DA:1091:U:H6	1.71	0.54
31:DA:1162:C:H2'	31:DA:1163:C:C6	2.41	0.54
31:DA:1332:A:O5'	31:DA:1332:A:H8	1.89	0.54
32:DB:24:TRP:CE3	32:DB:26:PRO:HA	2.42	0.54
1:AA:17:G:H2'	1:AA:18:C:C6	2.43	0.54
1:AA:244:A:C2	1:AA:255:A:C4	2.96	0.54
1:AA:1173:G:N1	1:AA:1176:G:OP2	2.26	0.54
1:AA:1201:C:H2'	1:AA:1202:C:H6	1.70	0.54
1:AA:1503:U:H2'	1:AA:1504:C:C6	2.42	0.54
1:AA:2137:C:O2	1:AA:2137:C:H2'	2.07	0.54
7:AH:70:THR:HA	7:AH:73:ALA:HB3	1.87	0.54
12:AQ:109:VAL:HG13	12:AQ:113:GLN:HB2	1.88	0.54
1:BA:1265:A:H8	1:BA:1265:A:OP1	1.89	0.54
1:BA:2129:C:N3	1:BA:2159:G:O6	2.40	0.54
1:BA:2318:G:N2	14:BS:3:ARG:HD2	2.22	0.54
3:BD:16:MET:HG2	3:BD:211:ARG:HH21	1.71	0.54
6:BG:9:ARG:NH1	6:BG:13:GLU:OE1	2.37	0.54
6:BG:108:ASN:HA	26:B4:37:SER:HB3	1.89	0.54
15:BT:95:ARG:HG2	15:BT:95:ARG:HH11	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:1001:A:N6	31:CA:1040:U:H3	2.02	0.54
31:CA:1063:C:H2'	31:CA:1064:G:C8	2.42	0.54
32:CB:87:ARG:HD2	32:CB:219:VAL:HG11	1.88	0.54
31:DA:254:G:OP1	47:DQ:66:SER:OG	2.21	0.54
31:DA:1027:C:H2'	31:DA:1028:C:C5	2.43	0.54
45:DO:23:GLY:O	45:DO:27:VAL:HB	2.07	0.54
1:AA:221:A:H4'	1:AA:222:A:O5'	2.07	0.54
1:AA:1608:A:H1'	1:AA:1610:A:OP2	2.08	0.54
2:AB:38:C:H2'	2:AB:39:A:O4'	2.07	0.54
3:AD:3:VAL:HG13	3:AD:17:THR:HB	1.87	0.54
1:BA:652(G):G:H2'	1:BA:652(H):C:H6	1.73	0.54
1:BA:2050:C:H2'	1:BA:2051:A:O4'	2.08	0.54
1:BA:2279:G:O6	22:B0:14:ARG:HD2	2.06	0.54
31:CA:101:A:H2'	31:CA:102:G:H8	1.72	0.54
31:CA:1504:G:OP1	31:CA:1507:A:H4'	2.07	0.54
31:DA:683:G:H2'	31:DA:684:A:C8	2.42	0.54
31:DA:944:G:N1	31:DA:1338:G:OP2	2.30	0.54
31:DA:1028:C:C4	31:DA:1033:G:N1	2.66	0.54
31:DA:1244:C:H2'	31:DA:1245:A:C8	2.41	0.54
35:DE:76:ILE:HG13	35:DE:77:PRO:HD2	1.90	0.54
11:AP:38:GLN:O	11:AP:39:LYS:HB2	2.08	0.54
14:AS:95:HIS:C	14:AS:99:LYS:HB3	2.28	0.54
1:BA:1332:G:N2	1:BA:1610:A:C8	2.75	0.54
11:BP:101:VAL:O	11:BP:103:ALA:N	2.40	0.54
21:BZ:111:VAL:HG12	21:BZ:112:ARG:H	1.72	0.54
21:BZ:156:LYS:HD2	21:BZ:158:PRO:HD3	1.89	0.54
31:CA:394:G:H2'	31:CA:395:C:C6	2.43	0.54
37:CG:139:GLU:HB3	37:CG:143:ARG:NH1	2.23	0.54
39:CI:28:VAL:HA	39:CI:63:ILE:O	2.08	0.54
40:CJ:44:VAL:HG12	40:CJ:46:ARG:HG3	1.89	0.54
42:CL:70:ILE:HG12	42:CL:100:ILE:HD12	1.89	0.54
45:CO:15:PHE:CE2	45:CO:84:LYS:HD2	2.42	0.54
49:CS:39:THR:O	49:CS:41:VAL:N	2.41	0.54
31:DA:859:A:H2'	31:DA:860:A:O4'	2.08	0.54
31:DA:989:C:H42	31:DA:1216:G:H1	1.55	0.54
31:DA:1128:C:H1'	31:DA:1146:A:H61	1.73	0.54
34:DD:108:LEU:HB3	34:DD:110:PHE:CE1	2.41	0.54
36:DF:22:GLU:OE2	36:DF:82:ARG:HG2	2.07	0.54
46:DP:70:ALA:O	46:DP:73:LEU:N	2.40	0.54
1:AA:76:C:O3'	24:A2:59:ARG:HG3	2.08	0.54
1:AA:492:A:H2'	1:AA:493:G:O4'	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:719:C:H2'	1:AA:720:C:H6	1.72	0.54
1:AA:1693:U:H4'	1:AA:1694:C:OP2	2.08	0.54
1:AA:1915:U:H3	52:AA:3001:T8B:C27	2.20	0.54
1:AA:2872:G:C2	1:AA:2873:A:N6	2.76	0.54
5:AF:129:PHE:CD2	5:AF:163:VAL:HG21	2.42	0.54
14:AS:14:VAL:O	14:AS:18:ILE:HG12	2.07	0.54
1:BA:2109:U:H3	1:BA:2180:U:H3	1.55	0.54
3:BD:274:ARG:HA	3:BD:275:LYS:HB2	1.90	0.54
26:B4:24:THR:OG1	26:B4:25:TYR:N	2.40	0.54
27:B5:33:CYS:SG	27:B5:46:CYS:SG	3.03	0.54
31:CA:737:A:H2'	31:CA:738:C:C6	2.43	0.54
31:CA:840:C:H4'	31:CA:841:U:OP1	2.08	0.54
31:CA:1355:G:H2'	31:CA:1356:G:C8	2.43	0.54
34:CD:9:CYS:O	34:CD:13:ARG:HG3	2.08	0.54
35:CE:71:LEU:O	35:CE:73:ASN:N	2.38	0.54
31:DA:364:A:H2'	31:DA:365:U:C6	2.42	0.54
36:DF:45:LEU:HD11	36:DF:57:GLN:HB3	1.88	0.54
1:AA:448:U:C4	1:AA:583:G:H1'	2.43	0.54
1:AA:702:G:C2	1:AA:731:C:C2	2.95	0.54
1:AA:1278:A:OP1	13:AR:36:THR:HG23	2.07	0.54
1:AA:1656:C:H2'	1:AA:1657:C:H6	1.72	0.54
1:AA:2363:C:O2	22:A0:39:ARG:NH2	2.41	0.54
13:AR:85:PRO:O	13:AR:87:TYR:N	2.41	0.54
26:A4:40:HIS:O	26:A4:43:TYR:HB3	2.07	0.54
1:BA:453:C:O2	1:BA:457:A:O2'	2.23	0.54
1:BA:2206:G:H5'	1:BA:2207:G:N7	2.22	0.54
3:BD:85:ASP:OD2	3:BD:88:ARG:NH1	2.37	0.54
6:BG:66:GLN:HG2	26:B4:1:MET:HE1	1.90	0.54
23:B1:86:SER:N	23:B1:89:GLU:OE1	2.36	0.54
31:CA:964:A:N3	31:CA:969:A:O2'	2.36	0.54
32:CB:104:ASN:O	32:CB:108:ILE:HG12	2.06	0.54
40:CJ:16:LEU:HD21	40:CJ:70:ARG:HG3	1.90	0.54
41:CK:48:ILE:HD11	41:CK:64:ALA:HA	1.89	0.54
31:DA:359:U:H2'	31:DA:360:A:H8	1.72	0.54
1:AA:64:A:O3'	19:AX:71:GLY:HA3	2.07	0.54
1:AA:141:A:H8	1:AA:1408:C:HO2'	1.49	0.54
1:AA:2165:G:H2'	1:AA:2166:G:C8	2.42	0.54
5:AF:101:LEU:O	5:AF:106:ARG:NH1	2.41	0.54
21:AZ:153:SER:OG	21:AZ:154:ASP:N	2.41	0.54
28:A6:11:LEU:HD12	28:A6:21:TYR:HB2	1.90	0.54
1:BA:363(B):G:H2'	1:BA:363(C):G:H8	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:796:C:H2'	1:BA:797:C:C6	2.43	0.54
1:BA:910:A:N1	1:BA:2277:G:H1'	2.23	0.54
1:BA:1688:U:O2	1:BA:1700:A:H5'	2.07	0.54
1:BA:2319:G:C2	14:BS:3:ARG:HA	2.42	0.54
1:BA:2572:A:N7	4:BE:145:LYS:HB2	2.22	0.54
31:CA:113:G:H2'	31:CA:114:U:H6	1.72	0.54
31:CA:674:G:H2'	31:CA:675:A:C8	2.39	0.54
31:CA:1030:C:C5	31:CA:1030(A):G:H1'	2.43	0.54
31:CA:1255:G:H3'	31:CA:1279:A:N6	2.23	0.54
32:CB:155:LEU:HD11	32:CB:159:PRO:HD3	1.89	0.54
37:CG:132:GLY:H	37:CG:135:VAL:HB	1.72	0.54
47:CQ:96:GLU:O	47:CQ:97:SER:HB2	2.08	0.54
31:DA:491:G:H2'	31:DA:492:G:H8	1.72	0.54
37:DG:15:ASP:O	37:DG:19:GLY:N	2.34	0.54
37:DG:26:PHE:CE2	37:DG:30:ILE:HD11	2.42	0.54
50:DT:38:LYS:O	50:DT:41:ILE:HG13	2.08	0.54
1:AA:1174:A:H5'	1:AA:1177:A:H61	1.71	0.54
1:AA:1394:U:H2'	1:AA:1395:A:O4'	2.07	0.54
1:AA:1766:U:H2'	1:AA:1767:C:C6	2.43	0.54
1:AA:2099:U:H3	1:AA:2190:G:H1	1.55	0.54
8:AI:92:VAL:CG1	8:AI:120:ILE:HB	2.37	0.54
1:BA:689:A:N3	1:BA:779:U:O2'	2.35	0.54
1:BA:1991:U:H2'	1:BA:1992:G:H5''	1.89	0.54
1:BA:2172:U:H4'	1:BA:2173:A:OP2	2.07	0.54
4:BE:11:MET:HG2	4:BE:24:THR:HB	1.89	0.54
7:BH:20:ALA:HB1	7:BH:21:PRO:HD2	1.89	0.54
31:CA:279:A:OP2	47:CQ:95:TYR:OH	2.20	0.54
31:CA:1458:G:H5''	50:CT:31:SER:HB2	1.90	0.54
34:CD:100:ARG:NH1	34:CD:137:SER:HB3	2.23	0.54
31:DA:1031:G:H2'	31:DA:1032:G:C8	2.42	0.54
32:DB:54:THR:HG23	32:DB:199:TYR:HB3	1.90	0.54
1:AA:249:C:O2	30:A8:12:LYS:NZ	2.31	0.54
1:AA:500:G:N2	1:AA:502:A:H3'	2.22	0.54
1:AA:2302:G:O2'	6:AG:126:ASP:O	2.22	0.54
2:AB:45:A:OP2	6:AG:96:ARG:NH2	2.37	0.54
6:AG:15:VAL:HG13	6:AG:175:LEU:HD23	1.90	0.54
6:AG:145:THR:OG1	6:AG:146:TYR:N	2.41	0.54
8:AI:93:THR:HG22	8:AI:119:PRO:HB3	1.90	0.54
14:AS:104:GLY:O	14:AS:107:GLU:HB2	2.08	0.54
16:AU:52:ARG:O	16:AU:55:ARG:HG3	2.08	0.54
1:BA:1188:U:H4'	17:BV:79:VAL:HG22	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:2079:U:O3'	23:B1:35:THR:OG1	2.24	0.54
1:BA:2319:G:N2	14:BS:3:ARG:HB2	2.23	0.54
20:BY:76:CYS:HG	20:BY:102:CYS:HG	1.54	0.54
22:B0:11:ARG:O	22:B0:14:ARG:NH2	2.40	0.54
27:B5:33:CYS:SG	27:B5:46:CYS:HB2	2.48	0.54
31:CA:712:A:H8	31:CA:712:A:O5'	1.91	0.54
31:CA:1142:G:H3'	31:CA:1143:G:H8	1.73	0.54
34:CD:3:ARG:O	34:CD:5:ILE:HG12	2.08	0.54
41:CK:20:TYR:HB2	41:CK:31:THR:HG23	1.90	0.54
31:DA:425:G:O3'	34:DD:45:GLN:NE2	2.41	0.54
31:DA:445:G:C6	31:DA:490:G:C6	2.96	0.54
31:DA:673:G:H5''	36:DF:87:ARG:NH1	2.23	0.54
32:DB:59:GLU:O	32:DB:63:MET:HG2	2.08	0.54
1:AA:321:G:C2	1:AA:341:G:H4'	2.43	0.54
1:AA:548:A:H61	17:AV:19:LYS:H	1.56	0.54
1:AA:1991:U:H2'	1:AA:1992:G:H5''	1.90	0.54
1:AA:2420:C:OP2	30:A8:33:ASN:HB2	2.08	0.54
5:AF:157:VAL:HB	5:AF:194:MET:HG2	1.88	0.54
7:AH:118:PRO:HD2	7:AH:121:ILE:HG21	1.89	0.54
9:AN:97:ARG:HA	9:AN:100:GLU:HB2	1.89	0.54
17:AV:5:VAL:HG11	17:AV:57:VAL:HG21	1.89	0.54
23:A1:51:VAL:HG11	23:A1:74:VAL:HG21	1.90	0.54
1:BA:271(F):C:H2'	1:BA:271(G):C:H6	1.72	0.54
1:BA:898:C:H2'	1:BA:899:A:O4'	2.08	0.54
1:BA:1570:A:H2'	1:BA:1571:A:C8	2.43	0.54
14:BS:11:LYS:HG3	14:BS:91:PRO:HD3	1.89	0.54
14:BS:101:LEU:O	14:BS:102:ALA:HB3	2.08	0.54
25:B3:44:ARG:O	25:B3:48:GLU:HG3	2.08	0.54
31:CA:922:G:C6	31:CA:923:A:C6	2.95	0.54
31:CA:1201:A:H1'	31:CA:1202:G:OP2	2.07	0.54
31:CA:1261:A:H3'	31:CA:1262:C:H6	1.70	0.54
36:CF:62:TRP:CH2	36:CF:64:GLN:HB2	2.43	0.54
31:DA:382:A:H2'	31:DA:383:A:C8	2.42	0.54
31:DA:936:C:H2'	31:DA:937:A:C8	2.43	0.54
1:AA:996:A:O3'	16:AU:91:ASP:HB2	2.09	0.53
1:AA:1798:U:C5'	3:AD:259:THR:HG22	2.37	0.53
1:AA:2172:U:H1'	1:AA:2173:A:OP1	2.08	0.53
3:AD:93:ALA:HB3	3:AD:105:ILE:HG13	1.90	0.53
4:AE:14:ILE:HB	15:AT:14:TYR:CE2	2.43	0.53
18:AW:56:ALA:O	18:AW:60:ASN:N	2.41	0.53
1:BA:271(D):G:H1	1:BA:271(T):C:H42	1.56	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:588:U:H1'	5:BF:90:PHE:HB3	1.90	0.53
1:BA:1298:C:H5''	1:BA:1299:G:OP2	2.08	0.53
1:BA:1365:A:OP1	23:B1:41:ARG:NH1	2.39	0.53
8:BI:27:ARG:HD2	23:B1:71:TYR:CE1	2.43	0.53
17:BV:98:GLU:OE2	17:BV:100:ARG:NH1	2.41	0.53
31:CA:31:G:O2'	31:CA:48:C:N4	2.41	0.53
31:CA:1090:U:H2'	31:CA:1091:U:C6	2.43	0.53
41:CK:31:THR:HA	41:CK:42:TRP:HA	1.89	0.53
46:CP:28:ARG:NH1	46:CP:29:ASP:OD2	2.42	0.53
31:DA:100:C:H2'	31:DA:101:A:C8	2.43	0.53
31:DA:1198:G:H2'	31:DA:1199:U:C6	2.43	0.53
39:DI:11:LYS:H	39:DI:104:ARG:HH22	1.56	0.53
1:AA:2152:G:H2'	1:AA:2153:G:C8	2.43	0.53
1:AA:2686:G:H2'	1:AA:2687:U:O4'	2.08	0.53
1:AA:2808:U:H5''	1:AA:2891:G:O6	2.09	0.53
3:AD:97:TYR:HB2	3:AD:101:GLU:O	2.07	0.53
10:AO:15:GLY:O	10:AO:47:ILE:HG12	2.08	0.53
17:AV:76:LYS:HD2	17:AV:81:TYR:CD1	2.44	0.53
26:A4:15:ILE:HB	26:A4:32:TYR:CD2	2.43	0.53
31:CA:748:C:H4'	31:CA:749:C:O5'	2.08	0.53
31:CA:857:C:H2'	31:CA:858:G:O4'	2.08	0.53
31:CA:1304:G:O6	31:CA:1331:G:O2'	2.24	0.53
33:CC:74:GLY:O	33:CC:76:VAL:N	2.41	0.53
34:CD:111:ALA:HB1	34:CD:116:GLN:HB3	1.89	0.53
31:DA:513:C:H2'	31:DA:514:C:C6	2.42	0.53
43:DM:65:LYS:CA	43:DM:66:LEU:HB2	2.28	0.53
1:AA:300:A:H2'	1:AA:334:C:O2'	2.07	0.53
1:AA:528:A:H3'	1:AA:528:A:C8	2.43	0.53
1:AA:698:C:O2'	1:AA:734:A:N6	2.40	0.53
1:AA:995:C:O2	9:AN:3:THR:OG1	2.23	0.53
16:AU:76:TYR:OH	16:AU:92:ARG:NH1	2.41	0.53
1:BA:26:G:C6	1:BA:27:G:N1	2.77	0.53
11:BP:83:VAL:HG13	11:BP:112:LEU:HD21	1.90	0.53
24:B2:32:LEU:HD22	24:B2:36:ARG:NH1	2.23	0.53
31:CA:1321:C:H5''	31:CA:1322:C:H5''	1.90	0.53
45:CO:3:ILE:HD13	45:CO:3:ILE:H	1.73	0.53
34:DD:170:VAL:HG22	34:DD:171:GLY:H	1.73	0.53
41:DK:65:ALA:HB1	41:DK:98:LEU:HD21	1.90	0.53
46:DP:58:TYR:O	46:DP:61:SER:OG	2.11	0.53
1:AA:100:G:OP2	1:AA:100:G:H4'	2.08	0.53
1:AA:1113:U:H2'	1:AA:1114:G:H8	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1547:C:H2'	1:AA:1548:C:H6	1.72	0.53
1:AA:2263:C:H2'	1:AA:2264:C:H6	1.74	0.53
1:AA:2376:A:H8	1:AA:2376:A:OP1	1.92	0.53
2:AB:63:G:C6	2:AB:64:C:C4	2.97	0.53
26:A4:26:SER:OG	26:A4:27:THR:N	2.41	0.53
1:BA:1268:A:H2'	1:BA:1269:A:O4'	2.08	0.53
1:BA:1430:C:H2'	1:BA:1431:U:C6	2.42	0.53
1:BA:2637:U:H5''	4:BE:82:ARG:HH21	1.73	0.53
1:BA:2833:G:H3'	1:BA:2834:G:H5''	1.90	0.53
4:BE:47:VAL:HG23	4:BE:84:PHE:O	2.09	0.53
13:BR:33:ARG:NH1	13:BR:115:GLU:OE2	2.32	0.53
31:CA:3:G:H2'	31:CA:3:G:N3	2.23	0.53
31:CA:391:G:OP1	46:CP:28:ARG:NH2	2.40	0.53
31:CA:1323:G:H2'	31:CA:1324:A:O4'	2.09	0.53
47:CQ:59:ILE:HG22	47:CQ:73:VAL:HA	1.90	0.53
48:CR:45:SER:OG	48:CR:47:THR:HG22	2.08	0.53
31:DA:858:G:O6	31:DA:869:G:H3'	2.08	0.53
31:DA:1011:G:H1	31:DA:1018:C:H42	1.56	0.53
43:DM:60:VAL:HA	43:DM:64:TRP:CZ3	2.42	0.53
1:AA:77:C:OP1	24:A2:59:ARG:HD3	2.07	0.53
1:AA:637:A:H4'	1:AA:638:G:O5'	2.09	0.53
1:AA:1337:G:H2'	1:AA:1338:G:H8	1.73	0.53
1:AA:2111:C:H42	1:AA:2147:G:H22	1.57	0.53
1:AA:2113:U:H2'	1:AA:2114:A:O4'	2.08	0.53
1:AA:2134:A:N3	1:AA:2159:G:O2'	2.32	0.53
10:AO:98:VAL:HG11	10:AO:114:ILE:HG23	1.91	0.53
20:AY:30:VAL:HG22	20:AY:37:VAL:HG12	1.91	0.53
1:BA:760:G:H2'	1:BA:761:A:O4'	2.08	0.53
1:BA:1913:A:C2'	52:BA:3001:T8B:O2	2.56	0.53
13:BR:83:ILE:O	13:BR:86:ARG:HG2	2.07	0.53
48:CR:71:LYS:O	48:CR:75:ILE:HG13	2.09	0.53
31:DA:1316:G:H4'	44:DN:18:VAL:HG13	1.89	0.53
32:DB:69:LEU:HB3	32:DB:162:ILE:HG22	1.91	0.53
40:DJ:5:ARG:N	40:DJ:73:ASP:OD1	2.42	0.53
1:AA:652(D):C:H2'	1:AA:652(E):G:O4'	2.07	0.53
1:AA:699:A:H2'	1:AA:700:G:O4'	2.08	0.53
1:AA:2881:C:H2'	1:AA:2882:A:O4'	2.09	0.53
6:AG:76:SER:HA	6:AG:83:ARG:HA	1.91	0.53
28:A6:16:CYS:HB2	28:A6:18:ARG:NH1	2.24	0.53
1:BA:84:A:H61	1:BA:102:G:H1'	1.74	0.53
1:BA:1106:A:H4'	1:BA:1107:G:OP2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:1877:A:H5'	1:BA:1878:G:OP2	2.09	0.53
1:BA:2009:G:OP1	18:BW:41:LYS:HE2	2.07	0.53
1:BA:2126:A:H2	1:BA:2127:G:N3	2.06	0.53
1:BA:2712:U:H1'	1:BA:2712(A):A:C8	2.44	0.53
22:B0:43:THR:OG1	22:B0:46:LYS:HG3	2.09	0.53
22:B0:55:ARG:HB2	22:B0:55:ARG:HH11	1.74	0.53
31:CA:158:G:N2	31:CA:163:C:O2	2.39	0.53
31:CA:186:C:H2'	31:CA:187:C:C6	2.44	0.53
31:CA:662:G:O2'	31:CA:836:G:OP1	2.26	0.53
31:CA:1025:U:C2	31:CA:1036:G:O6	2.61	0.53
31:CA:1032:G:H2'	31:CA:1033:G:C8	2.44	0.53
31:CA:1115:C:H1'	44:CN:61:TRP:O	2.09	0.53
31:CA:1129:C:N4	31:CA:1134:G:N7	2.56	0.53
31:DA:781:A:C8	31:DA:782:A:C8	2.97	0.53
35:DE:135:THR:O	35:DE:138:ALA:HB3	2.09	0.53
37:DG:48:LYS:O	37:DG:52:GLU:HG2	2.09	0.53
51:DU:12:LYS:HE2	51:DU:21:TYR:HB2	1.90	0.53
1:AA:686:G:H8	29:A7:7:PRO:HA	1.74	0.53
1:AA:715:G:C2	45:DO:56:LEU:HD21	2.43	0.53
7:AH:22:GLY:HA2	7:AH:37:VAL:O	2.08	0.53
7:AH:164:TYR:HB2	7:AH:167:GLU:HB2	1.89	0.53
23:A1:67:ILE:N	23:A1:68:PRO:HD2	2.23	0.53
1:BA:1815:A:OP2	3:BD:54:ARG:NH2	2.42	0.53
1:BA:2016:U:H1'	27:B5:6:VAL:HG13	1.90	0.53
5:BF:158:THR:OG1	5:BF:159:GLY:N	2.41	0.53
6:BG:109:VAL:C	6:BG:112:PRO:HD2	2.29	0.53
10:BO:2:ILE:HB	10:BO:33:ALA:HB3	1.90	0.53
28:B6:14:THR:OG1	28:B6:48:VAL:O	2.27	0.53
31:CA:141:A:H1'	31:CA:182:U:O2	2.08	0.53
31:CA:194:C:H2'	31:CA:195:A:H5''	1.91	0.53
31:CA:564:C:O2'	38:CH:91:ARG:NH2	2.41	0.53
31:CA:657:G:C2	31:CA:658:G:C8	2.96	0.53
31:CA:859:A:H2	38:CH:19:VAL:HG11	1.73	0.53
31:CA:983:A:H5'	31:CA:984:C:OP2	2.08	0.53
31:DA:539:A:H2'	31:DA:540:G:C8	2.43	0.53
31:DA:673:G:O3'	36:DF:87:ARG:NH2	2.42	0.53
31:DA:696:A:N1	31:DA:797:C:O2'	2.33	0.53
31:DA:708:C:OP1	41:DK:85:ARG:NH2	2.39	0.53
31:DA:724:G:H2'	31:DA:725:G:H8	1.73	0.53
31:DA:1151:A:O2'	31:DA:1152:A:H8	1.91	0.53
34:DD:200:GLU:O	34:DD:204:ILE:HG12	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DH:112:LEU:HB3	38:DH:133:LEU:HA	1.90	0.53
39:DI:71:SER:HA	39:DI:74:ILE:HD12	1.91	0.53
41:DK:52:GLY:H	41:DK:55:LYS:HE2	1.72	0.53
1:AA:184:C:H2'	1:AA:185:U:C6	2.43	0.53
1:AA:1109:C:H5'	1:AA:1110:G:OP2	2.09	0.53
1:AA:1486:A:H2'	1:AA:1487:G:H8	1.74	0.53
1:AA:1837:C:OP1	31:DA:784:C:H4'	2.09	0.53
1:AA:2286:A:H4'	1:AA:2287:A:O4'	2.08	0.53
1:AA:2641:G:P	9:AN:74:ARG:HH21	2.31	0.53
5:AF:135:LYS:O	5:AF:138:GLU:HB2	2.09	0.53
6:AG:125:PHE:CE1	6:AG:170:ARG:HG2	2.43	0.53
12:AQ:32:TYR:HB2	12:AQ:106:VAL:HG23	1.90	0.53
1:BA:184:C:H2'	1:BA:185:U:H6	1.72	0.53
1:BA:336:C:HO2'	20:BY:35:TYR:HH	1.45	0.53
1:BA:372:G:OP2	23:B1:69:LYS:NZ	2.39	0.53
1:BA:883:G:H1	1:BA:893:C:H42	1.57	0.53
3:BD:146:GLU:HB2	3:BD:189:CYS:HB3	1.90	0.53
4:BE:203:LYS:CB	4:BE:204:ALA:HA	2.39	0.53
8:BI:106:GLY:HA2	8:BI:107:VAL:HB	1.91	0.53
15:BT:90:GLN:OE1	15:BT:121:ILE:HD11	2.08	0.53
20:BY:76:CYS:CB	20:BY:99:CYS:HG	2.22	0.53
31:CA:110:C:O2'	46:CP:25:ARG:O	2.23	0.53
31:CA:132:C:H2'	31:CA:133:U:C6	2.44	0.53
31:CA:1092:A:H2'	31:CA:1093:A:C8	2.44	0.53
38:CH:21:LYS:O	38:CH:65:TYR:OH	2.16	0.53
46:CP:21:VAL:HG13	46:CP:33:ILE:HB	1.89	0.53
31:DA:36:C:OP1	42:DL:123:LYS:NZ	2.42	0.53
31:DA:186:C:H2'	31:DA:187:C:C6	2.43	0.53
31:DA:736:C:H2'	31:DA:737:A:C8	2.44	0.53
32:DB:97:TRP:CH2	32:DB:173:ALA:HA	2.44	0.53
42:DL:27:LEU:C	42:DL:29:GLY:H	2.12	0.53
42:DL:66:VAL:HG11	42:DL:98:TYR:HE1	1.73	0.53
48:DR:70:ILE:HG22	48:DR:74:ARG:HD2	1.90	0.53
50:DT:10:LEU:HD23	50:DT:12:ALA:H	1.74	0.53
1:AA:1045:A:H1'	1:AA:1047:G:C2	2.43	0.53
1:AA:1826:G:H4'	3:AD:242:ARG:CZ	2.39	0.53
4:AE:75:VAL:HG13	4:AE:77:ILE:H	1.74	0.53
8:AI:61:ARG:HB3	8:AI:133:HIS:CD2	2.43	0.53
1:BA:251:A:C5	1:BA:252:G:H1'	2.43	0.53
1:BA:334:C:P	1:BA:335:C:H41	2.32	0.53
1:BA:511:U:C5	1:BA:512:G:C5	2.97	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:615:G:OP1	5:BF:40:GLN:NE2	2.37	0.53
1:BA:1020:A:N1	1:BA:1141:U:O2'	2.38	0.53
1:BA:2023:G:H5'	1:BA:2617:C:H4'	1.91	0.53
21:BZ:63:ASP:OD1	21:BZ:65:GLN:HB2	2.09	0.53
31:CA:616:G:C2	31:CA:617:G:C8	2.97	0.53
31:CA:1322:C:OP1	31:CA:1322:C:H6	1.92	0.53
32:CB:24:TRP:CD2	32:CB:26:PRO:HD3	2.44	0.53
31:DA:110:C:H2'	31:DA:111:G:O4'	2.08	0.53
31:DA:573:A:N3	31:DA:883:C:O2'	2.40	0.53
31:DA:827:U:H5''	31:DA:828:A:OP2	2.09	0.53
31:DA:828:A:N6	31:DA:858:G:O2'	2.41	0.53
32:DB:100:GLY:O	32:DB:104:ASN:N	2.41	0.53
32:DB:149:LEU:HD22	32:DB:152:PHE:HB3	1.91	0.53
33:DC:19:GLU:O	33:DC:40:ARG:NH2	2.28	0.53
51:DU:6:ARG:O	51:DU:8:THR:N	2.38	0.53
1:AA:1359:A:H2'	1:AA:1360:A:H5'	1.90	0.53
1:AA:2593:U:H2'	1:AA:2594:C:C6	2.43	0.53
15:AT:94:ALA:HB1	15:AT:99:LEU:HD21	1.91	0.53
24:A2:4:SER:HA	24:A2:7:ARG:NH1	2.24	0.53
1:BA:1262:A:N3	27:B5:10:LYS:HE3	2.24	0.53
1:BA:1278:A:OP1	13:BR:36:THR:HG23	2.09	0.53
7:BH:154:PRO:HB3	7:BH:163:TYR:CE2	2.43	0.53
13:BR:59:ASP:N	13:BR:59:ASP:OD2	2.40	0.53
19:BX:92:LEU:C	19:BX:94:GLY:H	2.13	0.53
31:CA:15:G:H4'	35:CE:24:ARG:NH1	2.23	0.53
31:CA:184:G:H2'	31:CA:185:A:H8	1.74	0.53
31:CA:583:A:H2'	31:CA:584:G:O4'	2.08	0.53
31:CA:1124:G:H1	31:CA:1149:C:H42	1.56	0.53
31:CA:1288:A:O3'	51:CU:10:ARG:NH2	2.42	0.53
42:CL:49:ASN:ND2	42:CL:92:ASP:OD1	2.42	0.53
31:DA:297:G:N2	31:DA:300:A:OP2	2.42	0.53
31:DA:1005:A:O3'	31:DA:1037:C:O2'	2.26	0.53
31:DA:1390:U:H2'	31:DA:1391:U:C6	2.44	0.53
36:DF:82:ARG:HB3	36:DF:85:VAL:HG23	1.91	0.53
44:DN:23:ARG:HG3	44:DN:24:CYS:N	2.15	0.53
47:DQ:58:GLU:HG3	47:DQ:77:VAL:HG21	1.91	0.53
1:AA:631:A:H2'	1:AA:632:A:O4'	2.09	0.52
1:AA:1331:A:O2'	1:AA:1332:G:H8	1.91	0.52
1:AA:2261:C:O2'	1:AA:2262:U:H5'	2.10	0.52
17:AV:24:LYS:HA	17:AV:92:THR:OG1	2.10	0.52
24:A2:31:GLU:HB3	24:A2:53:LEU:HD11	1.89	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:299:A:N1	1:BA:322:A:O2'	2.38	0.52
1:BA:548:A:O2'	1:BA:549:G:OP1	2.23	0.52
1:BA:2273:A:H2'	1:BA:2274:A:C8	2.44	0.52
1:BA:2732:G:H3'	1:BA:2733:A:O4'	2.09	0.52
31:CA:250:A:H4'	31:CA:251:G:O5'	2.08	0.52
47:CQ:50:LYS:HE3	47:CQ:51:TYR:CE1	2.44	0.52
31:DA:936:C:H2'	31:DA:937:A:O4'	2.08	0.52
31:DA:1001:A:H2'	31:DA:1001(A):G:H8	1.74	0.52
31:DA:1246:C:H2'	31:DA:1247:U:H6	1.74	0.52
31:DA:1260:C:O5'	31:DA:1284:C:H4'	2.09	0.52
32:DB:34:ALA:O	32:DB:41:ILE:N	2.33	0.52
39:DI:89:ASN:O	39:DI:92:TYR:N	2.40	0.52
1:AA:102:G:HO2'	1:AA:103:A:P	2.27	0.52
1:AA:1410:G:H2'	1:AA:1411:C:C6	2.44	0.52
1:AA:1963:U:H4'	1:AA:1964:G:OP1	2.09	0.52
1:AA:2875:C:O2'	15:AT:2:ASN:OD1	2.25	0.52
17:AV:35:LEU:HB2	17:AV:57:VAL:CG2	2.40	0.52
1:BA:350:U:H2'	1:BA:351:G:O4'	2.09	0.52
1:BA:548:A:H61	17:BV:19:LYS:H	1.57	0.52
1:BA:587:C:O2	11:BP:33:ARG:NH2	2.41	0.52
1:BA:1009:A:OP1	9:BN:37:LYS:NZ	2.35	0.52
2:BB:24:G:N7	2:BB:56:G:H2'	2.24	0.52
4:BE:201:THR:OG1	4:BE:202:LYS:N	2.41	0.52
6:BG:50:ALA:C	6:BG:52:ILE:H	2.12	0.52
10:BO:102:VAL:HB	10:BO:106:LEU:HD12	1.91	0.52
24:B2:22:GLU:OE2	24:B2:68:ARG:NH2	2.39	0.52
31:CA:119:A:H4'	31:CA:120:A:O5'	2.09	0.52
31:CA:160:A:H3'	31:CA:161:A:H8	1.73	0.52
31:CA:266:G:H5'	31:CA:267:C:H5	1.74	0.52
31:CA:1025:U:HO2'	31:CA:1026:G:H8	1.54	0.52
31:CA:1274:G:H2'	31:CA:1275:A:H8	1.74	0.52
31:CA:1392:G:N2	31:CA:1502:A:H8	2.06	0.52
36:CF:68:PRO:HG2	36:CF:71:ARG:HD3	1.91	0.52
36:CF:69:GLU:O	36:CF:72:VAL:HG13	2.09	0.52
40:CJ:50:ILE:HB	44:CN:41:ARG:HE	1.74	0.52
47:CQ:13:ASP:O	47:CQ:15:MET:N	2.42	0.52
31:DA:108:G:C6	50:DT:15:ARG:HG3	2.44	0.52
31:DA:473:G:O2'	31:DA:474:G:H5'	2.10	0.52
31:DA:1135:U:H2'	31:DA:1137:C:O2	2.09	0.52
31:DA:1162:C:H2'	31:DA:1163:C:H6	1.74	0.52
40:DJ:38:ILE:HG12	40:DJ:71:LEU:O	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DL:24:VAL:HG13	42:DL:98:TYR:CE2	2.44	0.52
45:DO:40:SER:O	45:DO:44:LYS:HG3	2.09	0.52
1:AA:17:G:H2'	1:AA:18:C:H6	1.73	0.52
1:AA:1517:G:C6	1:AA:1518:U:C4	2.97	0.52
1:AA:2136:C:N4	1:AA:2156:G:C2	2.77	0.52
1:AA:2494:G:C5	1:AA:2495:G:N7	2.78	0.52
14:AS:62:LYS:HB3	14:AS:97:ARG:HD2	1.90	0.52
17:AV:89:GLN:OE1	17:AV:90:PRO:HD2	2.09	0.52
1:BA:587:C:OP2	11:BP:21:ARG:NH2	2.42	0.52
1:BA:2689:U:H5'	1:BA:2690:C:H5'	1.91	0.52
6:BG:7:LEU:HD13	6:BG:176:LEU:HD22	1.91	0.52
7:BH:83:TYR:CE1	7:BH:138:LYS:HD2	2.44	0.52
15:BT:125:ARG:O	15:BT:127:ALA:N	2.35	0.52
22:B0:72:ARG:HB2	22:B0:75:LEU:HB2	1.91	0.52
31:CA:1310:G:H5'	43:CM:77:ASN:HD22	1.75	0.52
31:CA:1359:C:O2'	31:CA:1362:C:N4	2.42	0.52
31:DA:193:C:H2'	31:DA:194:C:H6	1.73	0.52
31:DA:456:C:N3	31:DA:476:G:C2	2.77	0.52
31:DA:603:U:H2'	31:DA:604:G:C8	2.43	0.52
31:DA:1186:G:H21	44:DN:61:TRP:C	2.13	0.52
31:DA:1423:G:H2'	31:DA:1424:C:C6	2.45	0.52
33:DC:179:ARG:HH12	33:DC:206:GLU:CD	2.11	0.52
36:DF:69:GLU:O	36:DF:72:VAL:HG13	2.10	0.52
37:DG:41:ARG:O	37:DG:45:ASP:N	2.37	0.52
1:AA:375:C:H2'	1:AA:376:C:H6	1.73	0.52
1:AA:1265:A:C8	1:AA:1267:U:C2	2.97	0.52
1:AA:2178:C:H2'	1:AA:2179:C:O4'	2.09	0.52
1:AA:2319:G:N1	14:AS:3:ARG:HA	2.25	0.52
1:AA:2364:C:H2'	1:AA:2365:G:O4'	2.10	0.52
1:AA:2880:C:O3'	13:AR:90:ARG:NH1	2.43	0.52
9:AN:97:ARG:O	9:AN:100:GLU:N	2.43	0.52
21:AZ:144:LEU:HD21	21:AZ:150:LEU:HG	1.91	0.52
1:BA:86:C:H4'	1:BA:104:U:H1'	1.91	0.52
1:BA:657:U:H2'	1:BA:658:C:C6	2.44	0.52
1:BA:1110:G:H8	1:BA:1110:G:OP2	1.93	0.52
1:BA:1578:U:C2'	1:BA:1579:A:H5'	2.40	0.52
1:BA:2130:U:O2'	1:BA:2133:G:O2'	2.10	0.52
1:BA:2134:A:H1'	1:BA:2159:G:H1'	1.90	0.52
1:BA:2206:G:H3'	1:BA:2207:G:C8	2.45	0.52
5:BF:176:LEU:HD21	5:BF:180:GLY:O	2.08	0.52
6:BG:13:GLU:O	6:BG:14:GLU:HB2	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:BN:75:TYR:HA	9:BN:81:GLY:O	2.09	0.52
21:BZ:65:GLN:OE1	21:BZ:67:LEU:HD21	2.10	0.52
31:CA:17:U:H2'	31:CA:18:C:C6	2.44	0.52
31:CA:725:G:H2'	31:CA:726:C:H6	1.74	0.52
31:CA:872:A:C8	31:CA:874:G:C8	2.98	0.52
31:CA:1037:C:O5'	31:CA:1037:C:H6	1.92	0.52
32:CB:91:PRO:HD3	32:CB:154:LEU:HD22	1.90	0.52
45:CO:3:ILE:HA	45:CO:7:GLU:OE2	2.10	0.52
31:DA:575:G:OP1	31:DA:575:G:H4'	2.09	0.52
31:DA:662:G:H2'	31:DA:663:A:C8	2.44	0.52
31:DA:1129:C:N4	31:DA:1134:G:N7	2.57	0.52
31:DA:1132:C:H2'	31:DA:1133:G:C8	2.44	0.52
36:DF:22:GLU:OE1	36:DF:84:ASN:HB2	2.08	0.52
39:DI:26:VAL:O	39:DI:33:PHE:N	2.42	0.52
1:AA:139(A):G:O2'	1:AA:140:G:H5'	2.09	0.52
1:AA:271(I):G:H1	1:AA:271(O):C:H42	1.56	0.52
1:AA:673:C:H5''	5:AF:81:PRO:HD2	1.90	0.52
1:AA:1022:G:N2	1:AA:1142(A):A:C2	2.71	0.52
1:AA:1899:G:N3	1:AA:1899:G:H2'	2.24	0.52
1:AA:2543:G:HO2'	1:AA:2645:G:HO2'	1.57	0.52
4:AE:47:VAL:HG12	4:AE:49:LEU:HD13	1.92	0.52
13:AR:103:ARG:HD3	18:AW:40:ASN:ND2	2.25	0.52
27:A5:51:TYR:CE1	27:A5:56:LYS:HG2	2.44	0.52
1:BA:579:G:H2'	1:BA:580:C:C6	2.45	0.52
1:BA:652(S):C:H5'	1:BA:652(T):C:OP2	2.08	0.52
1:BA:829:A:N7	1:BA:2248:C:H5'	2.24	0.52
1:BA:1257:C:H4'	5:BF:83:PHE:CE2	2.45	0.52
10:BO:107:ARG:NE	15:BT:36:GLU:HG2	2.24	0.52
31:CA:402:G:C6	31:CA:403:C:C4	2.98	0.52
31:CA:445:G:H2'	31:CA:446:G:C8	2.45	0.52
31:CA:491:G:H2'	31:CA:492:G:C8	2.45	0.52
32:CB:81:VAL:HG22	32:CB:215:LEU:HD11	1.91	0.52
34:CD:28:SER:C	34:CD:30:LYS:H	2.13	0.52
31:DA:1097:C:H1'	31:DA:1169:A:H2	1.75	0.52
41:DK:22:HIS:O	41:DK:29:ILE:N	2.31	0.52
1:AA:93:G:H2'	1:AA:94:C:C6	2.44	0.52
1:AA:807:U:C2	1:AA:808:G:C8	2.97	0.52
1:AA:2133:G:N2	1:AA:2157:G:H2'	2.24	0.52
1:AA:2489:G:O2'	1:AA:2490:G:H5'	2.10	0.52
1:AA:2641:G:OP2	9:AN:74:ARG:NH2	2.39	0.52
1:BA:1366:A:OP1	23:B1:3:LYS:NZ	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:1667:G:N2	1:BA:1992:G:OP2	2.35	0.52
1:BA:1919:A:O3'	31:CA:1517:G:H1'	2.09	0.52
1:BA:2138:C:N4	1:BA:2153:G:H1	2.07	0.52
1:BA:2789:C:HO2'	1:BA:2790:A:HO2'	1.56	0.52
5:BF:107:LYS:HE3	5:BF:206:ILE:HA	1.91	0.52
19:BX:34:ALA:O	19:BX:77:LYS:NZ	2.41	0.52
31:CA:113:G:H2'	31:CA:114:U:C6	2.45	0.52
31:CA:599:C:H4'	38:CH:130:GLY:C	2.30	0.52
31:CA:769:G:H4'	31:CA:1513:A:H4'	1.91	0.52
31:CA:1298:C:H4'	31:CA:1299:A:O4'	2.08	0.52
36:DF:10:LEU:HD23	36:DF:61:LEU:HD13	1.90	0.52
44:DN:24:CYS:SG	44:DN:40:CYS:HB3	2.50	0.52
1:AA:637:A:OP1	11:AP:133:SER:OG	2.28	0.52
1:AA:1471:A:OP2	1:AA:1519:G:N1	2.33	0.52
1:AA:2126:A:H1'	1:AA:2127:G:OP2	2.10	0.52
1:AA:2405:G:HO2'	1:AA:2406:U:P	2.33	0.52
1:AA:2468:G:C2	1:AA:2481:G:N3	2.77	0.52
1:AA:2778:A:O2'	1:AA:2781:A:H5'	2.10	0.52
6:AG:114:ILE:HB	6:AG:117:PHE:HB2	1.91	0.52
1:BA:1799:G:O2'	1:BA:1800:C:OP2	2.23	0.52
1:BA:1833:U:H2'	1:BA:1834:U:C6	2.45	0.52
31:CA:266:G:H5''	31:CA:267:C:C5	2.44	0.52
31:CA:491:G:H2'	31:CA:492:G:H8	1.74	0.52
32:CB:134:GLU:O	32:CB:137:ARG:NH2	2.42	0.52
31:DA:1076:C:C2	31:DA:1082:G:N2	2.78	0.52
36:DF:11:ASN:O	36:DF:14:LEU:HB2	2.10	0.52
45:DO:82:ILE:HB	45:DO:87:ILE:HG22	1.91	0.52
1:AA:27:G:O2'	1:AA:28:A:OP2	2.28	0.52
1:AA:271(H):G:HO2'	1:AA:271(I):G:H8	1.58	0.52
1:AA:298:G:H5''	1:AA:299:A:OP1	2.09	0.52
1:AA:1027:A:C2	1:AA:2488:A:H5'	2.45	0.52
1:AA:1047:G:C2'	1:AA:1110:G:H22	2.21	0.52
1:AA:1187:G:H8	1:AA:1187:G:O5'	1.93	0.52
1:AA:1948:G:H1'	31:DA:1483:A:H1'	1.90	0.52
1:AA:2543:G:H2'	1:AA:2544:G:C8	2.45	0.52
1:AA:2575:C:H2'	1:AA:2578:G:O6	2.09	0.52
1:BA:1651:G:N7	13:BR:11:ASN:ND2	2.58	0.52
11:BP:59:LEU:HG	30:B8:58:ILE:HD13	1.92	0.52
12:BQ:31:ASP:HA	12:BQ:134:ARG:HG3	1.91	0.52
31:CA:96:U:O2'	31:CA:97:G:P	2.68	0.52
31:CA:333:G:H4'	50:CT:16:HIS:CE1	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:CB:15:VAL:H	32:CB:16:HIS:CE1	2.28	0.52
32:CB:97:TRP:CH2	32:CB:173:ALA:HA	2.45	0.52
34:CD:10:ARG:HG3	34:CD:11:LEU:HD23	1.91	0.52
34:CD:60:GLU:OE1	34:CD:199:ASN:N	2.36	0.52
43:CM:91:ARG:NE	43:CM:97:PRO:O	2.39	0.52
31:DA:675:A:H2'	31:DA:676:A:C8	2.45	0.52
31:DA:729:A:H2'	31:DA:730:G:H8	1.75	0.52
36:DF:2:ARG:CZ	36:DF:69:GLU:HG2	2.39	0.52
1:AA:1003:G:N2	1:AA:1153:C:C2	2.78	0.52
1:AA:1364:G:C8	23:A1:3:LYS:HD3	2.45	0.52
1:AA:1592:C:H2'	1:AA:1593:G:H8	1.75	0.52
2:AB:111:G:H2'	2:AB:112:U:C6	2.45	0.52
7:AH:98:LEU:HD12	7:AH:102:ALA:O	2.08	0.52
8:AI:77:LEU:HD12	8:AI:142:VAL:HG12	1.92	0.52
14:AS:25:ARG:NH1	14:AS:42:ASP:OD2	2.43	0.52
14:AS:85:VAL:O	14:AS:112:PHE:HB3	2.10	0.52
24:A2:22:GLU:OE2	24:A2:68:ARG:NH2	2.40	0.52
1:BA:764:A:H5'	3:BD:210:GLY:HA2	1.91	0.52
1:BA:2007:C:H5'	1:BA:2824:C:H1'	1.92	0.52
1:BA:2790:A:N3	1:BA:2790:A:H2'	2.24	0.52
8:BI:38:LEU:HB3	8:BI:40:THR:HG23	1.92	0.52
9:BN:47:ALA:HB2	9:BN:112:LEU:CD1	2.39	0.52
9:BN:128:HIS:O	9:BN:131:GLN:NE2	2.43	0.52
12:BQ:47:ILE:HD11	12:BQ:68:ILE:HD12	1.92	0.52
15:BT:125:ARG:C	15:BT:127:ALA:H	2.11	0.52
17:BV:5:VAL:HG11	17:BV:57:VAL:HG21	1.92	0.52
31:CA:21:G:H2'	31:CA:22:G:C8	2.45	0.52
31:CA:269:C:H2'	31:CA:270:A:H8	1.74	0.52
31:CA:436:C:H2'	31:CA:436:C:OP2	2.09	0.52
31:CA:538:G:OP2	42:CL:115:LYS:HB2	2.10	0.52
31:CA:1442:G:H1	31:CA:1461:G:H21	1.56	0.52
31:DA:130:A:N3	31:DA:263:A:O2'	2.34	0.52
31:DA:1143:G:H2'	31:DA:1144:G:H8	1.73	0.52
31:DA:1165:C:H2'	31:DA:1166:G:H8	1.75	0.52
31:DA:1504:G:H4'	31:DA:1505:G:O5'	2.09	0.52
37:DG:105:VAL:HG12	37:DG:109:ASN:ND2	2.24	0.52
1:AA:753:C:H2'	1:AA:754:C:H6	1.74	0.52
1:AA:2122:U:H2'	1:AA:2123:G:C8	2.44	0.52
1:AA:2252:G:H2'	1:AA:2253:G:O4'	2.09	0.52
1:AA:2298:A:H2'	1:AA:2299:G:O4'	2.10	0.52
1:AA:2391:G:O6	1:AA:2425:A:H8	1.93	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2473:U:H2'	1:AA:2473:U:O2	2.10	0.52
8:AI:124:GLY:H	8:AI:144:VAL:HG13	1.75	0.52
8:AI:132:PRO:HD3	8:AI:138:ILE:HD12	1.92	0.52
9:AN:18:ALA:HB1	9:AN:60:ILE:HD12	1.91	0.52
26:A4:15:ILE:HG13	26:A4:21:VAL:HG13	1.92	0.52
1:BA:9:U:O2'	1:BA:10:G:OP1	2.24	0.52
1:BA:2464:C:O2'	1:BA:2465:C:OP2	2.21	0.52
2:BB:52:A:HO2'	2:BB:53:A:H2	1.56	0.52
14:BS:78:LEU:HA	14:BS:82:ILE:O	2.10	0.52
28:B6:15:GLU:HG3	28:B6:47:THR:HG23	1.91	0.52
31:CA:500:G:N2	31:CA:546:G:H1'	2.25	0.52
33:CC:47:LEU:HD11	33:CC:68:VAL:HG11	1.92	0.52
31:DA:600:C:H2'	31:DA:601:C:H6	1.73	0.52
31:DA:1184:G:H8	31:DA:1184:G:OP1	1.92	0.52
34:DD:36:ARG:HG2	34:DD:36:ARG:HH11	1.74	0.52
46:DP:49:LEU:HD21	46:DP:77:ALA:HB2	1.90	0.52
1:AA:1201:C:H2'	1:AA:1202:C:C6	2.45	0.51
1:AA:1329:U:H5''	1:AA:1330:C:H5	1.75	0.51
1:AA:1826:G:H2'	1:AA:1827:C:C6	2.44	0.51
1:AA:2831:G:O2'	1:AA:2883:A:H2'	2.10	0.51
3:AD:71:ASP:CB	3:AD:103:ARG:HH22	2.22	0.51
4:AE:7:VAL:HG13	4:AE:27:LEU:HB3	1.92	0.51
7:AH:64:LEU:O	7:AH:67:LEU:HB3	2.10	0.51
8:AI:101:LEU:O	8:AI:107:VAL:HB	2.10	0.51
13:AR:85:PRO:C	13:AR:87:TYR:H	2.12	0.51
1:BA:55:G:O2'	1:BA:127:A:N1	2.36	0.51
1:BA:762:U:H4'	1:BA:763:G:O5'	2.10	0.51
1:BA:1113:U:H2'	1:BA:1114:G:C8	2.45	0.51
1:BA:1174:A:N3	1:BA:1175:U:H5''	2.25	0.51
5:BF:32:LEU:HD11	5:BF:105:VAL:HG13	1.92	0.51
11:BP:98:GLU:OE2	11:BP:99:LEU:HD23	2.10	0.51
16:BU:114:LYS:O	16:BU:117:GLN:HB2	2.11	0.51
31:CA:667:G:H4'	45:CO:51:HIS:CE1	2.45	0.51
31:CA:1051:C:H2'	31:CA:1052:U:C6	2.44	0.51
31:CA:1250:A:C2	31:CA:1370:G:H1'	2.45	0.51
31:CA:1315:U:O2'	31:CA:1360:A:N3	2.36	0.51
31:CA:1346:A:H5''	39:CI:120:ARG:HH12	1.75	0.51
41:CK:48:ILE:HG21	41:CK:63:LEU:HD22	1.92	0.51
46:CP:28:ARG:HH11	46:CP:28:ARG:HG2	1.74	0.51
31:DA:308:C:H2'	31:DA:309:G:C8	2.42	0.51
31:DA:513:C:H2'	31:DA:514:C:H6	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:1143:G:H2'	31:DA:1144:G:C8	2.45	0.51
31:DA:1207:G:H2'	31:DA:1208:C:C6	2.45	0.51
33:DC:9:GLY:HA2	33:DC:12:LEU:HG	1.92	0.51
43:DM:108:ARG:O	43:DM:112:GLY:N	2.41	0.51
1:AA:528:A:C2	1:AA:2042:A:H2'	2.44	0.51
1:AA:1028:A:N3	1:AA:2486:G:O2'	2.35	0.51
1:AA:1364:G:OP1	23:A1:2:SER:HA	2.11	0.51
1:AA:1839:G:N7	1:AA:1927:A:H1'	2.25	0.51
1:AA:2582:G:C2	1:AA:2583:G:C8	2.98	0.51
1:AA:2745:C:C4	1:AA:2746:U:C4	2.98	0.51
2:AB:24:G:N3	2:AB:27:C:N4	2.53	0.51
10:AO:75:SER:OG	10:AO:76:ALA:N	2.41	0.51
11:AP:86:LYS:HG2	11:AP:117:GLU:O	2.11	0.51
24:A2:28:LYS:HB3	24:A2:57:ILE:HG12	1.93	0.51
28:A6:14:THR:OG1	28:A6:48:VAL:O	2.29	0.51
1:BA:652(Q):G:O6	1:BA:652(R):C:N4	2.44	0.51
7:BH:106:THR:O	7:BH:106:THR:OG1	2.23	0.51
15:BT:5:ALA:O	15:BT:9:LEU:N	2.39	0.51
31:CA:721:G:H4'	31:CA:722:A:O4'	2.10	0.51
32:CB:210:SER:O	32:CB:214:ILE:HG12	2.09	0.51
37:CG:100:ALA:O	37:CG:104:LEU:HB2	2.09	0.51
41:CK:27:ASN:OD1	41:CK:28:THR:N	2.34	0.51
31:DA:616:G:C2	31:DA:617:G:C8	2.99	0.51
31:DA:1142:G:H3'	31:DA:1143:G:H8	1.74	0.51
34:DD:61:LYS:HD3	34:DD:206:PHE:CD2	2.46	0.51
1:AA:102:G:OP1	24:A2:7:ARG:NH2	2.43	0.51
1:AA:186:G:H2'	1:AA:187:G:C8	2.44	0.51
1:AA:315:G:H2'	1:AA:316:C:C6	2.45	0.51
1:AA:814:C:O2'	1:AA:815:C:H5'	2.11	0.51
1:AA:819:A:C4	1:AA:1189:A:C2	2.98	0.51
1:AA:2489:G:C2'	1:AA:2490:G:H5'	2.40	0.51
2:AB:11:C:OP2	2:AB:12:C:H5	1.92	0.51
7:AH:96:ALA:O	7:AH:125:VAL:HG11	2.10	0.51
11:AP:112:LEU:HD22	11:AP:113:LYS:N	2.26	0.51
14:AS:67:ARG:O	14:AS:71:ARG:HG3	2.10	0.51
1:BA:1131:G:OP2	1:BA:2515:C:H4'	2.10	0.51
1:BA:2267:A:H5''	1:BA:2268:A:H5'	1.91	0.51
2:BB:8:U:O3'	14:BS:25:ARG:NH2	2.43	0.51
32:CB:167:PRO:O	32:CB:171:ALA:N	2.43	0.51
35:CE:84:PHE:CE2	35:CE:133:TYR:HD2	2.28	0.51
47:CQ:78:GLU:HG2	47:CQ:79:SER:H	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:427:U:OP1	34:DD:13:ARG:NH2	2.42	0.51
31:DA:779:C:H5''	41:DK:122:LYS:HG2	1.92	0.51
31:DA:825:G:N2	38:DH:11:THR:HG21	2.25	0.51
42:DL:27:LEU:O	42:DL:29:GLY:N	2.38	0.51
1:AA:330:A:H2	1:AA:1210:A:HO2'	1.57	0.51
1:AA:548:A:O2'	1:AA:549:G:OP1	2.26	0.51
1:AA:1488:G:C6	1:AA:1489:U:N3	2.78	0.51
1:AA:1686:C:N4	1:AA:1687:G:C6	2.79	0.51
1:AA:2207:G:HO2'	1:AA:2208:A:P	2.32	0.51
2:AB:9:G:OP1	14:AS:15:ARG:HD3	2.09	0.51
11:AP:82:GLY:HA2	11:AP:113:LYS:O	2.10	0.51
12:AQ:42:ILE:HD13	12:AQ:97:VAL:HG21	1.91	0.51
25:A3:23:LEU:HD13	25:A3:50:VAL:HG11	1.92	0.51
1:BA:1109:C:H5'	1:BA:1110:G:OP2	2.11	0.51
1:BA:1173:G:O2'	1:BA:1174:A:O4'	2.27	0.51
1:BA:1547:C:H2'	1:BA:1548:C:C6	2.44	0.51
1:BA:1818:U:O4	3:BD:154:LYS:HE2	2.11	0.51
1:BA:2463:C:O2'	1:BA:2464:C:H5'	2.11	0.51
2:BB:32:C:C2	2:BB:51:G:N2	2.79	0.51
2:BB:48:A:P	14:BS:30:ARG:HH12	2.33	0.51
4:BE:14:ILE:HG13	4:BE:21:VAL:HG13	1.93	0.51
8:BI:130:TYR:HB3	8:BI:138:ILE:HB	1.92	0.51
31:CA:260:G:H2'	31:CA:261:U:C6	2.45	0.51
31:CA:872:A:C4	31:CA:874:G:N7	2.79	0.51
31:CA:920:U:H2'	31:CA:921:U:C6	2.44	0.51
31:CA:1028:C:C2	31:CA:1033:G:N2	2.78	0.51
32:CB:170:GLU:O	32:CB:174:VAL:HG23	2.10	0.51
35:CE:144:THR:O	35:CE:148:VAL:HG23	2.11	0.51
39:CI:10:ARG:HA	39:CI:104:ARG:NH1	2.26	0.51
43:CM:19:LEU:HB3	43:CM:25:ILE:HG21	1.92	0.51
31:DA:433:C:H2'	31:DA:434:U:C6	2.46	0.51
31:DA:491:G:H2'	31:DA:492:G:C8	2.45	0.51
31:DA:603:U:H2'	31:DA:604:G:H8	1.74	0.51
31:DA:620:C:H2'	31:DA:621:A:O4'	2.11	0.51
31:DA:1025:U:O2	31:DA:1036:G:O6	2.28	0.51
37:DG:88:PRO:HD2	37:DG:151:TYR:HB2	1.91	0.51
41:DK:82:VAL:HB	41:DK:108:ILE:HG12	1.92	0.51
1:AA:918:A:H5''	2:AB:98:G:O2'	2.11	0.51
1:AA:2687:U:H2'	1:AA:2688:U:O4'	2.10	0.51
2:AB:49:C:OP1	14:AS:97:ARG:N	2.42	0.51
25:A3:46:ASN:O	25:A3:50:VAL:HG22	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:616:G:H5'	5:BF:205:ARG:HD2	1.92	0.51
5:BF:28:ILE:HG12	5:BF:116:ASP:HB2	1.93	0.51
20:BY:90:LEU:O	20:BY:91:GLU:HB2	2.09	0.51
31:CA:1225:A:N3	31:CA:1225:A:H2'	2.25	0.51
31:CA:1249:C:O2	39:CI:69:GLY:HA2	2.10	0.51
47:CQ:3:LYS:HD2	47:CQ:60:ILE:HD11	1.93	0.51
31:DA:96:U:O2'	31:DA:97:G:P	2.69	0.51
31:DA:1111:A:H2'	31:DA:1112:C:C6	2.46	0.51
31:DA:1123:A:H4'	40:DJ:36:GLY:HA3	1.92	0.51
31:DA:1164:G:H1	31:DA:1172:C:H42	1.59	0.51
31:DA:1284:C:H2'	31:DA:1285:A:C8	2.45	0.51
41:DK:29:ILE:HA	41:DK:44:SER:HB3	1.93	0.51
1:AA:1283:G:N2	1:AA:1286:A:OP2	2.36	0.51
1:AA:2319:G:H22	14:AS:3:ARG:NE	2.09	0.51
1:AA:2320:A:H2'	1:AA:2320:A:N3	2.24	0.51
1:AA:2331:G:O2'	22:A0:43:THR:HG22	2.10	0.51
1:AA:2586:C:O5'	1:AA:2586:C:H6	1.93	0.51
2:AB:51:G:N7	14:AS:62:LYS:NZ	2.48	0.51
3:AD:16:MET:HG2	3:AD:211:ARG:HH21	1.75	0.51
6:AG:107:LEU:HD13	6:AG:177:GLY:O	2.11	0.51
24:A2:32:LEU:HD11	24:A2:54:LYS:HG3	1.91	0.51
1:BA:218:A:C2	1:BA:235:U:H4'	2.45	0.51
1:BA:528:A:O2'	1:BA:529:A:H5'	2.11	0.51
1:BA:2086:U:H2'	1:BA:2087:G:C8	2.45	0.51
1:BA:2577:A:H5''	1:BA:2578:G:H5'	1.93	0.51
1:BA:2834:G:H5''	1:BA:2834:G:H8	1.75	0.51
3:BD:142:VAL:HG23	3:BD:193:VAL:HA	1.91	0.51
6:BG:122:PRO:HG2	6:BG:182:LYS:O	2.10	0.51
14:BS:34:HIS:O	14:BS:97:ARG:NH2	2.42	0.51
27:B5:16:ARG:NH1	27:B5:16:ARG:HG2	2.25	0.51
31:CA:444:C:H2'	31:CA:445:G:H8	1.76	0.51
31:CA:957:U:O2'	31:CA:959:A:N7	2.37	0.51
31:CA:968:A:C8	31:CA:1062:U:H4'	2.45	0.51
31:CA:1373:G:H8	31:CA:1373:G:O5'	1.94	0.51
32:CB:185:ILE:HG12	32:CB:185:ILE:O	2.10	0.51
39:CI:3:GLN:HG3	39:CI:20:ARG:HG2	1.91	0.51
41:CK:34:ASP:OD2	41:CK:38:ASN:N	2.42	0.51
46:CP:56:ALA:O	46:CP:60:LEU:HD23	2.10	0.51
47:CQ:13:ASP:C	47:CQ:15:MET:H	2.14	0.51
31:DA:1034:G:H2'	31:DA:1035:A:C8	2.45	0.51
31:DA:1292:U:O2'	31:DA:1293:G:H5'	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:DI:5:TYR:HD1	39:DI:17:VAL:O	1.93	0.51
1:AA:195:A:H61	1:AA:198:C:H3'	1.75	0.51
1:AA:833:U:O2	11:AP:55:ARG:NH2	2.44	0.51
1:AA:1012:U:C5	9:AN:28:THR:HG21	2.46	0.51
1:AA:1210:A:H5'	1:AA:1210:A:C8	2.43	0.51
1:AA:1445:A:C8	1:AA:1460:A:C5	2.99	0.51
1:AA:1636:C:H2'	1:AA:1637:A:C8	2.45	0.51
1:AA:2274:A:C5	1:AA:2276:G:C8	2.99	0.51
3:AD:132:PRO:HD3	3:AD:190:TYR:CZ	2.46	0.51
4:AE:120:TRP:CE3	4:AE:155:LYS:HD3	2.46	0.51
6:AG:44:GLY:HA2	6:AG:88:ILE:HG22	1.92	0.51
6:AG:46:ALA:HB1	6:AG:51:ARG:HA	1.92	0.51
11:AP:52:GLU:OE2	30:A8:57:ARG:NH1	2.44	0.51
1:BA:14:A:N6	1:BA:15:G:C2	2.79	0.51
1:BA:573:G:O2'	1:BA:574:C:H3'	2.11	0.51
1:BA:805:G:N2	1:BA:829:A:OP1	2.43	0.51
1:BA:2139:C:N4	1:BA:2152:G:H1	2.08	0.51
1:BA:2571:C:H5''	1:BA:2572:A:H5''	1.92	0.51
12:BQ:12:GLN:HG2	12:BQ:73:PRO:HD2	1.90	0.51
32:CB:51:LEU:O	32:CB:55:PHE:HD2	1.94	0.51
43:CM:52:GLU:HG3	43:CM:55:ARG:HH22	1.75	0.51
50:CT:26:ASN:OD1	50:CT:71:THR:HG23	2.10	0.51
31:DA:1030:C:N4	31:DA:1031:G:H1	2.04	0.51
31:DA:1059:C:H2'	31:DA:1060:C:H6	1.75	0.51
47:DQ:57:VAL:HG12	47:DQ:76:LEU:HA	1.93	0.51
1:AA:570:G:H2'	1:AA:2030:A:C5	2.46	0.51
1:AA:1530:C:HO2'	1:AA:1531:C:P	2.29	0.51
1:AA:2136:C:N4	1:AA:2155:G:N1	2.59	0.51
1:AA:2278:A:O2'	21:AZ:198:LYS:HD3	2.11	0.51
6:AG:62:LEU:O	6:AG:143:GLU:HG3	2.10	0.51
13:AR:13:HIS:HE1	13:AR:15:SER:OG	1.94	0.51
1:BA:107:C:H2'	1:BA:108:U:H6	1.76	0.51
1:BA:365:C:H2'	1:BA:366:C:O4'	2.11	0.51
1:BA:1022:G:C5	1:BA:1140:C:C4	2.99	0.51
1:BA:1753:G:H5''	15:BT:95:ARG:HE	1.76	0.51
1:BA:1913:A:H4'	1:BA:1914:C:O5'	2.11	0.51
31:CA:171:A:H2'	31:CA:172:A:H8	1.75	0.51
31:CA:1014:A:H1'	49:CS:34:TRP:HB2	1.92	0.51
37:CG:57:GLU:O	37:CG:61:VAL:HG23	2.10	0.51
46:CP:20:VAL:HG23	46:CP:35:LYS:HA	1.93	0.51
31:DA:624:C:O3'	46:DP:10:GLY:HA2	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:790:A:C6	31:DA:791:G:C6	2.99	0.51
31:DA:1030(C):G:H2'	31:DA:1030(D):A:C8	2.43	0.51
31:DA:1150:U:O2	40:DJ:39:PRO:HG2	2.10	0.51
31:DA:1202:G:H2'	31:DA:1203:C:O4'	2.10	0.51
31:DA:1279:A:OP1	40:DJ:9:ARG:NH1	2.43	0.51
32:DB:189:ASP:N	32:DB:189:ASP:OD1	2.44	0.51
46:DP:20:VAL:HG23	46:DP:35:LYS:HA	1.93	0.51
1:AA:1165:U:H2'	1:AA:1166:C:C6	2.46	0.51
1:AA:1359:A:H5'	1:AA:1359:A:N3	2.26	0.51
1:AA:1657:C:OP1	4:AE:136:ARG:N	2.42	0.51
1:AA:2345:G:N3	1:AA:2381:C:H2'	2.26	0.51
1:AA:2712:U:H2'	1:AA:2714:G:H5''	1.92	0.51
14:AS:11:LYS:O	14:AS:15:ARG:HB2	2.11	0.51
15:AT:65:LYS:HE3	15:AT:67:SER:HB2	1.92	0.51
1:BA:271(H):G:HO2'	1:BA:271(I):G:H8	1.55	0.51
1:BA:813:U:H2'	1:BA:814:C:C6	2.46	0.51
1:BA:848:G:N3	1:BA:933:A:H1'	2.25	0.51
1:BA:2364:C:H2'	1:BA:2365:G:O4'	2.10	0.51
16:BU:113:ALA:O	16:BU:117:GLN:HG2	2.10	0.51
21:BZ:101:PRO:O	21:BZ:102:LEU:HD12	2.11	0.51
31:CA:578:C:O2'	31:CA:728:A:N3	2.36	0.51
39:CI:28:VAL:O	39:CI:31:GLN:NE2	2.44	0.51
31:DA:233:C:H2'	31:DA:234:C:H6	1.76	0.51
31:DA:769:G:H4'	31:DA:1513:A:H4'	1.92	0.51
31:DA:932:C:H5'	37:DG:4:ARG:HG3	1.93	0.51
31:DA:1107:C:C4	31:DA:1108:G:C8	2.99	0.51
33:DC:15:THR:HG21	33:DC:181:ASN:HA	1.93	0.51
33:DC:199:LYS:HD3	33:DC:201:TYR:HE1	1.76	0.51
38:DH:86:ILE:HG12	38:DH:135:CYS:HA	1.93	0.51
1:AA:271(P):C:C2'	1:AA:271(Q):G:H5'	2.41	0.51
1:AA:769:G:O2'	1:AA:770:G:H5'	2.11	0.51
1:AA:774:A:H2'	1:AA:774:A:N3	2.26	0.51
1:AA:1319:G:C6	1:AA:1320:C:N4	2.79	0.51
1:AA:2695:C:H2'	1:AA:2696:U:C6	2.45	0.51
1:AA:2870:C:H2'	1:AA:2871:C:O4'	2.11	0.51
13:AR:55:ALA:HB2	13:AR:79:LEU:HD13	1.93	0.51
21:AZ:158:PRO:O	21:AZ:161:VAL:HG11	2.11	0.51
1:BA:764:A:O4'	3:BD:213:ARG:HG3	2.11	0.51
1:BA:1486:A:H2'	1:BA:1487:G:C8	2.45	0.51
1:BA:1913:A:N6	31:CA:1493:A:O5'	2.44	0.51
1:BA:2115:G:H2'	1:BA:2117:A:N6	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:2294:C:OP2	14:BS:13:ARG:NH1	2.44	0.51
1:BA:2865:U:C4	1:BA:2866:U:C4	2.99	0.51
17:BV:85:LYS:HB2	17:BV:85:LYS:NZ	2.26	0.51
31:CA:22:G:H4'	31:CA:885:G:C8	2.46	0.51
31:CA:911:U:H2'	31:CA:912:C:C6	2.46	0.51
31:CA:1129:C:H5''	31:CA:1139:G:O6	2.11	0.51
31:CA:1187:G:H5'	39:CI:113:LYS:HE2	1.93	0.51
34:CD:14:ARG:HA	34:CD:39:PRO:HB3	1.93	0.51
39:CI:9:ARG:HA	39:CI:13:ALA:O	2.11	0.51
31:DA:256:U:H2'	31:DA:257:G:H8	1.75	0.51
31:DA:531:U:O3'	31:DA:532:A:H4'	2.11	0.51
31:DA:1072:G:H2'	31:DA:1073:U:C6	2.45	0.51
31:DA:1123:A:H4'	40:DJ:37:PRO:HD2	1.92	0.51
31:DA:1270:C:H2'	31:DA:1271:G:H8	1.76	0.51
31:DA:1310:G:H5'	43:DM:77:ASN:ND2	2.26	0.51
34:DD:57:ARG:HB3	34:DD:206:PHE:HB2	1.92	0.51
49:DS:53:ASN:HB2	49:DS:77:THR:HA	1.91	0.51
1:AA:2790:A:H2'	1:AA:2790:A:N3	2.27	0.50
11:AP:33:ARG:NH1	11:AP:39:LYS:O	2.44	0.50
12:AQ:32:TYR:CZ	12:AQ:133:ARG:HG3	2.47	0.50
15:AT:61:PHE:CZ	15:AT:76:PHE:HB2	2.46	0.50
23:A1:85:LEU:HB3	23:A1:89:GLU:HG3	1.93	0.50
1:BA:902:C:H2'	1:BA:903:C:C6	2.47	0.50
1:BA:996:A:H4'	16:BU:91:ASP:OD1	2.11	0.50
1:BA:2206:G:O2'	1:BA:2207:G:OP1	2.29	0.50
1:BA:2698:U:H2'	1:BA:2699:C:C6	2.46	0.50
4:BE:120:TRP:CE3	4:BE:155:LYS:HD3	2.46	0.50
10:BO:63:VAL:HG12	10:BO:106:LEU:HD11	1.93	0.50
21:BZ:111:VAL:C	21:BZ:113:ALA:H	2.14	0.50
26:B4:26:SER:OG	26:B4:27:THR:N	2.44	0.50
31:CA:160:A:H2'	31:CA:161:A:O4'	2.11	0.50
31:CA:1391:U:H2'	31:CA:1392:G:H8	1.76	0.50
31:DA:17:U:H2'	31:DA:18:C:C6	2.45	0.50
31:DA:1010:G:H2'	31:DA:1011:G:C8	2.47	0.50
31:DA:1287:A:C2	31:DA:1353:G:H1'	2.47	0.50
50:DT:73:HIS:HB3	50:DT:74:LYS:HE2	1.93	0.50
1:AA:829:A:N7	1:AA:2247:A:O2'	2.41	0.50
1:AA:928:G:O5'	1:AA:928:G:H8	1.94	0.50
1:AA:1231:G:H2'	1:AA:1232:G:C8	2.46	0.50
1:AA:1547:C:H2'	1:AA:1548:C:C6	2.46	0.50
1:AA:1783:A:C2	1:AA:2587:A:C5	2.99	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2286:A:OP1	28:A6:29:ASN:ND2	2.41	0.50
2:AB:27:C:N4	2:AB:56:G:O6	2.43	0.50
4:AE:11:MET:HG2	4:AE:24:THR:HB	1.94	0.50
4:AE:71:GLY:HA2	4:AE:72:VAL:O	2.12	0.50
16:AU:49:HIS:HA	16:AU:52:ARG:HB3	1.92	0.50
21:AZ:160:GLY:CA	21:AZ:161:VAL:HB	2.37	0.50
25:A3:4:LEU:O	25:A3:36:VAL:HA	2.10	0.50
27:A5:36:CYS:O	27:A5:37:LYS:HD3	2.11	0.50
28:A6:44:ARG:HB3	28:A6:44:ARG:HH11	1.75	0.50
1:BA:2151:G:H2'	1:BA:2152:G:O4'	2.12	0.50
1:BA:2158:A:H4'	1:BA:2159:G:H5'	1.92	0.50
1:BA:2751:G:C5	7:BH:2:SER:N	2.79	0.50
19:BX:31:HIS:HD2	19:BX:33:LYS:H	1.59	0.50
19:BX:35:THR:HG22	19:BX:37:THR:H	1.76	0.50
21:BZ:141:VAL:O	21:BZ:144:LEU:HB2	2.12	0.50
31:CA:452:A:H62	31:CA:480:U:H3	1.58	0.50
35:CE:77:PRO:HD2	35:CE:142:LEU:HD22	1.93	0.50
31:DA:266:G:H5'	31:DA:268:C:H41	1.75	0.50
31:DA:375:U:C4	31:DA:376:G:N7	2.79	0.50
31:DA:717:C:H6	31:DA:717:C:H5''	1.76	0.50
31:DA:1255:G:H2'	31:DA:1258:G:H21	1.76	0.50
31:DA:1308:U:H2'	31:DA:1309:G:H8	1.75	0.50
32:DB:47:THR:HG23	32:DB:202:PRO:HG2	1.93	0.50
38:DH:82:HIS:CE1	38:DH:84:ARG:HG2	2.46	0.50
39:DI:14:VAL:O	39:DI:65:VAL:HA	2.10	0.50
40:DJ:67:THR:O	40:DJ:67:THR:OG1	2.25	0.50
1:AA:1153:C:OP1	16:AU:92:ARG:NH1	2.43	0.50
1:AA:1173:G:O2'	1:AA:1174:A:O4'	2.29	0.50
1:AA:1914:C:OP1	1:AA:1914:C:H3'	2.11	0.50
1:AA:2030:A:H5''	1:AA:2031:A:OP1	2.10	0.50
1:AA:2318:G:N2	14:AS:3:ARG:HD2	2.26	0.50
1:AA:2751:G:C5	7:AH:2:SER:N	2.80	0.50
12:AQ:133:ARG:HG2	12:AQ:134:ARG:H	1.77	0.50
21:AZ:45:ASP:OD1	21:AZ:49:ARG:HD2	2.10	0.50
1:BA:1505:C:H2'	1:BA:1506:C:C6	2.43	0.50
1:BA:1823:G:OP1	3:BD:54:ARG:NH1	2.44	0.50
7:BH:25:LYS:NZ	7:BH:32:GLU:OE2	2.26	0.50
12:BQ:32:TYR:CE2	12:BQ:133:ARG:HG3	2.47	0.50
17:BV:42:GLY:O	17:BV:43:GLU:HG2	2.12	0.50
31:CA:236:G:H5''	47:CQ:42:TYR:OH	2.11	0.50
31:CA:390:C:O3'	46:CP:28:ARG:NH2	2.39	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:477:A:H2'	31:CA:479:C:H6	1.76	0.50
31:CA:598:U:H2'	31:CA:599:C:C6	2.46	0.50
37:CG:51:GLN:O	37:CG:55:GLY:HA2	2.11	0.50
31:DA:391:G:C6	31:DA:392:G:C5	3.00	0.50
34:DD:128:VAL:O	34:DD:130:GLY:N	2.44	0.50
42:DL:32:PHE:HE1	42:DL:86:ARG:HG3	1.75	0.50
1:AA:86:C:H4'	1:AA:104:U:H1'	1.92	0.50
1:AA:729:G:H2'	1:AA:1775:U:H1'	1.93	0.50
1:AA:782:A:N7	3:AD:221:VAL:HG11	2.27	0.50
1:AA:1332:G:N2	1:AA:1610:A:H8	2.07	0.50
1:AA:1342:A:H2	1:AA:1396:U:HO2'	1.58	0.50
1:AA:1637:A:H4'	1:AA:2711:A:O2'	2.11	0.50
1:AA:2115:G:H2'	1:AA:2117:A:N6	2.25	0.50
1:BA:296:C:H2'	1:BA:297:C:H6	1.75	0.50
1:BA:1405:U:H2'	1:BA:1406:U:C6	2.47	0.50
1:BA:1480:G:C6	1:BA:1481:U:N3	2.80	0.50
3:BD:108:PRO:HG2	3:BD:111:LEU:HG	1.93	0.50
9:BN:20:GLY:HA2	9:BN:61:ARG:HH11	1.77	0.50
21:BZ:25:PRO:O	21:BZ:85:HIS:HA	2.10	0.50
28:B6:8:LYS:HD3	30:B8:34:TRP:CD2	2.46	0.50
31:CA:175:C:H2'	31:CA:176:C:C6	2.46	0.50
31:CA:391:G:C6	31:CA:392:G:C5	3.00	0.50
31:CA:663:A:H2'	31:CA:664:G:O4'	2.11	0.50
31:CA:677:U:H2'	31:CA:678:U:C6	2.45	0.50
31:CA:973:G:H3'	31:CA:974:A:H5''	1.93	0.50
31:CA:1219:U:OP1	44:CN:19:ARG:NH1	2.37	0.50
31:CA:1251:A:O2'	31:CA:1369:C:O2'	2.23	0.50
33:CC:20:SER:OG	33:CC:36:ASP:OD1	2.29	0.50
33:CC:34:LEU:O	33:CC:38:ARG:HG3	2.12	0.50
48:CR:52:PRO:O	48:CR:56:THR:HG23	2.12	0.50
31:DA:1512:U:H2'	31:DA:1513:A:C8	2.46	0.50
51:DU:12:LYS:HB3	51:DU:17:THR:O	2.11	0.50
1:AA:530:G:C5	1:AA:2022:U:H5''	2.47	0.50
1:AA:1198:U:H2'	1:AA:1199:U:C6	2.46	0.50
1:AA:1223:G:N2	1:AA:1226:A:OP2	2.44	0.50
4:AE:96:PHE:O	4:AE:175:VAL:HG11	2.12	0.50
8:AI:116:LEU:CD1	8:AI:119:PRO:HA	2.41	0.50
15:AT:41:ARG:CZ	31:DA:346:G:H5'	2.42	0.50
7:BH:11:VAL:HG13	7:BH:15:VAL:HG22	1.94	0.50
8:BI:110:GLU:N	8:BI:130:TYR:OH	2.36	0.50
14:BS:46:VAL:HG12	14:BS:48:LEU:HD12	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:187:C:H2'	31:CA:188:C:H6	1.76	0.50
31:CA:985:C:C2	31:CA:1221:G:N2	2.79	0.50
31:CA:1446:U:H4'	31:CA:1447:A:C5	2.46	0.50
32:CB:55:PHE:CE1	32:CB:218:ALA:HA	2.46	0.50
33:CC:134:ILE:HD11	33:CC:153:VAL:HB	1.93	0.50
39:CI:8:GLY:HA3	39:CI:76:ALA:O	2.11	0.50
45:CO:78:TYR:O	45:CO:82:ILE:HG12	2.12	0.50
31:DA:674:G:H2'	31:DA:675:A:C8	2.44	0.50
31:DA:1060:C:O2'	31:DA:1061:G:H5'	2.11	0.50
31:DA:1346:A:O2'	37:DG:10:ARG:NH1	2.40	0.50
32:DB:212:GLN:NE2	32:DB:235:SER:HB3	2.27	0.50
46:DP:20:VAL:HG21	46:DP:32:TYR:CG	2.46	0.50
1:AA:61:G:H5'	24:A2:50:ILE:HD12	1.93	0.50
1:AA:107:C:H2'	1:AA:108:U:C6	2.46	0.50
1:AA:222:A:N3	1:AA:224:G:H1'	2.27	0.50
1:AA:511:U:H5''	1:AA:512:G:OP2	2.11	0.50
1:AA:764:A:O4'	3:AD:213:ARG:HG3	2.12	0.50
1:AA:815:C:C2	1:AA:1193:G:C2	2.99	0.50
1:AA:1049:C:H4'	1:AA:1050:A:OP1	2.12	0.50
1:AA:1665:A:H2'	1:AA:1666:G:O4'	2.11	0.50
1:AA:2738:A:C2	1:AA:2739:U:H1'	2.46	0.50
1:AA:2784:C:H1'	4:AE:37:ARG:NH1	2.26	0.50
25:A3:50:VAL:HB	25:A3:53:LEU:HD12	1.93	0.50
1:BA:8:A:H2'	1:BA:9:U:H6	1.76	0.50
1:BA:459:U:H4'	29:B7:40:TRP:CZ3	2.47	0.50
1:BA:2178:C:H2'	1:BA:2179:C:O4'	2.12	0.50
1:BA:2533:A:H2'	1:BA:2534:A:O4'	2.12	0.50
12:BQ:42:ILE:HD13	12:BQ:97:VAL:HG21	1.94	0.50
31:CA:91:C:O2'	31:CA:92:C:H5'	2.11	0.50
31:CA:391:G:O3'	46:CP:8:ARG:NH2	2.44	0.50
31:CA:1169:A:H2'	31:CA:1170:A:C8	2.47	0.50
31:CA:1258:G:H2'	31:CA:1259:C:C6	2.46	0.50
31:CA:1378:C:C2	31:CA:1379:G:H1'	2.47	0.50
50:CT:53:LEU:O	50:CT:57:ARG:HG3	2.11	0.50
31:DA:1490:C:H2'	31:DA:1491:G:C8	2.45	0.50
1:AA:642:G:H21	1:AA:646:A:H2	1.58	0.50
1:AA:1109:C:H5	1:AA:1110:G:C5	2.30	0.50
1:AA:1131:G:O6	1:AA:2040:C:H1'	2.11	0.50
1:AA:1876:A:H2'	1:AA:1877:A:C8	2.46	0.50
1:AA:1927:A:H2'	1:AA:1928:A:C8	2.47	0.50
1:AA:2505:G:O6	1:AA:2576:G:H2'	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2740:A:C6	1:AA:2741:A:C6	3.00	0.50
2:AB:11:C:H3'	2:AB:12:C:C6	2.46	0.50
3:AD:242:ARG:HG2	3:AD:246:PRO:HG3	1.93	0.50
13:AR:38:VAL:O	13:AR:42:LYS:HG3	2.11	0.50
16:AU:105:VAL:O	16:AU:108:GLU:HB2	2.12	0.50
18:AW:76:VAL:HG22	18:AW:103:ILE:HG23	1.93	0.50
21:AZ:104:PHE:HB3	21:AZ:141:VAL:HG21	1.93	0.50
27:A5:53:ALA:N	27:A5:54:GLY:HA2	2.27	0.50
30:A8:26:LYS:HE3	30:A8:48:PHE:CD1	2.47	0.50
1:BA:107:C:H2'	1:BA:108:U:C6	2.46	0.50
1:BA:414:C:O2'	1:BA:415:A:H5'	2.12	0.50
1:BA:1568:G:H5'	3:BD:60:ARG:HA	1.93	0.50
1:BA:1914:C:O4'	1:BA:1914:C:O2	2.29	0.50
1:BA:2492:U:H2'	1:BA:2493:U:H6	1.76	0.50
1:BA:2502:G:H5''	1:BA:2503:A:H5''	1.94	0.50
1:BA:2627:G:O2'	1:BA:2781:A:N1	2.33	0.50
4:BE:115:GLY:O	4:BE:119:ARG:HB2	2.12	0.50
6:BG:122:PRO:HG3	6:BG:180:PHE:HB3	1.94	0.50
31:CA:1468:A:H2'	31:CA:1469:G:O4'	2.11	0.50
36:CF:36:ARG:HD3	36:CF:37:VAL:O	2.12	0.50
36:CF:49:ALA:N	48:CR:77:GLY:O	2.34	0.50
39:CI:31:GLN:HG2	39:CI:36:TYR:HB2	1.94	0.50
40:CJ:38:ILE:HG13	40:CJ:38:ILE:O	2.11	0.50
40:CJ:54:PHE:CD2	40:CJ:55:LYS:HD2	2.44	0.50
39:DI:44:VAL:N	39:DI:45:ALA:HB2	2.26	0.50
41:DK:95:ILE:HD12	41:DK:108:ILE:HD13	1.92	0.50
44:DN:24:CYS:HG	44:DN:40:CYS:CB	2.23	0.50
1:AA:1009:A:P	9:AN:37:LYS:HZ1	2.34	0.50
1:AA:2280:G:O6	22:A0:14:ARG:HD3	2.11	0.50
1:AA:2394:C:OP1	30:A8:30:ARG:NH1	2.42	0.50
1:AA:2408:U:H2'	1:AA:2409:G:C8	2.46	0.50
1:AA:2572:A:C8	4:AE:144:ARG:HD2	2.47	0.50
2:AB:13:A:N1	2:AB:69:G:O2'	2.37	0.50
5:AF:88:VAL:HG21	5:AF:91:GLY:HA3	1.93	0.50
1:BA:334:C:OP1	1:BA:335:C:N4	2.44	0.50
1:BA:652(R):C:H2'	1:BA:652(R):C:OP2	2.12	0.50
1:BA:2773:C:H2'	1:BA:2774:C:H6	1.75	0.50
5:BF:150:GLY:HA2	5:BF:172:TRP:CE3	2.46	0.50
11:BP:38:GLN:O	11:BP:38:GLN:HG3	2.11	0.50
29:B7:22:MET:HA	29:B7:28:ARG:HG2	1.94	0.50
31:CA:397:A:C6	31:CA:548:G:N7	2.80	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:628:G:H2'	31:CA:629:G:C8	2.47	0.50
31:CA:1034:G:H2'	31:CA:1035:A:C8	2.47	0.50
31:CA:1065:U:H4'	31:CA:1066:C:O5'	2.12	0.50
31:CA:1288:A:N1	31:CA:1371:G:H1'	2.26	0.50
38:CH:6:ILE:O	38:CH:10:LEU:HG	2.12	0.50
38:CH:120:THR:H	38:CH:123:GLU:HB2	1.77	0.50
31:DA:1233:G:H2'	31:DA:1234:C:C6	2.46	0.50
32:DB:210:SER:O	32:DB:214:ILE:HG12	2.11	0.50
33:DC:47:LEU:HB3	33:DC:52:LEU:HD13	1.93	0.50
37:DG:44:TYR:HA	37:DG:47:CYS:HB2	1.93	0.50
38:DH:20:TYR:CE1	38:DH:76:PRO:HG2	2.47	0.50
43:DM:5:ALA:HA	43:DM:61:GLU:HG2	1.94	0.50
1:AA:23:G:OP1	1:AA:504:U:N3	2.34	0.50
1:AA:1827:C:H2'	1:AA:1828:G:H5'	1.93	0.50
1:AA:2572:A:N7	4:AE:144:ARG:HD2	2.27	0.50
2:AB:48:A:C2	2:AB:49:C:C2	3.00	0.50
13:AR:84:ALA:HB3	13:AR:85:PRO:HD3	1.94	0.50
20:AY:5:MET:HE1	20:AY:32:PRO:HA	1.94	0.50
21:AZ:70:LEU:HG	21:AZ:91:LEU:HD21	1.94	0.50
1:BA:2100:G:C6	1:BA:2190:G:C6	3.00	0.50
15:BT:20:PRO:HD2	15:BT:86:ILE:HB	1.93	0.50
17:BV:22:VAL:HG23	17:BV:23:GLU:O	2.12	0.50
31:CA:44:G:H2'	31:CA:45:U:O4'	2.12	0.50
31:CA:565:U:H3'	31:CA:566:G:H2'	1.94	0.50
31:CA:1131:G:H2'	31:CA:1132:C:H6	1.76	0.50
31:CA:1361:G:H2'	31:CA:1362:C:O4'	2.12	0.50
32:CB:47:THR:HA	32:CB:202:PRO:HG2	1.94	0.50
48:CR:73:ALA:HB3	48:CR:79:LEU:HD12	1.93	0.50
31:DA:1014:A:H1'	49:DS:34:TRP:HB2	1.94	0.50
31:DA:1468:A:H2'	31:DA:1469:G:O4'	2.11	0.50
1:AA:208:C:H2'	1:AA:209:C:C6	2.47	0.49
1:AA:2319:G:C2	14:AS:3:ARG:HA	2.47	0.49
1:AA:2328:A:H2'	1:AA:2329:G:C8	2.47	0.49
1:AA:2468:G:N2	1:AA:2481:G:H1'	2.27	0.49
5:AF:33:LEU:HB3	11:AP:6:LEU:HD21	1.93	0.49
1:BA:271(Q):G:O2'	1:BA:271(R):G:OP2	2.21	0.49
5:BF:154:VAL:HB	5:BF:173:VAL:HG22	1.94	0.49
7:BH:164:TYR:HB2	7:BH:167:GLU:HB2	1.93	0.49
28:B6:25:LYS:NZ	28:B6:51:GLU:OE2	2.44	0.49
31:CA:240:C:H2'	31:CA:241:C:C6	2.47	0.49
31:CA:246:A:N1	31:CA:278:G:O2'	2.40	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:501:C:H2'	31:CA:502:G:C8	2.47	0.49
31:CA:649:G:H2'	31:CA:650:G:H8	1.76	0.49
31:CA:1124:G:H2'	31:CA:1126:U:O4	2.12	0.49
31:CA:1316:G:N2	31:CA:1318:A:H3'	2.27	0.49
32:CB:208:ILE:HA	32:CB:211:ILE:HD12	1.93	0.49
37:CG:108:ALA:HB1	37:CG:120:ILE:HD13	1.92	0.49
40:CJ:51:ARG:CZ	40:CJ:61:GLU:HB3	2.42	0.49
31:DA:658:G:C6	31:DA:659:U:C4	3.00	0.49
31:DA:1090:U:H2'	31:DA:1091:U:C6	2.47	0.49
31:DA:1097:C:H2'	31:DA:1098:C:H6	1.77	0.49
43:DM:33:ALA:HA	43:DM:59:TYR:HE1	1.77	0.49
1:AA:19:C:H2'	1:AA:20:C:C6	2.47	0.49
1:AA:176:G:O2'	1:AA:177:G:H5'	2.12	0.49
1:AA:265:A:H1'	1:AA:266:G:O4'	2.13	0.49
1:AA:311:A:C6	1:AA:328:U:C4	3.01	0.49
1:AA:411:G:OP2	1:AA:2406:U:O2'	2.27	0.49
1:AA:708:C:H42	1:AA:723:G:H1	1.58	0.49
1:AA:2390:U:O2'	1:AA:2391:G:H5'	2.12	0.49
1:AA:2728:U:H2'	1:AA:2729:G:C8	2.47	0.49
5:AF:109:GLY:O	5:AF:113:ALA:N	2.37	0.49
5:AF:179:GLU:H	5:AF:179:GLU:CD	2.15	0.49
21:AZ:72:ARG:NH2	21:AZ:97:GLU:O	2.46	0.49
1:BA:271(E):U:H2'	1:BA:271(F):C:C6	2.47	0.49
1:BA:581:C:OP1	16:BU:33:ARG:HG3	2.12	0.49
1:BA:2307:G:H4'	1:BA:2308:G:C4	2.47	0.49
11:BP:50:ARG:HD3	30:B8:7:HIS:CD2	2.47	0.49
22:B0:51:VAL:HG22	22:B0:81:VAL:HG23	1.92	0.49
31:CA:382:A:H2'	31:CA:383:A:H8	1.76	0.49
31:CA:730:G:C5	31:CA:731:G:H1'	2.47	0.49
31:CA:1084:G:H2'	31:CA:1085:U:C6	2.47	0.49
31:CA:1399:C:H4'	31:CA:1400:C:H5''	1.93	0.49
34:CD:88:VAL:O	34:CD:92:VAL:HG23	2.13	0.49
36:CF:75:LEU:O	36:CF:79:LEU:HG	2.12	0.49
31:DA:4:U:H5	38:DH:105:ARG:HD3	1.77	0.49
31:DA:857:C:H2'	31:DA:858:G:O4'	2.12	0.49
31:DA:1144:G:N2	31:DA:1146:A:H62	2.10	0.49
31:DA:1513:A:H2'	31:DA:1514:C:C6	2.47	0.49
33:DC:150:LYS:HD3	33:DC:169:ALA:HB2	1.94	0.49
39:DI:37:PHE:HB3	39:DI:43:ALA:HB1	1.93	0.49
1:AA:296:C:H2'	1:AA:297:C:C6	2.47	0.49
7:AH:87:LEU:HB2	7:AH:131:VAL:HB	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AN:96:GLU:HB2	9:AN:122:VAL:HG12	1.94	0.49
20:AY:12:THR:OG1	20:AY:26:LYS:HG2	2.12	0.49
1:BA:672:C:C2'	1:BA:673:C:H5'	2.42	0.49
1:BA:819:A:C4	1:BA:1189:A:C2	3.01	0.49
8:BI:104:GLN:CB	8:BI:105:HIS:HD2	2.25	0.49
31:CA:618:C:H5''	31:CA:619:U:H5''	1.93	0.49
31:CA:1012:U:H2'	31:CA:1013:G:C8	2.48	0.49
31:CA:1218:C:H2'	31:CA:1219:U:C6	2.47	0.49
31:CA:1271:G:H5'	31:CA:1314:C:H5'	1.94	0.49
33:CC:155:GLY:O	33:CC:157:ILE:N	2.43	0.49
31:DA:819:A:H4'	31:DA:820:U:OP2	2.11	0.49
31:DA:1315:U:O2'	31:DA:1360:A:N3	2.31	0.49
39:DI:40:LEU:HD11	39:DI:70:LYS:HB3	1.95	0.49
1:AA:26:G:H1'	1:AA:514:A:N6	2.28	0.49
1:AA:469:G:O6	29:A7:37:LYS:NZ	2.38	0.49
1:AA:590:A:H2'	1:AA:591:C:C6	2.46	0.49
1:AA:768:G:O2'	1:AA:1379:A:N1	2.33	0.49
1:AA:830:G:H4'	1:AA:831:G:OP2	2.12	0.49
1:AA:858:U:O2	1:AA:2268:A:H2'	2.13	0.49
1:AA:2433:A:H5''	1:AA:2434:A:OP1	2.12	0.49
1:AA:2748:A:C6	1:AA:2749:A:C5	3.00	0.49
10:AO:7:TYR:CE1	10:AO:20:MET:HB2	2.48	0.49
13:AR:21:TYR:CZ	13:AR:43:GLU:HG2	2.47	0.49
1:BA:143:G:H4'	19:BX:35:THR:HG21	1.95	0.49
1:BA:484:C:H2'	1:BA:485:C:C6	2.47	0.49
1:BA:1801:G:OP2	3:BD:154:LYS:NZ	2.41	0.49
1:BA:2315:G:H2'	1:BA:2316:C:C6	2.47	0.49
4:BE:9:VAL:HG22	4:BE:25:VAL:HB	1.93	0.49
8:BI:79:ILE:N	8:BI:143:SER:O	2.39	0.49
31:CA:22:G:H2'	31:CA:23:C:C6	2.47	0.49
31:CA:262:A:C6	31:CA:263:A:C6	3.00	0.49
31:CA:374:A:C6	31:CA:375:U:C4	3.00	0.49
31:CA:850:U:H2'	31:CA:851:G:H5''	1.94	0.49
33:CC:156:ARG:HG2	33:CC:160:ALA:O	2.12	0.49
31:DA:96:U:HO2'	31:DA:97:G:P	2.34	0.49
31:DA:545:C:O2'	31:DA:549:C:OP1	2.28	0.49
31:DA:709:G:H2'	31:DA:710:G:H8	1.77	0.49
31:DA:737:A:H2'	31:DA:738:C:H6	1.77	0.49
31:DA:741:G:H2'	31:DA:742:G:O4'	2.12	0.49
31:DA:1165:C:H2'	31:DA:1166:G:C8	2.47	0.49
1:AA:342:G:H2'	1:AA:343:C:H6	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:612:C:H2'	1:AA:613:G:O4'	2.12	0.49
1:AA:1668:A:H4'	1:AA:1669:A:O5'	2.12	0.49
1:AA:2097:C:H2'	1:AA:2098:U:C6	2.48	0.49
1:AA:2228:G:C6	1:AA:2229:C:C4	3.00	0.49
1:AA:2263:C:H2'	1:AA:2264:C:C6	2.47	0.49
1:AA:2273:A:H2'	1:AA:2274:A:C8	2.47	0.49
2:AB:11:C:H3'	2:AB:12:C:H6	1.78	0.49
2:AB:48:A:C2	2:AB:49:C:N3	2.81	0.49
7:AH:149:ARG:HH22	7:AH:154:PRO:HG2	1.78	0.49
8:AI:111:PRO:C	8:AI:114:LEU:HB2	2.33	0.49
15:AT:110:ILE:O	15:AT:114:LEU:N	2.44	0.49
1:BA:1429:G:H2'	1:BA:1430:C:H6	1.75	0.49
1:BA:1949:G:C6	1:BA:1950:G:C6	3.00	0.49
1:BA:2207:G:HO2'	1:BA:2208:A:P	2.34	0.49
1:BA:2790:A:OP1	1:BA:2790:A:H4'	2.10	0.49
6:BG:138:GLN:O	6:BG:144:ILE:HD12	2.13	0.49
15:BT:66:VAL:HA	15:BT:71:GLY:HA2	1.94	0.49
31:CA:41:G:C6	31:CA:402:G:C6	3.00	0.49
31:CA:185:A:H2'	31:CA:186:C:H6	1.78	0.49
31:CA:328:C:H4'	31:CA:329:A:H5'	1.92	0.49
31:CA:666:G:H5'	31:CA:726:C:H1'	1.94	0.49
31:CA:1001:A:N6	31:CA:1001(A):G:O6	2.46	0.49
31:CA:1030(D):A:H62	31:CA:1031:G:N2	2.09	0.49
31:CA:1237:C:H4'	31:CA:1334:G:N2	2.28	0.49
33:CC:150:LYS:HB2	33:CC:169:ALA:HB2	1.95	0.49
34:CD:127:THR:OG1	34:CD:128:VAL:N	2.45	0.49
39:CI:5:TYR:OH	39:CI:16:ARG:HG2	2.12	0.49
31:DA:102:G:H2'	31:DA:103:C:C6	2.46	0.49
31:DA:503:C:OP2	42:DL:116:SER:HB3	2.12	0.49
31:DA:529:G:HO2'	31:DA:533:A:N6	2.10	0.49
31:DA:716:A:C6	31:DA:717:C:N3	2.81	0.49
31:DA:991:U:O4	31:DA:1212:U:O2'	2.16	0.49
33:DC:36:ASP:O	33:DC:40:ARG:HG3	2.12	0.49
45:DO:18:PHE:HB2	45:DO:19:PRO:HD2	1.94	0.49
1:AA:207:A:H2'	1:AA:208:C:O4'	2.11	0.49
1:AA:248:G:O2'	1:AA:2432:A:OP1	2.22	0.49
1:AA:300:A:H3'	20:AY:84:ARG:HH21	1.78	0.49
1:AA:963:U:H1'	1:AA:2250:G:O6	2.12	0.49
1:AA:1420:U:HO2'	1:AA:1421:G:P	2.34	0.49
1:AA:1479:G:C6	1:AA:1480:G:C5	3.00	0.49
1:AA:1491:G:O2'	3:AD:101:GLU:HB2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2697:G:H2'	1:AA:2698:U:O4'	2.13	0.49
7:AH:24:VAL:HG13	7:AH:37:VAL:HG21	1.94	0.49
1:BA:1530:C:HO2'	1:BA:1531:C:P	2.32	0.49
1:BA:2663:G:C6	1:BA:2664:G:C4	3.01	0.49
7:BH:88:LEU:CD2	7:BH:165:ALA:HA	2.43	0.49
14:BS:96:GLY:N	14:BS:99:LYS:H	2.05	0.49
31:CA:137:C:O2'	31:CA:138:G:H5'	2.13	0.49
31:CA:561:U:HO2'	31:CA:562:C:P	2.36	0.49
31:CA:1157:A:C6	31:CA:1180:A:C6	3.01	0.49
31:CA:1237:C:H2'	31:CA:1336:C:H5	1.76	0.49
31:CA:1318:A:H4'	49:CS:10:PHE:CE2	2.48	0.49
32:CB:79:ASP:O	32:CB:82:ARG:N	2.45	0.49
33:CC:11:ARG:HD3	33:CC:15:THR:HB	1.94	0.49
43:CM:66:LEU:H	43:CM:70:LEU:HD12	1.77	0.49
50:CT:36:LEU:HD12	50:CT:55:ILE:HG23	1.95	0.49
31:DA:109:A:H2'	31:DA:326:G:N2	2.27	0.49
31:DA:240:C:H2'	31:DA:241:C:C6	2.46	0.49
31:DA:968:A:C8	31:DA:1062:U:H4'	2.48	0.49
31:DA:1070:U:H2'	31:DA:1071:C:C6	2.48	0.49
31:DA:1174:G:C2	31:DA:1175:G:C8	3.00	0.49
31:DA:1308:U:H2'	31:DA:1309:G:C8	2.47	0.49
50:DT:26:ASN:OD1	50:DT:71:THR:HG23	2.12	0.49
1:AA:565:C:H4'	1:AA:1253:A:C6	2.47	0.49
1:AA:603:A:H4'	1:AA:604:G:H5'	1.95	0.49
1:AA:620:G:N3	1:AA:620:G:H5'	2.27	0.49
1:AA:688:U:O2'	1:AA:786:C:O2'	2.26	0.49
1:AA:833:U:H2'	1:AA:834:C:C6	2.47	0.49
1:AA:2282:G:H4'	1:AA:2283:C:O5'	2.11	0.49
1:AA:2659:G:OP1	7:AH:158:HIS:NE2	2.42	0.49
10:AO:64:ARG:HB2	10:AO:83:ALA:HB3	1.94	0.49
13:AR:54:LEU:O	13:AR:57:ARG:HB2	2.12	0.49
17:AV:2:PHE:CZ	17:AV:41:GLY:HA3	2.48	0.49
1:BA:603:A:H4'	1:BA:604:G:H5'	1.95	0.49
1:BA:2164:C:H3'	1:BA:2165:G:C8	2.44	0.49
3:BD:238:GLY:O	3:BD:239:ARG:HB2	2.11	0.49
4:BE:101:ARG:NH1	4:BE:169:ASN:O	2.44	0.49
5:BF:107:LYS:HE2	5:BF:208:GLY:N	2.28	0.49
31:CA:232:G:H1'	31:CA:262:A:N1	2.27	0.49
31:CA:1027:C:H5	31:CA:1029:C:N4	2.11	0.49
31:CA:1096:C:H2'	31:CA:1097:C:C6	2.47	0.49
40:CJ:11:PHE:HB3	44:CN:55:GLY:HA3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:CS:32:LYS:HD2	49:CS:57:HIS:HD2	1.78	0.49
31:DA:45:U:H2'	31:DA:46:G:C8	2.48	0.49
31:DA:60:A:P	31:DA:60:A:H8	2.36	0.49
31:DA:583:A:H2'	31:DA:584:G:O4'	2.12	0.49
31:DA:671:G:H2'	31:DA:672:U:O4'	2.12	0.49
31:DA:677:U:H3	31:DA:713:G:N2	1.95	0.49
31:DA:788:U:H2'	31:DA:789:U:O4'	2.12	0.49
31:DA:815:A:N7	31:DA:1509:C:O2'	2.37	0.49
31:DA:1030(B):C:H2'	31:DA:1030(C):G:H5'	1.95	0.49
31:DA:1233:G:H2'	31:DA:1234:C:H6	1.77	0.49
32:DB:54:THR:O	32:DB:58:ILE:HG13	2.13	0.49
33:DC:5:ILE:HD12	33:DC:6:HIS:H	1.78	0.49
38:DH:25:ASP:OD2	38:DH:60:ARG:HG2	2.12	0.49
42:DL:119:LYS:HB2	42:DL:120:TYR:CD1	2.48	0.49
1:AA:311:A:C8	1:AA:332:A:N7	2.81	0.49
1:AA:475:U:C4	1:AA:481:G:O6	2.66	0.49
1:AA:652(E):G:H2'	1:AA:652(F):G:C8	2.48	0.49
1:AA:1191:G:H2'	1:AA:1192:G:O4'	2.12	0.49
4:AE:176:ILE:HB	4:AE:181:LEU:HB2	1.95	0.49
7:AH:40:GLU:OE2	7:AH:60:ARG:NH1	2.46	0.49
14:AS:59:LYS:H	14:AS:60:GLY:HA3	1.77	0.49
1:BA:363(B):G:H2'	1:BA:363(C):G:C8	2.47	0.49
1:BA:534:U:H2'	1:BA:535:C:C6	2.48	0.49
1:BA:620:G:N3	1:BA:620:G:H2'	2.27	0.49
1:BA:620:G:N3	1:BA:620:G:H5'	2.27	0.49
1:BA:1165:U:H2'	1:BA:1166:C:C6	2.47	0.49
1:BA:2305:A:H5''	6:BG:134:GLY:HA3	1.95	0.49
31:CA:64:G:H4'	31:CA:65:U:H3'	1.95	0.49
31:CA:78:G:N2	31:CA:91:C:N3	2.52	0.49
31:CA:601:C:H2'	31:CA:602:A:H8	1.76	0.49
46:CP:49:LEU:HD21	46:CP:77:ALA:HB2	1.94	0.49
46:CP:72:ARG:HH21	46:CP:73:LEU:HD21	1.77	0.49
48:CR:31:LEU:H	48:CR:31:LEU:HD23	1.78	0.49
49:CS:23:ASN:HA	49:CS:27:GLU:HB2	1.94	0.49
31:DA:3:G:H2'	31:DA:3:G:N3	2.27	0.49
31:DA:743:U:H2'	31:DA:744:C:C6	2.48	0.49
31:DA:1001:A:H2'	31:DA:1001(A):G:C8	2.48	0.49
31:DA:1182:G:H5'	31:DA:1184:G:H5''	1.94	0.49
31:DA:1236:A:O2'	31:DA:1304:G:H4'	2.12	0.49
32:DB:179:LYS:HA	38:DH:72:PRO:HG3	1.95	0.49
33:DC:134:ILE:HG22	33:DC:168:ALA:HB3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DJ:16:LEU:HD21	40:DJ:70:ARG:HG3	1.94	0.49
44:DN:23:ARG:CG	44:DN:24:CYS:H	2.19	0.49
1:AA:628:G:H2'	1:AA:629:G:C8	2.47	0.49
1:AA:892:G:H2'	1:AA:893:C:H6	1.77	0.49
1:AA:1689:A:OP2	1:AA:1698:A:N6	2.46	0.49
1:AA:1913:A:H4'	1:AA:1914:C:H5''	1.95	0.49
1:AA:1948:G:N3	31:DA:1418:A:H2	2.11	0.49
2:AB:90:A:N7	2:AB:91:C:H1'	2.27	0.49
7:AH:8:PRO:HB3	7:AH:51:ARG:HD3	1.95	0.49
9:AN:71:ILE:HG21	9:AN:84:LYS:HB3	1.93	0.49
12:AQ:30:GLY:HA2	12:AQ:107:ALA:HB2	1.94	0.49
29:A7:25:PRO:HA	29:A7:28:ARG:CZ	2.43	0.49
1:BA:84:A:H5''	20:BY:8:LYS:HE3	1.94	0.49
1:BA:203:C:H3'	1:BA:204:A:H5''	1.94	0.49
1:BA:435:C:H2'	1:BA:436:C:H5'	1.95	0.49
1:BA:1914:C:O2'	1:BA:1915:U:OP2	2.30	0.49
1:BA:1918:A:O2'	1:BA:1920:C:N4	2.46	0.49
1:BA:2689:U:H4'	1:BA:2690:C:OP2	2.12	0.49
5:BF:123:LEU:HD13	5:BF:192:LEU:HB3	1.95	0.49
24:B2:32:LEU:HD22	24:B2:36:ARG:HH11	1.76	0.49
31:CA:277:C:H2'	31:CA:278:G:H8	1.77	0.49
31:CA:448:A:H2'	31:CA:449:C:C6	2.48	0.49
31:CA:1005:A:H1'	31:CA:1036:G:H22	1.78	0.49
31:CA:1250:A:H2	31:CA:1370:G:H1'	1.77	0.49
33:CC:22:TRP:CD1	33:CC:59:ARG:HB2	2.47	0.49
35:CE:8:GLU:OE2	35:CE:63:ARG:NH2	2.46	0.49
45:CO:23:GLY:O	45:CO:27:VAL:HB	2.13	0.49
31:DA:189(B):C:H2'	31:DA:189(C):C:C6	2.48	0.49
31:DA:370:C:H2'	31:DA:371:G:C8	2.48	0.49
31:DA:542:G:P	34:DD:10:ARG:HH22	2.36	0.49
34:DD:128:VAL:HA	34:DD:145:GLU:O	2.12	0.49
45:DO:33:THR:HG21	45:DO:85:LEU:HD22	1.94	0.49
46:DP:51:VAL:HG12	46:DP:53:VAL:H	1.78	0.49
1:AA:1014:U:H2'	1:AA:1015:G:H8	1.77	0.49
1:AA:1449:A:H5'	1:AA:1450:G:OP2	2.13	0.49
1:AA:2199:A:H3'	1:AA:2200:C:H6	1.78	0.49
6:AG:106:LEU:HA	6:AG:110:ALA:HB3	1.94	0.49
12:AQ:16:ARG:O	12:AQ:17:LEU:HD23	2.13	0.49
17:AV:37:VAL:HG11	17:AV:40:LEU:HD21	1.94	0.49
21:AZ:99:TYR:HA	21:AZ:124:ILE:O	2.13	0.49
27:A5:36:CYS:HB3	27:A5:49:CYS:HB3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:298:G:O5'	1:BA:298:G:H8	1.96	0.49
1:BA:889:C:HO2'	1:BA:890:A:H8	1.60	0.49
1:BA:2273:A:O2'	1:BA:2274:A:H5'	2.13	0.49
1:BA:2845:G:O2'	1:BA:2846:G:H5'	2.13	0.49
3:BD:93:ALA:HB3	3:BD:105:ILE:HG13	1.95	0.49
4:BE:12:THR:HG21	15:BT:11:GLU:HG2	1.94	0.49
8:BI:83:ALA:HA	8:BI:89:TYR:CD1	2.47	0.49
9:BN:108:PRO:O	9:BN:113:GLY:HA3	2.13	0.49
31:CA:60:A:H4'	31:CA:61:G:O5'	2.13	0.49
31:CA:60:A:N3	31:CA:61:G:H1'	2.28	0.49
31:CA:626:U:C2	31:CA:627:G:C8	3.01	0.49
31:CA:1007:C:O5'	31:CA:1007:C:H6	1.96	0.49
31:CA:1038:C:H2'	31:CA:1039:C:O4'	2.13	0.49
31:CA:1237:C:H2'	31:CA:1336:C:C5	2.48	0.49
35:CE:11:ILE:HG22	35:CE:12:LEU:HB2	1.94	0.49
35:CE:71:LEU:HD13	35:CE:114:GLY:O	2.12	0.49
49:CS:7:LYS:O	49:CS:7:LYS:HG2	2.13	0.49
31:DA:713:G:H2'	31:DA:714:G:C8	2.48	0.49
31:DA:858:G:N1	31:DA:870:U:OP2	2.38	0.49
31:DA:1065:U:H4'	31:DA:1066:C:O5'	2.13	0.49
44:DN:3:ARG:C	44:DN:3:ARG:HH21	2.16	0.49
1:AA:922:U:H2'	1:AA:923:C:C6	2.47	0.48
1:AA:2445:G:C2'	1:AA:2446:G:H5'	2.43	0.48
1:AA:2502:G:H5''	1:AA:2503:A:H5''	1.95	0.48
1:AA:2815:C:H2'	1:AA:2816:C:H6	1.78	0.48
4:AE:50:GLY:CA	4:AE:75:VAL:HG11	2.42	0.48
6:AG:9:ARG:NH1	6:AG:13:GLU:OE1	2.45	0.48
1:BA:195:A:H61	1:BA:198:C:H3'	1.77	0.48
1:BA:630:G:OP1	30:B8:47:LYS:NZ	2.43	0.48
24:B2:50:ILE:O	24:B2:51:ARG:HB3	2.13	0.48
31:CA:93:G:O2'	31:CA:96:U:P	2.70	0.48
31:CA:413:G:O2'	31:CA:428:G:N2	2.46	0.48
31:CA:456:C:H2'	31:CA:457:C:H6	1.78	0.48
31:CA:987:G:H2'	31:CA:988:G:C8	2.48	0.48
34:CD:111:ALA:HB2	34:CD:120:LEU:HD12	1.95	0.48
38:CH:96:GLY:N	38:CH:99:GLU:OE2	2.40	0.48
31:DA:452:A:C2	31:DA:453:A:C4	3.00	0.48
31:DA:1117:G:H21	31:DA:1180:A:H1'	1.78	0.48
31:DA:1129:C:H5''	31:DA:1139:G:O6	2.13	0.48
33:DC:43:LEU:O	33:DC:47:LEU:HB2	2.12	0.48
50:DT:72:LEU:HD11	50:DT:80:ARG:HD2	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:821:A:H5''	1:AA:822:U:O5'	2.13	0.48
1:AA:897:C:H2'	1:AA:898:C:H6	1.78	0.48
1:AA:954:G:O2'	1:AA:2274:A:N1	2.37	0.48
1:AA:975(A):G:C2	1:AA:990:A:C8	3.01	0.48
1:AA:1019:U:O2'	1:AA:1021:A:H2	1.96	0.48
1:AA:1331:A:H2'	1:AA:1333:C:H5	1.78	0.48
1:AA:1929:G:H4'	1:AA:1930:G:OP1	2.13	0.48
1:AA:2347:C:H2'	1:AA:2348:U:C6	2.48	0.48
2:AB:44:G:C2	2:AB:48:A:C2	3.01	0.48
2:AB:55:U:C1'	6:AG:29:TRP:HE1	2.26	0.48
4:AE:147:PRO:HB2	4:AE:149:ARG:HG2	1.95	0.48
5:AF:155:LEU:HD11	5:AF:176:LEU:HD22	1.94	0.48
15:AT:39:ARG:HH21	31:DA:345:C:H5'	1.78	0.48
21:AZ:111:VAL:HG12	21:AZ:112:ARG:H	1.78	0.48
1:BA:493:G:H2'	1:BA:494:G:O4'	2.13	0.48
1:BA:1153:C:H5'	16:BU:76:TYR:HE2	1.78	0.48
1:BA:1759:A:H1'	1:BA:2711:A:C2	2.48	0.48
1:BA:2115:G:O2'	1:BA:2167:U:H1'	2.12	0.48
31:CA:487:A:H2'	31:CA:488:C:O4'	2.13	0.48
31:CA:913:A:O2'	31:CA:914:A:OP2	2.25	0.48
31:CA:1228:C:H2'	31:CA:1229:A:H8	1.78	0.48
31:CA:1256:A:H5'	31:CA:1258:G:C1'	2.43	0.48
31:CA:1325:C:H4'	51:CU:17:THR:HG21	1.94	0.48
38:CH:112:LEU:HB3	38:CH:133:LEU:HA	1.94	0.48
31:DA:335:C:H2'	31:DA:336:C:C6	2.48	0.48
31:DA:1070:U:H2'	31:DA:1071:C:H6	1.77	0.48
31:DA:1120:G:H2'	31:DA:1121:U:C6	2.49	0.48
33:DC:35:GLU:O	33:DC:39:ILE:HG13	2.13	0.48
43:DM:40:ASN:HA	43:DM:41:PRO:HD2	1.69	0.48
1:AA:71:A:H2	19:AX:31:HIS:HE1	1.58	0.48
1:AA:515:A:H1'	1:AA:581:C:H1'	1.94	0.48
1:AA:914:C:H2'	1:AA:915:C:H5'	1.95	0.48
1:AA:1292:U:H2'	1:AA:1293:C:C6	2.47	0.48
1:AA:1638:C:H2'	1:AA:1639:U:O4'	2.14	0.48
1:AA:2299:G:C2	1:AA:2318:G:H8	2.31	0.48
1:AA:2405:G:O2'	1:AA:2406:U:P	2.71	0.48
4:AE:181:LEU:HD21	15:AT:6:LEU:HD12	1.96	0.48
5:AF:34:TRP:CE3	5:AF:35:GLU:HG2	2.48	0.48
7:AH:7:LEU:O	7:AH:69:ARG:HD3	2.13	0.48
17:AV:43:GLU:HG2	17:AV:44:LYS:H	1.78	0.48
1:BA:656:G:H2'	1:BA:657:U:O4'	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:1257:C:H4'	5:BF:83:PHE:CD2	2.49	0.48
1:BA:1364:G:C8	23:B1:3:LYS:HD3	2.47	0.48
1:BA:1530:C:O2'	1:BA:1531:C:P	2.71	0.48
1:BA:2314:C:H2'	1:BA:2315:G:H8	1.78	0.48
5:BF:123:LEU:HD12	5:BF:124:LEU:N	2.29	0.48
14:BS:96:GLY:H	14:BS:99:LYS:N	2.05	0.48
20:BY:86:ARG:O	20:BY:87:LYS:HD3	2.13	0.48
29:B7:25:PRO:HA	29:B7:28:ARG:CZ	2.43	0.48
31:CA:663:A:H5'	31:CA:836:G:OP1	2.13	0.48
31:CA:1048:G:OP1	44:CN:4:LYS:N	2.42	0.48
31:CA:1320:C:H2'	31:CA:1321:C:O4'	2.12	0.48
32:CB:211:ILE:O	32:CB:215:LEU:HB2	2.13	0.48
40:CJ:68:HIS:H	40:CJ:68:HIS:CD2	2.31	0.48
42:CL:119:LYS:HB2	42:CL:120:TYR:CD1	2.48	0.48
31:DA:436:C:O2'	31:DA:437:U:OP2	2.27	0.48
31:DA:942:G:N2	39:DI:124:GLN:HE22	2.08	0.48
32:DB:178:ARG:HH22	38:DH:68:ARG:HH22	1.59	0.48
32:DB:185:ILE:HG22	32:DB:199:TYR:HB2	1.94	0.48
34:DD:65:ARG:HG2	34:DD:75:PHE:CD1	2.49	0.48
1:AA:226:G:H21	1:AA:228:A:H62	1.61	0.48
1:AA:553:G:H2'	1:AA:554:U:O4'	2.14	0.48
1:AA:848:G:N9	1:AA:933:A:H8	2.12	0.48
1:AA:1014:U:H2'	1:AA:1015:G:C8	2.48	0.48
1:AA:1656:C:H2'	1:AA:1657:C:C6	2.48	0.48
1:AA:1839:G:H5'	1:AA:1839:G:H8	1.78	0.48
2:AB:66:A:H61	2:AB:108:U:H2'	1.77	0.48
5:AF:64:ILE:HG21	5:AF:78:ILE:HG23	1.94	0.48
5:AF:132:VAL:CG2	5:AF:163:VAL:HG22	2.44	0.48
7:AH:107:VAL:HG21	7:AH:152:ARG:HB2	1.94	0.48
9:AN:42:TRP:HA	9:AN:48:MET:SD	2.53	0.48
25:A3:26:LEU:HD21	25:A3:46:ASN:HB3	1.95	0.48
1:BA:1673:U:C5	4:BE:129:HIS:CE1	3.01	0.48
1:BA:1814:G:O3'	3:BD:54:ARG:NH2	2.46	0.48
1:BA:1925:C:O2'	1:BA:1926:U:H5'	2.13	0.48
1:BA:2152:G:H2'	1:BA:2153:G:C8	2.49	0.48
4:BE:147:PRO:HB2	4:BE:149:ARG:HG2	1.96	0.48
6:BG:61:ALA:HB2	6:BG:68:PRO:HD3	1.94	0.48
8:BI:105:HIS:N	8:BI:105:HIS:CD2	2.82	0.48
15:BT:33:LYS:O	15:BT:82:LEU:HD23	2.12	0.48
31:CA:708:C:H2'	31:CA:709:G:H8	1.77	0.48
31:CA:1394:A:C5	31:CA:1501:C:H4'	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:CC:188:LEU:HD21	33:CC:195:VAL:HG22	1.95	0.48
31:DA:1291:G:C6	31:DA:1292:U:C4	3.01	0.48
41:DK:82:VAL:O	41:DK:109:VAL:HG23	2.12	0.48
48:DR:31:LEU:HD11	48:DR:62:GLU:HB3	1.96	0.48
1:AA:1141:U:OP2	9:AN:63:THR:OG1	2.18	0.48
1:AA:2291:U:H2'	1:AA:2292:C:C6	2.48	0.48
1:AA:2695:C:H2'	1:AA:2696:U:H6	1.78	0.48
2:AB:14:U:OP2	2:AB:71:C:H5'	2.13	0.48
18:AW:28:SER:O	18:AW:31:GLU:N	2.46	0.48
24:A2:63:VAL:HA	24:A2:66:GLU:HB2	1.95	0.48
1:BA:286:C:H2'	1:BA:287:C:H6	1.78	0.48
1:BA:1784:A:H4'	1:BA:1785:A:O5'	2.13	0.48
1:BA:2208:A:H1'	1:BA:2219:G:C4	2.48	0.48
3:BD:148:GLU:HB2	3:BD:151:LYS:HD2	1.95	0.48
4:BE:54:GLN:HB2	4:BE:76:ARG:HB3	1.95	0.48
4:BE:119:ARG:HG2	4:BE:160:TYR:HB2	1.96	0.48
8:BI:29:TYR:O	8:BI:32:PRO:HD2	2.13	0.48
14:BS:96:GLY:HA2	14:BS:97:ARG:C	2.34	0.48
31:CA:148:G:C2	31:CA:149:A:N7	2.82	0.48
31:CA:674:G:N2	31:CA:717:C:O2	2.46	0.48
31:CA:1305:G:H5''	51:CU:4:GLY:HA3	1.94	0.48
45:CO:3:ILE:O	45:CO:3:ILE:HG12	2.13	0.48
31:DA:73:G:H2'	31:DA:76:C:O4'	2.14	0.48
31:DA:503:C:H2'	31:DA:504:C:H6	1.78	0.48
31:DA:1023:G:C2	31:DA:1024:G:H1'	2.49	0.48
40:DJ:42:THR:OG1	40:DJ:68:HIS:HB3	2.14	0.48
48:DR:53:ARG:HE	48:DR:59:SER:C	2.16	0.48
1:AA:93:G:H2'	1:AA:94:C:H6	1.79	0.48
1:AA:652(S):C:H5'	1:AA:652(T):C:OP2	2.14	0.48
1:AA:980:A:C4	1:AA:1136:G:O4'	2.67	0.48
1:AA:1205:U:C4	5:AF:171:PRO:HA	2.47	0.48
1:AA:1231:G:H2'	1:AA:1232:G:H8	1.79	0.48
1:AA:1419:A:C8	1:AA:1421:G:C6	3.02	0.48
1:AA:1701:A:H5''	1:AA:1702:G:OP2	2.14	0.48
1:AA:1827:C:C2'	1:AA:1828:G:H5'	2.44	0.48
1:AA:2758:A:C2	1:AA:2759:G:H1'	2.48	0.48
7:AH:54:ARG:NH2	7:AH:57:ASP:OD1	2.46	0.48
14:AS:11:LYS:HD3	14:AS:15:ARG:NH1	2.29	0.48
14:AS:87:PHE:HB2	14:AS:112:PHE:CE1	2.48	0.48
15:AT:121:ILE:O	15:AT:124:ASP:HB2	2.13	0.48
19:AX:18:TYR:C	19:AX:20:GLY:H	2.17	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:A0:72:ARG:HB2	22:A0:75:LEU:HB2	1.95	0.48
1:BA:29:U:H6	1:BA:29:U:O5'	1.96	0.48
1:BA:861:A:N3	2:BB:79:C:O2'	2.44	0.48
1:BA:2133:G:H1'	1:BA:2158:A:H61	1.79	0.48
4:BE:31:CYS:HB3	4:BE:49:LEU:HG	1.95	0.48
6:BG:111:LEU:HB3	6:BG:117:PHE:CZ	2.48	0.48
12:BQ:37:LEU:HD21	12:BQ:130:LYS:HE3	1.96	0.48
14:BS:58:LEU:HG	14:BS:65:VAL:HG13	1.96	0.48
17:BV:35:LEU:HB2	17:BV:57:VAL:CG2	2.43	0.48
22:B0:51:VAL:CG2	22:B0:81:VAL:HG23	2.44	0.48
31:CA:376:G:H2'	31:CA:377:G:H8	1.78	0.48
31:CA:531:U:O3'	31:CA:532:A:H4'	2.13	0.48
31:CA:738:C:H2'	31:CA:739:C:H6	1.79	0.48
31:CA:838:G:H2'	31:CA:839:U:H5''	1.94	0.48
31:CA:1343:G:H2'	31:CA:1344:C:C6	2.48	0.48
31:CA:1409:C:H2'	31:CA:1410:G:C8	2.49	0.48
34:CD:18:LYS:HD3	34:CD:20:TYR:OH	2.13	0.48
36:CF:62:TRP:CZ3	36:CF:64:GLN:HB2	2.49	0.48
39:CI:29:ASN:OD1	39:CI:64:THR:HA	2.13	0.48
46:CP:51:VAL:HG12	46:CP:53:VAL:N	2.22	0.48
47:CQ:67:LYS:HA	47:CQ:70:ARG:NH1	2.25	0.48
31:DA:9:G:H2'	31:DA:10:A:H8	1.77	0.48
31:DA:189(K):U:H2'	31:DA:189(L):G:C8	2.48	0.48
31:DA:425:G:H4'	34:DD:45:GLN:OE1	2.14	0.48
36:DF:3:ARG:HB3	36:DF:93:SER:CB	2.43	0.48
37:DG:16:LEU:H	37:DG:16:LEU:HD12	1.78	0.48
38:DH:124:ALA:O	38:DH:128:GLY:N	2.44	0.48
1:AA:118:A:N3	1:AA:178:G:H1'	2.29	0.48
1:AA:528:A:H3'	1:AA:528:A:H8	1.77	0.48
1:AA:2438:U:H5''	1:AA:2600:A:OP1	2.12	0.48
7:AH:70:THR:O	7:AH:71:LEU:HB3	2.14	0.48
10:AO:73:ASP:OD1	15:AT:32:TYR:OH	2.09	0.48
23:A1:15:ALA:O	23:A1:40:ARG:HG3	2.14	0.48
1:BA:740:U:H2'	1:BA:741:G:C8	2.48	0.48
1:BA:2114:A:H3'	1:BA:2115:G:H8	1.73	0.48
1:BA:2687:U:H2'	1:BA:2688:U:O4'	2.12	0.48
16:BU:28:ARG:HD3	16:BU:38:THR:OG1	2.13	0.48
18:BW:14:PRO:HG2	18:BW:78:GLU:CG	2.37	0.48
27:B5:53:ALA:N	27:B5:54:GLY:HA2	2.29	0.48
31:CA:59:A:H5''	31:CA:60:A:C5'	2.43	0.48
31:CA:156:G:N1	31:CA:166:G:C6	2.82	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:445:G:C2	31:CA:446:G:C4	3.01	0.48
31:CA:607:A:H2'	31:CA:608:A:O4'	2.14	0.48
31:CA:741:G:H2'	31:CA:742:G:O4'	2.14	0.48
31:CA:833:U:H3	31:CA:853:G:H1	1.60	0.48
37:CG:139:GLU:O	37:CG:143:ARG:HG3	2.14	0.48
50:CT:90:GLN:O	50:CT:93:GLU:HB3	2.13	0.48
51:CU:17:THR:O	51:CU:22:ARG:HD3	2.13	0.48
31:DA:413:G:N2	31:DA:428:G:H1'	2.29	0.48
31:DA:691:G:H2'	31:DA:692:U:C6	2.49	0.48
31:DA:1159:U:O4'	31:DA:1182:G:N2	2.47	0.48
39:DI:11:LYS:H	39:DI:104:ARG:NH2	2.11	0.48
45:DO:15:PHE:CE2	45:DO:84:LYS:HD2	2.48	0.48
50:DT:10:LEU:HD23	50:DT:11:SER:N	2.28	0.48
1:AA:325:G:H2'	1:AA:326:G:O4'	2.13	0.48
1:AA:548:A:N6	17:AV:19:LYS:H	2.12	0.48
1:AA:1463:C:H2'	1:AA:1464:C:C6	2.48	0.48
1:AA:2729:G:H2'	1:AA:2730:C:C6	2.48	0.48
4:AE:108:SER:HB3	4:AE:165:VAL:HG21	1.96	0.48
6:AG:173:LEU:HB3	6:AG:178:PHE:CG	2.48	0.48
18:AW:56:ALA:O	18:AW:60:ASN:HB2	2.12	0.48
19:AX:18:TYR:O	19:AX:20:GLY:N	2.47	0.48
1:BA:36:G:N3	1:BA:450:G:O2'	2.45	0.48
1:BA:797:C:OP2	5:BF:62:ARG:HB2	2.14	0.48
1:BA:830:G:H4'	1:BA:831:G:OP2	2.12	0.48
1:BA:928:G:H8	1:BA:928:G:O5'	1.96	0.48
1:BA:1312:U:H4'	1:BA:1313:U:O5'	2.13	0.48
1:BA:1514:U:H2'	1:BA:1515:G:H8	1.78	0.48
1:BA:1914:C:HO2'	1:BA:1915:U:C5'	2.27	0.48
1:BA:2438:U:O2'	1:BA:2440:C:OP1	2.29	0.48
7:BH:7:LEU:HG	7:BH:69:ARG:HH12	1.78	0.48
8:BI:101:LEU:HD12	8:BI:140:LEU:HD11	1.96	0.48
17:BV:52:VAL:HG22	17:BV:55:ALA:HB3	1.96	0.48
18:BW:83:LYS:O	18:BW:84:ARG:NH1	2.43	0.48
25:B3:19:GLN:OE1	25:B3:52:HIS:NE2	2.45	0.48
31:CA:655:A:N1	31:CA:656:C:C2	2.82	0.48
31:CA:1168:A:C6	31:CA:1169:A:C6	3.01	0.48
31:CA:1324:A:H2'	31:CA:1325:C:H6	1.79	0.48
32:CB:181:PHE:O	32:CB:183:PRO:HD3	2.14	0.48
35:CE:136:MET:O	35:CE:139:LEU:N	2.47	0.48
42:CL:6:THR:HG23	42:CL:9:GLN:OE1	2.13	0.48
31:DA:240:C:H2'	31:DA:241:C:H6	1.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:666:G:H5'	31:DA:726:C:H1'	1.96	0.48
31:DA:1246:C:H2'	31:DA:1247:U:C6	2.48	0.48
31:DA:1309:G:OP1	43:DM:88:ARG:NH1	2.47	0.48
37:DG:20:ASP:HB3	37:DG:23:VAL:HG23	1.94	0.48
38:DH:92:ARG:HB3	38:DH:94:TYR:CE2	2.49	0.48
39:DI:24:GLY:HA2	39:DI:59:PHE:O	2.14	0.48
1:AA:71:A:H2	19:AX:31:HIS:CE1	2.31	0.48
1:AA:1005:C:H2'	1:AA:1006:C:C6	2.49	0.48
1:AA:2355:C:H5''	1:AA:2356:C:OP2	2.13	0.48
1:AA:2884:U:H2'	1:AA:2885:C:O4'	2.13	0.48
2:AB:111:G:H2'	2:AB:112:U:H6	1.79	0.48
7:AH:41:MET:SD	7:AH:64:LEU:HB3	2.53	0.48
8:AI:82:ARG:HB3	8:AI:89:TYR:CD1	2.49	0.48
9:AN:23:LEU:HB2	9:AN:60:ILE:HG12	1.95	0.48
1:BA:1173:G:O2'	1:BA:1174:A:O5'	2.32	0.48
1:BA:1174:A:H1'	1:BA:1175:U:C5'	2.41	0.48
1:BA:1514:U:H2'	1:BA:1515:G:C8	2.49	0.48
1:BA:1756:G:H4'	1:BA:1758:G:O4'	2.14	0.48
1:BA:1784:A:H4'	1:BA:1785:A:C5'	2.44	0.48
1:BA:1876:A:H2'	1:BA:1877:A:C8	2.48	0.48
1:BA:2287:A:H2	1:BA:2346:A:H62	1.59	0.48
1:BA:2492:U:H2'	1:BA:2493:U:C6	2.49	0.48
4:BE:143:ASN:HD22	4:BE:147:PRO:CD	2.27	0.48
16:BU:47:TYR:HA	16:BU:50:ARG:NH2	2.29	0.48
17:BV:14:VAL:HB	17:BV:96:ILE:HG13	1.95	0.48
31:CA:397:A:H3'	31:CA:397:A:N3	2.28	0.48
31:CA:405:U:O4	34:CD:2:GLY:N	2.47	0.48
35:CE:143:ARG:NH1	38:CH:77:GLU:OE1	2.47	0.48
31:DA:266:G:O3'	47:DQ:67:LYS:HB2	2.14	0.48
31:DA:568:G:N3	31:DA:574:A:H2	2.12	0.48
31:DA:742:G:P	45:DO:35:ARG:HH21	2.36	0.48
31:DA:942:G:H21	39:DI:124:GLN:NE2	2.11	0.48
33:DC:7:PRO:HG3	33:DC:201:TYR:HE2	1.78	0.48
42:DL:83:VAL:HG13	42:DL:100:ILE:HG23	1.96	0.48
1:AA:786:C:C2'	1:AA:787:U:H5'	2.44	0.48
1:AA:876:C:H2'	1:AA:877:U:O4'	2.14	0.48
1:AA:1047:G:H2'	1:AA:1110:G:N2	2.26	0.48
1:AA:1141:U:H2'	9:AN:63:THR:HG21	1.96	0.48
1:AA:1282:U:H2'	1:AA:1283:G:O4'	2.14	0.48
1:AA:1445:A:H8	1:AA:1460:A:C5	2.32	0.48
1:AA:1826:G:H2'	1:AA:1827:C:H6	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1946:U:H2'	1:AA:1947:C:C6	2.48	0.48
1:AA:2056:G:C4	1:AA:2577:A:C2	3.02	0.48
1:AA:2115:G:H21	1:AA:2171:A:N6	2.12	0.48
1:AA:2305:A:H2'	1:AA:2306:C:C6	2.49	0.48
1:AA:2406:U:H4'	1:AA:2407:G:H5''	1.96	0.48
7:AH:9:ILE:N	7:AH:50:VAL:O	2.31	0.48
24:A2:48:HIS:O	24:A2:52:ASP:HB2	2.14	0.48
1:BA:154:G:C6	1:BA:154(A):C:N4	2.82	0.48
1:BA:1126:A:H8	1:BA:1126:A:OP1	1.97	0.48
1:BA:2316:C:H2'	1:BA:2317:C:C6	2.49	0.48
21:BZ:31:ARG:NH1	21:BZ:94:GLU:OE2	2.42	0.48
23:B1:3:LYS:HB2	23:B1:61:ARG:HH12	1.78	0.48
26:B4:42:PHE:CB	26:B4:43:TYR:HB2	2.43	0.48
31:CA:56:U:H2'	31:CA:57:G:C8	2.48	0.48
31:CA:277:C:H5''	47:CQ:68:ARG:NH2	2.28	0.48
31:CA:598:U:H4'	38:CH:94:TYR:CD2	2.48	0.48
31:CA:649:G:H2'	31:CA:650:G:C8	2.49	0.48
31:CA:1067:A:N3	31:CA:1068:G:H1'	2.29	0.48
31:CA:1349:A:OP1	39:CI:121:ARG:N	2.46	0.48
31:CA:1418:A:C2	31:CA:1483:A:C2	3.02	0.48
34:CD:19:LEU:HD23	34:CD:21:LEU:HD11	1.95	0.48
43:CM:65:LYS:HA	43:CM:66:LEU:CB	2.41	0.48
33:DC:40:ARG:O	33:DC:44:GLU:HB2	2.14	0.48
33:DC:125:GLU:HA	33:DC:191:THR:HG22	1.95	0.48
45:DO:71:GLN:HA	45:DO:78:TYR:HB2	1.95	0.48
50:DT:72:LEU:HD23	50:DT:73:HIS:N	2.29	0.48
1:AA:289:A:H2'	1:AA:290:G:O4'	2.12	0.47
1:AA:952:G:C6	1:AA:966:G:C6	3.02	0.47
1:AA:966:G:C6	1:AA:967:C:N4	2.82	0.47
1:AA:2207:G:O2'	1:AA:2208:A:P	2.72	0.47
6:AG:11:TYR:HB2	6:AG:176:LEU:HD21	1.96	0.47
1:BA:628:G:H2'	1:BA:629:G:C8	2.49	0.47
1:BA:848:G:O6	1:BA:928:G:H2'	2.14	0.47
1:BA:1625:C:H2'	1:BA:1626:G:O4'	2.14	0.47
1:BA:1833:U:H2'	1:BA:1834:U:H6	1.79	0.47
1:BA:2362:G:OP1	30:B8:44:LYS:NZ	2.43	0.47
2:BB:90:A:N7	2:BB:91:C:H1'	2.28	0.47
7:BH:5:GLY:HA2	7:BH:69:ARG:HB3	1.96	0.47
16:BU:61:TRP:O	16:BU:65:ILE:HG13	2.14	0.47
31:CA:174:C:H2'	31:CA:175:C:H6	1.79	0.47
31:CA:363:A:OP1	42:CL:34:ARG:N	2.35	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:1072:G:C5	31:CA:1073:U:C4	3.02	0.47
31:CA:1438:G:H2'	31:CA:1439:C:H6	1.79	0.47
32:CB:116:GLU:HA	32:CB:119:GLU:CB	2.41	0.47
34:CD:148:VAL:HG12	34:CD:149:ALA:H	1.79	0.47
31:DA:4:U:C5	38:DH:105:ARG:HD3	2.49	0.47
31:DA:236:G:C6	31:DA:237:C:C4	3.02	0.47
31:DA:1074:G:C2	31:DA:1075:C:C2	3.02	0.47
31:DA:1330:U:H4'	43:DM:23:TYR:CE2	2.49	0.47
31:DA:1362:C:H2'	31:DA:1363:C:H5''	1.96	0.47
40:DJ:24:VAL:O	40:DJ:34:VAL:HG11	2.14	0.47
48:DR:21:LYS:HD3	48:DR:21:LYS:HA	1.58	0.47
1:AA:96:G:H4'	24:A2:48:HIS:CD2	2.49	0.47
1:AA:221:A:C4	1:AA:266:G:N7	2.83	0.47
1:AA:391:G:C2	1:AA:411:G:C5	3.02	0.47
1:AA:1448:G:O2'	1:AA:1528(A):A:N1	2.37	0.47
1:AA:1919:A:N1	31:DA:1495:U:O2'	2.38	0.47
6:AG:3:LEU:HD22	26:A4:25:TYR:CE2	2.49	0.47
8:AI:126:TYR:HB2	8:AI:142:VAL:HG23	1.95	0.47
15:AT:51:ARG:HB3	15:AT:62:THR:HB	1.96	0.47
1:BA:300:A:P	20:BY:86:ARG:HH22	2.37	0.47
1:BA:511:U:C5	1:BA:512:G:C4	3.02	0.47
1:BA:637:A:H8	11:BP:117:GLU:HG3	1.79	0.47
1:BA:2106:G:C2	1:BA:2184:G:C2	3.01	0.47
3:BD:246:PRO:O	3:BD:254:THR:HG22	2.14	0.47
6:BG:132:ASN:HA	6:BG:157:ILE:O	2.13	0.47
31:CA:184:G:O2'	31:CA:185:A:H5'	2.14	0.47
31:CA:601:C:H2'	31:CA:602:A:C8	2.49	0.47
31:CA:735:C:H2'	31:CA:736:C:C6	2.46	0.47
31:CA:952:U:O2'	31:CA:953:G:H5'	2.14	0.47
31:CA:1375:A:C6	31:CA:1376:U:N3	2.82	0.47
32:CB:102:LEU:HB3	32:CB:180:LEU:CD1	2.44	0.47
33:CC:51:GLY:HA3	33:CC:71:ALA:HB3	1.95	0.47
36:CF:3:ARG:NE	36:CF:38:GLU:OE1	2.48	0.47
39:CI:5:TYR:HA	39:CI:17:VAL:O	2.14	0.47
43:CM:48:LEU:O	43:CM:52:GLU:HB2	2.14	0.47
44:CN:39:LEU:HD22	44:CN:43:CYS:HB3	1.96	0.47
50:CT:73:HIS:HB3	50:CT:74:LYS:HE2	1.96	0.47
31:DA:299:G:H2'	31:DA:300:A:C8	2.49	0.47
31:DA:475:G:C2'	31:DA:476:G:H5'	2.43	0.47
31:DA:833:U:H3	31:DA:853:G:H1	1.61	0.47
31:DA:1059:C:H2'	31:DA:1060:C:C6	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:1165:C:H2'	31:DA:1166:G:O4'	2.13	0.47
31:DA:1510:U:H2'	31:DA:1511:G:C8	2.49	0.47
44:DN:24:CYS:SG	44:DN:40:CYS:CB	3.02	0.47
44:DN:40:CYS:H	44:DN:43:CYS:HB2	1.80	0.47
1:AA:1009:A:H8	1:AA:1009:A:O5'	1.96	0.47
1:AA:1131:G:C2	1:AA:1132:A:C4	3.01	0.47
1:AA:1161:C:H2'	1:AA:1162:G:H8	1.80	0.47
1:AA:1356:G:C6	1:AA:1357:U:C4	3.03	0.47
1:AA:1405:U:H2'	1:AA:1406:U:H6	1.77	0.47
1:AA:1524:G:N2	1:AA:1525:G:H1'	2.30	0.47
1:AA:1721:G:H8	1:AA:1741:A:H62	1.62	0.47
1:AA:1740:G:H2'	1:AA:1741:A:H8	1.79	0.47
1:AA:1784:A:H4'	1:AA:1785:A:H5''	1.94	0.47
1:AA:1833:U:H2'	1:AA:1834:U:C6	2.50	0.47
1:AA:2051:A:H5'	1:AA:2578:G:O4'	2.13	0.47
1:AA:2472:G:H2'	1:AA:2475:C:H42	1.79	0.47
3:AD:223:GLY:HA3	3:AD:231:HIS:CE1	2.49	0.47
6:AG:56:ALA:HB2	6:AG:153:ARG:NE	2.29	0.47
7:AH:25:LYS:NZ	7:AH:32:GLU:OE2	2.38	0.47
12:AQ:35:VAL:CG1	12:AQ:130:LYS:HB3	2.44	0.47
14:AS:58:LEU:HD12	14:AS:59:LYS:HD3	1.96	0.47
21:AZ:102:LEU:HD23	21:AZ:104:PHE:CZ	2.49	0.47
1:BA:1019:U:H3	1:BA:1142(A):A:H62	1.62	0.47
2:BB:6:C:C2	2:BB:116:G:N2	2.82	0.47
31:CA:397:A:N6	31:CA:548:G:C5	2.82	0.47
31:CA:787:A:N1	31:CA:795:C:N4	2.57	0.47
31:CA:1077:G:N2	31:CA:1080:A:OP2	2.34	0.47
31:CA:1412:C:H2'	31:CA:1413:A:C8	2.49	0.47
32:CB:185:ILE:HA	32:CB:199:TYR:O	2.14	0.47
34:CD:33:MET:SD	34:CD:37:PRO:HA	2.54	0.47
48:CR:74:ARG:HG2	48:CR:79:LEU:HB2	1.96	0.47
50:CT:33:ILE:O	50:CT:37:SER:OG	2.22	0.47
31:DA:78:G:N2	31:DA:91:C:N3	2.53	0.47
31:DA:1137:C:H5''	31:DA:1138:G:OP1	2.14	0.47
31:DA:1252:A:H2'	31:DA:1253:G:O4'	2.13	0.47
31:DA:1289:A:OP1	51:DU:9:ARG:NH2	2.47	0.47
31:DA:1320:C:N4	49:DS:36:ARG:HG3	2.30	0.47
34:DD:25:ARG:C	34:DD:27:TYR:H	2.16	0.47
1:AA:836:G:C5	1:AA:837:C:C4	3.02	0.47
1:AA:902:C:H2'	1:AA:903:C:C6	2.49	0.47
1:AA:1411:C:H2'	1:AA:1412:A:C8	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1434:A:H61	1:AA:1558:A:N6	2.12	0.47
1:AA:1494:A:O2'	1:AA:1495:A:H5'	2.14	0.47
1:AA:1591:G:C6	1:AA:1592:C:C4	3.03	0.47
1:AA:1793:C:H2'	1:AA:1794:U:H6	1.79	0.47
4:AE:9:VAL:HG13	4:AE:25:VAL:O	2.14	0.47
4:AE:31:CYS:HB3	4:AE:49:LEU:HG	1.96	0.47
4:AE:36:ARG:NH2	4:AE:88:GLY:O	2.46	0.47
6:AG:43:LEU:HB2	6:AG:89:GLY:HA2	1.97	0.47
9:AN:15:LEU:HB2	9:AN:135:PRO:HB2	1.97	0.47
21:AZ:102:LEU:HB3	21:AZ:104:PHE:CE2	2.50	0.47
1:BA:828:U:C5	1:BA:2247:A:H4'	2.49	0.47
1:BA:879:G:H2'	1:BA:880:G:O4'	2.13	0.47
1:BA:1177:A:P	1:BA:1177:A:H2'	2.54	0.47
1:BA:1324:G:C4	1:BA:1328:G:O6	2.67	0.47
3:BD:5:LYS:HB3	3:BD:5:LYS:HE3	1.67	0.47
30:B8:34:TRP:CE2	30:B8:35:GLN:HB3	2.50	0.47
31:CA:303:A:H2'	31:CA:304:U:O4'	2.15	0.47
31:CA:316:G:OP2	31:CA:351:G:O2'	2.32	0.47
31:CA:382:A:H2'	31:CA:383:A:C8	2.50	0.47
31:CA:778:G:H1'	41:CK:119:CYS:HB3	1.96	0.47
31:CA:952:U:H5'	31:CA:972:C:N4	2.29	0.47
31:CA:1026:G:N3	31:CA:1026:G:H2'	2.30	0.47
31:CA:1165:C:H2'	31:CA:1166:G:O4'	2.14	0.47
32:CB:132:LYS:O	32:CB:135:GLN:HG2	2.14	0.47
34:CD:53:ASP:O	34:CD:57:ARG:NH1	2.47	0.47
40:CJ:43:ARG:O	40:CJ:67:THR:HG23	2.14	0.47
43:CM:108:ARG:O	43:CM:112:GLY:N	2.47	0.47
31:DA:153:C:O5'	31:DA:153:C:H6	1.97	0.47
31:DA:688:G:H2'	31:DA:689:C:H6	1.78	0.47
31:DA:1113:C:N4	31:DA:1187:G:H1	2.11	0.47
31:DA:1305:G:H22	31:DA:1331:G:HO2'	1.57	0.47
31:DA:1333:A:H2'	31:DA:1334:G:O4'	2.14	0.47
32:DB:213:LEU:HD22	32:DB:214:ILE:HD13	1.96	0.47
35:DE:53:LEU:HD12	35:DE:53:LEU:H	1.79	0.47
42:DL:66:VAL:HG11	42:DL:98:TYR:CE1	2.49	0.47
44:DN:24:CYS:HB3	44:DN:27:CYS:O	2.14	0.47
1:AA:271(G):C:O2'	1:AA:271(H):G:H5'	2.14	0.47
1:AA:1109:C:C5	1:AA:1110:G:C5	3.03	0.47
1:AA:1287:A:N1	1:AA:1648:C:O2'	2.43	0.47
1:AA:1545:A:H2'	1:AA:1546:C:O4'	2.15	0.47
1:AA:1575:C:H2'	1:AA:1576:U:C6	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1671:U:O2'	1:AA:1673:U:H5	1.96	0.47
1:AA:2249:U:H4'	1:AA:2250:G:OP2	2.14	0.47
1:AA:2789:C:H4'	1:AA:2790:A:OP1	2.15	0.47
52:AA:3001:T8B:C13	31:DA:1492:A:O5'	2.62	0.47
19:AX:11:PRO:HG2	19:AX:13:LEU:HD21	1.96	0.47
1:BA:747:U:O2	1:BA:2014:A:H1'	2.13	0.47
1:BA:858:U:O2	1:BA:2268:A:H2'	2.13	0.47
1:BA:955:C:OP1	12:BQ:87:LYS:HE2	2.15	0.47
1:BA:1021:A:C8	1:BA:1021:A:C3'	2.98	0.47
1:BA:2335:A:O2'	1:BA:2336:A:H5''	2.14	0.47
1:BA:2786:U:O2'	4:BE:62:PRO:O	2.20	0.47
3:BD:13:ARG:NH1	3:BD:16:MET:SD	2.87	0.47
15:BT:13:ARG:HG3	15:BT:14:TYR:CE1	2.50	0.47
31:CA:110:C:H2'	31:CA:111:G:O4'	2.15	0.47
31:CA:189(C):C:H2'	31:CA:189(D):C:O4'	2.14	0.47
31:CA:191:G:N3	50:CT:103:GLY:HA2	2.29	0.47
31:CA:427:U:C4	31:CA:428:G:C6	3.03	0.47
31:CA:731:G:OP1	31:CA:766:A:H1'	2.14	0.47
31:CA:738:C:H2'	31:CA:739:C:C6	2.50	0.47
31:CA:881:G:H2'	31:CA:882:C:O4'	2.15	0.47
35:CE:36:ASP:OD2	35:CE:37:ARG:N	2.48	0.47
44:CN:4:LYS:O	44:CN:7:ILE:HG23	2.14	0.47
31:DA:922:G:H2'	31:DA:923:A:C8	2.49	0.47
33:DC:113:ALA:HB2	33:DC:202:ILE:HG13	1.96	0.47
45:DO:53:HIS:O	45:DO:56:LEU:HB3	2.15	0.47
1:AA:363(B):G:H2'	1:AA:363(C):G:H8	1.79	0.47
1:AA:1141:U:H4'	1:AA:1142(A):A:O4'	2.14	0.47
1:AA:1689:A:N6	1:AA:1698:A:H2	1.95	0.47
2:AB:72:G:O2'	2:AB:105:A:N6	2.48	0.47
4:AE:52:LEU:O	4:AE:76:ARG:N	2.45	0.47
15:AT:16:ARG:NH1	15:AT:18:ASP:OD2	2.48	0.47
22:A0:43:THR:O	22:A0:45:PHE:N	2.47	0.47
1:BA:225:A:H2'	1:BA:226:G:H5'	1.96	0.47
1:BA:1289:C:H2'	1:BA:1290:C:H6	1.80	0.47
1:BA:1418:G:H8	1:BA:1418:G:O5'	1.98	0.47
1:BA:1657:C:H4'	4:BE:133:LYS:HB3	1.96	0.47
1:BA:2699:C:H2'	1:BA:2700:C:O4'	2.13	0.47
8:BI:65:ALA:HB1	8:BI:136:VAL:HG11	1.97	0.47
14:BS:59:LYS:H	14:BS:60:GLY:HA3	1.78	0.47
31:CA:575:G:O2'	31:CA:821:G:H5'	2.15	0.47
31:CA:955:U:H2'	31:CA:956:U:H6	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:991:U:H2'	31:CA:1212:U:O2	2.15	0.47
31:CA:1086:U:H2'	31:CA:1087:G:C8	2.44	0.47
31:CA:1118:C:H2'	31:CA:1119:C:C6	2.50	0.47
31:CA:1250:A:H4'	39:CI:68:GLY:H	1.79	0.47
32:CB:115:LEU:O	32:CB:119:GLU:N	2.46	0.47
39:CI:44:VAL:H	39:CI:45:ALA:HA	1.78	0.47
31:DA:187:C:H2'	31:DA:188:C:H6	1.78	0.47
31:DA:1291:G:H4'	39:DI:39:GLY:HA3	1.96	0.47
31:DA:1392:G:N2	31:DA:1502:A:H8	2.13	0.47
1:AA:760:G:H2'	1:AA:761:A:O4'	2.14	0.47
1:AA:1198:U:H2'	1:AA:1199:U:H6	1.78	0.47
1:AA:1331:A:O2'	1:AA:1332:G:H5''	2.14	0.47
1:AA:1354:A:H2'	1:AA:1355:G:O4'	2.14	0.47
1:AA:1472:A:H2'	1:AA:1473:G:O4'	2.14	0.47
1:AA:1660:C:H2'	1:AA:1661:G:H8	1.79	0.47
1:AA:1803:A:O2'	3:AD:259:THR:HG21	2.13	0.47
1:AA:1849:G:N3	1:AA:1850:G:C8	2.82	0.47
1:AA:2405:G:O2'	1:AA:2411:A:N6	2.47	0.47
52:AA:3001:T8B:H10	31:DA:1492:A:C4'	2.43	0.47
2:AB:28:C:C2	2:AB:29:A:C8	3.02	0.47
2:AB:38:C:O2	2:AB:48:A:H1'	2.14	0.47
2:AB:45:A:H2'	2:AB:46:A:H5'	1.96	0.47
3:AD:221:VAL:HG22	3:AD:226:MET:CE	2.44	0.47
6:AG:3:LEU:HD22	26:A4:25:TYR:CZ	2.50	0.47
11:AP:121:LYS:HD3	11:AP:123:LEU:HD11	1.96	0.47
16:AU:113:ALA:O	16:AU:117:GLN:HG2	2.15	0.47
21:AZ:40:ASP:HB3	21:AZ:43:GLU:HB2	1.97	0.47
24:A2:17:SER:HB3	24:A2:20:GLU:OE2	2.14	0.47
26:A4:15:ILE:HB	26:A4:32:TYR:HD2	1.80	0.47
28:A6:26:ASN:HB3	28:A6:29:ASN:HB2	1.97	0.47
1:BA:51:G:N2	1:BA:120:U:O2	2.45	0.47
1:BA:244:A:H4'	11:BP:74:GLU:HB2	1.95	0.47
1:BA:566:U:H5''	11:BP:29:LYS:HE3	1.97	0.47
1:BA:645:C:O2	1:BA:645:C:H2'	2.15	0.47
1:BA:652(F):G:H2'	1:BA:652(G):G:C8	2.49	0.47
1:BA:1173:G:N1	1:BA:1176:G:OP2	2.25	0.47
1:BA:1282:U:H2'	1:BA:1283:G:O4'	2.15	0.47
1:BA:1668:A:H4'	1:BA:1669:A:O5'	2.15	0.47
1:BA:1797:C:H4'	3:BD:257:LEU:O	2.15	0.47
1:BA:1803:A:O2'	3:BD:259:THR:HG21	2.15	0.47
1:BA:1839:G:H5'	1:BA:1839:G:C8	2.46	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:2128:C:N4	1:BA:2160:G:H1	2.11	0.47
1:BA:2199:A:OP2	1:BA:2200:C:H5	1.96	0.47
1:BA:2282:G:H5''	1:BA:2283:C:O4'	2.14	0.47
3:BD:71:ASP:HB3	3:BD:103:ARG:HH22	1.78	0.47
5:BF:53:THR:C	5:BF:55:GLY:H	2.18	0.47
7:BH:124:GLU:OE1	7:BH:132:ARG:HD2	2.14	0.47
12:BQ:26:TYR:CE1	12:BQ:28:ALA:HB2	2.50	0.47
12:BQ:103:MET:HE2	12:BQ:104:PHE:CE1	2.50	0.47
13:BR:36:THR:HG22	13:BR:37:THR:H	1.78	0.47
15:BT:127:ALA:HA	15:BT:128:GLU:C	2.35	0.47
31:CA:96:U:HO2'	31:CA:97:G:P	2.35	0.47
31:CA:624:C:H2'	31:CA:625:G:C8	2.50	0.47
31:CA:826:C:H2'	31:CA:827:U:C6	2.50	0.47
31:CA:944:G:O6	31:CA:1337:G:H8	1.98	0.47
31:CA:945:G:N1	31:CA:1337:G:C2	2.82	0.47
31:CA:1161:C:H2'	31:CA:1162:C:C6	2.50	0.47
31:CA:1226:C:C5	43:CM:104:ARG:HB2	2.49	0.47
33:CC:189:ALA:O	33:CC:191:THR:N	2.44	0.47
49:CS:31:ILE:O	49:CS:50:ALA:N	2.41	0.47
31:DA:115:G:H4'	31:DA:116:A:O5'	2.14	0.47
31:DA:344:A:H4'	31:DA:345:C:OP2	2.15	0.47
31:DA:355:C:H5''	31:DA:389:A:OP2	2.15	0.47
31:DA:417:C:H2'	31:DA:418:C:C6	2.49	0.47
31:DA:460:G:C6	31:DA:470:C:H5''	2.50	0.47
31:DA:487:A:H2'	31:DA:488:C:O4'	2.15	0.47
31:DA:521:G:O5'	42:DL:73:GLU:HG2	2.14	0.47
31:DA:542:G:OP1	34:DD:10:ARG:NH1	2.45	0.47
31:DA:791:G:C6	31:DA:792:A:N7	2.83	0.47
31:DA:991:U:C4	31:DA:1212:U:H1'	2.49	0.47
31:DA:993:G:O2'	31:DA:994:A:N7	2.46	0.47
31:DA:1030(A):G:H3'	31:DA:1030(B):C:C5'	2.45	0.47
31:DA:1238:A:H62	31:DA:1299:A:N6	2.12	0.47
31:DA:1285:A:H4'	31:DA:1286:A:O5'	2.15	0.47
31:DA:1504:G:H3'	31:DA:1504:G:P	2.54	0.47
32:DB:28:PHE:CE2	32:DB:31:TYR:HD1	2.32	0.47
33:DC:113:ALA:HB1	33:DC:200:ALA:HB1	1.96	0.47
37:DG:113:GLU:H	37:DG:113:GLU:HG2	1.50	0.47
43:DM:65:LYS:O	43:DM:70:LEU:HG	2.14	0.47
46:DP:29:ASP:OD2	46:DP:29:ASP:N	2.48	0.47
51:DU:6:ARG:C	51:DU:8:THR:H	2.17	0.47
1:AA:102:G:O2'	1:AA:103:A:P	2.72	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1161:C:H2'	1:AA:1162:G:C8	2.49	0.47
1:AA:1564:C:O2'	1:AA:1565:C:H5'	2.15	0.47
1:AA:1792:G:C5'	3:AD:205:VAL:HG13	2.45	0.47
1:AA:2301:C:H2'	1:AA:2302:G:H8	1.80	0.47
1:AA:2396:G:OP1	23:A1:25:LYS:NZ	2.35	0.47
1:AA:2736:G:N2	1:AA:2737:G:H1'	2.30	0.47
2:AB:50:G:H2'	2:AB:51:G:O4'	2.15	0.47
8:AI:12:LEU:HD22	8:AI:19:VAL:HG21	1.96	0.47
8:AI:76:THR:O	8:AI:105:HIS:NE2	2.47	0.47
1:BA:668:G:H5'	1:BA:669:G:OP2	2.14	0.47
1:BA:779:U:H5''	3:BD:49:ILE:HD11	1.95	0.47
1:BA:1701:A:H5''	1:BA:1702:G:OP2	2.15	0.47
3:BD:35:LYS:HE2	3:BD:64:ILE:HD11	1.95	0.47
14:BS:65:VAL:O	14:BS:69:VAL:HG12	2.15	0.47
14:BS:95:HIS:C	14:BS:99:LYS:HB3	2.35	0.47
18:BW:4:LYS:HB2	18:BW:106:ILE:HG12	1.97	0.47
18:BW:28:SER:O	18:BW:31:GLU:N	2.46	0.47
18:BW:31:GLU:O	18:BW:35:ILE:HG13	2.14	0.47
31:CA:189(K):U:H2'	31:CA:189(L):G:C8	2.50	0.47
31:CA:564:C:H5'	47:CQ:32:TYR:CE2	2.49	0.47
31:CA:1215:G:H2'	31:CA:1216:G:C8	2.46	0.47
31:CA:1318:A:N6	44:CN:18:VAL:HG11	2.30	0.47
32:CB:60:ASP:OD2	32:CB:64:ARG:NE	2.43	0.47
36:CF:61:LEU:HD23	36:CF:63:TYR:OH	2.14	0.47
38:CH:25:ASP:OD2	38:CH:25:ASP:N	2.48	0.47
38:CH:58:TYR:O	38:CH:59:LEU:HD23	2.14	0.47
46:CP:28:ARG:HG2	46:CP:29:ASP:OD2	2.15	0.47
50:CT:16:HIS:O	50:CT:19:SER:OG	2.20	0.47
50:CT:72:LEU:HD23	50:CT:73:HIS:N	2.30	0.47
31:DA:25:C:H2'	31:DA:26:A:C8	2.50	0.47
31:DA:193:C:H2'	31:DA:194:C:C6	2.50	0.47
31:DA:512:U:H2'	31:DA:513:C:C6	2.49	0.47
31:DA:948:C:H2'	31:DA:949:A:H8	1.78	0.47
31:DA:1084:G:H5'	31:DA:1102:A:OP2	2.14	0.47
32:DB:27:LYS:HD2	32:DB:193:ASP:OD1	2.15	0.47
32:DB:145:LEU:O	32:DB:149:LEU:HB2	2.14	0.47
33:DC:6:HIS:CE1	33:DC:8:ILE:HB	2.50	0.47
33:DC:30:ARG:HH21	44:DN:38:GLY:HA2	1.80	0.47
34:DD:10:ARG:HB2	34:DD:40:PRO:HG3	1.97	0.47
34:DD:135:LEU:O	34:DD:137:SER:N	2.48	0.47
42:DL:49:ASN:ND2	42:DL:92:ASP:OD1	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:27:G:N2	1:AA:512:G:H1'	2.30	0.47
1:AA:581:C:OP1	16:AU:33:ARG:HG3	2.14	0.47
1:AA:646:A:H2'	1:AA:647:G:O4'	2.15	0.47
1:AA:748:G:O6	18:AW:90:ARG:NH1	2.48	0.47
1:AA:777:A:H2'	1:AA:778:G:H8	1.79	0.47
1:AA:1174:A:H1'	1:AA:1175:U:C5'	2.42	0.47
1:AA:1721:G:H2'	1:AA:1740:G:O6	2.15	0.47
1:AA:2136:C:N3	1:AA:2155:G:N2	2.61	0.47
1:AA:2349:G:H3'	1:AA:2350:C:H5''	1.97	0.47
1:AA:2536:G:C6	1:AA:2537:U:C4	3.03	0.47
1:AA:2784:C:H2'	1:AA:2785:C:C6	2.50	0.47
1:AA:2838:G:C6	1:AA:2839:G:C5	3.03	0.47
2:AB:33:G:N2	2:AB:50:G:C4	2.83	0.47
2:AB:61:G:C6	2:AB:62:C:C4	3.02	0.47
5:AF:28:ILE:HG12	5:AF:116:ASP:HB2	1.97	0.47
13:AR:92:GLY:HA2	13:AR:94:TYR:CE1	2.50	0.47
24:A2:66:GLU:HA	24:A2:69:ARG:NH1	2.30	0.47
25:A3:8:LEU:HD13	25:A3:31:LEU:HD23	1.97	0.47
1:BA:596:G:H2'	1:BA:597:U:O4'	2.15	0.47
1:BA:1531:C:N4	1:BA:1538:G:H1	2.06	0.47
1:BA:1782:C:H2'	1:BA:2608:G:O2'	2.15	0.47
1:BA:2199:A:N3	1:BA:2199:A:H2'	2.30	0.47
1:BA:2207:G:O2'	1:BA:2208:A:P	2.71	0.47
1:BA:2320:A:N3	1:BA:2320:A:H2'	2.30	0.47
1:BA:2377:A:H2'	1:BA:2378:A:C8	2.49	0.47
6:BG:138:GLN:HE22	6:BG:153:ARG:HB2	1.79	0.47
6:BG:145:THR:OG1	6:BG:146:TYR:N	2.47	0.47
7:BH:69:ARG:C	7:BH:70:THR:O	2.52	0.47
9:BN:76:SER:HB2	9:BN:77:GLY:H	1.49	0.47
11:BP:132:LYS:HE2	11:BP:132:LYS:HB2	1.52	0.47
24:B2:1:MET:N	24:B2:52:ASP:OD1	2.43	0.47
31:CA:34:C:H2'	31:CA:35:G:C8	2.50	0.47
31:CA:124:G:C6	31:CA:125:U:N3	2.83	0.47
31:CA:189(F):U:C4	47:CQ:72:ARG:NH2	2.83	0.47
31:CA:460:G:C6	31:CA:470:C:H5''	2.50	0.47
31:CA:1027:C:H2'	31:CA:1028:C:C4	2.50	0.47
31:CA:1063:C:H3'	31:CA:1064:G:H2'	1.96	0.47
33:CC:32:LEU:HD13	33:CC:59:ARG:HG2	1.97	0.47
33:CC:40:ARG:O	33:CC:44:GLU:HB2	2.15	0.47
42:CL:119:LYS:H	42:CL:119:LYS:HG3	1.49	0.47
31:DA:109:A:C6	31:DA:326:G:C6	3.03	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:617:G:H4'	46:DP:44:THR:O	2.14	0.47
31:DA:977:A:N6	31:DA:1224:G:O5'	2.48	0.47
31:DA:1092:A:H2'	31:DA:1093:A:C8	2.50	0.47
31:DA:1216:G:H5''	44:DN:5:ALA:HB3	1.97	0.47
34:DD:88:VAL:O	34:DD:92:VAL:HG23	2.15	0.47
40:DJ:51:ARG:HD3	44:DN:45:ARG:HH21	1.79	0.47
46:DP:23:ASP:O	46:DP:25:ARG:N	2.48	0.47
47:DQ:22:LEU:HD11	47:DQ:39:SER:HB3	1.97	0.47
1:AA:24:G:O2'	18:AW:77:ASP:HB3	2.15	0.47
1:AA:900:A:H2'	1:AA:901:A:H8	1.79	0.47
1:AA:1177:A:P	1:AA:1177:A:H2'	2.55	0.47
1:AA:1754:C:H2'	1:AA:1755:A:O4'	2.15	0.47
1:AA:1796:U:H2'	1:AA:1797:C:H6	1.79	0.47
1:AA:1934:C:H2'	1:AA:1935:G:O4'	2.15	0.47
1:AA:2019:A:OP2	27:A5:9:LYS:NZ	2.42	0.47
1:AA:2139:C:H42	1:AA:2152:G:H1	1.62	0.47
1:AA:2291:U:H5''	1:AA:2380:C:O2'	2.15	0.47
1:AA:2784:C:H2'	1:AA:2785:C:H6	1.80	0.47
2:AB:29:A:P	14:AS:32:LEU:HD12	2.55	0.47
9:AN:111:PRO:HA	9:AN:114:ARG:NH1	2.29	0.47
13:AR:80:PHE:O	13:AR:84:ALA:HB3	2.15	0.47
1:BA:1385:G:H4'	1:BA:1386:C:OP1	2.15	0.47
1:BA:1475:G:C2	1:BA:1517:G:C2	3.03	0.47
1:BA:1495:A:H2'	1:BA:1496:A:H8	1.78	0.47
1:BA:2171:A:H4'	1:BA:2172:U:OP1	2.14	0.47
1:BA:2749:A:H4'	7:BH:62:LYS:HB3	1.97	0.47
14:BS:32:LEU:O	14:BS:62:LYS:HE2	2.15	0.47
21:BZ:151:HIS:N	21:BZ:171:ILE:HG12	2.30	0.47
31:CA:323:U:H2'	31:CA:324:G:O4'	2.14	0.47
31:CA:633:G:H2'	31:CA:634:C:H6	1.80	0.47
31:CA:788:U:H2'	31:CA:789:U:O4'	2.15	0.47
31:CA:975:A:C2	31:CA:1357:A:H1'	2.49	0.47
31:CA:1433:A:C4	31:CA:1468:A:C2	3.04	0.47
35:CE:20:GLN:NE2	35:CE:21:ALA:O	2.48	0.47
42:CL:66:VAL:HG11	42:CL:98:TYR:CE1	2.50	0.47
50:CT:63:ILE:HG22	50:CT:77:ALA:HB1	1.97	0.47
31:DA:1202:G:H1'	44:DN:29:ARG:HD2	1.96	0.47
31:DA:1221:G:H4'	49:DS:77:THR:HG21	1.98	0.47
41:DK:58:PRO:HB2	41:DK:93:GLN:HG3	1.96	0.47
42:DL:38:THR:OG1	42:DL:57:LYS:HB3	2.15	0.47
45:DO:39:LEU:HD13	45:DO:56:LEU:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:DS:12:ASP:OD2	49:DS:37:ARG:NH1	2.48	0.47
50:DT:76:ALA:HA	50:DT:79:ARG:NH1	2.30	0.47
1:AA:898:C:H2'	1:AA:899:A:O4'	2.16	0.46
1:AA:1188:U:O2'	1:AA:1189:A:H5'	2.15	0.46
1:AA:1297:C:OP1	1:AA:2710:C:H4'	2.16	0.46
1:AA:1441:G:H2'	1:AA:1442:G:H8	1.79	0.46
1:AA:2251:G:OP2	12:AQ:82:ARG:NH1	2.48	0.46
1:AA:2327:A:H2'	1:AA:2328:A:H8	1.74	0.46
1:AA:2740:A:N6	1:AA:2764:A:C8	2.83	0.46
12:AQ:137:TYR:O	12:AQ:141:GLN:HG2	2.15	0.46
16:AU:47:TYR:HA	16:AU:50:ARG:NH2	2.30	0.46
26:A4:14:ILE:HA	26:A4:31:ILE:O	2.16	0.46
1:BA:248:G:H5'	1:BA:250:G:N7	2.30	0.46
1:BA:528:A:C8	1:BA:528:A:H3'	2.50	0.46
1:BA:674:G:H1'	5:BF:74:ARG:HD3	1.97	0.46
1:BA:862:G:H2'	1:BA:863:A:O4'	2.15	0.46
1:BA:987:G:O2'	1:BA:1000:A:N3	2.40	0.46
1:BA:1587:A:H2'	1:BA:1588:C:C6	2.50	0.46
1:BA:1721:G:H2'	1:BA:1740:G:O6	2.15	0.46
1:BA:2190:G:C2'	1:BA:2191:G:H5''	2.41	0.46
1:BA:2747:G:O6	1:BA:2755:C:H5''	2.15	0.46
5:BF:53:THR:CG2	5:BF:55:GLY:H	2.28	0.46
5:BF:65:TRP:CZ2	5:BF:75:HIS:HD2	2.32	0.46
7:BH:17:VAL:HG21	7:BH:50:VAL:HG21	1.97	0.46
13:BR:72:ASP:OD1	13:BR:75:LEU:HB2	2.14	0.46
31:CA:189(F):U:O2	47:CQ:63:ARG:NH2	2.47	0.46
31:CA:192:U:H2'	31:CA:193:C:C6	2.49	0.46
31:CA:559:A:H4'	31:CA:560:U:C3'	2.45	0.46
33:CC:184:TYR:OH	33:CC:186:PHE:HB2	2.15	0.46
45:CO:52:SER:O	45:CO:55:GLY:N	2.48	0.46
31:DA:64:G:H4'	31:DA:65:U:H3'	1.96	0.46
31:DA:167:G:H2'	31:DA:168:G:H8	1.80	0.46
31:DA:657:G:H4'	45:DO:28:GLN:HG2	1.97	0.46
31:DA:1359:C:O2'	31:DA:1362:C:N4	2.45	0.46
31:DA:1501:C:N4	31:DA:1504:G:C2	2.83	0.46
32:DB:24:TRP:CZ3	32:DB:29:ALA:HB2	2.45	0.46
32:DB:55:PHE:CE1	32:DB:218:ALA:HA	2.51	0.46
1:AA:566:U:H5''	11:AP:29:LYS:HE3	1.98	0.46
1:AA:619:G:H5''	1:AA:620:G:OP2	2.15	0.46
1:AA:1582:C:O2'	1:AA:1586:A:N3	2.40	0.46
1:AA:1892:C:O5'	1:AA:1892:C:H6	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2094:G:OP1	8:AI:22:LYS:HD2	2.15	0.46
1:AA:2440:C:N3	1:AA:2441:C:H1'	2.30	0.46
1:AA:2749:A:O4'	7:AH:63:SER:HA	2.15	0.46
1:AA:2869:G:H2'	1:AA:2870:C:O4'	2.15	0.46
2:AB:54:G:H21	6:AG:29:TRP:HZ2	1.64	0.46
2:AB:94:C:H2'	2:AB:95:C:C6	2.48	0.46
4:AE:37:ARG:HB2	4:AE:46:ALA:N	2.30	0.46
11:AP:50:ARG:HD3	30:A8:7:HIS:CD2	2.50	0.46
15:AT:50:ILE:HA	15:AT:99:LEU:HD12	1.96	0.46
21:AZ:23:LYS:O	21:AZ:25:PRO:HD3	2.14	0.46
1:BA:1806:C:O2'	3:BD:46:GLN:OE1	2.14	0.46
1:BA:2074:U:H2'	1:BA:2075:U:C6	2.50	0.46
1:BA:2304:G:H22	1:BA:2312:U:H3	1.63	0.46
5:BF:102:PRO:HB2	5:BF:105:VAL:HG23	1.96	0.46
15:BT:16:ARG:HG3	15:BT:79:HIS:HA	1.97	0.46
18:BW:13:SER:HB3	18:BW:16:LYS:HD2	1.96	0.46
24:B2:13:ALA:HA	24:B2:16:LEU:HD12	1.96	0.46
31:CA:237:C:H5''	47:CQ:25:ARG:CZ	2.45	0.46
31:CA:1157:A:C6	31:CA:1181:G:C5	3.04	0.46
31:CA:1414:U:H2'	31:CA:1415:G:H8	1.80	0.46
35:CE:78:HIS:CE1	35:CE:143:ARG:H	2.32	0.46
42:CL:66:VAL:HG11	42:CL:98:TYR:HE1	1.79	0.46
45:CO:20:GLY:O	45:CO:22:THR:HG23	2.16	0.46
31:DA:9:G:OP1	35:DE:122:GLU:HB2	2.16	0.46
31:DA:15:G:H2'	31:DA:16:A:H8	1.80	0.46
31:DA:59:A:H2'	31:DA:59:A:N3	2.31	0.46
31:DA:668:G:O4'	45:DO:49:ASP:HB2	2.16	0.46
31:DA:1055:A:H62	31:DA:1200:C:H42	1.63	0.46
32:DB:187:LEU:HD23	32:DB:201:ILE:CG2	2.45	0.46
32:DB:208:ILE:HA	32:DB:211:ILE:HD12	1.97	0.46
1:AA:64:A:H1'	19:AX:66:LEU:HB2	1.98	0.46
1:AA:518:G:H4'	18:AW:18:ARG:NH1	2.31	0.46
1:AA:706:A:H2'	1:AA:707:G:O4'	2.16	0.46
1:AA:723:G:H2'	1:AA:724:U:O4'	2.15	0.46
1:AA:921:G:C6	1:AA:922:U:C4	3.04	0.46
1:AA:1789:A:H2'	1:AA:1790:C:O4'	2.15	0.46
1:AA:2419:U:H2'	1:AA:2420:C:C6	2.50	0.46
7:AH:13:LYS:HA	7:AH:14:GLY:HA2	1.53	0.46
23:A1:54:ALA:HB1	23:A1:83:GLU:HG3	1.97	0.46
1:BA:1332:G:N2	1:BA:1609:A:HO2'	2.10	0.46
1:BA:1359:A:N6	1:BA:1372:U:H3	2.09	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:1790:C:H2'	1:BA:1791:A:C5	2.50	0.46
8:BI:3:VAL:HG12	8:BI:38:LEU:HA	1.97	0.46
15:BT:41:ARG:HH11	15:BT:41:ARG:HG3	1.80	0.46
31:CA:14:U:O2	31:CA:17:U:H5	1.98	0.46
31:CA:407:G:C2	31:CA:436:C:N3	2.83	0.46
31:CA:1082:G:H2'	31:CA:1083:U:O4'	2.15	0.46
31:CA:1500:A:H5''	31:CA:1508:G:H5''	1.97	0.46
31:CA:1510:U:H2'	31:CA:1511:G:C8	2.51	0.46
32:CB:32:ILE:HD13	32:CB:40:HIS:ND1	2.29	0.46
32:CB:54:THR:O	32:CB:58:ILE:HG13	2.15	0.46
34:CD:59:ARG:NH2	34:CD:66:ARG:HH12	2.13	0.46
31:DA:955:U:H2'	31:DA:956:U:C6	2.49	0.46
31:DA:1084:G:H2'	31:DA:1085:U:C6	2.50	0.46
33:DC:121:ALA:HB1	33:DC:189:ALA:HB2	1.97	0.46
1:AA:674:G:O2'	5:AF:74:ARG:HD3	2.15	0.46
1:AA:1740:G:H2'	1:AA:1741:A:C8	2.51	0.46
1:AA:1788:C:H5''	3:AD:225:ALA:HB1	1.98	0.46
1:AA:1913:A:O2'	52:AA:3001:T8B:C23	2.63	0.46
1:AA:2319:G:H22	14:AS:3:ARG:HE	1.61	0.46
1:AA:2463:C:O2'	1:AA:2464:C:H5'	2.14	0.46
2:AB:17:C:H2'	2:AB:18:G:O4'	2.15	0.46
3:AD:175:LEU:HD12	3:AD:185:VAL:HG21	1.98	0.46
4:AE:14:ILE:HB	15:AT:14:TYR:CZ	2.51	0.46
4:AE:101:ARG:HB2	4:AE:201:THR:HG21	1.98	0.46
20:AY:9:LYS:HA	20:AY:10:GLY:HA2	1.59	0.46
25:A3:6:VAL:O	25:A3:34:GLU:HA	2.15	0.46
1:BA:225:A:C2'	1:BA:226:G:H5'	2.46	0.46
1:BA:588:U:H2'	1:BA:589:C:C6	2.51	0.46
1:BA:1406:U:H2'	1:BA:1407:C:C6	2.51	0.46
1:BA:2305:A:H5''	6:BG:134:GLY:CA	2.46	0.46
3:BD:248:SER:C	3:BD:250:TRP:H	2.19	0.46
3:BD:274:ARG:H	3:BD:274:ARG:HG2	1.46	0.46
4:BE:30:PRO:HB3	4:BE:92:THR:HG22	1.97	0.46
4:BE:119:ARG:HA	4:BE:160:TYR:CD1	2.50	0.46
6:BG:165:THR:OG1	6:BG:168:GLU:HG3	2.16	0.46
11:BP:82:GLY:HA2	11:BP:113:LYS:O	2.14	0.46
14:BS:3:ARG:HG3	14:BS:4:LEU:N	2.25	0.46
31:CA:370:C:H2'	31:CA:371:G:C8	2.50	0.46
31:CA:568:G:N7	42:CL:5:PRO:HD3	2.31	0.46
31:CA:952:U:H2'	31:CA:953:G:H8	1.81	0.46
31:CA:984:C:H2'	31:CA:985:C:C6	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:1206:G:O4'	33:CC:194:GLY:HA2	2.16	0.46
32:CB:87:ARG:NH1	32:CB:220:ASP:OD1	2.37	0.46
42:CL:32:PHE:CE1	42:CL:86:ARG:HG3	2.49	0.46
43:CM:5:ALA:HA	43:CM:61:GLU:CG	2.46	0.46
31:DA:682:G:C6	31:DA:709:G:C6	3.04	0.46
31:DA:1147:C:C4	31:DA:1148:U:C4	3.03	0.46
31:DA:1274:G:N2	31:DA:1275:A:H62	2.13	0.46
32:DB:77:ALA:HB1	32:DB:211:ILE:HG21	1.96	0.46
41:DK:31:THR:HA	41:DK:42:TRP:HA	1.98	0.46
46:DP:6:LEU:HD11	46:DP:73:LEU:HD12	1.97	0.46
1:AA:2144:U:O2	1:AA:2148:G:N1	2.48	0.46
1:AA:2274:A:N1	1:AA:2276:G:H1'	2.30	0.46
2:AB:2:C:C2	2:AB:3:C:C5	3.03	0.46
2:AB:87:G:N2	2:AB:89:G:H3'	2.31	0.46
4:AE:110:GLY:O	13:AR:3:HIS:CE1	2.69	0.46
8:AI:44:LEU:HD13	8:AI:44:LEU:HA	1.76	0.46
10:AO:14:THR:O	10:AO:52:VAL:HG22	2.16	0.46
10:AO:20:MET:HE3	10:AO:44:LYS:HE3	1.98	0.46
13:AR:101:ALA:HA	27:A5:44:THR:CG2	2.46	0.46
1:BA:247:G:H4'	1:BA:386:G:C5	2.50	0.46
1:BA:885:C:N4	1:BA:890:A:C6	2.62	0.46
1:BA:2736:G:N2	1:BA:2769:C:H1'	2.31	0.46
31:CA:33:A:H2'	31:CA:34:C:C6	2.50	0.46
31:CA:579:G:O5'	31:CA:728:A:H1'	2.16	0.46
31:CA:1210:C:H2'	31:CA:1211:U:H5''	1.98	0.46
31:CA:1288:A:H2'	31:CA:1289:A:O4'	2.16	0.46
36:CF:19:LEU:HD23	36:CF:19:LEU:O	2.15	0.46
31:DA:15:G:H2'	31:DA:16:A:C8	2.51	0.46
31:DA:630:G:H2'	31:DA:631:G:C8	2.49	0.46
31:DA:782:A:H4'	31:DA:1514:C:O2'	2.16	0.46
31:DA:921:U:O2'	35:DE:19:MET:O	2.27	0.46
31:DA:991:U:O2'	31:DA:992:U:OP2	2.32	0.46
31:DA:1237:C:HO2'	31:DA:1300:G:H22	1.56	0.46
31:DA:1496:C:H2'	31:DA:1497:G:O4'	2.15	0.46
50:DT:21:LYS:O	50:DT:25:ARG:HG3	2.15	0.46
1:AA:39:C:H2'	1:AA:40:C:C6	2.50	0.46
1:AA:441:U:H2'	1:AA:442:G:C8	2.50	0.46
1:AA:907:U:O2'	12:AQ:101:ARG:NH2	2.36	0.46
1:AA:950:G:H2'	1:AA:951:C:O4'	2.15	0.46
1:AA:1227:G:OP1	16:AU:13:LYS:HG2	2.16	0.46
1:AA:1248:G:O2'	16:AU:3:ARG:HA	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2543:G:N3	1:AA:2765:A:H2'	2.31	0.46
1:AA:2630:G:H2'	1:AA:2631:G:C8	2.51	0.46
7:AH:11:VAL:CG2	7:AH:50:VAL:HG23	2.45	0.46
8:AI:105:HIS:H	8:AI:105:HIS:HD2	1.57	0.46
18:AW:4:LYS:HB2	18:AW:106:ILE:HG12	1.97	0.46
23:A1:7:ILE:HG21	23:A1:69:LYS:HG2	1.97	0.46
1:BA:1999:C:H5''	1:BA:2723:C:O2'	2.15	0.46
5:BF:37:VAL:HG21	11:BP:6:LEU:CD1	2.45	0.46
6:BG:121:ASN:HA	6:BG:122:PRO:HD3	1.74	0.46
31:CA:174:C:H2'	31:CA:175:C:C6	2.50	0.46
31:CA:184:G:N2	31:CA:194:C:C2	2.83	0.46
31:CA:968:A:H8	31:CA:968:A:OP1	1.99	0.46
31:CA:1492:A:H8	42:CL:47:LYS:HD2	1.79	0.46
31:DA:266:G:H5''	31:DA:267:C:C5	2.51	0.46
33:DC:111:LEU:HD23	33:DC:141:VAL:HG13	1.98	0.46
35:DE:129:ILE:O	35:DE:132:ALA:HB3	2.14	0.46
36:DF:14:LEU:HB3	36:DF:15:ASP:H	1.49	0.46
39:DI:26:VAL:HG13	39:DI:61:ALA:HB3	1.98	0.46
42:DL:69:TYR:HD2	42:DL:99:HIS:CD2	2.34	0.46
43:DM:44:ARG:HB2	43:DM:47:ASP:OD2	2.15	0.46
1:AA:8:A:H2'	1:AA:9:U:H6	1.80	0.46
1:AA:53:A:H2'	1:AA:54:G:O4'	2.15	0.46
1:AA:330:A:H2	1:AA:1210:A:C2'	2.26	0.46
1:AA:795:C:O2'	1:AA:796:C:H5'	2.15	0.46
1:AA:1000:A:C6	1:AA:1001:A:C6	3.03	0.46
1:AA:1142(A):A:C8	1:AA:1144:G:N7	2.84	0.46
1:AA:1270:C:H5''	1:AA:1271:G:O5'	2.16	0.46
1:AA:1913:A:C4'	1:AA:1914:C:H5''	2.46	0.46
1:AA:2223:G:H2'	1:AA:2224:G:H5'	1.97	0.46
3:AD:75:ILE:HG21	3:AD:99:ASP:HB2	1.97	0.46
13:AR:50:HIS:CE1	13:AR:54:LEU:HD13	2.51	0.46
1:BA:41:C:H2'	1:BA:42:G:O4'	2.16	0.46
1:BA:95:G:H4'	24:B2:46:GLN:H	1.81	0.46
1:BA:335:C:H4'	20:BY:73:ARG:HD3	1.97	0.46
1:BA:443:A:N6	5:BF:41:LEU:O	2.40	0.46
1:BA:489:G:N7	18:BW:49:LYS:NZ	2.64	0.46
1:BA:738:G:H1'	1:BA:759:G:N2	2.30	0.46
1:BA:2205:C:O2	1:BA:2220:G:C2	2.69	0.46
1:BA:2287:A:C4	1:BA:2289:G:C8	3.04	0.46
4:BE:24:THR:HG22	4:BE:186:GLY:O	2.15	0.46
5:BF:104:LYS:O	5:BF:108:LYS:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:BN:138:LEU:HA	9:BN:138:LEU:HD23	1.62	0.46
10:BO:70:LYS:HB3	10:BO:70:LYS:HE2	1.71	0.46
11:BP:90:ARG:HG2	11:BP:91:PHE:CD2	2.51	0.46
19:BX:72:LYS:HB3	19:BX:72:LYS:HE2	1.66	0.46
27:B5:36:CYS:SG	27:B5:46:CYS:SG	3.13	0.46
31:CA:834:C:C2	31:CA:853:G:C2	3.04	0.46
31:CA:949:A:OP1	43:CM:101:GLN:HB3	2.16	0.46
31:CA:1014:A:C2	31:CA:1219:U:H1'	2.51	0.46
31:CA:1070:U:H2'	31:CA:1071:C:H6	1.80	0.46
31:CA:1285:A:H4'	31:CA:1286:A:O5'	2.16	0.46
31:CA:1298:C:C4	37:CG:114:ARG:HD3	2.51	0.46
38:CH:6:ILE:HB	38:CH:85:ARG:NH1	2.31	0.46
38:CH:44:PHE:HB3	38:CH:80:ILE:HG12	1.96	0.46
31:DA:238:G:C6	31:DA:239:U:C4	3.04	0.46
31:DA:445:G:C4	31:DA:446:G:C8	3.04	0.46
31:DA:477:A:H2'	31:DA:479:C:H6	1.80	0.46
31:DA:949:A:H1'	31:DA:1364:U:H3	1.81	0.46
31:DA:992:U:O2	31:DA:992:U:H2'	2.16	0.46
31:DA:1097:C:H2'	31:DA:1098:C:C6	2.51	0.46
33:DC:8:ILE:O	33:DC:11:ARG:N	2.43	0.46
33:DC:11:ARG:O	33:DC:13:GLY:N	2.48	0.46
35:DE:78:HIS:CE1	35:DE:143:ARG:H	2.33	0.46
43:DM:11:ARG:HA	43:DM:45:VAL:HB	1.97	0.46
50:DT:16:HIS:O	50:DT:19:SER:OG	2.20	0.46
50:DT:53:LEU:O	50:DT:57:ARG:HG3	2.16	0.46
50:DT:77:ALA:O	50:DT:81:LYS:HG3	2.15	0.46
1:AA:70:G:H5''	1:AA:112:U:O2	2.15	0.46
1:AA:280:C:C2	1:AA:361:G:C2	3.04	0.46
1:AA:300:A:H3'	20:AY:84:ARG:NH2	2.31	0.46
1:AA:1313:U:H2'	1:AA:1610:A:C2	2.51	0.46
1:AA:2144:U:H2'	1:AA:2146:C:C4	2.50	0.46
1:AA:2280:G:C2'	1:AA:2281:C:H5'	2.46	0.46
1:AA:2293:C:H2'	1:AA:2294:C:C6	2.50	0.46
1:AA:2755:C:O2'	1:AA:2756:U:H6	1.98	0.46
6:AG:16:ARG:NH2	6:AG:28:VAL:O	2.49	0.46
6:AG:174:GLU:HG3	6:AG:178:PHE:O	2.16	0.46
9:AN:99:LEU:O	9:AN:103:VAL:HG23	2.16	0.46
21:AZ:166:SER:HA	21:AZ:167:PRO:HD3	1.78	0.46
1:BA:1412:A:H2'	1:BA:1413:G:C8	2.51	0.46
1:BA:2031:A:C6	1:BA:2498:C:H1'	2.51	0.46
1:BA:2181:G:H2'	1:BA:2182:G:C8	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:2261:C:C6	22:B0:16:SER:HB3	2.51	0.46
9:BN:12:ARG:HD3	9:BN:14:VAL:HG23	1.97	0.46
12:BQ:110:THR:HG23	12:BQ:113:GLN:OE1	2.16	0.46
13:BR:18:LEU:HD23	13:BR:18:LEU:HA	1.52	0.46
14:BS:61:ASN:O	14:BS:64:GLU:N	2.48	0.46
31:CA:377:G:H2'	31:CA:378:G:C8	2.51	0.46
31:CA:598:U:H4'	38:CH:94:TYR:CG	2.50	0.46
31:CA:654:G:C5	31:CA:655:A:C8	3.04	0.46
31:CA:868:C:H2'	31:CA:869:G:O4'	2.16	0.46
31:DA:229:U:O2'	46:DP:23:ASP:OD2	2.29	0.46
31:DA:380:G:N2	31:DA:384:G:C5	2.84	0.46
31:DA:683:G:H2'	31:DA:684:A:H8	1.80	0.46
31:DA:929:G:C6	31:DA:930:C:C4	3.04	0.46
31:DA:1030(B):C:C2'	31:DA:1030(C):G:H5'	2.46	0.46
31:DA:1086:U:H2'	31:DA:1087:G:C8	2.51	0.46
31:DA:1109:C:H2'	31:DA:1110:A:O4'	2.16	0.46
31:DA:1277:C:H2'	31:DA:1278:U:H5'	1.96	0.46
31:DA:1387:G:H2'	31:DA:1388:C:C6	2.51	0.46
34:DD:13:ARG:O	34:DD:15:GLU:N	2.48	0.46
34:DD:159:ARG:O	34:DD:163:GLU:N	2.41	0.46
42:DL:33:ARG:O	42:DL:85:ILE:HB	2.16	0.46
1:AA:323:G:C2	1:AA:333:G:H1'	2.50	0.46
1:AA:370:G:H4'	1:AA:371:A:OP2	2.15	0.46
1:AA:593:G:H4'	30:A8:63:PRO:HB3	1.97	0.46
1:AA:1479:G:H5''	1:AA:1560:G:H4'	1.98	0.46
1:AA:1913:A:H3'	52:AA:3001:T8B:O2	2.16	0.46
1:AA:2133:G:C2	1:AA:2157:G:H2'	2.51	0.46
1:AA:2164:C:H3'	1:AA:2165:G:C8	2.48	0.46
1:AA:2306:C:H3'	1:AA:2307:G:C8	2.50	0.46
1:AA:2628:C:O2'	1:AA:2781:A:H2'	2.16	0.46
1:AA:2789:C:O3'	1:AA:2790:A:H4'	2.15	0.46
5:AF:17:ARG:O	5:AF:18:ARG:HG2	2.16	0.46
8:AI:86:THR:HA	8:AI:123:LEU:HD11	1.98	0.46
20:AY:9:LYS:HA	20:AY:27:VAL:HB	1.98	0.46
1:BA:536:A:H2'	1:BA:537:C:C6	2.51	0.46
1:BA:1045:A:O2'	1:BA:1047:G:C4	2.62	0.46
1:BA:1047:G:C2'	1:BA:1110:G:H22	2.23	0.46
1:BA:1051:G:H5''	1:BA:1052:C:OP2	2.15	0.46
1:BA:1825:A:OP1	3:BD:249:PRO:HD3	2.16	0.46
1:BA:2884:U:H1'	27:B5:53:ALA:HB2	1.97	0.46
5:BF:117:ARG:HD3	5:BF:117:ARG:HA	1.65	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BG:11:TYR:CZ	6:BG:16:ARG:HD3	2.51	0.46
14:BS:58:LEU:HB2	14:BS:59:LYS:CB	2.44	0.46
21:BZ:151:HIS:C	21:BZ:153:SER:H	2.12	0.46
27:B5:36:CYS:O	27:B5:38:ALA:N	2.49	0.46
31:CA:194:C:C2'	31:CA:195:A:H5''	2.46	0.46
31:CA:414:A:H2'	31:CA:415:A:H8	1.81	0.46
31:CA:1358:U:H5''	44:CN:33:VAL:O	2.16	0.46
47:CQ:22:LEU:HD11	47:CQ:39:SER:HB3	1.98	0.46
31:DA:294:U:H2'	31:DA:295:C:C6	2.50	0.46
31:DA:799:G:O6	31:DA:800:G:C2	2.69	0.46
31:DA:1056:U:H2'	31:DA:1057:G:H8	1.80	0.46
31:DA:1187:G:H2'	31:DA:1188:A:C8	2.51	0.46
31:DA:1279:A:H5''	31:DA:1280:A:OP1	2.16	0.46
31:DA:1290:G:H2'	31:DA:1290:G:N3	2.30	0.46
39:DI:102:LEU:O	39:DI:103:THR:HG22	2.15	0.46
45:DO:36:ILE:O	45:DO:39:LEU:N	2.49	0.46
46:DP:26:ARG:CZ	46:DP:31:LYS:HB3	2.45	0.46
48:DR:66:LEU:HG	48:DR:70:ILE:HD11	1.97	0.46
1:AA:2531:A:N3	1:AA:2658:C:O2'	2.34	0.46
2:AB:29:A:OP2	14:AS:32:LEU:N	2.41	0.46
5:AF:117:ARG:HD3	5:AF:117:ARG:HA	1.67	0.46
12:AQ:72:LYS:HA	12:AQ:73:PRO:HD3	1.73	0.46
14:AS:14:VAL:HG21	14:AS:90:GLY:O	2.16	0.46
14:AS:59:LYS:HB3	14:AS:60:GLY:HA2	1.97	0.46
18:AW:37:ARG:HD3	18:AW:38:TYR:CE2	2.51	0.46
1:BA:616:G:H2'	1:BA:618:C:O4'	2.16	0.46
1:BA:1002:G:H2'	1:BA:1003:G:O4'	2.16	0.46
1:BA:1846:G:N2	1:BA:1895:C:C2	2.84	0.46
1:BA:2138:C:H42	1:BA:2153:G:H1	1.63	0.46
1:BA:2464:C:O2'	1:BA:2465:C:H6	1.98	0.46
1:BA:2760:C:H2'	1:BA:2761:G:H5''	1.97	0.46
5:BF:129:PHE:HB3	5:BF:132:VAL:HG22	1.98	0.46
6:BG:107:LEU:HD11	6:BG:178:PHE:CE1	2.50	0.46
8:BI:29:TYR:C	8:BI:32:PRO:HD2	2.36	0.46
12:BQ:7:MET:HE3	12:BQ:7:MET:HB2	1.89	0.46
20:BY:90:LEU:HD12	20:BY:90:LEU:HA	1.82	0.46
21:BZ:111:VAL:HG12	21:BZ:112:ARG:N	2.31	0.46
24:B2:10:LEU:HD23	24:B2:10:LEU:HA	1.83	0.46
28:B6:21:TYR:CE2	28:B6:38:LYS:HG2	2.51	0.46
31:CA:202:U:O2'	31:CA:203:U:O5'	2.26	0.46
31:CA:203:U:H2'	31:CA:203:U:OP2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:1074:G:C2	31:CA:1102:A:C6	3.04	0.46
31:CA:1300:G:HO2'	31:CA:1301:U:P	2.37	0.46
36:CF:11:ASN:HB3	36:CF:14:LEU:HG	1.98	0.46
36:CF:86:ARG:O	36:CF:87:ARG:HG2	2.16	0.46
42:CL:8:ASN:O	42:CL:12:ARG:HG3	2.16	0.46
44:CN:40:CYS:HB2	44:CN:43:CYS:HB2	1.97	0.46
46:CP:68:ASP:C	46:CP:70:ALA:H	2.20	0.46
31:DA:192:U:H4'	50:DT:57:ARG:HD2	1.97	0.46
31:DA:303:A:H2'	31:DA:304:U:O4'	2.16	0.46
31:DA:433:C:H2'	31:DA:434:U:H6	1.79	0.46
31:DA:938:A:C6	31:DA:939:G:C5	3.04	0.46
31:DA:1034:G:C2	31:DA:1035:A:C4	3.04	0.46
31:DA:1168:A:H2'	31:DA:1169:A:O4'	2.15	0.46
31:DA:1220:G:H2'	31:DA:1221:G:O4'	2.16	0.46
31:DA:1324:A:H2'	31:DA:1325:C:C6	2.51	0.46
38:DH:124:ALA:HB1	38:DH:129:VAL:O	2.16	0.46
1:AA:12:U:O2	1:AA:12:U:H2'	2.17	0.45
1:AA:1109:C:C5	1:AA:1110:G:C6	3.00	0.45
1:AA:1913:A:C3'	1:AA:1914:C:H5''	2.46	0.45
1:AA:2012:G:H8	1:AA:2012:G:O5'	1.98	0.45
1:AA:2127:G:O6	1:AA:2161:C:N3	2.49	0.45
1:AA:2730:C:H4'	4:AE:168:MET:O	2.16	0.45
7:AH:86:GLU:HB2	7:AH:165:ALA:HB2	1.98	0.45
7:AH:149:ARG:HA	7:AH:162:ILE:CG2	2.46	0.45
8:AI:127:VAL:HA	8:AI:140:LEU:O	2.16	0.45
9:AN:28:THR:HG22	9:AN:29:LYS:N	2.31	0.45
9:AN:56:ASN:H	9:AN:125:GLY:H	1.64	0.45
14:AS:53:SER:O	14:AS:57:LYS:N	2.49	0.45
25:A3:43:ILE:O	25:A3:47:VAL:HG23	2.17	0.45
1:BA:675:A:C4	1:BA:804:A:C2	3.03	0.45
1:BA:1275:A:N1	1:BA:1295:C:O2'	2.39	0.45
1:BA:1453:U:O2'	1:BA:1455:G:N7	2.42	0.45
1:BA:2467:C:H4'	12:BQ:123:HIS:CD2	2.51	0.45
1:BA:2881:C:H2'	1:BA:2882:A:O4'	2.16	0.45
3:BD:26:LYS:HE2	3:BD:28:GLU:O	2.16	0.45
31:CA:579:G:C6	31:CA:580:U:C4	3.03	0.45
31:CA:655:A:N1	31:CA:754:C:N4	2.64	0.45
31:CA:657:G:C2	31:CA:750:G:C5	3.05	0.45
33:CC:127:ARG:HB3	33:CC:127:ARG:NH1	2.32	0.45
49:CS:41:VAL:O	49:CS:43:GLU:N	2.49	0.45
31:DA:130:A:O2'	31:DA:131:C:O5'	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:142:G:H2'	31:DA:143:A:H8	1.80	0.45
31:DA:262:A:C6	31:DA:263:A:C6	3.04	0.45
31:DA:352:C:O2'	31:DA:354:G:OP1	2.23	0.45
31:DA:540:G:H2'	31:DA:541:G:O4'	2.15	0.45
31:DA:1148:U:H2'	31:DA:1149:C:O4'	2.16	0.45
31:DA:1266:G:N2	31:DA:1268:A:H3'	2.31	0.45
31:DA:1511:G:H2'	31:DA:1512:U:O4'	2.17	0.45
34:DD:173:TRP:HB2	34:DD:187:ARG:O	2.16	0.45
43:DM:102:ARG:NE	43:DM:105:THR:OG1	2.49	0.45
46:DP:21:VAL:HG13	46:DP:33:ILE:HB	1.97	0.45
1:AA:342:G:H2'	1:AA:343:C:C6	2.51	0.45
1:AA:863:A:H2'	1:AA:864:G:C8	2.50	0.45
1:AA:1384:A:N3	1:AA:1405:U:H1'	2.31	0.45
1:AA:1417:C:H2'	1:AA:1418:G:O4'	2.15	0.45
7:AH:116:GLU:HA	7:AH:117:PRO:HD2	1.79	0.45
7:AH:124:GLU:HB2	7:AH:132:ARG:HB3	1.98	0.45
8:AI:45:LYS:O	8:AI:48:GLU:N	2.49	0.45
8:AI:77:LEU:HD12	8:AI:142:VAL:CG1	2.46	0.45
8:AI:79:ILE:HA	8:AI:80:PRO:HD2	1.73	0.45
14:AS:101:LEU:O	14:AS:102:ALA:HB3	2.16	0.45
15:AT:118:ARG:HA	15:AT:118:ARG:HD2	1.64	0.45
1:BA:102:G:O2'	1:BA:103:A:P	2.73	0.45
1:BA:280:C:C2	1:BA:361:G:C2	3.04	0.45
1:BA:783:A:O2'	1:BA:785:G:OP1	2.16	0.45
1:BA:1331:A:O2'	1:BA:1332:G:C8	2.69	0.45
1:BA:1541:G:H3'	1:BA:1542:A:H2'	1.98	0.45
1:BA:2469:A:H5''	1:BA:2470:G:OP2	2.16	0.45
3:BD:83:GLU:OE1	3:BD:104:TYR:OH	2.26	0.45
8:BI:93:THR:O	8:BI:97:ILE:N	2.28	0.45
9:BN:104:LYS:HB2	9:BN:117:PHE:CE1	2.50	0.45
12:BQ:109:VAL:HG13	12:BQ:113:GLN:HB2	1.99	0.45
31:CA:417:C:H2'	31:CA:418:C:H6	1.80	0.45
31:CA:655:A:C2	31:CA:754:C:C4	3.04	0.45
31:CA:939:G:H2'	31:CA:940:C:C6	2.50	0.45
31:CA:1085:U:H3'	31:CA:1086:U:C5	2.51	0.45
31:CA:1096:C:H2'	31:CA:1097:C:H6	1.81	0.45
32:CB:97:TRP:CZ2	32:CB:173:ALA:HA	2.52	0.45
35:CE:78:HIS:ND1	35:CE:142:LEU:HD23	2.30	0.45
45:CO:54:ARG:O	45:CO:58:MET:HG3	2.16	0.45
46:CP:75:ARG:O	46:CP:78:GLY:N	2.49	0.45
47:CQ:68:ARG:H	47:CQ:70:ARG:NH1	2.14	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
50:CT:36:LEU:HD13	50:CT:36:LEU:HA	1.81	0.45
50:CT:41:ILE:O	50:CT:45:GLN:N	2.40	0.45
31:DA:186:C:O2'	50:DT:85:MET:SD	2.64	0.45
31:DA:1256:A:H2	31:DA:1277:C:N3	2.14	0.45
31:DA:1392:G:N2	31:DA:1502:A:C8	2.83	0.45
31:DA:1446:U:H4'	31:DA:1447:A:C5	2.51	0.45
33:DC:132:ARG:O	33:DC:136:GLN:HB2	2.17	0.45
35:DE:68:GLU:CG	35:DE:70:PRO:HG3	2.46	0.45
47:DQ:20:THR:HG21	47:DQ:41:LYS:HD2	1.98	0.45
1:AA:116:C:H2'	1:AA:117:G:O4'	2.16	0.45
1:AA:325:G:N2	1:AA:337:C:O2	2.49	0.45
1:AA:510:C:H2'	1:AA:511:U:O4'	2.16	0.45
1:AA:1250:G:N7	11:AP:18:ARG:NH2	2.65	0.45
1:AA:1328:G:H2'	1:AA:1330:C:C5	2.50	0.45
1:AA:2055:C:H5'	1:AA:2056:G:O5'	2.16	0.45
1:AA:2096:U:H3	1:AA:2193:G:H1	1.65	0.45
1:AA:2115:G:H21	1:AA:2171:A:H61	1.63	0.45
1:AA:2256:G:H2'	1:AA:2257:U:H6	1.81	0.45
1:AA:2283:C:C2	1:AA:2389:G:C2	3.04	0.45
4:AE:152:LYS:HB2	9:AN:77:GLY:O	2.15	0.45
7:AH:7:LEU:HA	7:AH:8:PRO:HD3	1.67	0.45
1:BA:90:U:O2'	1:BA:92:A:C8	2.65	0.45
31:CA:7:G:H5'	31:CA:298:A:O4'	2.16	0.45
31:CA:655:A:C2	31:CA:656:C:C2	3.04	0.45
31:CA:1361:G:H8	31:CA:1361:G:O5'	2.00	0.45
34:CD:165:MET:O	34:CD:167:GLY:N	2.49	0.45
36:CF:30:LEU:HB3	36:CF:35:ALA:HB3	1.98	0.45
42:CL:53:ARG:CB	42:CL:93:LEU:HD11	2.46	0.45
46:CP:72:ARG:HE	46:CP:73:LEU:CD2	2.30	0.45
47:CQ:45:HIS:HB3	47:CQ:72:ARG:HB3	1.98	0.45
50:CT:38:LYS:O	50:CT:41:ILE:HG13	2.16	0.45
31:DA:60:A:H4'	31:DA:61:G:O5'	2.17	0.45
31:DA:973:G:H4'	40:DJ:54:PHE:O	2.17	0.45
31:DA:1010:G:H2'	31:DA:1011:G:H8	1.79	0.45
31:DA:1015:A:N6	31:DA:1016:A:C6	2.85	0.45
32:DB:24:TRP:CZ3	32:DB:26:PRO:HA	2.51	0.45
35:DE:33:VAL:HG21	35:DE:109:ILE:HA	1.99	0.45
1:AA:296:C:O3'	20:AY:95:LYS:NZ	2.49	0.45
1:AA:322:A:H3'	5:AF:169:ASN:ND2	2.31	0.45
1:AA:372:G:O2'	1:AA:373:U:OP2	2.31	0.45
1:AA:448:U:O4	1:AA:583:G:H1'	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1614:A:H8	1:AA:1614:A:O5'	2.00	0.45
1:AA:2082:A:H2'	1:AA:2083:G:O4'	2.16	0.45
1:AA:2821:A:H2'	1:AA:2822:G:O4'	2.16	0.45
1:AA:2850:A:N7	1:AA:2868:A:O2'	2.33	0.45
52:AA:3001:T8B:H8	31:DA:1492:A:O5'	2.15	0.45
2:AB:31:C:O2'	2:AB:53:A:N6	2.49	0.45
3:AD:172:TYR:CD1	3:AD:186:HIS:HA	2.52	0.45
4:AE:2:LYS:HA	4:AE:84:PHE:CE2	2.52	0.45
8:AI:94:ALA:HA	8:AI:97:ILE:HG13	1.99	0.45
10:AO:103:ALA:HB1	10:AO:105:GLU:OE1	2.16	0.45
26:A4:14:ILE:HG23	26:A4:31:ILE:HB	1.98	0.45
1:BA:300:A:H3'	20:BY:84:ARG:NH2	2.32	0.45
1:BA:597:U:H2'	1:BA:598:G:C8	2.51	0.45
1:BA:742:G:H4'	1:BA:1676:A:H5'	1.98	0.45
1:BA:1357:U:H2'	1:BA:1358:G:O4'	2.16	0.45
1:BA:1638:C:H5''	1:BA:2710:C:O2'	2.16	0.45
1:BA:1711:C:H2'	1:BA:1712:C:C6	2.52	0.45
1:BA:1913:A:H3'	52:BA:3001:T8B:O10	2.13	0.45
1:BA:1914:C:O2'	1:BA:1915:U:O5'	2.29	0.45
1:BA:2052:G:H4'	4:BE:143:ASN:O	2.17	0.45
1:BA:2302:G:N1	1:BA:2315:G:C6	2.84	0.45
1:BA:2392:A:N3	11:BP:61:ARG:HG2	2.32	0.45
1:BA:2463:C:H2'	1:BA:2464:C:H5'	1.99	0.45
5:BF:140:LEU:HD13	5:BF:140:LEU:HA	1.78	0.45
6:BG:11:TYR:O	6:BG:15:VAL:HB	2.17	0.45
26:B4:15:ILE:O	26:B4:32:TYR:HA	2.17	0.45
31:CA:258:G:H2'	31:CA:259:G:H8	1.82	0.45
31:CA:263:A:OP1	50:CT:79:ARG:NH1	2.49	0.45
31:CA:718:G:H5'	41:CK:117:ASN:HB2	1.97	0.45
31:CA:750:G:N2	31:CA:751:U:C2	2.85	0.45
31:CA:1005:A:C6	31:CA:1025:U:H1'	2.51	0.45
31:CA:1250:A:H4'	39:CI:68:GLY:N	2.32	0.45
31:CA:1409:C:H2'	31:CA:1410:G:H8	1.80	0.45
35:CE:105:VAL:O	35:CE:109:ILE:HG13	2.17	0.45
37:CG:23:VAL:HG13	37:CG:43:PHE:CZ	2.52	0.45
31:DA:1036:G:H2'	31:DA:1036:G:N3	2.29	0.45
31:DA:1232:U:O5'	31:DA:1232:U:H6	1.99	0.45
33:DC:181:ASN:ND2	33:DC:204:LEU:HB2	2.32	0.45
35:DE:93:PRO:HG2	38:DH:105:ARG:NE	2.31	0.45
38:DH:92:ARG:HD3	38:DH:92:ARG:HA	1.61	0.45
39:DI:4:TYR:CE2	39:DI:88:TYR:HD1	2.34	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:DM:15:VAL:O	43:DM:19:LEU:HD22	2.16	0.45
1:AA:45:C:H2'	1:AA:47:C:C6	2.52	0.45
1:AA:107:C:C2	1:AA:108:U:C5	3.04	0.45
1:AA:610:G:N2	1:AA:619:G:H1'	2.31	0.45
1:AA:683:C:O5'	1:AA:683:C:H6	2.00	0.45
1:AA:861:A:C2	1:AA:917:A:C4	3.04	0.45
1:AA:1359:A:N6	1:AA:1372:U:H3	2.10	0.45
1:AA:1412:A:C2	1:AA:1591:G:C2	3.05	0.45
1:AA:1779:U:C2	1:AA:1783:A:N7	2.85	0.45
1:AA:1839:G:C8	1:AA:1927:A:H1'	2.52	0.45
1:AA:2156:G:O6	1:AA:2157:G:N2	2.40	0.45
1:AA:2174:C:H2'	1:AA:2175:C:C6	2.46	0.45
1:AA:2627:G:N2	1:AA:2777:G:OP2	2.41	0.45
1:AA:2727:G:O3'	10:AO:70:LYS:HD3	2.16	0.45
7:AH:149:ARG:NH2	7:AH:154:PRO:HG2	2.31	0.45
9:AN:4:TYR:CD2	16:AU:100:VAL:HG11	2.50	0.45
15:AT:95:ARG:HG2	15:AT:95:ARG:NH1	2.31	0.45
16:AU:20:LEU:HD23	16:AU:20:LEU:HA	1.87	0.45
1:BA:885:C:OP1	1:BA:885:C:H4'	2.16	0.45
1:BA:1009:A:O4'	16:BU:59:ARG:HG2	2.17	0.45
1:BA:1176:G:H4'	1:BA:1177:A:OP2	2.17	0.45
1:BA:2781:A:H5''	1:BA:2782:G:H5'	1.98	0.45
2:BB:2:C:H2'	2:BB:3:C:C6	2.52	0.45
8:BI:77:LEU:HD23	8:BI:77:LEU:HA	1.79	0.45
11:BP:126:VAL:CG1	11:BP:148:LEU:HD13	2.47	0.45
14:BS:74:ALA:O	14:BS:77:ALA:HB3	2.17	0.45
31:CA:175:C:H2'	31:CA:176:C:H6	1.80	0.45
31:CA:947:G:O2'	31:CA:1306:A:H4'	2.17	0.45
32:CB:17:PHE:HB3	32:CB:44:LEU:HD11	1.98	0.45
33:CC:139:GLN:HG3	33:CC:143:GLU:OE1	2.15	0.45
37:CG:146:GLU:C	37:CG:148:ASN:H	2.20	0.45
46:CP:72:ARG:HE	46:CP:73:LEU:HD23	1.81	0.45
31:DA:578:C:O2'	31:DA:728:A:N3	2.38	0.45
31:DA:889:A:H4'	31:DA:890:G:OP1	2.17	0.45
31:DA:1060:C:C5	33:DC:2:GLY:HA3	2.51	0.45
31:DA:1235:U:O2'	31:DA:1305:G:OP1	2.28	0.45
36:DF:2:ARG:NH1	36:DF:69:GLU:HB3	2.32	0.45
1:AA:51:G:N3	1:AA:119:A:C2	2.85	0.45
1:AA:52:A:OP2	1:AA:117:G:N1	2.30	0.45
1:AA:686:G:N2	1:AA:788:A:H61	2.15	0.45
1:AA:1026:U:H4'	1:AA:1027:A:OP1	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1212:G:C2	1:AA:1236:G:C4	3.04	0.45
1:AA:1798:U:OP2	3:AD:274:ARG:NH2	2.50	0.45
1:AA:1986:A:H2'	1:AA:1987:G:H8	1.81	0.45
1:AA:2041:U:H2'	1:AA:2042:A:C8	2.52	0.45
1:AA:2483:C:H2'	1:AA:2484:G:O4'	2.17	0.45
1:AA:2884:U:O2	27:A5:53:ALA:HB2	2.17	0.45
2:AB:44:G:N2	2:AB:48:A:C4	2.85	0.45
2:AB:55:U:O4'	6:AG:29:TRP:NE1	2.45	0.45
5:AF:12:LEU:HB2	5:AF:124:LEU:HD11	1.98	0.45
12:AQ:38:GLU:HB2	12:AQ:127:ILE:HG22	1.97	0.45
22:A0:23:VAL:HA	22:A0:38:VAL:HG22	1.98	0.45
1:BA:108:U:C2	1:BA:109:G:C8	3.05	0.45
1:BA:528:A:N1	1:BA:2042:A:H2'	2.31	0.45
1:BA:600:G:H2'	1:BA:601:C:O4'	2.15	0.45
1:BA:1175:U:H1'	1:BA:1176:G:OP1	2.16	0.45
1:BA:1177:A:O2'	1:BA:1178:C:C6	2.70	0.45
1:BA:1263:U:H2'	1:BA:1264:G:C8	2.52	0.45
12:BQ:84:GLY:O	12:BQ:85:LYS:HB2	2.17	0.45
31:CA:499:A:H4'	31:CA:500:G:OP1	2.17	0.45
31:CA:515:G:N1	31:CA:537:G:C6	2.84	0.45
31:CA:782:A:H4'	31:CA:1514:C:O2'	2.17	0.45
31:CA:1217:C:H2'	31:CA:1218:C:C6	2.51	0.45
31:CA:1228:C:P	43:CM:108:ARG:HH12	2.38	0.45
31:CA:1273:G:H3'	31:CA:1274:G:H8	1.82	0.45
34:CD:111:ALA:HB2	34:CD:120:LEU:CD1	2.46	0.45
34:CD:117:ALA:HA	34:CD:120:LEU:HB2	1.99	0.45
37:CG:20:ASP:OD2	37:CG:23:VAL:HG23	2.16	0.45
40:CJ:56:HIS:O	40:CJ:58:ASP:N	2.44	0.45
43:CM:105:THR:OG1	43:CM:106:ASN:N	2.49	0.45
31:DA:60:A:N6	31:DA:110:C:N3	2.64	0.45
31:DA:337:C:H2'	31:DA:338:A:C8	2.52	0.45
31:DA:538:G:H2'	31:DA:539:A:H8	1.81	0.45
31:DA:840:C:H4'	31:DA:841:U:OP1	2.16	0.45
34:DD:149:ALA:O	34:DD:152:SER:N	2.42	0.45
37:DG:111:ARG:HB2	37:DG:119:ARG:HD2	1.99	0.45
42:DL:7:ILE:O	42:DL:11:VAL:HG23	2.16	0.45
1:AA:95:G:H4'	24:A2:46:GLN:N	2.27	0.45
1:AA:300:A:N3	1:AA:319:C:H1'	2.32	0.45
1:AA:597:U:H2'	1:AA:598:G:C8	2.52	0.45
1:AA:652(P):G:N3	1:AA:652(Q):G:N7	2.64	0.45
1:AA:1171:G:H3'	1:AA:1173:G:H5'	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1721:G:H5''	1:AA:1721:G:N3	2.32	0.45
1:AA:2016:U:H2'	1:AA:2017:U:C6	2.52	0.45
1:AA:2056:G:C2	1:AA:2057:A:C8	3.04	0.45
1:AA:2165:G:H2'	1:AA:2166:G:H8	1.81	0.45
1:AA:2272:U:H5''	1:AA:2273:A:OP1	2.16	0.45
1:AA:2337:G:C2	1:AA:2338:G:C8	3.04	0.45
1:AA:2730:C:H2'	1:AA:2731:G:H8	1.82	0.45
3:AD:144:ALA:HB3	3:AD:192:THR:HG23	1.99	0.45
4:AE:116:VAL:HG13	4:AE:122:PHE:HB2	1.99	0.45
5:AF:184:TYR:CD2	5:AF:188:ARG:HD2	2.50	0.45
7:AH:86:GLU:OE2	7:AH:132:ARG:NH2	2.46	0.45
8:AI:63:ALA:O	8:AI:66:GLU:N	2.50	0.45
8:AI:73:GLU:OE2	8:AI:139:GLN:N	2.49	0.45
27:A5:37:LYS:HG2	27:A5:37:LYS:O	2.15	0.45
30:A8:16:ILE:HD13	30:A8:59:LYS:HD2	1.97	0.45
1:BA:24:G:H2'	1:BA:25:U:O4'	2.17	0.45
1:BA:83:G:H1	1:BA:102:G:H2'	1.82	0.45
1:BA:637:A:P	11:BP:116:GLY:HA3	2.57	0.45
1:BA:1558:A:N3	1:BA:1558:A:O4'	2.49	0.45
1:BA:2134:A:N6	1:BA:2157:G:H4'	2.32	0.45
1:BA:2228:G:C6	1:BA:2229:C:C4	3.05	0.45
1:BA:2469:A:C2	1:BA:2482:G:C8	3.05	0.45
2:BB:21:G:H2'	2:BB:22:U:O4'	2.17	0.45
6:BG:43:LEU:HB2	6:BG:89:GLY:HA2	1.98	0.45
21:BZ:106:GLY:O	21:BZ:108:PRO:HD3	2.16	0.45
26:B4:16:CYS:HA	26:B4:33:VAL:O	2.16	0.45
31:CA:68:G:O4'	31:CA:171:A:H1'	2.17	0.45
31:CA:987:G:H2'	31:CA:988:G:H8	1.81	0.45
31:CA:1236:A:O2'	31:CA:1304:G:H4'	2.16	0.45
33:CC:23:TYR:CG	40:CJ:10:GLY:HA2	2.51	0.45
38:CH:51:VAL:HG21	38:CH:60:ARG:HG3	1.98	0.45
38:CH:84:ARG:HD3	38:CH:85:ARG:O	2.16	0.45
42:CL:117:ARG:HB3	42:CL:122:THR:HB	1.99	0.45
50:CT:76:ALA:HA	50:CT:79:ARG:NH1	2.32	0.45
31:DA:91:C:O2'	31:DA:92:C:H5'	2.17	0.45
31:DA:238:G:C2	31:DA:239:U:C2	3.04	0.45
31:DA:999:C:H2'	31:DA:1000:U:O4'	2.16	0.45
32:DB:185:ILE:HG22	32:DB:199:TYR:CD1	2.52	0.45
33:DC:54:ARG:HG2	33:DC:55:VAL:N	2.32	0.45
39:DI:31:GLN:N	39:DI:31:GLN:HE21	2.14	0.45
41:DK:58:PRO:HA	41:DK:90:GLY:HA2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:DS:32:LYS:HD2	49:DS:57:HIS:CD2	2.52	0.45
1:AA:13:A:N1	1:AA:525:U:H2'	2.32	0.45
1:AA:223:A:O2'	1:AA:420:C:O2	2.31	0.45
1:AA:797:C:H2'	1:AA:798:G:C8	2.52	0.45
1:AA:1045:A:H2'	1:AA:1045:A:N3	2.31	0.45
1:AA:1252:G:N1	16:AU:37:GLU:OE1	2.40	0.45
1:AA:1494:A:H2'	1:AA:1495:A:H8	1.81	0.45
2:AB:59:A:H2'	2:AB:60:C:O4'	2.17	0.45
9:AN:123:TYR:CZ	9:AN:129:PRO:HD2	2.52	0.45
18:AW:17:VAL:O	18:AW:20:VAL:N	2.41	0.45
22:A0:43:THR:C	22:A0:45:PHE:H	2.19	0.45
1:BA:36:G:H4'	1:BA:451:C:C2	2.52	0.45
1:BA:375:C:H2'	1:BA:376:C:C6	2.52	0.45
1:BA:1028:A:N6	1:BA:1125:G:H2'	2.31	0.45
1:BA:1665:A:C4'	10:BO:67:LYS:HB2	2.46	0.45
2:BB:38:C:O4'	14:BS:95:HIS:NE2	2.48	0.45
3:BD:71:ASP:HB3	3:BD:103:ARG:NH2	2.31	0.45
31:CA:277:C:H2'	31:CA:278:G:C8	2.52	0.45
31:CA:646:U:H2'	31:CA:647:C:C6	2.51	0.45
31:CA:954:G:H21	31:CA:1227:A:N6	2.08	0.45
31:CA:994:A:O2'	44:CN:8:GLU:HG2	2.16	0.45
31:CA:1206:G:H2'	31:CA:1207:G:O4'	2.17	0.45
34:CD:60:GLU:HG3	34:CD:198:VAL:HG13	1.98	0.45
36:CF:97:PHE:O	48:CR:31:LEU:HD23	2.17	0.45
37:CG:106:GLN:O	37:CG:110:GLN:HG3	2.17	0.45
49:CS:44:MET:O	49:CS:47:HIS:HB2	2.17	0.45
31:DA:19:C:H4'	31:DA:864:A:O4'	2.17	0.45
31:DA:277:C:OP2	47:DQ:41:LYS:NZ	2.42	0.45
31:DA:430:A:OP1	34:DD:9:CYS:N	2.42	0.45
31:DA:606:G:H2'	31:DA:631:G:H1	1.81	0.45
31:DA:642:A:H2'	31:DA:643:C:C6	2.52	0.45
31:DA:673:G:C4	31:DA:734:G:C2	3.05	0.45
31:DA:763:G:H2'	31:DA:764:C:H6	1.81	0.45
31:DA:1030(B):C:C3'	31:DA:1030(C):G:H5'	2.46	0.45
31:DA:1291:G:H5''	39:DI:39:GLY:HA3	1.98	0.45
35:DE:79:GLU:HG3	35:DE:93:PRO:HD2	1.99	0.45
39:DI:113:LYS:H	39:DI:119:ALA:HA	1.81	0.45
42:DL:66:VAL:HG21	42:DL:98:TYR:CD1	2.52	0.45
46:DP:22:THR:HA	46:DP:33:ILE:HG12	1.99	0.45
48:DR:53:ARG:HH21	48:DR:60:ALA:N	2.15	0.45
1:AA:126:A:O5'	29:A7:19:ARG:HG3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:954:G:H5''	12:AQ:13:GLN:HB3	1.99	0.45
1:AA:1176:G:H4'	1:AA:1177:A:OP2	2.17	0.45
1:AA:1275:A:O2'	1:AA:1645:G:N3	2.49	0.45
1:AA:1463:C:H2'	1:AA:1464:C:H6	1.82	0.45
1:AA:1494:A:C6	1:AA:1495:A:C6	3.05	0.45
1:AA:1612:C:O3'	29:A7:5:TRP:HB3	2.17	0.45
1:AA:2122:U:O2'	1:AA:2123:G:OP1	2.28	0.45
1:AA:2139:C:H2'	1:AA:2140:C:O4'	2.17	0.45
1:AA:2438:U:O2'	1:AA:2440:C:OP1	2.30	0.45
2:AB:95:C:H2'	2:AB:96:U:C6	2.52	0.45
4:AE:68:ALA:C	4:AE:70:ALA:H	2.19	0.45
5:AF:101:LEU:HB3	5:AF:106:ARG:HD3	1.98	0.45
15:AT:39:ARG:NH2	31:DA:345:C:H5'	2.31	0.45
21:AZ:151:HIS:HD2	21:AZ:168:GLU:O	2.00	0.45
22:A0:43:THR:C	22:A0:45:PHE:N	2.70	0.45
1:BA:69:C:O2'	1:BA:70:G:H5'	2.17	0.45
1:BA:470:A:H2'	1:BA:471:A:O4'	2.17	0.45
1:BA:588:U:O4	1:BA:670:A:H1'	2.16	0.45
1:BA:1482:G:C6	1:BA:1507:A:C6	3.05	0.45
1:BA:2282:G:OP1	1:BA:2283:C:H1'	2.17	0.45
4:BE:179:GLU:O	4:BE:181:LEU:HD13	2.17	0.45
5:BF:7:TYR:HB2	5:BF:22:ALA:HB3	1.97	0.45
31:CA:103:C:H1'	31:CA:171:A:N1	2.32	0.45
31:CA:377:G:OP2	46:CP:3:LYS:NZ	2.46	0.45
31:CA:963:G:HO2'	40:CJ:54:PHE:HZ	1.63	0.45
31:CA:1492:A:C8	42:CL:47:LYS:HB2	2.52	0.45
31:CA:1494:G:C6	31:CA:1495:U:C4	3.04	0.45
32:CB:167:PRO:HG3	32:CB:188:ALA:HB2	1.98	0.45
43:CM:102:ARG:HH21	43:CM:105:THR:HG23	1.81	0.45
45:CO:8:LYS:O	45:CO:12:ILE:HG13	2.17	0.45
45:CO:40:SER:O	45:CO:44:LYS:HG3	2.16	0.45
46:CP:74:LEU:HB3	46:CP:79:VAL:HG11	1.99	0.45
50:CT:92:LEU:HD23	50:CT:92:LEU:HA	1.72	0.45
31:DA:137:C:O2'	31:DA:138:G:H5'	2.17	0.45
31:DA:475:G:C4	31:DA:476:G:C8	3.05	0.45
31:DA:858:G:H1	31:DA:870:U:P	2.39	0.45
33:DC:149:ALA:HA	33:DC:201:TYR:O	2.16	0.45
40:DJ:12:ASP:O	40:DJ:16:LEU:HB3	2.17	0.45
45:DO:3:ILE:H	45:DO:3:ILE:HD13	1.81	0.45
49:DS:37:ARG:H	49:DS:37:ARG:HG3	1.46	0.45
49:DS:53:ASN:HB3	49:DS:75:ALA:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:271(D):G:H2'	1:AA:271(E):U:C6	2.51	0.45
1:AA:868:U:H2'	1:AA:869:G:O4'	2.18	0.45
1:AA:1015:G:O2'	1:AA:1016:G:H5'	2.16	0.45
1:AA:1316:U:H2'	1:AA:1317:A:C8	2.52	0.45
1:AA:1684:C:C2	1:AA:1705:G:N2	2.85	0.45
1:AA:1913:A:C2'	52:AA:3001:T8B:O2	2.65	0.45
1:AA:2163:C:O5'	1:AA:2163:C:H6	2.00	0.45
1:AA:2399:G:H2'	1:AA:2400:G:O4'	2.17	0.45
1:AA:2489:G:C6	1:AA:2490:G:N1	2.85	0.45
1:AA:2525:G:C2	1:AA:2539:C:C2	3.04	0.45
1:AA:2531:A:H2	1:AA:2658:C:O2	2.00	0.45
4:AE:172:VAL:HG13	4:AE:182:LEU:HD11	1.98	0.45
5:AF:80:ALA:HB3	5:AF:83:PHE:HD1	1.82	0.45
15:AT:55:ASN:N	15:AT:59:THR:HG22	2.31	0.45
20:AY:38:ILE:HD13	20:AY:66:PRO:HA	1.99	0.45
21:AZ:63:ASP:OD1	21:AZ:65:GLN:HB2	2.17	0.45
1:BA:72:U:OP1	19:BX:1:MET:N	2.49	0.45
1:BA:548:A:N6	17:BV:19:LYS:H	2.14	0.45
1:BA:848:G:N9	1:BA:933:A:H8	2.14	0.45
1:BA:2808:U:H5''	1:BA:2891:G:O6	2.17	0.45
5:BF:157:VAL:HB	5:BF:194:MET:HG2	1.99	0.45
5:BF:160:ASN:CG	5:BF:163:VAL:HG23	2.38	0.45
10:BO:17:ARG:NH1	10:BO:47:ILE:HD13	2.31	0.45
12:BQ:16:ARG:HG2	12:BQ:18:LYS:HE2	1.98	0.45
13:BR:96:ARG:HD2	13:BR:115:GLU:OE1	2.17	0.45
15:BT:106:SER:O	15:BT:110:ILE:HG13	2.17	0.45
18:BW:14:PRO:CG	18:BW:78:GLU:HG2	2.39	0.45
21:BZ:6:LYS:HB3	21:BZ:6:LYS:HE2	1.81	0.45
21:BZ:100:VAL:O	21:BZ:124:ILE:N	2.50	0.45
31:CA:189:G:C5	31:CA:189(A):C:C5	3.05	0.45
31:CA:189:G:C6	31:CA:189(L):G:N1	2.85	0.45
31:CA:270:A:H2'	31:CA:271:C:C6	2.52	0.45
31:CA:295:C:H2'	31:CA:296:U:O4'	2.17	0.45
31:CA:436:C:O2'	31:CA:437:U:OP2	2.26	0.45
31:CA:948:C:H2'	31:CA:949:A:H8	1.82	0.45
31:CA:950:U:H3	31:CA:1231:G:H1	1.65	0.45
31:CA:1227:A:OP2	43:CM:96:LEU:HD21	2.16	0.45
31:CA:1288:A:H1'	31:CA:1353:G:O4'	2.17	0.45
32:CB:54:THR:HG23	32:CB:199:TYR:HB3	1.99	0.45
37:CG:22:LEU:HG	37:CG:62:PHE:HE2	1.81	0.45
44:CN:3:ARG:HD2	44:CN:3:ARG:HA	1.59	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:22:G:H2'	31:DA:23:C:C6	2.52	0.45
31:DA:38:G:H22	31:DA:397:A:H5''	1.82	0.45
31:DA:542:G:OP1	34:DD:10:ARG:NH2	2.48	0.45
31:DA:949:A:H1'	31:DA:1364:U:O2	2.17	0.45
31:DA:971:G:H1'	31:DA:1365:G:O2'	2.17	0.45
31:DA:1079:G:H2'	31:DA:1080:A:C8	2.52	0.45
32:DB:178:ARG:HH21	38:DH:74:PRO:CB	2.30	0.45
43:DM:16:ASP:HB3	43:DM:34:LEU:HD11	1.99	0.45
43:DM:114:ARG:H	43:DM:114:ARG:HG2	1.44	0.45
1:AA:587:C:C5	1:AA:671:C:H1'	2.52	0.44
1:AA:664:C:H4'	1:AA:941:A:OP1	2.16	0.44
1:AA:672:C:H2'	1:AA:673:C:H6	1.82	0.44
1:AA:700:G:H2'	1:AA:701:G:O4'	2.17	0.44
1:AA:1336:A:H2'	1:AA:1337:G:H8	1.77	0.44
1:AA:1344:G:O2'	1:AA:1385:G:H2'	2.17	0.44
1:AA:2206:G:HO2'	1:AA:2207:G:P	2.39	0.44
1:AA:2787:C:H2'	1:AA:2788:C:H6	1.82	0.44
3:AD:53:PHE:C	3:AD:218:ARG:HB2	2.37	0.44
7:AH:11:VAL:HG13	7:AH:15:VAL:HG22	1.99	0.44
10:AO:104:ARG:HB2	10:AO:104:ARG:CZ	2.47	0.44
12:AQ:74:TYR:O	12:AQ:90:VAL:HA	2.18	0.44
22:A0:27:GLU:HB2	22:A0:69:PHE:HD1	1.82	0.44
1:BA:100:G:H4'	1:BA:100:G:OP2	2.17	0.44
1:BA:515:A:H1'	1:BA:581:C:H1'	2.00	0.44
1:BA:1837:C:OP1	31:CA:784:C:H4'	2.17	0.44
1:BA:2007:C:H4'	1:BA:2824:C:O2'	2.17	0.44
1:BA:2093:G:O5'	8:BI:24:GLY:HA3	2.16	0.44
1:BA:2113:U:H2'	1:BA:2114:A:C8	2.52	0.44
1:BA:2139:C:H2'	1:BA:2140:C:O4'	2.18	0.44
1:BA:2778:A:H4'	1:BA:2779:U:OP2	2.17	0.44
1:BA:2853:C:H2'	1:BA:2854:G:H8	1.82	0.44
1:BA:2875:C:O2'	15:BT:4:GLY:HA3	2.17	0.44
8:BI:72:LEU:HA	8:BI:75:LEU:HD11	1.98	0.44
29:B7:16:HIS:HB2	29:B7:44:PRO:HG2	1.98	0.44
31:CA:1146:A:C5	31:CA:1147:C:C6	3.06	0.44
31:CA:1151:A:O2'	31:CA:1152:A:H8	1.99	0.44
32:CB:12:GLU:C	32:CB:16:HIS:HD1	2.21	0.44
33:CC:11:ARG:HB3	33:CC:15:THR:HB	1.98	0.44
36:CF:4:TYR:CD1	36:CF:92:LYS:HA	2.52	0.44
36:CF:19:LEU:HD11	36:CF:59:TYR:CZ	2.52	0.44
41:CK:32:ILE:HG21	41:CK:72:ALA:CB	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:270:A:H2'	31:DA:271:C:C6	2.51	0.44
31:DA:277:C:H5''	47:DQ:68:ARG:NH2	2.32	0.44
31:DA:328:C:H4'	31:DA:329:A:H5'	1.98	0.44
31:DA:623:C:C4	31:DA:624:C:C4	3.05	0.44
31:DA:728:A:H2'	31:DA:729:A:C8	2.51	0.44
31:DA:825:G:C6	31:DA:826:C:C4	3.05	0.44
31:DA:841:U:H6	31:DA:841:U:P	2.40	0.44
31:DA:991:U:O2'	31:DA:992:U:P	2.75	0.44
31:DA:1332:A:O5'	31:DA:1332:A:C8	2.70	0.44
31:DA:1386:G:C2	31:DA:1387:G:C8	3.05	0.44
33:DC:182:ILE:HA	33:DC:202:ILE:O	2.17	0.44
39:DI:21:PRO:HA	39:DI:59:PHE:HA	1.99	0.44
45:DO:87:ILE:O	45:DO:88:ARG:HB3	2.17	0.44
50:DT:67:ALA:HA	50:DT:72:LEU:O	2.17	0.44
1:AA:71:A:H3'	1:AA:71:A:OP2	2.16	0.44
1:AA:300:A:P	20:AY:86:ARG:NH2	2.89	0.44
1:AA:459:U:OP2	29:A7:39:ARG:NH1	2.47	0.44
1:AA:825:C:H2'	1:AA:826:U:O4'	2.17	0.44
1:AA:870:A:C2	1:AA:908:C:C2	3.05	0.44
1:AA:1915:U:O2	52:AA:3001:T8B:H20	2.17	0.44
1:AA:2638:G:OP1	4:AE:82:ARG:NH2	2.49	0.44
3:AD:253:GLN:HE21	3:AD:253:GLN:HB3	1.55	0.44
8:AI:88:ILE:HG22	8:AI:90:GLY:H	1.83	0.44
12:AQ:37:LEU:HB2	12:AQ:128:LYS:O	2.16	0.44
12:AQ:41:TRP:HB3	12:AQ:94:VAL:HG11	1.99	0.44
14:AS:23:ARG:NH2	14:AS:84:GLN:HB3	2.32	0.44
21:AZ:137:ILE:HA	21:AZ:156:LYS:HZ2	1.83	0.44
23:A1:25:LYS:O	23:A1:28:GLY:N	2.49	0.44
1:BA:71:A:OP2	1:BA:71:A:H3'	2.17	0.44
1:BA:93:G:H2'	1:BA:94:C:C6	2.53	0.44
1:BA:128:C:H2'	1:BA:129:C:C6	2.53	0.44
1:BA:181:A:H1'	1:BA:435:C:H5'	1.98	0.44
1:BA:278:A:H4'	1:BA:279:C:OP1	2.17	0.44
1:BA:663:G:H2'	1:BA:664:C:O4'	2.17	0.44
1:BA:1178:C:H2'	1:BA:1179:C:H6	1.81	0.44
1:BA:1952:A:C6	1:BA:1953:A:N1	2.85	0.44
1:BA:2778:A:O2'	1:BA:2781:A:H5'	2.17	0.44
1:BA:2839:G:H5'	13:BR:46:GLY:CA	2.47	0.44
3:BD:3:VAL:HG13	3:BD:17:THR:HB	1.98	0.44
6:BG:96:ARG:O	6:BG:99:MET:HB3	2.18	0.44
19:BX:3:THR:O	19:BX:6:ASP:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:129(A):G:H1'	31:CA:189(F):U:H2'	1.99	0.44
31:CA:157:G:C2	31:CA:158:G:C8	3.05	0.44
31:CA:589:C:H2'	31:CA:590:C:H6	1.81	0.44
31:CA:814:A:N7	31:CA:816:A:C4	2.86	0.44
31:CA:1020:U:H2'	31:CA:1021:G:C8	2.52	0.44
34:CD:17:VAL:HG11	34:CD:197:PRO:HG3	2.00	0.44
43:CM:14:ARG:HA	43:CM:44:ARG:HA	2.00	0.44
44:CN:23:ARG:NH1	44:CN:30:ALA:HB2	2.32	0.44
45:CO:78:TYR:CE2	45:CO:82:ILE:HD13	2.51	0.44
31:DA:160:A:H3'	31:DA:161:A:H8	1.83	0.44
31:DA:517:G:H4'	31:DA:519:C:C6	2.52	0.44
31:DA:532:A:H61	33:DC:193:TYR:HB2	1.82	0.44
31:DA:1033:G:H3'	31:DA:1034:G:H8	1.83	0.44
31:DA:1267:C:H5	31:DA:1268:A:C4	2.35	0.44
32:DB:95:GLN:HG3	32:DB:147:LYS:HG2	1.98	0.44
35:DE:92:LYS:HA	35:DE:93:PRO:HD2	1.88	0.44
37:DG:43:PHE:HD2	37:DG:44:TYR:CE2	2.35	0.44
37:DG:69:VAL:HG11	37:DG:134:ALA:HB1	1.98	0.44
1:AA:243:U:O2'	1:AA:244:A:H5'	2.17	0.44
1:AA:775:G:N2	1:AA:793:A:O3'	2.50	0.44
1:AA:883:G:H1	1:AA:893:C:H42	1.65	0.44
1:AA:1002:G:C6	1:AA:1003:G:C5	3.05	0.44
1:AA:1416:G:O2'	1:AA:1417:C:H5	2.00	0.44
1:AA:1495:A:H2'	1:AA:1496:A:H8	1.79	0.44
1:AA:1526:G:C6	1:AA:1527:G:C2	3.06	0.44
1:AA:2315:G:H2'	1:AA:2316:C:C6	2.52	0.44
1:AA:2351:G:HO2'	1:AA:2352:A:H8	1.65	0.44
2:AB:61:G:H2'	2:AB:62:C:C6	2.52	0.44
21:AZ:101:PRO:O	21:AZ:102:LEU:HD12	2.17	0.44
27:A5:16:ARG:HG2	27:A5:16:ARG:HH11	1.81	0.44
29:A7:22:MET:HA	29:A7:28:ARG:HG2	1.99	0.44
1:BA:1430:C:H2'	1:BA:1431:U:H6	1.82	0.44
1:BA:1803:A:N6	1:BA:1814:G:H1'	2.32	0.44
1:BA:1887:C:H2'	1:BA:1888:G:H5'	1.99	0.44
4:BE:71:GLY:HA2	4:BE:72:VAL:C	2.38	0.44
6:BG:149:VAL:HG22	6:BG:150:ASP:O	2.18	0.44
10:BO:88:ASN:HB3	10:BO:94:ARG:HD3	1.98	0.44
12:BQ:54:MET:HG2	12:BQ:117:ALA:HB1	1.99	0.44
13:BR:26:LYS:HE2	13:BR:70:LEU:O	2.18	0.44
22:B0:25:ARG:HD2	22:B0:29:GLN:NE2	2.32	0.44
31:CA:142:G:H2'	31:CA:143:A:C8	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:198:G:C5	31:CA:220:G:C2	3.05	0.44
31:CA:414:A:H2'	31:CA:415:A:C8	2.52	0.44
31:CA:1265:G:H2'	31:CA:1266:G:O4'	2.17	0.44
33:CC:113:ALA:HB3	33:CC:114:PRO:HD3	2.00	0.44
31:DA:604:G:C5	31:DA:605:U:C5	3.05	0.44
31:DA:1072:G:C6	31:DA:1073:U:C4	3.06	0.44
31:DA:1187:G:H5'	39:DI:113:LYS:HE2	1.98	0.44
31:DA:1327:C:H2'	31:DA:1328:C:C6	2.52	0.44
32:DB:185:ILE:HA	32:DB:199:TYR:O	2.16	0.44
34:DD:158:ILE:O	34:DD:162:LEU:HB2	2.17	0.44
37:DG:62:PHE:O	37:DG:66:VAL:HG23	2.17	0.44
38:DH:97:VAL:HA	38:DH:100:ILE:HD11	1.98	0.44
41:DK:59:TYR:O	41:DK:63:LEU:HB2	2.17	0.44
46:DP:43:LYS:HA	46:DP:48:TRP:HB3	1.98	0.44
1:AA:773:U:O2'	3:AD:48:ARG:HD3	2.17	0.44
1:AA:813:U:H2'	1:AA:814:C:C6	2.52	0.44
1:AA:849:A:C8	1:AA:850:C:C5	3.06	0.44
1:AA:947:G:N2	1:AA:971:C:C2	2.85	0.44
1:AA:1386:C:H2'	1:AA:1387:C:C6	2.52	0.44
1:AA:2350:C:H2'	1:AA:2351:G:O4'	2.17	0.44
1:AA:2394:C:P	30:A8:30:ARG:NH1	2.91	0.44
1:AA:2505:G:H2'	1:AA:2576:G:O6	2.17	0.44
3:AD:182:LEU:HD23	3:AD:182:LEU:HA	1.71	0.44
4:AE:110:GLY:O	13:AR:3:HIS:HE1	2.00	0.44
5:AF:123:LEU:HD12	5:AF:124:LEU:H	1.83	0.44
8:AI:140:LEU:HD22	8:AI:141:LYS:N	2.33	0.44
15:AT:26:ASP:OD1	15:AT:120:ARG:NH2	2.45	0.44
18:AW:19:LEU:O	27:A5:25:LEU:HD12	2.18	0.44
1:BA:7:G:H2'	1:BA:8:A:C8	2.52	0.44
1:BA:125:G:C6	29:B7:10:ARG:HG3	2.52	0.44
1:BA:1187:G:H8	1:BA:1187:G:O5'	2.00	0.44
1:BA:2119:A:C2	1:BA:2170:A:H2'	2.53	0.44
6:BG:15:VAL:HG13	6:BG:175:LEU:HB3	1.99	0.44
8:BI:65:ALA:CB	8:BI:136:VAL:HG11	2.47	0.44
8:BI:145:VAL:HG12	8:BI:146:ALA:N	2.32	0.44
13:BR:57:ARG:HB3	13:BR:59:ASP:OD2	2.17	0.44
20:BY:47:LYS:HE3	20:BY:48:ALA:O	2.17	0.44
31:CA:189:G:C6	31:CA:189(A):C:C4	3.05	0.44
31:CA:524:G:H2'	31:CA:525:C:C6	2.53	0.44
31:CA:975:A:H61	31:CA:1367:C:C1'	2.30	0.44
32:CB:189:ASP:HB3	32:CB:203:GLY:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:CB:212:GLN:NE2	32:CB:235:SER:HB3	2.32	0.44
33:CC:22:TRP:CH2	44:CN:54:PRO:HG2	2.52	0.44
34:CD:189:PRO:CB	34:CD:194:LEU:HD11	2.34	0.44
35:CE:43:LEU:HD21	35:CE:132:ALA:HB1	1.98	0.44
45:CO:76:GLU:OE1	45:CO:76:GLU:HA	2.18	0.44
46:CP:6:LEU:HG	46:CP:17:TYR:CB	2.47	0.44
31:DA:277:C:OP1	47:DQ:41:LYS:HE3	2.18	0.44
31:DA:444:C:H2'	31:DA:445:G:H8	1.83	0.44
31:DA:729:A:H2'	31:DA:730:G:O4'	2.17	0.44
31:DA:765:G:N2	31:DA:813:U:H5	2.16	0.44
31:DA:1055:A:C6	31:DA:1206:G:C5	3.05	0.44
31:DA:1106:G:H2'	31:DA:1107:C:C6	2.53	0.44
31:DA:1123:A:O2'	40:DJ:37:PRO:O	2.32	0.44
31:DA:1126:U:H1'	31:DA:1280:A:C5	2.52	0.44
31:DA:1376:U:H2'	31:DA:1377:A:C8	2.53	0.44
32:DB:163:PHE:CD2	32:DB:185:ILE:HG13	2.52	0.44
32:DB:168:THR:O	32:DB:169:LYS:HB2	2.18	0.44
34:DD:201:GLN:O	34:DD:204:ILE:HB	2.18	0.44
35:DE:51:VAL:O	35:DE:55:VAL:HG23	2.18	0.44
1:AA:125:G:C6	29:A7:10:ARG:HG3	2.53	0.44
1:AA:251:A:C5	1:AA:252:G:H1'	2.53	0.44
1:AA:644:A:H4'	1:AA:645:C:H5	1.80	0.44
1:AA:729:G:C6	3:AD:208:LYS:HB2	2.52	0.44
1:AA:863:A:C2	1:AA:864:G:C4	3.06	0.44
1:AA:885:C:N4	1:AA:890:A:C6	2.69	0.44
1:AA:1139:G:O2'	1:AA:1143:A:N1	2.38	0.44
1:AA:1488:G:H5''	1:AA:1488:G:H8	1.82	0.44
1:AA:1539:G:H2'	1:AA:1540:U:O4'	2.16	0.44
1:AA:2199:A:H2'	1:AA:2199:A:N3	2.32	0.44
1:AA:2690:C:N4	1:AA:2713:A:H1'	2.33	0.44
1:AA:2728:U:H2'	1:AA:2729:G:H8	1.82	0.44
2:AB:45:A:C2'	2:AB:46:A:H5'	2.47	0.44
8:AI:105:HIS:CD2	8:AI:105:HIS:N	2.82	0.44
12:AQ:7:MET:HB3	12:AQ:9:TYR:O	2.18	0.44
24:A2:32:LEU:HD12	24:A2:57:ILE:HD12	2.00	0.44
1:BA:438:G:H2'	1:BA:440:G:C8	2.53	0.44
1:BA:783:A:H2'	1:BA:783:A:N3	2.33	0.44
1:BA:1045:A:N3	1:BA:1045:A:C2'	2.80	0.44
1:BA:1649:G:C6	1:BA:2009:G:C6	3.06	0.44
1:BA:2184:G:H8	1:BA:2184:G:OP2	2.01	0.44
1:BA:2350:C:H2'	1:BA:2351:G:O4'	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:BI:84:GLY:O	8:BI:85:GLU:HB3	2.16	0.44
9:BN:19:GLU:HA	9:BN:59:LYS:HB2	1.98	0.44
18:BW:12:ILE:HD13	18:BW:17:VAL:HG22	1.98	0.44
21:BZ:26:GLY:HA2	21:BZ:85:HIS:CD2	2.53	0.44
31:CA:153:C:H42	31:CA:169:C:N4	2.15	0.44
31:CA:355:C:N3	31:CA:356:A:C5	2.85	0.44
31:CA:418:C:H1'	31:CA:540:G:O2'	2.17	0.44
31:CA:748:C:H1'	31:CA:749:C:OP2	2.18	0.44
31:CA:947:G:H2'	31:CA:948:C:O4'	2.17	0.44
31:CA:949:A:C1'	31:CA:1364:U:H3	2.30	0.44
31:CA:1256:A:H5'	31:CA:1258:G:H1'	1.99	0.44
31:CA:1261:A:H3'	31:CA:1262:C:C6	2.51	0.44
34:CD:94:LEU:HA	34:CD:94:LEU:HD23	1.69	0.44
35:CE:70:PRO:O	35:CE:72:GLN:N	2.50	0.44
36:CF:37:VAL:HA	36:CF:65:VAL:HG12	1.99	0.44
39:CI:42:ARG:O	39:CI:45:ALA:HA	2.18	0.44
42:CL:70:ILE:HD13	42:CL:77:LEU:HD12	1.99	0.44
31:DA:872:A:C4	31:DA:874:G:N7	2.86	0.44
31:DA:1176:A:H2'	31:DA:1177:G:H8	1.83	0.44
34:DD:13:ARG:HB2	34:DD:14:ARG:H	1.36	0.44
34:DD:111:ALA:HB2	34:DD:120:LEU:HD12	2.00	0.44
40:DJ:44:VAL:HG22	40:DJ:66:ARG:HD3	1.99	0.44
50:DT:35:THR:O	50:DT:39:LYS:N	2.41	0.44
1:AA:208:C:H2'	1:AA:209:C:H6	1.82	0.44
1:AA:271(R):G:H5''	23:A1:97:LEU:HD21	2.00	0.44
1:AA:600:G:N2	1:AA:605:C:O3'	2.51	0.44
1:AA:1027:A:H2	1:AA:2487:G:HO2'	1.65	0.44
1:AA:1665:A:H4'	10:AO:67:LYS:HB2	2.00	0.44
1:AA:2173:A:H2'	1:AA:2174:C:H5'	1.99	0.44
1:AA:2203:U:O4'	3:AD:151:LYS:HE2	2.18	0.44
1:AA:2218:U:H6	1:AA:2218:U:H2'	1.27	0.44
1:AA:2303:G:O2'	6:AG:132:ASN:HB2	2.17	0.44
1:AA:2514:U:H2'	1:AA:2515:C:C6	2.53	0.44
6:AG:126:ASP:HB2	6:AG:130:ASN:O	2.18	0.44
9:AN:3:THR:HG22	9:AN:4:TYR:H	1.82	0.44
12:AQ:30:GLY:CA	12:AQ:107:ALA:HB2	2.48	0.44
1:BA:103:A:H8	1:BA:103:A:O5'	2.00	0.44
1:BA:319:C:H2'	1:BA:320:A:O4'	2.18	0.44
1:BA:1143:A:OP1	9:BN:25:ARG:NH2	2.51	0.44
1:BA:1210:A:H5''	1:BA:1212:G:O4'	2.18	0.44
1:BA:1697:G:OP2	1:BA:1698:A:O2'	2.31	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:1854:A:H2'	1:BA:1855:G:O4'	2.17	0.44
1:BA:2461:C:H42	1:BA:2489:G:H1	1.66	0.44
20:BY:68:HIS:ND1	20:BY:70:SER:HB3	2.32	0.44
31:CA:352:C:O2'	31:CA:354:G:OP1	2.23	0.44
31:CA:670:G:C6	31:CA:671:G:C5	3.06	0.44
31:CA:1160:G:OP1	32:CB:133:LYS:HD3	2.17	0.44
31:CA:1237:C:H4'	31:CA:1334:G:H21	1.82	0.44
32:CB:18:GLY:HA2	32:CB:42:ILE:H	1.82	0.44
34:CD:52:SER:O	34:CD:55:ALA:N	2.51	0.44
41:CK:32:ILE:HG21	41:CK:72:ALA:HB2	1.99	0.44
42:CL:24:VAL:HG13	42:CL:98:TYR:HE2	1.82	0.44
42:CL:85:ILE:HD13	42:CL:85:ILE:HA	1.74	0.44
47:CQ:7:THR:OG1	47:CQ:58:GLU:HG2	2.17	0.44
31:DA:343:U:C2	31:DA:347:G:N1	2.85	0.44
31:DA:838:G:H2'	31:DA:839:U:H5''	2.00	0.44
31:DA:1244:C:OP2	51:DU:9:ARG:HB2	2.17	0.44
31:DA:1250:A:H2'	31:DA:1251:A:O4'	2.17	0.44
32:DB:185:ILE:HG12	32:DB:185:ILE:O	2.18	0.44
38:DH:37:ARG:HH21	38:DH:38:ILE:HG12	1.82	0.44
40:DJ:13:HIS:O	40:DJ:17:ASP:N	2.43	0.44
43:DM:30:ALA:O	43:DM:34:LEU:HG	2.16	0.44
1:AA:96:G:H4'	24:A2:48:HIS:NE2	2.32	0.44
1:AA:141:A:C8	1:AA:1408:C:O2'	2.62	0.44
1:AA:146:G:H2'	1:AA:147:U:O4'	2.17	0.44
1:AA:271(K):U:HO2'	1:AA:271(M):G:H22	1.64	0.44
1:AA:902:C:H2'	1:AA:903:C:H6	1.83	0.44
1:AA:993:G:OP1	16:AU:50:ARG:NH2	2.50	0.44
1:AA:1000:A:C6	1:AA:1001:A:N1	2.86	0.44
1:AA:1636:C:H2'	1:AA:1637:A:H8	1.83	0.44
1:AA:2031:A:H1'	1:AA:2455:G:O2'	2.18	0.44
2:AB:29:A:O5'	14:AS:32:LEU:HD12	2.17	0.44
8:AI:111:PRO:HB2	8:AI:112:LYS:HE3	1.99	0.44
12:AQ:29:PHE:HB2	12:AQ:105:GLU:OE2	2.18	0.44
12:AQ:72:LYS:HB3	12:AQ:94:VAL:HG23	1.99	0.44
16:AU:91:ASP:N	17:AV:11:GLN:OE1	2.35	0.44
20:AY:5:MET:HE2	20:AY:5:MET:HB2	1.90	0.44
20:AY:91:GLU:H	20:AY:93:GLY:HA2	1.82	0.44
23:A1:62:VAL:HG13	23:A1:63:ALA:O	2.18	0.44
1:BA:1434:A:H61	1:BA:1558:A:H62	1.65	0.44
1:BA:2568:C:H2'	1:BA:2569:G:O4'	2.18	0.44
1:BA:2789:C:H4'	1:BA:2790:A:OP1	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BE:170:LEU:HD12	4:BE:170:LEU:HA	1.76	0.44
9:BN:42:TRP:CH2	9:BN:44:PRO:HB3	2.52	0.44
12:BQ:32:TYR:HB2	12:BQ:106:VAL:HG23	1.99	0.44
26:B4:15:ILE:HG13	26:B4:21:VAL:HG22	1.99	0.44
26:B4:42:PHE:CA	26:B4:43:TYR:HB2	2.47	0.44
31:CA:317:G:C6	31:CA:318:G:C5	3.06	0.44
31:CA:919:A:O2'	31:CA:920:U:H5'	2.18	0.44
31:CA:1097:C:H1'	31:CA:1170:A:H1'	1.99	0.44
31:CA:1380:U:C4	37:CG:3:ARG:HG2	2.52	0.44
32:CB:178:ARG:HG2	38:CH:71:GLY:O	2.18	0.44
38:CH:9:MET:HG2	38:CH:10:LEU:HD23	1.99	0.44
31:DA:21:G:C2	31:DA:22:G:C6	3.05	0.44
31:DA:509:A:N3	31:DA:543:C:O2'	2.50	0.44
31:DA:1028:C:H3'	31:DA:1028:C:H6	1.82	0.44
31:DA:1046:A:H2'	31:DA:1047:G:O4'	2.18	0.44
31:DA:1255:G:H1	31:DA:1282:C:H42	1.63	0.44
31:DA:1287:A:H2'	31:DA:1288:A:C8	2.53	0.44
32:DB:19:HIS:CE1	32:DB:206:ASP:HB2	2.53	0.44
34:DD:26:CYS:HB3	34:DD:31:CYS:SG	2.54	0.44
34:DD:64:LEU:HD22	34:DD:198:VAL:HG11	1.99	0.44
37:DG:68:ASN:ND2	37:DG:128:ALA:HA	2.32	0.44
38:DH:20:TYR:HA	38:DH:65:TYR:CZ	2.52	0.44
47:DQ:10:VAL:HG13	47:DQ:19:VAL:HB	1.98	0.44
1:AA:57:C:H2'	1:AA:58:G:O4'	2.17	0.44
1:AA:271(P):C:H2'	1:AA:271(Q):G:H5'	1.99	0.44
1:AA:445:C:O2'	1:AA:446:G:H5'	2.18	0.44
1:AA:709:U:C2	1:AA:723:G:N2	2.85	0.44
1:AA:997:G:O2'	1:AA:998:C:H5'	2.17	0.44
1:AA:1323:U:OP1	18:AW:98:LYS:NZ	2.41	0.44
1:AA:1578:U:C2'	1:AA:1579:A:H5'	2.48	0.44
1:AA:1613:G:C2	1:AA:1619:G:C5	3.06	0.44
1:AA:1939:U:OP1	1:AA:2604:U:O2'	2.36	0.44
1:AA:1992:G:H5'	1:AA:1994:C:H41	1.82	0.44
1:AA:2203:U:O2'	1:AA:2205:C:H5'	2.17	0.44
1:AA:2378:A:C5	1:AA:2379:G:H1'	2.52	0.44
1:AA:2464:C:O2'	1:AA:2465:C:OP2	2.32	0.44
1:AA:2694:G:C6	1:AA:2695:C:C4	3.05	0.44
1:AA:2787:C:H1'	4:AE:62:PRO:HG3	2.00	0.44
2:AB:19:G:H5'	2:AB:20:C:OP2	2.18	0.44
4:AE:151:TYR:O	4:AE:154:LYS:HB2	2.17	0.44
8:AI:82:ARG:HB3	8:AI:89:TYR:CG	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:AR:92:GLY:HA2	13:AR:94:TYR:CZ	2.53	0.44
17:AV:77:ALA:O	17:AV:79:VAL:HB	2.17	0.44
20:AY:43:ASN:OD1	20:AY:65:ALA:HB3	2.18	0.44
20:AY:76:CYS:SG	20:AY:99:CYS:HB2	2.57	0.44
24:A2:22:GLU:HG2	24:A2:64:LEU:HD11	2.00	0.44
26:A4:16:CYS:HA	26:A4:33:VAL:O	2.18	0.44
1:BA:428:A:H8	1:BA:428:A:OP2	2.01	0.44
1:BA:729:G:H5'	1:BA:730:C:H5''	2.00	0.44
1:BA:2331:G:O2'	1:BA:2336:A:N1	2.40	0.44
5:BF:29:ASN:HB3	5:BF:112:MET:HE1	2.00	0.44
12:BQ:27:VAL:O	12:BQ:67:ARG:NH1	2.51	0.44
21:BZ:138:GLU:H	21:BZ:156:LYS:HZ1	1.66	0.44
31:CA:178:C:H2'	31:CA:179:A:H8	1.81	0.44
31:CA:708:C:P	41:CK:85:ARG:HH12	2.41	0.44
31:CA:826:C:H2'	31:CA:827:U:H6	1.83	0.44
31:CA:1030(D):A:H2'	31:CA:1031:G:O4'	2.18	0.44
31:CA:1095:U:P	31:CA:1108:G:H1	2.40	0.44
31:CA:1217:C:H2'	31:CA:1218:C:O4'	2.18	0.44
31:CA:1347:G:H22	31:CA:1374:A:P	2.41	0.44
31:CA:1347:G:N2	31:CA:1373:G:H2'	2.32	0.44
31:CA:1399:C:C2	31:CA:1401:G:C5	3.06	0.44
31:CA:1456:G:O2'	50:CT:39:LYS:HG2	2.17	0.44
37:CG:64:GLN:O	37:CG:68:ASN:HB2	2.18	0.44
38:CH:10:LEU:O	38:CH:13:ILE:HB	2.18	0.44
42:CL:33:ARG:O	42:CL:85:ILE:HB	2.17	0.44
42:CL:44:THR:HG22	42:CL:51:ALA:O	2.18	0.44
43:CM:86:CYS:H	43:CM:89:GLY:HA3	1.82	0.44
50:CT:80:ARG:O	50:CT:84:LEU:HB2	2.18	0.44
31:DA:375:U:H5''	46:DP:6:LEU:HD23	1.99	0.44
31:DA:410:G:C2	31:DA:429:U:C2	3.06	0.44
31:DA:600:C:OP1	38:DH:97:VAL:N	2.40	0.44
31:DA:1490:C:H2'	31:DA:1491:G:H8	1.82	0.44
33:DC:67:THR:HA	33:DC:102:ASN:HB2	1.99	0.44
36:DF:30:LEU:HB3	36:DF:35:ALA:HB3	2.00	0.44
40:DJ:96:ILE:H	40:DJ:96:ILE:HG13	1.56	0.44
1:AA:53:A:H61	1:AA:117:G:C2'	2.30	0.44
1:AA:274:G:H3'	1:AA:275:G:C8	2.53	0.44
1:AA:435:C:H2'	1:AA:436:C:H5'	2.00	0.44
1:AA:806:C:OP2	11:AP:37:GLY:HA2	2.18	0.44
1:AA:857:C:C2	1:AA:921:G:N2	2.86	0.44
1:AA:873:G:N2	1:AA:905:U:C2	2.86	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1709:U:H2'	1:AA:1710:C:C6	2.53	0.44
1:AA:2115:G:O2'	1:AA:2167:U:H1'	2.18	0.44
1:AA:2243:U:H2'	1:AA:2244:U:C6	2.53	0.44
2:AB:117:G:H2'	2:AB:118:G:O4'	2.18	0.44
11:AP:135:LEU:HD23	11:AP:135:LEU:HA	1.91	0.44
19:AX:29:TRP:CZ3	19:AX:78:LYS:HB3	2.53	0.44
21:AZ:153:SER:HB3	21:AZ:167:PRO:O	2.17	0.44
23:A1:50:ARG:HG2	23:A1:59:THR:HB	2.00	0.44
30:A8:34:TRP:NE1	30:A8:35:GLN:HB3	2.32	0.44
1:BA:990:A:OP2	1:BA:991:C:OP2	2.36	0.44
1:BA:1638:C:O2	1:BA:2698:U:O2'	2.32	0.44
1:BA:1963:U:H4'	1:BA:1964:G:OP1	2.17	0.44
31:CA:59:A:H3'	31:CA:331:G:H22	1.83	0.44
31:CA:489:C:C4	31:CA:490:G:N7	2.86	0.44
31:CA:818:G:HO2'	31:CA:820:U:H6	1.59	0.44
31:CA:828:A:N6	31:CA:858:G:O2'	2.49	0.44
40:CJ:43:ARG:HD2	40:CJ:43:ARG:HA	1.72	0.44
42:CL:79:GLU:HB3	42:CL:80:HIS:ND1	2.32	0.44
31:DA:292:G:N2	31:DA:309:G:C4	2.86	0.44
31:DA:324:G:H8	31:DA:324:G:O5'	2.01	0.44
31:DA:643:C:H5'	38:DH:31:PHE:CD1	2.53	0.44
31:DA:1074:G:O2'	31:DA:1101:A:N1	2.31	0.44
31:DA:1089:G:C6	31:DA:1090:U:C4	3.05	0.44
33:DC:77:ILE:HA	33:DC:84:ILE:N	2.31	0.44
39:DI:28:VAL:HG22	39:DI:63:ILE:HD12	1.99	0.44
43:DM:38:GLY:O	43:DM:55:ARG:NH1	2.51	0.44
1:AA:95:G:O5'	24:A2:45:SER:HB2	2.17	0.43
1:AA:1916:A:H2'	1:AA:1917:U:O4'	2.17	0.43
2:AB:64:C:H2'	2:AB:65:C:C6	2.53	0.43
2:AB:117:G:O5'	2:AB:117:G:H8	2.00	0.43
6:AG:36:LYS:HB3	6:AG:95:ARG:HG2	2.00	0.43
7:AH:3:ARG:HG2	7:AH:6:ARG:NE	2.32	0.43
10:AO:73:ASP:OD2	10:AO:75:SER:HB3	2.18	0.43
14:AS:32:LEU:O	14:AS:62:LYS:HE2	2.18	0.43
15:AT:6:LEU:HA	15:AT:6:LEU:HD13	1.64	0.43
15:AT:91:ARG:HD2	15:AT:120:ARG:NH1	2.33	0.43
20:AY:90:LEU:O	20:AY:91:GLU:HB2	2.18	0.43
1:BA:568:U:H5'	1:BA:945:A:C6	2.54	0.43
1:BA:1545:A:H2'	1:BA:1546:C:O4'	2.18	0.43
1:BA:2076:U:OP2	1:BA:2238:G:N2	2.44	0.43
1:BA:2544:G:H1'	1:BA:2646:C:H4'	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BB:9:G:OP1	14:BS:15:ARG:HD3	2.17	0.43
2:BB:31:C:H4'	6:BG:29:TRP:HH2	1.83	0.43
2:BB:31:C:O5'	2:BB:31:C:H6	2.01	0.43
7:BH:154:PRO:HB3	7:BH:163:TYR:CZ	2.53	0.43
21:BZ:45:ASP:O	21:BZ:49:ARG:HG3	2.18	0.43
31:CA:66:G:N2	31:CA:172:A:N3	2.66	0.43
31:CA:684:A:C6	31:CA:685:G:C6	3.06	0.43
31:CA:1051:C:H2'	31:CA:1052:U:H6	1.83	0.43
35:CE:48:ALA:O	35:CE:50:GLU:N	2.51	0.43
37:CG:99:LEU:HD22	37:CG:103:TRP:CZ2	2.53	0.43
42:CL:119:LYS:HB2	42:CL:120:TYR:HD1	1.82	0.43
43:CM:69:GLU:HG3	43:CM:70:LEU:H	1.82	0.43
49:CS:37:ARG:H	49:CS:37:ARG:HG3	1.57	0.43
50:CT:56:MET:HE3	50:CT:88:VAL:HG11	1.99	0.43
31:DA:93:G:O2'	31:DA:96:U:P	2.75	0.43
31:DA:986:A:H2'	31:DA:987:G:C8	2.52	0.43
31:DA:1002:G:H2'	31:DA:1003:G:O4'	2.18	0.43
31:DA:1004:A:H5'	31:DA:1024:G:N2	2.33	0.43
37:DG:70:LYS:HA	37:DG:71:PRO:HD3	1.91	0.43
40:DJ:61:GLU:OE2	44:DN:45:ARG:NE	2.49	0.43
42:DL:85:ILE:HD13	42:DL:85:ILE:HA	1.80	0.43
1:AA:793:A:OP2	1:AA:2071:A:O2'	2.35	0.43
1:AA:1967:C:H2'	1:AA:1968:G:O4'	2.17	0.43
1:AA:2027:G:H2'	1:AA:2028:U:O4'	2.18	0.43
1:AA:2171:A:H1'	1:AA:2172:U:C6	2.53	0.43
1:AA:2227:A:H4'	3:AD:265:PRO:HD3	2.00	0.43
1:AA:2494:G:C6	1:AA:2495:G:C5	3.06	0.43
1:AA:2590:A:C2	1:AA:2605:U:C2	3.06	0.43
1:AA:2698:U:H2'	1:AA:2699:C:C6	2.53	0.43
1:AA:2834:G:H5''	1:AA:2834:G:H8	1.82	0.43
2:AB:29:A:H5''	2:AB:30:C:OP2	2.19	0.43
11:AP:100:LEU:O	11:AP:103:ALA:HB3	2.17	0.43
13:AR:21:TYR:OH	13:AR:43:GLU:HG2	2.19	0.43
17:AV:28:GLU:HA	17:AV:29:PRO:HD3	1.84	0.43
25:A3:52:HIS:CD2	25:A3:53:LEU:HD21	2.53	0.43
28:A6:40:CYS:SG	28:A6:42:TRP:HB2	2.58	0.43
1:BA:12:U:H2'	1:BA:12:U:O2	2.18	0.43
1:BA:271(D):G:C6	1:BA:271(E):U:C4	3.05	0.43
1:BA:886:C:H2'	1:BA:887:A:H5''	2.00	0.43
1:BA:1773:A:H2'	1:BA:1774:C:H5'	1.99	0.43
1:BA:1983:C:H4'	1:BA:2606:C:H4'	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:1990:C:H2'	1:BA:1991:U:O4'	2.17	0.43
5:BF:89:VAL:HG12	5:BF:90:PHE:N	2.33	0.43
13:BR:1:MET:H1	13:BR:1:MET:HG2	1.70	0.43
21:BZ:144:LEU:HA	21:BZ:144:LEU:HD12	1.70	0.43
31:CA:160:A:H8	31:CA:160:A:OP1	2.01	0.43
31:CA:377:G:H2'	31:CA:378:G:H8	1.82	0.43
31:CA:456:C:H2'	31:CA:457:C:C6	2.54	0.43
31:CA:1318:A:O2'	49:CS:37:ARG:HB3	2.17	0.43
35:CE:30:ALA:N	35:CE:46:GLY:O	2.47	0.43
40:CJ:50:ILE:HG13	40:CJ:60:ARG:HE	1.82	0.43
44:CN:47:LEU:HD23	44:CN:50:LYS:HD3	2.00	0.43
31:DA:804:U:H5''	31:DA:805:C:OP2	2.18	0.43
31:DA:1104:G:C4	31:DA:1105:A:C8	3.06	0.43
31:DA:1412:C:H2'	31:DA:1413:A:H8	1.84	0.43
34:DD:60:GLU:HG2	34:DD:202:LEU:HB2	2.00	0.43
38:DH:21:LYS:O	38:DH:65:TYR:OH	2.26	0.43
42:DL:74:GLY:O	42:DL:102:ARG:NH1	2.51	0.43
1:AA:83:G:N1	1:AA:102:G:H2'	2.34	0.43
1:AA:255:A:C6	1:AA:256:A:C5	3.06	0.43
1:AA:271(Q):G:C2	1:AA:271(R):G:C5	3.07	0.43
1:AA:1016:G:C4	1:AA:1017:G:C8	3.06	0.43
1:AA:1179:C:O2'	1:AA:1180:C:H5'	2.19	0.43
1:AA:1482:G:C6	1:AA:1507:A:C6	3.06	0.43
2:AB:24:G:H4'	2:AB:25:A:H8	1.81	0.43
5:AF:96:ASP:OD1	5:AF:98:SER:HB3	2.18	0.43
6:AG:11:TYR:O	6:AG:16:ARG:HG2	2.19	0.43
6:AG:109:VAL:C	6:AG:112:PRO:HD2	2.38	0.43
6:AG:124:SER:HB3	6:AG:132:ASN:O	2.18	0.43
12:AQ:134:ARG:HG3	12:AQ:134:ARG:HH11	1.82	0.43
15:AT:28:VAL:HG12	15:AT:30:VAL:HG23	2.00	0.43
25:A3:44:ARG:O	25:A3:48:GLU:HG3	2.19	0.43
1:BA:185:U:H2'	1:BA:186:G:C8	2.54	0.43
1:BA:371:A:O3'	1:BA:372:G:H4'	2.19	0.43
1:BA:554:U:C4	1:BA:555:U:C4	3.06	0.43
1:BA:614:U:H5'	1:BA:614(C):A:N6	2.33	0.43
1:BA:699:A:H2'	1:BA:700:G:O4'	2.19	0.43
1:BA:1042:G:C6	1:BA:1043:C:C4	3.07	0.43
1:BA:1799:G:H4'	1:BA:1800:C:O5'	2.18	0.43
1:BA:2164:C:H2'	1:BA:2165:G:O4'	2.18	0.43
1:BA:2579:C:H2'	1:BA:2580:U:O4'	2.18	0.43
4:BE:111:ARG:HB3	13:BR:1:MET:CE	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:BQ:72:LYS:HA	12:BQ:73:PRO:HD3	1.84	0.43
21:BZ:152:ALA:HA	21:BZ:155:LEU:HD22	1.99	0.43
21:BZ:154:ASP:N	21:BZ:154:ASP:OD1	2.46	0.43
21:BZ:159:PRO:HA	21:BZ:160:GLY:HA2	1.75	0.43
31:CA:196:A:N3	31:CA:222:U:H1'	2.33	0.43
31:CA:584:G:H5'	47:CQ:91:ARG:NH2	2.30	0.43
31:CA:630:G:H2'	31:CA:631:G:C8	2.51	0.43
31:CA:1126:U:O2	31:CA:1126:U:H3'	2.17	0.43
31:CA:1431:C:H2'	31:CA:1432:G:O4'	2.18	0.43
35:CE:122:GLU:OE1	35:CE:131:ILE:HG13	2.18	0.43
38:CH:120:THR:OG1	38:CH:123:GLU:HG3	2.18	0.43
45:CO:74:ASP:CG	45:CO:77:ARG:H	2.22	0.43
31:DA:5:U:H5'	31:DA:6:G:C5	2.53	0.43
31:DA:186:C:O2'	31:DA:187:C:H5'	2.17	0.43
31:DA:361:G:H2'	31:DA:362:G:O4'	2.18	0.43
31:DA:412:A:C6	34:DD:35:ARG:HB3	2.54	0.43
31:DA:501:C:H1'	31:DA:549:C:H1'	2.01	0.43
31:DA:683:G:C4	31:DA:684:A:C8	3.06	0.43
31:DA:719:C:N4	48:DR:71:LYS:HE2	2.33	0.43
31:DA:778:G:H2'	31:DA:779:C:O4'	2.17	0.43
31:DA:1288:A:H2'	31:DA:1289:A:O4'	2.18	0.43
31:DA:1306:A:H1'	31:DA:1332:A:C2	2.53	0.43
31:DA:1375:A:H4'	37:DG:29:LYS:NZ	2.32	0.43
32:DB:185:ILE:HG22	32:DB:199:TYR:HD1	1.83	0.43
34:DD:116:GLN:O	34:DD:120:LEU:HG	2.19	0.43
35:DE:84:PHE:CE2	35:DE:133:TYR:HD2	2.36	0.43
39:DI:89:ASN:O	39:DI:91:ASP:N	2.51	0.43
40:DJ:32:ALA:HA	40:DJ:33:GLN:HA	1.84	0.43
49:DS:33:THR:O	49:DS:57:HIS:HE1	2.01	0.43
1:AA:228:A:H2'	1:AA:230:U:O4'	2.17	0.43
1:AA:614(C):A:C4	5:AF:180:GLY:HA2	2.52	0.43
1:AA:784:A:C8	1:AA:792:G:C5	3.06	0.43
1:AA:1047:G:H2'	1:AA:1110:G:N1	2.31	0.43
1:AA:1174:A:H5'	1:AA:1177:A:H62	1.78	0.43
1:AA:1494:A:C2'	1:AA:1495:A:H5'	2.48	0.43
1:AA:2469:A:H4'	12:AQ:56:ARG:HG2	1.99	0.43
1:AA:2552:U:H2'	1:AA:2554:U:OP2	2.18	0.43
1:AA:2600:A:C6	1:AA:2601:C:N4	2.86	0.43
1:AA:2789:C:H5''	1:AA:2790:A:OP2	2.18	0.43
14:AS:96:GLY:N	14:AS:99:LYS:HB3	2.32	0.43
18:AW:1:MET:HG3	18:AW:2:GLU:N	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:AX:40:LYS:HG3	19:AX:51:VAL:HB	2.00	0.43
20:AY:76:CYS:HG	20:AY:102:CYS:HG	1.60	0.43
21:AZ:35:ARG:HA	21:AZ:35:ARG:HD2	1.85	0.43
21:AZ:151:HIS:O	21:AZ:153:SER:N	2.50	0.43
24:A2:45:SER:O	24:A2:46:GLN:HB2	2.18	0.43
26:A4:18:CYS:SG	26:A4:39:CYS:CB	3.06	0.43
28:A6:4:GLU:HG3	28:A6:5:VAL:N	2.34	0.43
1:BA:322:A:OP1	5:BF:168:ARG:HD2	2.18	0.43
1:BA:351:G:H5''	1:BA:352:G:OP1	2.18	0.43
1:BA:468:G:O6	1:BA:469:G:C2	2.72	0.43
1:BA:998:C:H2'	1:BA:999:U:O4'	2.18	0.43
1:BA:1796:U:H2'	1:BA:1797:C:C6	2.53	0.43
1:BA:2111:C:H42	1:BA:2147:G:H22	1.66	0.43
1:BA:2198:A:O2'	1:BA:2224:G:N2	2.50	0.43
1:BA:2658:C:O5'	1:BA:2658:C:H6	2.01	0.43
1:BA:2684:U:H1'	10:BO:70:LYS:HD2	2.00	0.43
1:BA:2730:C:H4'	4:BE:168:MET:O	2.18	0.43
3:BD:68:LYS:HD2	3:BD:70:TRP:CZ2	2.53	0.43
4:BE:50:GLY:CA	4:BE:75:VAL:HG11	2.49	0.43
4:BE:144:ARG:HB3	4:BE:145:LYS:H	1.33	0.43
15:BT:7:ILE:O	15:BT:11:GLU:HG3	2.18	0.43
19:BX:12:VAL:HG22	19:BX:29:TRP:CE2	2.54	0.43
27:B5:51:TYR:CE1	27:B5:56:LYS:HG2	2.53	0.43
31:CA:10:A:H2'	31:CA:11:G:C8	2.53	0.43
31:CA:735:C:O2'	31:CA:736:C:H5'	2.19	0.43
31:CA:1084:G:H2'	31:CA:1085:U:C5	2.54	0.43
31:CA:1084:G:C5	31:CA:1085:U:C4	3.07	0.43
31:CA:1289:A:H2	31:CA:1372:U:O4'	2.01	0.43
31:CA:1399:C:C2	31:CA:1502:A:N6	2.86	0.43
31:CA:1511:G:H2'	31:CA:1512:U:O4'	2.17	0.43
32:CB:55:PHE:CD1	32:CB:58:ILE:HD12	2.53	0.43
32:CB:74:LYS:O	32:CB:78:GLN:N	2.51	0.43
32:CB:87:ARG:CZ	32:CB:233:SER:HB2	2.48	0.43
37:CG:22:LEU:HG	37:CG:62:PHE:CE2	2.52	0.43
38:CH:82:HIS:CE1	38:CH:84:ARG:HG2	2.53	0.43
40:CJ:45:ARG:HH21	44:CN:36:PHE:HB2	1.82	0.43
31:DA:115:G:C2	31:DA:289:G:N7	2.87	0.43
31:DA:117:G:H2'	31:DA:118:U:O4'	2.19	0.43
31:DA:191:G:C6	31:DA:192:U:N3	2.87	0.43
31:DA:1010:G:N1	31:DA:1020:U:O2	2.52	0.43
31:DA:1240:U:OP2	37:DG:116:ALA:N	2.50	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:1338:G:H2'	31:DA:1339:A:C8	2.54	0.43
32:DB:17:PHE:HB3	32:DB:44:LEU:HD11	2.00	0.43
32:DB:81:VAL:HB	32:DB:94:ASN:HD21	1.84	0.43
40:DJ:6:ILE:HA	40:DJ:97:GLU:O	2.18	0.43
41:DK:18:ARG:HB2	41:DK:33:THR:OG1	2.18	0.43
49:DS:60:VAL:O	49:DS:62:ILE:HD12	2.18	0.43
1:AA:108:U:C2	1:AA:109:G:C8	3.06	0.43
1:AA:1022:G:N2	1:AA:1142(A):A:H2	2.03	0.43
1:AA:1025:G:C4	1:AA:1135:C:H1'	2.53	0.43
1:AA:1686:C:C4	1:AA:1687:G:C5	3.06	0.43
1:AA:2186:G:H2'	1:AA:2187:G:H5''	2.00	0.43
1:AA:2292:C:O2'	1:AA:2293:C:H5'	2.18	0.43
6:AG:50:ALA:O	6:AG:52:ILE:N	2.51	0.43
8:AI:133:HIS:HA	8:AI:134:PRO:HD2	1.67	0.43
9:AN:46:VAL:HG12	9:AN:48:MET:HG2	2.00	0.43
9:AN:138:LEU:HD23	9:AN:138:LEU:HA	1.62	0.43
14:AS:102:ALA:HB1	14:AS:112:PHE:CZ	2.53	0.43
15:AT:16:ARG:HG3	15:AT:79:HIS:HA	2.01	0.43
16:AU:80:ILE:HD13	16:AU:80:ILE:HA	1.75	0.43
22:A0:78:TYR:HB3	22:A0:80:HIS:NE2	2.33	0.43
23:A1:3:LYS:HB3	23:A1:4:VAL:H	1.39	0.43
1:BA:443:A:C5	5:BF:45:ARG:HD2	2.54	0.43
1:BA:484:C:H2'	1:BA:485:C:H6	1.84	0.43
1:BA:807:U:OP2	11:BP:41:ARG:NH2	2.51	0.43
1:BA:821:A:O2'	1:BA:946:G:OP2	2.20	0.43
1:BA:994:C:H5''	1:BA:995:C:OP1	2.17	0.43
1:BA:1022:G:C6	1:BA:1140:C:C4	3.07	0.43
1:BA:1283:G:N2	1:BA:1286:A:C8	2.87	0.43
1:BA:1641:A:H2'	1:BA:1642:G:O4'	2.18	0.43
1:BA:2055:C:H5'	1:BA:2056:G:O5'	2.17	0.43
1:BA:2136:C:N4	1:BA:2155:G:N1	2.65	0.43
1:BA:2404:C:O3'	11:BP:77:ARG:NH2	2.52	0.43
2:BB:6:C:C2	2:BB:116:G:C2	3.06	0.43
5:BF:63:LYS:HA	5:BF:76:GLY:O	2.18	0.43
6:BG:106:LEU:HD12	6:BG:110:ALA:HB3	2.00	0.43
6:BG:131:TYR:HB3	6:BG:159:VAL:HG13	2.00	0.43
12:BQ:134:ARG:O	12:BQ:138:ASP:HB2	2.18	0.43
15:BT:22:PHE:HA	15:BT:91:ARG:HH12	1.84	0.43
23:B1:12:PRO:HG3	23:B1:43:TYR:HD2	1.83	0.43
24:B2:48:HIS:O	24:B2:52:ASP:HB2	2.19	0.43
31:CA:154:C:N3	31:CA:168:G:C2	2.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:258:G:N3	31:CA:259:G:C8	2.86	0.43
31:CA:445:G:H2'	31:CA:446:G:H8	1.84	0.43
31:CA:683:G:C2	31:CA:708:C:N3	2.87	0.43
31:CA:1078:U:C5	31:CA:1079:G:C5	3.07	0.43
31:CA:1106:G:H2'	31:CA:1107:C:C6	2.53	0.43
31:CA:1179:A:H4'	39:CI:103:THR:HA	1.99	0.43
31:CA:1202:G:O4'	44:CN:29:ARG:NH1	2.45	0.43
31:CA:1256:A:H62	31:CA:1278:U:H1'	1.84	0.43
32:CB:73:THR:HB	32:CB:95:GLN:O	2.19	0.43
35:CE:33:VAL:HG23	35:CE:42:GLY:O	2.19	0.43
39:CI:112:LYS:HA	39:CI:119:ALA:HB2	2.00	0.43
31:DA:298:A:H5''	31:DA:299:G:OP2	2.19	0.43
31:DA:579:G:H2'	31:DA:580:U:C6	2.53	0.43
31:DA:1030(A):G:C8	31:DA:1030(B):C:H5''	2.53	0.43
31:DA:1189:C:P	40:DJ:51:ARG:HH22	2.42	0.43
49:DS:53:ASN:O	49:DS:77:THR:HG22	2.18	0.43
1:AA:7:G:H2'	1:AA:8:A:H8	1.80	0.43
1:AA:83:G:H1	1:AA:102:G:H2'	1.83	0.43
1:AA:193:U:O3'	1:AA:803:U:H4'	2.19	0.43
1:AA:721:C:H2'	1:AA:722:A:C8	2.54	0.43
1:AA:1470:G:N2	1:AA:1523:U:C4	2.86	0.43
1:AA:1593:G:C2	1:AA:1594:G:C4	3.07	0.43
1:AA:2227:A:OP1	3:AD:263:ARG:HD2	2.19	0.43
1:AA:2313:C:H4'	6:AG:91:ARG:HG3	2.00	0.43
3:AD:43:ARG:HG2	3:AD:47:GLY:O	2.19	0.43
3:AD:102:LYS:C	3:AD:103:ARG:HG2	2.38	0.43
4:AE:37:ARG:HG3	4:AE:46:ALA:HB3	2.01	0.43
7:AH:9:ILE:HD13	7:AH:72:ILE:HG22	2.00	0.43
8:AI:77:LEU:HD21	8:AI:101:LEU:HG	2.00	0.43
9:AN:120:LEU:HD22	9:AN:122:VAL:HG23	2.01	0.43
21:AZ:67:LEU:HA	21:AZ:68:PRO:HD3	1.90	0.43
22:A0:31:VAL:HG11	22:A0:37:LEU:CD2	2.47	0.43
24:A2:32:LEU:HD11	24:A2:54:LYS:CG	2.48	0.43
1:BA:32:C:O2'	1:BA:33:U:H5'	2.19	0.43
1:BA:330:A:HO2'	1:BA:331:A:H8	1.61	0.43
1:BA:340:A:H2'	1:BA:341:G:O4'	2.18	0.43
1:BA:652(R):C:HO2'	1:BA:652(S):C:H6	1.61	0.43
1:BA:1673:U:C5	4:BE:129:HIS:HE1	2.36	0.43
1:BA:2286:A:OP1	28:B6:29:ASN:ND2	2.51	0.43
1:BA:2336:A:H61	22:B0:43:THR:HG22	1.84	0.43
1:BA:2670:A:O2'	1:BA:2671:A:H5'	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BG:43:LEU:HB3	6:BG:44:GLY:H	1.70	0.43
6:BG:46:ALA:HB1	6:BG:51:ARG:HA	2.00	0.43
7:BH:5:GLY:HA2	7:BH:69:ARG:CB	2.48	0.43
7:BH:91:GLY:HA3	7:BH:160:LYS:HG3	1.99	0.43
7:BH:158:HIS:O	7:BH:160:LYS:N	2.48	0.43
14:BS:69:VAL:HA	14:BS:72:ALA:HB3	2.00	0.43
19:BX:8:ILE:HD11	19:BX:43:VAL:HG23	2.01	0.43
21:BZ:144:LEU:HD21	21:BZ:150:LEU:HG	2.00	0.43
31:CA:602:A:C6	31:CA:603:U:C4	3.06	0.43
31:CA:833:U:H2'	31:CA:834:C:C6	2.53	0.43
31:CA:981:U:N3	31:CA:982:U:C4	2.86	0.43
31:CA:1128:C:N4	31:CA:1139:G:N3	2.67	0.43
35:CE:33:VAL:HG21	35:CE:109:ILE:HA	2.01	0.43
36:CF:11:ASN:O	36:CF:14:LEU:HB2	2.18	0.43
31:DA:298:A:C6	31:DA:299:G:C2	3.07	0.43
31:DA:447:G:H8	31:DA:447:G:O5'	2.01	0.43
31:DA:538:G:H5''	42:DL:114:LYS:HB2	2.01	0.43
31:DA:684:A:C2	31:DA:685:G:C4	3.07	0.43
31:DA:990:C:C2	31:DA:1216:G:N2	2.86	0.43
31:DA:1057:G:H2'	31:DA:1058:G:O4'	2.18	0.43
31:DA:1151:A:O2'	31:DA:1152:A:C8	2.70	0.43
34:DD:38:TYR:HA	34:DD:39:PRO:HD3	1.77	0.43
41:DK:21:ILE:HD12	41:DK:84:VAL:HG22	2.01	0.43
1:AA:35:G:H2'	1:AA:36:G:O4'	2.19	0.43
1:AA:340:A:N6	1:AA:341:G:C2	2.87	0.43
1:AA:631:A:OP2	30:A8:47:LYS:NZ	2.38	0.43
1:AA:862:G:H4'	2:AB:79:C:H5'	2.01	0.43
1:AA:1337:G:H2'	1:AA:1338:G:C8	2.53	0.43
1:AA:1711:C:H2'	1:AA:1712:C:C6	2.54	0.43
1:AA:1827:C:O2'	1:AA:1970:A:N3	2.32	0.43
1:AA:1997:G:O2'	1:AA:1998:G:H5'	2.19	0.43
1:AA:2050:C:O2	4:AE:156:MET:HE1	2.19	0.43
2:AB:15:A:OP2	2:AB:69:G:N2	2.50	0.43
5:AF:110:LEU:HD23	5:AF:110:LEU:HA	1.69	0.43
6:AG:36:LYS:HD3	6:AG:95:ARG:NH1	2.34	0.43
9:AN:55:VAL:HG22	9:AN:126:PRO:HA	2.00	0.43
11:AP:63:PRO:O	30:A8:13:ARG:HB2	2.19	0.43
19:AX:50:LYS:HG2	19:AX:84:ALA:HB2	2.01	0.43
22:A0:50:ASN:HB3	22:A0:63:VAL:HG22	2.00	0.43
1:BA:221:A:H4'	1:BA:222:A:O5'	2.16	0.43
1:BA:263:C:H2'	1:BA:264:C:O4'	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:271(H):G:O2'	1:BA:271(I):G:H8	2.02	0.43
1:BA:841:A:C2	1:BA:938:G:C2	3.07	0.43
1:BA:1582:C:C2	1:BA:1583:A:C8	3.06	0.43
1:BA:2165:G:H2'	1:BA:2166:G:H8	1.81	0.43
1:BA:2405:G:O2'	1:BA:2406:U:P	2.77	0.43
11:BP:3:LEU:HA	11:BP:3:LEU:HD12	1.82	0.43
12:BQ:2:LEU:HD12	12:BQ:2:LEU:HA	1.92	0.43
12:BQ:14:ARG:HG2	12:BQ:41:TRP:HH2	1.83	0.43
12:BQ:32:TYR:CZ	12:BQ:133:ARG:HG3	2.54	0.43
31:CA:355:C:O4'	31:CA:388:G:O2'	2.32	0.43
31:CA:1086:U:H2'	31:CA:1087:G:O4'	2.18	0.43
32:CB:71:VAL:HA	32:CB:93:VAL:HG23	2.01	0.43
32:CB:185:ILE:HG22	32:CB:199:TYR:HD1	1.84	0.43
37:CG:72:ARG:NH1	37:CG:142:GLU:OE1	2.52	0.43
37:CG:111:ARG:HB2	37:CG:119:ARG:HD2	2.01	0.43
45:CO:26:GLU:HA	45:CO:81:LEU:HD22	2.00	0.43
31:DA:130:A:H1'	31:DA:263:A:O2'	2.19	0.43
31:DA:626:U:C2	31:DA:627:G:C8	3.07	0.43
31:DA:1133:G:H2'	31:DA:1134:G:O4'	2.19	0.43
37:DG:149:ARG:HD2	41:DK:59:TYR:CE1	2.53	0.43
49:DS:38:SER:O	49:DS:70:LYS:HG3	2.18	0.43
1:AA:154:G:C6	1:AA:154(A):C:N4	2.86	0.43
1:AA:345:A:H1'	1:AA:346:A:N7	2.33	0.43
1:AA:573:G:O2'	1:AA:574:C:H3'	2.19	0.43
1:AA:619:G:H8	1:AA:619:G:O5'	2.01	0.43
1:AA:770:G:OP1	29:A7:8:ASN:ND2	2.47	0.43
1:AA:780:G:C2	1:AA:782:A:C2	3.06	0.43
1:AA:909:A:H2'	1:AA:912:C:H5	1.82	0.43
1:AA:1288:U:O4	13:AR:106:GLY:HA3	2.19	0.43
1:AA:1587:A:H2'	1:AA:1588:C:C6	2.53	0.43
1:AA:1689:A:N7	1:AA:1698:A:N1	2.67	0.43
1:AA:2335:A:C8	1:AA:2337:G:C5	3.06	0.43
1:AA:2699:C:H2'	1:AA:2700:C:O4'	2.18	0.43
1:AA:2751:G:H3'	1:AA:2752:C:C6	2.53	0.43
4:AE:101:ARG:HA	4:AE:170:LEU:O	2.18	0.43
13:AR:38:VAL:CG1	13:AR:42:LYS:HE3	2.49	0.43
17:AV:89:GLN:OE1	17:AV:89:GLN:HA	2.18	0.43
18:AW:78:GLU:OE1	18:AW:99:ARG:NH1	2.51	0.43
21:AZ:72:ARG:O	21:AZ:72:ARG:HD3	2.19	0.43
25:A3:19:GLN:O	25:A3:23:LEU:HD22	2.19	0.43
1:BA:222:A:C2	1:BA:233:A:H5''	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:725:G:C6	1:BA:726:G:N1	2.86	0.43
1:BA:994:C:OP1	16:BU:53:ARG:NH2	2.51	0.43
1:BA:1903:G:P	3:BD:241:PRO:HB2	2.58	0.43
1:BA:1907:G:C6	1:BA:1908:C:C4	3.06	0.43
1:BA:2445:G:C2'	1:BA:2446:G:H5'	2.49	0.43
1:BA:2803:C:H2'	1:BA:2804:C:O4'	2.18	0.43
2:BB:78:A:H2'	2:BB:79:C:O4'	2.18	0.43
19:BX:52:VAL:HG12	19:BX:82:GLN:HG3	2.00	0.43
20:BY:79:CYS:CB	20:BY:81:LYS:H	2.31	0.43
26:B4:35:VAL:O	26:B4:36:CYS:HB3	2.18	0.43
28:B6:9:LEU:HD21	28:B6:25:LYS:HB3	2.01	0.43
31:CA:708:C:H2'	31:CA:709:G:C8	2.52	0.43
31:CA:996:A:H2'	31:CA:997:U:O4'	2.19	0.43
31:CA:1232:U:H2'	31:CA:1233:G:O4'	2.19	0.43
31:CA:1347:G:N7	39:CI:11:LYS:HD2	2.34	0.43
31:CA:1352:C:H1'	31:CA:1371:G:N2	2.34	0.43
32:CB:17:PHE:CB	32:CB:44:LEU:HD11	2.49	0.43
32:CB:185:ILE:HG22	32:CB:199:TYR:CD1	2.52	0.43
33:CC:74:GLY:C	33:CC:76:VAL:N	2.72	0.43
35:CE:15:ARG:HD2	35:CE:26:PHE:CD2	2.54	0.43
38:CH:77:GLU:HG3	38:CH:78:GLN:N	2.34	0.43
40:CJ:75:LEU:O	40:CJ:77:PRO:HD3	2.19	0.43
31:DA:16:A:O2'	31:DA:17:U:H5'	2.18	0.43
31:DA:25:C:O2'	31:DA:26:A:H5'	2.19	0.43
31:DA:134:A:H61	46:DP:25:ARG:NH1	2.16	0.43
31:DA:149:A:O2'	31:DA:150:C:P	2.77	0.43
31:DA:367:U:C2	31:DA:369:C:C5	3.06	0.43
31:DA:373:A:N1	31:DA:391:G:O2'	2.45	0.43
31:DA:612:C:O2	31:DA:629:G:N2	2.52	0.43
31:DA:638:G:O2'	31:DA:639:G:H5'	2.19	0.43
31:DA:734:G:O2'	48:DR:71:LYS:HD3	2.19	0.43
31:DA:825:G:O2'	31:DA:826:C:H5'	2.19	0.43
31:DA:949:A:C1'	31:DA:1364:U:H3	2.32	0.43
31:DA:1106:G:C6	31:DA:1107:C:C4	3.07	0.43
31:DA:1372:U:OP1	39:DI:72:GLY:N	2.52	0.43
32:DB:87:ARG:HD2	32:DB:219:VAL:HG11	2.01	0.43
32:DB:133:LYS:C	32:DB:135:GLN:H	2.21	0.43
33:DC:39:ILE:HD11	33:DC:59:ARG:HH21	1.84	0.43
33:DC:148:GLY:HA3	33:DC:172:ARG:O	2.18	0.43
35:DE:31:LEU:HD23	35:DE:31:LEU:HA	1.76	0.43
37:DG:152:ALA:O	37:DG:155:ARG:HB3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:DI:3:GLN:HE21	39:DI:3:GLN:HB2	1.67	0.43
41:DK:73:MET:HE2	41:DK:103:LEU:HD23	2.00	0.43
1:AA:320:A:H4'	1:AA:322:A:N7	2.33	0.43
1:AA:1204:A:N1	1:AA:1241:A:N7	2.67	0.43
1:AA:1238:G:N2	1:AA:1239:G:H1'	2.34	0.43
1:AA:1285:G:N2	1:AA:1328:G:H5''	2.33	0.43
1:AA:1986:A:H2'	1:AA:1987:G:C8	2.54	0.43
1:AA:2472:G:O6	1:AA:2476:A:H4'	2.19	0.43
2:AB:75:G:N1	2:AB:103:G:N2	2.67	0.43
3:AD:132:PRO:HG3	3:AD:190:TYR:CE1	2.54	0.43
6:AG:86:MET:HA	6:AG:87:PRO:HD2	1.93	0.43
8:AI:145:VAL:HG12	8:AI:146:ALA:N	2.34	0.43
21:AZ:111:VAL:C	21:AZ:113:ALA:H	2.21	0.43
30:A8:6:THR:HG22	30:A8:63:PRO:HD2	1.99	0.43
1:BA:562:U:C4	1:BA:2036:C:O4'	2.72	0.43
1:BA:570:G:H2'	1:BA:2030:A:N7	2.33	0.43
1:BA:833:U:H2'	1:BA:834:C:C6	2.53	0.43
1:BA:1685:C:O2'	1:BA:1686:C:H5'	2.18	0.43
1:BA:2307:G:N1	6:BG:43:LEU:O	2.33	0.43
1:BA:2428:G:H5''	1:BA:2429:G:O5'	2.19	0.43
3:BD:26:LYS:HB3	3:BD:83:GLU:HG2	2.00	0.43
6:BG:60:LEU:HB3	6:BG:68:PRO:HG3	1.99	0.43
11:BP:59:LEU:HD21	30:B8:10:ALA:HA	2.01	0.43
12:BQ:109:VAL:HG22	12:BQ:113:GLN:OE1	2.19	0.43
13:BR:33:ARG:HA	13:BR:114:VAL:O	2.19	0.43
16:BU:58:ARG:HA	16:BU:61:TRP:CE3	2.53	0.43
17:BV:24:LYS:HA	17:BV:92:THR:OG1	2.18	0.43
22:B0:72:ARG:CB	22:B0:75:LEU:HB2	2.49	0.43
26:B4:40:HIS:O	26:B4:43:TYR:HB3	2.18	0.43
31:CA:49:U:H3	31:CA:362:G:H1'	1.84	0.43
31:CA:513:C:H2'	31:CA:514:C:C6	2.54	0.43
31:CA:749:C:O5'	31:CA:749:C:H6	2.01	0.43
31:CA:837:G:C2	31:CA:850:U:O2	2.72	0.43
31:CA:1015:A:H2'	31:CA:1016:A:C8	2.54	0.43
31:CA:1270:C:H2'	31:CA:1271:G:H8	1.83	0.43
31:CA:1385:G:C6	31:CA:1386:G:N7	2.86	0.43
31:CA:1438:G:H2'	31:CA:1439:C:C6	2.53	0.43
40:CJ:13:HIS:HB3	40:CJ:68:HIS:CE1	2.54	0.43
40:CJ:19:SER:OG	40:CJ:91:PRO:HD3	2.19	0.43
40:CJ:79:ARG:C	40:CJ:81:THR:H	2.21	0.43
48:CR:45:SER:C	48:CR:47:THR:H	2.21	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:509:A:H4'	31:DA:510:A:OP1	2.19	0.43
31:DA:542:G:H5'	34:DD:41:GLY:HA3	2.01	0.43
31:DA:585:G:N3	31:DA:879:C:H4'	2.34	0.43
31:DA:663:A:H2'	31:DA:664:G:O4'	2.18	0.43
31:DA:680:C:H6	31:DA:680:C:O5'	2.01	0.43
31:DA:867:G:N2	31:DA:868:C:C2	2.87	0.43
31:DA:1037:C:H3'	31:DA:1037:C:H6	1.84	0.43
31:DA:1288:A:N3	31:DA:1352:C:O2'	2.44	0.43
31:DA:1288:A:C6	31:DA:1289:A:C5	3.07	0.43
31:DA:1423:G:C2	31:DA:1424:C:C2	3.06	0.43
32:DB:145:LEU:HD13	32:DB:145:LEU:HA	1.94	0.43
33:DC:131:ARG:NE	33:DC:166:GLU:OE2	2.41	0.43
34:DD:123:HIS:O	34:DD:125:HIS:N	2.52	0.43
36:DF:67:MET:HG3	36:DF:68:PRO:O	2.18	0.43
1:AA:328:U:H4'	20:AY:68:HIS:CG	2.54	0.43
1:AA:422:A:H2'	1:AA:423:A:C8	2.54	0.43
1:AA:528:A:C2	1:AA:2043:C:H4'	2.53	0.43
1:AA:674:G:H1'	5:AF:74:ARG:CD	2.49	0.43
1:AA:897:C:H2'	1:AA:898:C:C6	2.54	0.43
1:AA:977:G:N3	1:AA:1001:A:H2	2.16	0.43
1:AA:1184:G:C6	1:AA:1185:C:C4	3.07	0.43
1:AA:1250:G:OP2	11:AP:21:ARG:HD3	2.19	0.43
1:AA:1786:A:H1'	1:AA:1938:A:H61	1.82	0.43
1:AA:1853:A:H2'	1:AA:1854:A:C8	2.54	0.43
1:AA:2063:C:O2	1:AA:2450:A:N1	2.52	0.43
1:AA:2173:A:C3'	1:AA:2174:C:H5'	2.49	0.43
1:AA:2299:G:N2	1:AA:2300:G:C4	2.87	0.43
1:AA:2360:A:H2'	1:AA:2361:A:O4'	2.19	0.43
1:AA:2661:G:H2'	1:AA:2662:A:C8	2.53	0.43
1:AA:2815:C:H2'	1:AA:2816:C:O4'	2.19	0.43
8:AI:99:GLU:HA	8:AI:102:SER:OG	2.19	0.43
9:AN:112:LEU:O	9:AN:115:ARG:HB2	2.19	0.43
12:AQ:58:PHE:O	12:AQ:60:ARG:N	2.52	0.43
14:AS:62:LYS:HB3	14:AS:97:ARG:CD	2.48	0.43
17:AV:38:LEU:O	17:AV:39:LEU:HB2	2.18	0.43
25:A3:26:LEU:HD21	25:A3:46:ASN:CB	2.49	0.43
1:BA:271(U):G:H2'	1:BA:271(V):G:H8	1.83	0.43
1:BA:543:C:H2'	1:BA:545:G:O4'	2.19	0.43
1:BA:637:A:C8	11:BP:117:GLU:HG3	2.54	0.43
1:BA:1110:G:OP2	1:BA:1110:G:C8	2.72	0.43
1:BA:1494:A:C6	1:BA:1495:A:C6	3.07	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:2037:G:C6	1:BA:2038:G:C6	3.07	0.43
1:BA:2127:G:O6	1:BA:2161:C:N3	2.51	0.43
1:BA:2370:G:C6	1:BA:2371:G:C6	3.07	0.43
1:BA:2653:U:O2'	7:BH:110:SER:HB2	2.18	0.43
1:BA:2788:C:N4	1:BA:2789:C:N4	2.67	0.43
2:BB:50:G:H2'	2:BB:51:G:O4'	2.19	0.43
3:BD:95:LEU:O	3:BD:102:LYS:HA	2.19	0.43
5:BF:132:VAL:HG23	5:BF:163:VAL:HG22	2.01	0.43
5:BF:157:VAL:HG21	5:BF:181:LEU:HD13	2.01	0.43
6:BG:47:LYS:O	6:BG:82:LEU:HD21	2.19	0.43
13:BR:53:HIS:O	13:BR:56:LYS:HB2	2.19	0.43
31:CA:530:G:OP1	31:CA:530:G:H3'	2.19	0.43
31:CA:545:C:O2'	31:CA:546:G:H5'	2.19	0.43
31:CA:606:G:H2'	31:CA:631:G:H1	1.83	0.43
31:CA:742:G:H5'	45:CO:58:MET:HE3	2.00	0.43
31:CA:763:G:H2'	31:CA:764:C:H6	1.84	0.43
31:CA:827:U:H5''	31:CA:828:A:OP2	2.19	0.43
31:CA:1162:C:H2'	31:CA:1163:C:O4'	2.18	0.43
31:CA:1171:G:H2'	31:CA:1172:C:C6	2.54	0.43
31:CA:1365:G:C6	31:CA:1366:C:C4	3.07	0.43
35:CE:90:VAL:O	35:CE:120:THR:HA	2.19	0.43
39:CI:17:VAL:HA	39:CI:63:ILE:HG12	2.01	0.43
42:CL:77:LEU:HD21	42:CL:107:ALA:HA	2.00	0.43
42:CL:124:LYS:HA	42:CL:125:PRO:HD3	1.90	0.43
49:CS:73:GLU:H	49:CS:73:GLU:HG2	1.70	0.43
50:CT:14:LYS:HG3	50:CT:17:ARG:NH2	2.34	0.43
31:DA:167:G:C4	31:DA:168:G:C8	3.07	0.43
31:DA:300:A:H1'	31:DA:565:U:O2	2.19	0.43
31:DA:473:G:C4	31:DA:474:G:C8	3.07	0.43
31:DA:622:A:C8	31:DA:623:C:C6	3.07	0.43
31:DA:953:G:C6	31:DA:954:G:C4	3.06	0.43
31:DA:1016:A:H4'	31:DA:1218:C:H4'	2.00	0.43
31:DA:1151:A:H5'	40:DJ:41:PRO:HA	1.99	0.43
33:DC:24:ALA:HB1	33:DC:28:GLN:O	2.19	0.43
35:DE:139:LEU:HA	35:DE:142:LEU:CD1	2.49	0.43
42:DL:75:HIS:ND1	42:DL:77:LEU:HB2	2.34	0.43
49:DS:36:ARG:HB2	49:DS:72:GLY:HA3	2.01	0.43
1:AA:817:C:H2'	1:AA:818:G:C8	2.54	0.42
1:AA:1007:C:P	9:AN:37:LYS:HZ2	2.42	0.42
1:AA:1167:U:C2	1:AA:1183:G:N2	2.87	0.42
1:AA:1799:G:H4'	1:AA:1800:C:O5'	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1915:U:N3	52:AA:3001:T8B:C32	2.82	0.42
1:AA:2109:U:H2'	1:AA:2110:G:C8	2.54	0.42
4:AE:2:LYS:HA	4:AE:84:PHE:CD2	2.54	0.42
6:AG:50:ALA:C	6:AG:52:ILE:N	2.72	0.42
9:AN:28:THR:HG22	9:AN:29:LYS:HG3	2.01	0.42
14:AS:67:ARG:HE	14:AS:67:ARG:HB2	1.32	0.42
16:AU:65:ILE:HD11	16:AU:95:LEU:HB3	2.01	0.42
16:AU:97:ASP:OD1	16:AU:101:ARG:HD2	2.19	0.42
17:AV:59:ALA:HB2	17:AV:96:ILE:HD13	2.00	0.42
20:AY:23:ARG:NH1	20:AY:23:ARG:HB2	2.34	0.42
20:AY:76:CYS:HA	20:AY:77:PRO:HD3	1.76	0.42
21:AZ:45:ASP:OD1	21:AZ:49:ARG:NH1	2.51	0.42
21:AZ:154:ASP:N	21:AZ:154:ASP:OD1	2.49	0.42
27:A5:46:CYS:HA	27:A5:47:PRO:HD3	1.87	0.42
1:BA:271(X):G:C2	1:BA:271(Y):U:O4	2.72	0.42
1:BA:775:G:C4	1:BA:794:G:C8	3.07	0.42
1:BA:1614:A:P	1:BA:1614:A:H8	2.41	0.42
1:BA:1719:G:C6	1:BA:1720:U:C4	3.07	0.42
1:BA:1846:G:H5''	1:BA:1847:A:OP2	2.18	0.42
1:BA:2012:G:O3'	18:BW:96:ILE:HG23	2.19	0.42
1:BA:2023:G:H4'	1:BA:2617:C:O3'	2.19	0.42
1:BA:2123:G:H2'	1:BA:2124:G:H8	1.84	0.42
1:BA:2394:C:P	30:B8:30:ARG:NH1	2.91	0.42
5:BF:102:PRO:HB2	5:BF:105:VAL:CG2	2.49	0.42
6:BG:123:ASN:C	6:BG:125:PHE:H	2.23	0.42
8:BI:111:PRO:HB2	8:BI:112:LYS:HE3	2.01	0.42
10:BO:104:ARG:NH1	15:BT:34:VAL:HG21	2.34	0.42
12:BQ:79:LEU:HD23	12:BQ:79:LEU:HA	1.78	0.42
31:CA:310:G:H5''	46:CP:31:LYS:HB2	2.01	0.42
31:CA:375:U:N3	31:CA:376:G:N7	2.67	0.42
31:CA:540:G:C4	31:CA:541:G:C8	3.07	0.42
31:CA:690:G:C6	31:CA:691:G:N1	2.87	0.42
31:CA:1005:A:H1'	31:CA:1036:G:N2	2.33	0.42
31:CA:1014:A:N3	31:CA:1219:U:H1'	2.34	0.42
32:CB:93:VAL:HG21	32:CB:97:TRP:HD1	1.84	0.42
34:CD:78:LEU:O	34:CD:81:GLU:N	2.52	0.42
34:CD:123:HIS:O	34:CD:125:HIS:N	2.52	0.42
38:CH:5:PRO:O	38:CH:8:ASP:HB3	2.19	0.42
31:DA:203:U:H2'	31:DA:203:U:OP2	2.19	0.42
31:DA:832:C:O2'	31:DA:833:U:P	2.77	0.42
31:DA:832:C:N4	31:DA:855:G:C6	2.86	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:892:A:H2'	31:DA:893:C:C6	2.54	0.42
31:DA:1226:C:H4'	49:DS:80:TYR:CZ	2.54	0.42
31:DA:1416:G:H2'	31:DA:1417:G:O4'	2.18	0.42
35:DE:78:HIS:CE1	35:DE:142:LEU:HD23	2.54	0.42
35:DE:127:ASN:HA	35:DE:128:PRO:HD3	1.84	0.42
47:DQ:81:ARG:HA	47:DQ:81:ARG:HD2	1.78	0.42
1:AA:247:G:H4'	1:AA:386:G:C6	2.54	0.42
1:AA:672:C:H2'	1:AA:673:C:C6	2.55	0.42
1:AA:910:A:C5	12:AQ:13:GLN:HG3	2.53	0.42
1:AA:1514:U:H2'	1:AA:1515:G:C8	2.54	0.42
1:AA:2052:G:C8	4:AE:141:ILE:HD11	2.54	0.42
1:AA:2299:G:N1	1:AA:2318:G:H8	2.17	0.42
1:AA:2322:A:H2'	1:AA:2323:G:O4'	2.19	0.42
1:AA:2680:C:H1'	4:AE:187:ALA:HB1	2.01	0.42
1:AA:2751:G:H3'	1:AA:2752:C:H6	1.84	0.42
3:AD:206:LEU:HD23	3:AD:206:LEU:HA	1.62	0.42
6:AG:19:LEU:HD13	6:AG:32:PRO:HD2	1.99	0.42
6:AG:122:PRO:HG2	6:AG:182:LYS:O	2.18	0.42
10:AO:66:LYS:N	10:AO:82:ASN:OD1	2.44	0.42
16:AU:59:ARG:HH11	16:AU:59:ARG:HB3	1.85	0.42
18:AW:7:ALA:O	18:AW:103:ILE:N	2.52	0.42
21:AZ:54:HIS:CG	21:AZ:101:PRO:HG3	2.54	0.42
23:A1:83:GLU:HA	23:A1:84:GLY:HA2	1.59	0.42
1:BA:484:C:H6	1:BA:484:C:O5'	2.02	0.42
1:BA:910:A:N7	12:BQ:13:GLN:HG3	2.34	0.42
1:BA:1329:U:H5'	1:BA:1330:C:H5	1.84	0.42
1:BA:2292:C:H4'	1:BA:2375:G:H4'	2.01	0.42
7:BH:11:VAL:HA	7:BH:12:PRO:HD3	1.79	0.42
8:BI:112:LYS:HA	8:BI:114:LEU:O	2.18	0.42
12:BQ:21:THR:HG21	12:BQ:101:ARG:HD2	2.01	0.42
12:BQ:51:ARG:HG3	12:BQ:66:ILE:HD11	2.01	0.42
15:BT:37:GLY:HA2	15:BT:38:ASN:HA	1.74	0.42
20:BY:51:VAL:HG22	20:BY:58:GLY:N	2.34	0.42
23:B1:83:GLU:HA	23:B1:84:GLY:HA2	1.65	0.42
24:B2:25:VAL:O	24:B2:29:LYS:N	2.48	0.42
24:B2:25:VAL:O	24:B2:28:LYS:N	2.51	0.42
25:B3:17:LYS:H	25:B3:17:LYS:HG2	1.61	0.42
31:CA:27:G:H2'	31:CA:28:G:O4'	2.19	0.42
31:CA:118:U:C5	31:CA:288:A:C6	3.07	0.42
31:CA:189:G:C6	31:CA:189(L):G:C6	3.07	0.42
31:CA:690:G:C6	31:CA:691:G:C2	3.07	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:949:A:H1'	31:CA:1364:U:H3	1.84	0.42
38:CH:44:PHE:HA	38:CH:79:VAL:CG1	2.50	0.42
40:CJ:47:PHE:HD2	44:CN:34:TYR:CE2	2.37	0.42
40:CJ:51:ARG:NH2	40:CJ:61:GLU:HB3	2.34	0.42
44:CN:53:LEU:HA	44:CN:54:PRO:HD3	1.84	0.42
48:CR:74:ARG:H	48:CR:74:ARG:HG3	1.60	0.42
31:DA:159:G:N2	31:DA:163:C:C4	2.87	0.42
31:DA:189(E):U:O2'	31:DA:189(F):U:H5'	2.19	0.42
31:DA:612:C:H2'	31:DA:613:C:C6	2.53	0.42
31:DA:724:G:C2	31:DA:725:G:C8	3.07	0.42
31:DA:1074:G:C4	31:DA:1102:A:C2	3.08	0.42
31:DA:1278:U:H5''	31:DA:1279:A:O4'	2.19	0.42
31:DA:1301:U:H2'	31:DA:1303:C:C5	2.54	0.42
33:DC:159:GLY:HA2	33:DC:193:TYR:CD1	2.53	0.42
37:DG:89:MET:HE2	37:DG:155:ARG:HG3	2.01	0.42
40:DJ:13:HIS:HA	40:DJ:16:LEU:HD22	2.00	0.42
42:DL:70:ILE:HD13	42:DL:77:LEU:HD12	2.01	0.42
46:DP:7:ALA:O	46:DP:9:PHE:HD2	2.01	0.42
46:DP:52:ASP:O	46:DP:54:GLU:N	2.52	0.42
1:AA:236:C:H2'	1:AA:237:C:C6	2.55	0.42
1:AA:247:G:H5''	1:AA:386:G:O2'	2.19	0.42
1:AA:284:U:H2'	1:AA:285:C:H6	1.84	0.42
1:AA:637:A:C8	11:AP:117:GLU:HG3	2.47	0.42
1:AA:652(R):C:O2'	1:AA:652(S):C:H6	2.01	0.42
1:AA:719:C:H2'	1:AA:720:C:C6	2.51	0.42
1:AA:817:C:H2'	1:AA:818:G:O4'	2.19	0.42
1:AA:886:C:H2'	1:AA:887:A:H5''	2.01	0.42
1:AA:1341:U:C5	1:AA:1395:A:H2	2.36	0.42
1:AA:1343:G:O2'	1:AA:1384:A:N1	2.37	0.42
1:AA:1814:G:H2'	1:AA:1815:A:C8	2.54	0.42
1:AA:2262:U:N3	1:AA:2279:G:C2	2.88	0.42
1:AA:2732:G:H3'	1:AA:2733:A:O4'	2.19	0.42
1:AA:2838:G:C2	1:AA:2881:C:C2	3.08	0.42
3:AD:26:LYS:O	3:AD:83:GLU:HG2	2.19	0.42
6:AG:138:GLN:NE2	6:AG:153:ARG:HB2	2.33	0.42
10:AO:80:ASP:OD2	15:AT:71:GLY:HA3	2.19	0.42
12:AQ:21:THR:HG21	12:AQ:101:ARG:HB2	2.01	0.42
14:AS:26:LEU:HD22	14:AS:87:PHE:HE1	1.82	0.42
15:AT:37:GLY:HA2	15:AT:38:ASN:HA	1.73	0.42
1:BA:185:U:H2'	1:BA:186:G:H8	1.84	0.42
1:BA:271(Q):G:O2'	1:BA:271(R):G:P	2.77	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:272(C):G:H2'	1:BA:272(D):G:O4'	2.19	0.42
1:BA:1047:G:O2'	1:BA:1048:A:P	2.77	0.42
1:BA:1256:G:H5'	1:BA:1257:C:OP2	2.19	0.42
1:BA:1489:U:O2'	1:BA:1490:A:H8	2.03	0.42
1:BA:1929:G:H4'	1:BA:1930:G:OP1	2.20	0.42
1:BA:2336:A:H61	22:B0:43:THR:CG2	2.32	0.42
5:BF:53:THR:HG22	5:BF:55:GLY:H	1.84	0.42
9:BN:22:THR:HB	9:BN:25:ARG:HB2	2.02	0.42
9:BN:125:GLY:HA3	9:BN:126:PRO:HA	1.88	0.42
11:BP:112:LEU:HD22	11:BP:113:LYS:N	2.33	0.42
15:BT:20:PRO:HG2	15:BT:86:ILE:O	2.19	0.42
15:BT:121:ILE:O	15:BT:124:ASP:HB2	2.18	0.42
31:CA:50:A:H2	31:CA:52:G:N3	2.16	0.42
31:CA:256:U:H2'	31:CA:257:G:C8	2.54	0.42
31:CA:834:C:C2	31:CA:853:G:N2	2.87	0.42
31:CA:995:C:H4'	44:CN:8:GLU:OE2	2.19	0.42
31:CA:1130:A:O2'	39:CI:3:GLN:NE2	2.52	0.42
31:CA:1226:C:C4	43:CM:104:ARG:HB2	2.54	0.42
33:CC:51:GLY:O	33:CC:52:LEU:HB3	2.19	0.42
34:CD:155:LEU:HD23	34:CD:156:GLU:N	2.34	0.42
42:CL:60:LEU:HD13	42:CL:60:LEU:HA	1.76	0.42
46:CP:55:ARG:NE	46:CP:55:ARG:HA	2.33	0.42
50:CT:47:GLY:HA2	50:CT:48:LYS:CB	2.49	0.42
31:DA:511:C:HO2'	31:DA:512:U:H6	1.65	0.42
31:DA:1115:C:C2	31:DA:1186:G:C2	3.08	0.42
31:DA:1292:U:C2'	31:DA:1293:G:H5'	2.50	0.42
31:DA:1300:G:O2'	31:DA:1301:U:P	2.78	0.42
31:DA:1408:A:C2	31:DA:1494:G:C2	3.08	0.42
31:DA:1452:C:H3'	31:DA:1456:G:N7	2.34	0.42
31:DA:1478:C:H2'	31:DA:1479:C:H6	1.84	0.42
36:DF:75:LEU:O	36:DF:79:LEU:HG	2.20	0.42
37:DG:108:ALA:O	37:DG:111:ARG:HB2	2.20	0.42
39:DI:44:VAL:HA	39:DI:45:ALA:HB2	2.02	0.42
44:DN:37:PHE:HZ	44:DN:56:VAL:HG21	1.84	0.42
1:AA:817:C:O2'	1:AA:839:U:H5''	2.20	0.42
1:AA:828:U:H4'	1:AA:831:G:N1	2.35	0.42
1:AA:1046:A:HO2'	1:AA:1047:G:P	2.42	0.42
1:AA:2331:G:H4'	22:A0:43:THR:H	1.83	0.42
1:AA:2813:A:C6	1:AA:2814:C:C4	3.07	0.42
1:AA:2855:C:H2'	1:AA:2856:C:H6	1.83	0.42
52:AA:3001:T8B:H10	31:DA:1492:A:H4'	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:87:G:H22	2:AB:89:G:H3'	1.85	0.42
2:AB:90:A:C5	2:AB:91:C:H1'	2.55	0.42
3:AD:159:ALA:HB1	3:AD:198:ASN:HB3	2.01	0.42
5:AF:39:TRP:CE2	5:AF:43:LYS:HD3	2.55	0.42
8:AI:25:TYR:CE2	8:AI:29:TYR:CD2	3.07	0.42
8:AI:29:TYR:C	8:AI:32:PRO:HD2	2.39	0.42
9:AN:96:GLU:H	9:AN:96:GLU:CD	2.23	0.42
14:AS:10:ARG:HH21	14:AS:91:PRO:HB2	1.85	0.42
21:AZ:151:HIS:C	21:AZ:153:SER:N	2.72	0.42
26:A4:42:PHE:CA	26:A4:43:TYR:HB2	2.49	0.42
1:BA:271(K):U:H1'	1:BA:271(L):U:OP1	2.20	0.42
1:BA:1329:U:C5'	1:BA:1330:C:H5	2.32	0.42
1:BA:1341:U:O2	19:BX:80:ILE:HD12	2.19	0.42
1:BA:1503:U:H2'	1:BA:1504:C:H6	1.82	0.42
1:BA:1971:A:C5	3:BD:241:PRO:HD3	2.55	0.42
1:BA:2126:A:H1'	1:BA:2127:G:OP2	2.18	0.42
1:BA:2126:A:N3	1:BA:2162:G:N2	2.68	0.42
1:BA:2223:G:OP1	3:BD:172:TYR:OH	2.24	0.42
1:BA:2290:G:O2'	1:BA:2381:C:H1'	2.19	0.42
1:BA:2413:G:H2'	1:BA:2414:G:O4'	2.19	0.42
1:BA:2418:A:H2'	1:BA:2419:U:O4'	2.19	0.42
1:BA:2638:G:O2'	1:BA:2775:A:N1	2.48	0.42
1:BA:2660:A:H2'	1:BA:2661:G:O4'	2.19	0.42
1:BA:2773:C:H2'	1:BA:2774:C:C6	2.52	0.42
6:BG:137:GLU:HB3	6:BG:138:GLN:H	1.54	0.42
8:BI:60:GLU:C	8:BI:61:ARG:HH11	2.23	0.42
13:BR:54:LEU:HD12	13:BR:54:LEU:HA	1.86	0.42
13:BR:100:LEU:HD11	13:BR:113:LEU:HD23	2.00	0.42
21:BZ:111:VAL:C	21:BZ:113:ALA:N	2.73	0.42
31:CA:226:G:N2	31:CA:227:G:H1'	2.34	0.42
31:CA:509:A:N3	31:CA:543:C:O2'	2.44	0.42
31:CA:1080:A:H5''	35:CE:16:THR:HG21	2.02	0.42
31:CA:1242:C:H4'	31:CA:1304:G:OP1	2.19	0.42
31:CA:1269:A:H1'	31:CA:1326:C:H1'	2.01	0.42
31:CA:1348:U:C4	31:CA:1349:A:N7	2.88	0.42
38:CH:83:ILE:HB	38:CH:137:VAL:HG13	2.01	0.42
49:CS:57:HIS:O	49:CS:59:PRO:HD3	2.20	0.42
31:DA:931:C:H2'	31:DA:932:C:H6	1.84	0.42
31:DA:1479:C:H2'	31:DA:1480:G:C8	2.54	0.42
32:DB:78:GLN:O	32:DB:94:ASN:ND2	2.53	0.42
40:DJ:79:ARG:C	40:DJ:81:THR:H	2.22	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:271(Q):G:O2'	1:AA:271(R):G:P	2.77	0.42
1:AA:278:A:H4'	1:AA:279:C:OP1	2.19	0.42
1:AA:478:A:N1	1:AA:500:G:H4'	2.35	0.42
1:AA:481:G:H1'	1:AA:507:A:N1	2.35	0.42
1:AA:795:C:C2'	1:AA:796:C:H5'	2.49	0.42
1:AA:1252:G:C2	1:AA:1253:A:C2	3.07	0.42
1:AA:1530:C:O2'	1:AA:1531:C:P	2.76	0.42
1:AA:1531:C:N4	1:AA:1538:G:H1	2.14	0.42
1:AA:1804:C:H2'	1:AA:1805:U:H6	1.85	0.42
1:AA:1976:U:H2'	1:AA:1977:A:C8	2.55	0.42
1:AA:2085:C:H2'	1:AA:2086:U:O4'	2.19	0.42
1:AA:2521:C:O2'	1:AA:2564:A:N3	2.45	0.42
1:AA:2575:C:O2'	1:AA:2578:G:N7	2.32	0.42
6:AG:43:LEU:HB3	6:AG:44:GLY:H	1.58	0.42
7:AH:8:PRO:O	7:AH:69:ARG:NH1	2.52	0.42
9:AN:94:HIS:HB3	9:AN:97:ARG:HD3	2.01	0.42
13:AR:72:ASP:O	13:AR:76:VAL:HG23	2.19	0.42
26:A4:14:ILE:HG22	26:A4:33:VAL:HG22	2.01	0.42
1:BA:118:A:N3	1:BA:178:G:H1'	2.34	0.42
1:BA:218:A:N1	1:BA:235:U:H4'	2.34	0.42
1:BA:1653:G:H3'	13:BR:2:ARG:HD3	2.00	0.42
1:BA:2012:G:H8	1:BA:2012:G:O5'	2.03	0.42
1:BA:2029:G:H2'	1:BA:2031:A:OP1	2.19	0.42
1:BA:2114:A:H1'	1:BA:2168:G:H5'	2.00	0.42
1:BA:2837:G:C2	1:BA:2882:A:C2	3.07	0.42
2:BB:55:U:H2'	2:BB:56:G:O4'	2.20	0.42
3:BD:133:LEU:HD23	3:BD:136:ILE:HD12	2.00	0.42
3:BD:210:GLY:O	3:BD:213:ARG:HB2	2.18	0.42
8:BI:25:TYR:CD1	8:BI:30:LEU:HD11	2.54	0.42
14:BS:10:ARG:O	14:BS:14:VAL:HG13	2.18	0.42
20:BY:92:ASN:HB3	20:BY:94:LYS:H	1.84	0.42
31:CA:107:G:H2'	31:CA:108:G:O4'	2.19	0.42
31:CA:112:G:H5'	31:CA:389:A:O2'	2.19	0.42
31:CA:223:U:H2'	31:CA:224:C:H6	1.84	0.42
31:CA:254:G:O2'	47:CQ:16:GLN:O	2.37	0.42
31:CA:327:A:C4	31:CA:329:A:C8	3.07	0.42
31:CA:335:C:H2'	31:CA:336:C:C6	2.54	0.42
31:CA:404:U:H2'	31:CA:405:U:C6	2.53	0.42
31:CA:624:C:H2'	31:CA:625:G:H8	1.84	0.42
31:CA:734:G:C2	31:CA:735:C:C2	3.08	0.42
31:CA:746:A:H4'	31:CA:837:G:O2'	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:1002:G:H2'	31:CA:1003:G:O4'	2.19	0.42
38:CH:21:LYS:O	38:CH:63:LEU:HD23	2.20	0.42
42:CL:27:LEU:C	42:CL:29:GLY:N	2.68	0.42
50:CT:63:ILE:HD13	50:CT:80:ARG:HB2	2.01	0.42
31:DA:101:A:H2'	31:DA:102:G:H8	1.84	0.42
31:DA:113:G:H2'	31:DA:114:U:C6	2.55	0.42
31:DA:269:C:H2'	31:DA:270:A:C8	2.55	0.42
31:DA:875:C:O2'	38:DH:14:ARG:HD2	2.19	0.42
31:DA:935:A:N6	37:DG:3:ARG:HG3	2.34	0.42
32:DB:87:ARG:HH11	32:DB:219:VAL:HG12	1.85	0.42
32:DB:187:LEU:HD23	32:DB:201:ILE:HG22	2.01	0.42
38:DH:11:THR:HG22	38:DH:15:ASN:ND2	2.34	0.42
38:DH:78:GLN:HG2	38:DH:80:ILE:O	2.20	0.42
38:DH:87:SER:HB2	38:DH:93:VAL:HB	2.00	0.42
43:DM:88:ARG:HG3	43:DM:98:VAL:HG13	2.01	0.42
50:DT:54:LYS:HA	50:DT:57:ARG:NH1	2.34	0.42
1:AA:528:A:N1	1:AA:2042:A:H2'	2.35	0.42
1:AA:598:G:C6	1:AA:599:G:C5	3.07	0.42
1:AA:634:C:H6	1:AA:634:C:O5'	2.03	0.42
1:AA:708:C:N4	1:AA:723:G:H1	2.17	0.42
1:AA:867:C:O2	1:AA:913:U:H5'	2.19	0.42
1:AA:900:A:H2'	1:AA:901:A:C8	2.54	0.42
1:AA:952:G:C6	1:AA:953:A:N7	2.88	0.42
1:AA:953:A:C2	1:AA:954:G:C8	3.07	0.42
1:AA:1005:C:O2	1:AA:1143:A:C6	2.72	0.42
1:AA:1016:G:C5	1:AA:1017:G:N7	2.87	0.42
1:AA:1797:C:H4'	3:AD:257:LEU:O	2.19	0.42
1:AA:2787:C:H2'	1:AA:2788:C:C6	2.55	0.42
2:AB:22:U:H3	2:AB:61:G:H1	1.67	0.42
4:AE:77:ILE:HG21	4:AE:195:LEU:HD11	2.01	0.42
6:AG:98:ARG:O	6:AG:102:PHE:N	2.48	0.42
9:AN:103:VAL:O	9:AN:107:LEU:HG	2.20	0.42
25:A3:10:LYS:O	25:A3:53:LEU:HB3	2.19	0.42
1:BA:945:A:C4	1:BA:2448:A:C2	3.07	0.42
1:BA:994:C:OP2	16:BU:54:LYS:NZ	2.45	0.42
1:BA:1426:G:H8	1:BA:1426:G:O5'	2.03	0.42
1:BA:2117:A:N6	1:BA:2171:A:N6	2.68	0.42
1:BA:2136:C:N3	1:BA:2155:G:N2	2.66	0.42
1:BA:2872:G:O2'	1:BA:2873:A:H5'	2.19	0.42
5:BF:123:LEU:HD12	5:BF:124:LEU:H	1.83	0.42
7:BH:8:PRO:O	7:BH:69:ARG:NH1	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:BN:39:ARG:HA	9:BN:40:PRO:HD3	1.90	0.42
12:BQ:18:LYS:HB2	12:BQ:18:LYS:HE3	1.75	0.42
14:BS:90:GLY:C	14:BS:92:TYR:H	2.23	0.42
19:BX:5:TYR:CE1	24:B2:30:ARG:HB2	2.54	0.42
21:BZ:152:ALA:O	21:BZ:155:LEU:HB2	2.19	0.42
31:CA:149:A:HO2'	31:CA:150:C:H6	1.67	0.42
31:CA:397:A:N6	31:CA:548:G:N7	2.68	0.42
31:CA:803:G:H2'	31:CA:804:U:O4'	2.19	0.42
31:CA:1240:U:H4'	31:CA:1241:G:OP2	2.20	0.42
31:CA:1522:U:H2'	31:CA:1523:G:H8	1.85	0.42
34:CD:58:LEU:O	34:CD:62:GLN:HG2	2.19	0.42
35:CE:68:GLU:HG2	35:CE:70:PRO:HG3	2.01	0.42
44:CN:50:LYS:HE2	44:CN:52:GLN:NE2	2.35	0.42
46:CP:39:TYR:CE2	46:CP:41:PRO:HB3	2.53	0.42
50:CT:14:LYS:HA	50:CT:17:ARG:CZ	2.49	0.42
31:DA:458:C:C2'	31:DA:460:G:H8	2.30	0.42
31:DA:559:A:H4'	31:DA:560:U:H3'	2.00	0.42
31:DA:945:G:C2	31:DA:946:A:C8	3.08	0.42
31:DA:981:U:O5'	31:DA:981:U:H6	2.03	0.42
31:DA:1277:C:O2'	31:DA:1279:A:H8	2.01	0.42
31:DA:1471:G:O2'	31:DA:1472:U:H5'	2.20	0.42
31:DA:1514:C:H2'	31:DA:1515:C:C6	2.54	0.42
34:DD:30:LYS:HA	34:DD:35:ARG:HD2	2.01	0.42
36:DF:10:LEU:HD21	36:DF:61:LEU:HD22	2.02	0.42
43:DM:52:GLU:HA	43:DM:55:ARG:HB3	2.01	0.42
44:DN:13:THR:HA	44:DN:14:PRO:HD3	1.86	0.42
1:AA:17:G:C4	1:AA:18:C:C5	3.08	0.42
1:AA:171:G:H2'	1:AA:172:C:C6	2.55	0.42
1:AA:829:A:N7	1:AA:2248:C:H5'	2.34	0.42
1:AA:918:A:O2'	2:AB:97:G:N2	2.51	0.42
1:AA:968:G:H2'	1:AA:969:U:O4'	2.20	0.42
1:AA:1159:U:H2'	1:AA:1160:G:C8	2.55	0.42
1:AA:1761:C:H2'	1:AA:1762:A:H5''	2.01	0.42
1:AA:1819:A:H3'	3:AD:178:PRO:HB2	2.02	0.42
1:AA:2168:G:H2'	1:AA:2170:A:OP2	2.20	0.42
1:AA:2190:G:H2'	1:AA:2191:G:H5''	2.01	0.42
1:AA:2256:G:H2'	1:AA:2257:U:C6	2.54	0.42
3:AD:5:LYS:HE3	3:AD:5:LYS:HB3	1.82	0.42
3:AD:16:MET:CG	3:AD:211:ARG:HH21	2.32	0.42
3:AD:159:ALA:HB1	3:AD:198:ASN:O	2.18	0.42
3:AD:221:VAL:HG22	3:AD:226:MET:HE3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AN:15:LEU:HD13	9:AN:16:ILE:N	2.34	0.42
15:AT:6:LEU:O	15:AT:9:LEU:HB3	2.20	0.42
25:A3:7:LYS:O	25:A3:54:VAL:HA	2.20	0.42
1:BA:143:G:H2'	1:BA:143(A):C:C6	2.54	0.42
1:BA:226:G:H21	1:BA:228:A:H62	1.67	0.42
1:BA:478:A:N1	1:BA:500:G:H4'	2.35	0.42
1:BA:1432:C:H2'	1:BA:1433:U:O4'	2.19	0.42
1:BA:1913:A:C2'	52:BA:3001:T8B:C23	2.95	0.42
1:BA:1913:A:C2	52:BA:3001:T8B:O8	2.73	0.42
1:BA:2131:G:H5''	1:BA:2132:U:C5'	2.49	0.42
1:BA:2262:U:O2'	1:BA:2263:C:H5'	2.19	0.42
1:BA:2563:U:H4'	10:BO:28:SER:HA	2.02	0.42
4:BE:6:GLY:HA2	4:BE:51:PHE:CE1	2.55	0.42
8:BI:145:VAL:HG12	8:BI:146:ALA:H	1.85	0.42
22:B0:53:MET:HA	22:B0:58:THR:O	2.20	0.42
24:B2:17:SER:OG	24:B2:20:GLU:HG3	2.19	0.42
25:B3:31:LEU:C	25:B3:33:GLN:H	2.23	0.42
30:B8:35:GLN:O	30:B8:35:GLN:HG3	2.20	0.42
31:CA:8:A:H5'	35:CE:101:ILE:HG22	2.00	0.42
31:CA:475:G:C2'	31:CA:476:G:H5'	2.49	0.42
31:CA:521:G:O5'	42:CL:73:GLU:HG2	2.20	0.42
31:CA:581:G:N2	31:CA:582:U:C4	2.87	0.42
31:CA:981:U:H5	31:CA:982:U:HO2'	1.66	0.42
31:CA:1030:C:N3	31:CA:1031:G:C2	2.88	0.42
31:CA:1277:C:H1'	31:CA:1282:C:C2	2.55	0.42
31:CA:1329:A:N7	51:CU:7:ARG:NH2	2.59	0.42
31:CA:1415:G:C4	31:CA:1486:G:C2	3.07	0.42
31:DA:59:A:H3'	31:DA:331:G:H22	1.84	0.42
31:DA:142:G:H2'	31:DA:143:A:C8	2.54	0.42
31:DA:667:G:H4'	45:DO:51:HIS:CE1	2.54	0.42
31:DA:841:U:OP1	31:DA:841:U:H6	2.03	0.42
31:DA:935:A:H61	37:DG:3:ARG:HG3	1.84	0.42
31:DA:1480:G:H2'	31:DA:1481:U:O4'	2.19	0.42
35:DE:126:ARG:HH11	35:DE:126:ARG:CG	2.31	0.42
42:DL:46:LYS:HE2	42:DL:47:LYS:O	2.20	0.42
49:DS:36:ARG:H	49:DS:36:ARG:HG2	1.74	0.42
1:AA:330:A:C2	1:AA:1210:A:H2'	2.46	0.42
1:AA:571:A:N6	1:AA:2499:C:O3'	2.49	0.42
1:AA:797:C:H2'	1:AA:798:G:H8	1.84	0.42
1:AA:911:A:H2'	12:AQ:9:TYR:CZ	2.54	0.42
1:AA:1833:U:H2'	1:AA:1834:U:H6	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2163:C:OP2	1:AA:2164:C:H5	2.03	0.42
1:AA:2206:G:H3'	1:AA:2207:G:N7	2.34	0.42
1:AA:2263:C:H42	1:AA:2278:A:N6	2.18	0.42
1:AA:2345:G:OP2	28:A6:38:LYS:HG3	2.20	0.42
1:AA:2735:G:N3	1:AA:2736:G:C8	2.88	0.42
52:AA:3001:T8B:O5	31:DA:1491:G:H2'	2.20	0.42
4:AE:143:ASN:ND2	4:AE:147:PRO:HD2	2.18	0.42
6:AG:128:ARG:HE	6:AG:128:ARG:HB2	1.31	0.42
8:AI:3:VAL:HG12	8:AI:38:LEU:HA	2.01	0.42
10:AO:26:LYS:HD2	10:AO:37:ASP:OD1	2.20	0.42
13:AR:44:LEU:HD23	13:AR:44:LEU:HA	1.92	0.42
21:AZ:126:VAL:HG13	21:AZ:127:LYS:O	2.19	0.42
28:A6:15:GLU:HG3	28:A6:47:THR:HG23	2.02	0.42
1:BA:195:A:OP1	11:BP:46:LYS:HE2	2.20	0.42
1:BA:312:G:H4'	1:BA:331:A:N3	2.34	0.42
1:BA:405:U:H4'	1:BA:406:G:OP2	2.20	0.42
1:BA:528:A:C2	1:BA:2042:A:H2'	2.54	0.42
1:BA:631:A:H1'	11:BP:66:GLY:HA2	2.02	0.42
1:BA:1022:G:H22	1:BA:1142(A):A:H2	1.60	0.42
1:BA:1169:G:N2	1:BA:1181:C:C2	2.88	0.42
1:BA:1815:A:P	3:BD:54:ARG:NH2	2.92	0.42
1:BA:2285:C:C5	28:B6:6:ARG:NH2	2.88	0.42
1:BA:2328:A:H2'	1:BA:2329:G:H8	1.82	0.42
1:BA:2360:A:H8	1:BA:2360:A:O5'	2.02	0.42
1:BA:2516:G:C5	1:BA:2517:C:C4	3.07	0.42
1:BA:2889:C:H2'	1:BA:2891:G:O4'	2.20	0.42
2:BB:29:A:H2'	2:BB:30:C:O4'	2.20	0.42
6:BG:25:TYR:HB3	6:BG:30:GLU:HB2	2.00	0.42
9:BN:56:ASN:CA	9:BN:125:GLY:H	2.29	0.42
11:BP:19:VAL:HG23	11:BP:31:ALA:HB1	2.02	0.42
19:BX:31:HIS:HA	19:BX:32:PRO:HD3	1.86	0.42
31:CA:93:G:HO2'	31:CA:96:U:P	2.37	0.42
31:CA:293:G:H2'	31:CA:294:U:H6	1.84	0.42
31:CA:979:C:C2	44:CN:19:ARG:HG2	2.55	0.42
31:CA:1256:A:N6	31:CA:1278:U:O4'	2.52	0.42
31:CA:1307:U:H2'	31:CA:1308:U:C6	2.55	0.42
31:CA:1416:G:C5	31:CA:1417:G:C5	3.08	0.42
31:CA:1443:G:N2	31:CA:1460:A:H1'	2.35	0.42
36:CF:14:LEU:HB3	36:CF:15:ASP:H	1.49	0.42
47:CQ:87:LYS:HD3	47:CQ:87:LYS:HA	1.60	0.42
31:DA:721:G:H8	31:DA:721:G:OP1	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:949:A:H1'	31:DA:1364:U:C2	2.54	0.42
34:DD:119:GLN:O	34:DD:123:HIS:CD2	2.73	0.42
34:DD:120:LEU:HD23	34:DD:120:LEU:HA	1.87	0.42
35:DE:93:PRO:HG2	38:DH:105:ARG:CZ	2.50	0.42
40:DJ:47:PHE:CZ	44:DN:37:PHE:HE2	2.38	0.42
42:DL:70:ILE:HG12	42:DL:100:ILE:HD12	2.02	0.42
45:DO:3:ILE:HA	45:DO:7:GLU:OE2	2.20	0.42
46:DP:28:ARG:HG2	46:DP:29:ASP:OD2	2.20	0.42
50:DT:22:ARG:O	50:DT:26:ASN:N	2.49	0.42
1:AA:140:G:N3	1:AA:142:A:N6	2.64	0.42
1:AA:481:G:C4	1:AA:507:A:C2	3.08	0.42
1:AA:583:G:C5	1:AA:584:C:C5	3.08	0.42
1:AA:1666:G:C2'	1:AA:1667:G:H5'	2.50	0.42
1:AA:2019:A:C6	1:AA:2020:A:N7	2.88	0.42
1:AA:2036:C:H5'	1:AA:2036:C:C6	2.53	0.42
1:AA:2173:A:C2'	1:AA:2174:C:H5'	2.50	0.42
1:AA:2294:C:P	14:AS:89:ARG:HH22	2.42	0.42
1:AA:2519:U:C2	1:AA:2542:A:C6	3.07	0.42
1:AA:2745:C:H2'	1:AA:2746:U:O4'	2.20	0.42
52:AA:3001:T8B:O4	31:DA:1492:A:C5'	2.65	0.42
3:AD:238:GLY:O	3:AD:239:ARG:HB2	2.20	0.42
6:AG:106:LEU:HD12	6:AG:110:ALA:HB3	2.01	0.42
6:AG:150:ASP:CG	6:AG:151:ALA:H	2.23	0.42
7:AH:83:TYR:CE2	7:AH:138:LYS:HB2	2.54	0.42
7:AH:149:ARG:HA	7:AH:162:ILE:HG21	2.00	0.42
14:AS:34:HIS:HB3	14:AS:35:ILE:H	1.55	0.42
14:AS:35:ILE:CD1	14:AS:101:LEU:HD12	2.50	0.42
18:AW:38:TYR:O	27:A5:28:PRO:HB3	2.20	0.42
21:AZ:65:GLN:OE1	21:AZ:67:LEU:HD21	2.20	0.42
24:A2:51:ARG:HA	24:A2:54:LYS:HB2	2.01	0.42
25:A3:8:LEU:HA	25:A3:8:LEU:HD23	1.72	0.42
1:BA:524:U:H4'	1:BA:555:U:H4'	2.02	0.42
1:BA:817:C:O2'	1:BA:839:U:H5''	2.20	0.42
1:BA:1262:A:C6	1:BA:1263:U:C4	3.08	0.42
1:BA:1793:C:H2'	1:BA:1794:U:H6	1.83	0.42
2:BB:111:G:H2'	2:BB:112:U:C6	2.54	0.42
6:BG:7:LEU:HD22	6:BG:7:LEU:HA	1.76	0.42
8:BI:9:LEU:HD21	8:BI:35:LEU:HD22	2.01	0.42
8:BI:108:THR:O	8:BI:108:THR:OG1	2.29	0.42
8:BI:123:LEU:H	8:BI:123:LEU:HG	1.64	0.42
20:BY:92:ASN:N	20:BY:93:GLY:HA2	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:B1:82:LEU:HA	23:B1:85:LEU:HD23	2.02	0.42
31:CA:20:U:C2'	31:CA:21:G:H5'	2.50	0.42
32:CB:215:LEU:HD23	32:CB:215:LEU:HA	1.70	0.42
34:CD:13:ARG:HD2	34:CD:38:TYR:O	2.19	0.42
34:CD:18:LYS:HE3	34:CD:26:CYS:O	2.20	0.42
35:CE:122:GLU:O	35:CE:123:LEU:HD23	2.19	0.42
44:CN:2:ALA:HB1	44:CN:6:LEU:HD13	2.01	0.42
45:CO:4:THR:OG1	45:CO:7:GLU:HG3	2.20	0.42
45:CO:74:ASP:OD2	45:CO:77:ARG:HB2	2.20	0.42
31:DA:992:U:O2'	31:DA:993:G:O4'	2.38	0.42
31:DA:1010:G:H22	31:DA:1020:U:H1'	1.84	0.42
31:DA:1401:G:H2'	31:DA:1402:C:O4'	2.19	0.42
34:DD:13:ARG:HD2	34:DD:38:TYR:O	2.20	0.42
35:DE:16:THR:OG1	35:DE:17:ALA:N	2.52	0.42
35:DE:139:LEU:HA	35:DE:142:LEU:HD12	2.02	0.42
38:DH:36:LEU:HD12	38:DH:59:LEU:HD13	2.01	0.42
1:AA:30:G:H2'	1:AA:31:C:H6	1.81	0.42
1:AA:271(C):C:C2	1:AA:271(V):G:C2	3.07	0.42
1:AA:458:G:O2'	29:A7:39:ARG:HD3	2.20	0.42
1:AA:647:G:H2'	1:AA:648:G:O4'	2.20	0.42
1:AA:734:A:O2'	1:AA:1635:G:H5'	2.20	0.42
1:AA:921:G:C5	1:AA:922:U:C4	3.08	0.42
1:AA:1418:G:OP1	1:AA:1588:C:O2'	2.32	0.42
1:AA:1434:A:H61	1:AA:1558:A:H62	1.67	0.42
1:AA:1690:A:H3'	1:AA:1691:C:H6	1.84	0.42
1:AA:1999:C:H5''	1:AA:2723:C:O2'	2.20	0.42
1:AA:2136:C:C5	1:AA:2137:C:H5	2.38	0.42
1:AA:2302:G:N1	1:AA:2315:G:C6	2.88	0.42
1:AA:2403:C:N3	1:AA:2415:G:C2	2.88	0.42
1:AA:2833:G:H3'	1:AA:2834:G:H5''	2.02	0.42
6:AG:15:VAL:HG22	6:AG:175:LEU:O	2.20	0.42
7:AH:167:GLU:HA	7:AH:168:PRO:HD3	1.90	0.42
10:AO:118:ALA:HA	10:AO:119:PRO:HD2	1.82	0.42
13:AR:26:LYS:HE2	13:AR:70:LEU:O	2.20	0.42
14:AS:99:LYS:HE3	14:AS:103:GLU:OE2	2.20	0.42
14:AS:103:GLU:O	14:AS:107:GLU:HG3	2.19	0.42
16:AU:21:ALA:HA	16:AU:24:TYR:CE1	2.55	0.42
20:AY:7:VAL:HG21	20:AY:72:VAL:HG12	2.02	0.42
25:A3:4:LEU:N	25:A3:37:LEU:O	2.46	0.42
1:BA:65:C:H2'	1:BA:66:C:C6	2.52	0.42
1:BA:171:G:H2'	1:BA:172:C:C6	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:311:A:C6	1:BA:328:U:C4	3.08	0.42
1:BA:312:G:H5'	1:BA:331:A:O2'	2.20	0.42
1:BA:612:C:H2'	1:BA:613:G:O4'	2.20	0.42
1:BA:652(F):G:H2'	1:BA:652(G):G:H8	1.85	0.42
1:BA:727:A:C6	1:BA:728:G:C6	3.07	0.42
1:BA:839:U:H1'	1:BA:1191:G:H1'	2.02	0.42
1:BA:1025:G:C4	1:BA:1135:C:H1'	2.55	0.42
1:BA:1141:U:C5	9:BN:64:GLY:HA3	2.55	0.42
1:BA:1202:C:N4	1:BA:1203:G:C6	2.88	0.42
1:BA:1289:C:H2'	1:BA:1290:C:C6	2.54	0.42
1:BA:1509(B):A:H2'	1:BA:1510:G:C8	2.54	0.42
1:BA:1538:G:H8	1:BA:1538:G:O5'	2.03	0.42
1:BA:2101:G:H2'	1:BA:2102:U:O4'	2.20	0.42
1:BA:2679:A:H2'	1:BA:2680:C:O4'	2.20	0.42
52:BA:3001:T8B:C12	31:CA:1491:G:C5	3.03	0.42
2:BB:60:C:C2	2:BB:61:G:C8	3.08	0.42
4:BE:72:VAL:CA	4:BE:73:GLU:HB3	2.49	0.42
8:BI:123:LEU:HA	8:BI:144:VAL:HG13	2.02	0.42
9:BN:112:LEU:HD12	9:BN:112:LEU:O	2.20	0.42
23:B1:73:LEU:O	23:B1:77:ALA:N	2.53	0.42
31:CA:38:G:C2	31:CA:397:A:C2	3.07	0.42
31:CA:741:G:H2'	31:CA:742:G:C8	2.55	0.42
31:CA:985:C:H2'	31:CA:986:A:C8	2.55	0.42
31:CA:1146:A:N6	31:CA:1147:C:C4	2.88	0.42
31:CA:1244:C:H42	31:CA:1293:G:H1	1.68	0.42
31:CA:1323:G:H2'	31:CA:1324:A:C8	2.55	0.42
31:CA:1373:G:H5''	37:CG:36:LYS:HB2	2.02	0.42
31:CA:1443:G:C2	31:CA:1460:A:N3	2.88	0.42
31:CA:1493:A:N3	31:CA:1493:A:C2'	2.83	0.42
32:CB:28:PHE:CZ	32:CB:189:ASP:HA	2.55	0.42
32:CB:162:ILE:HD11	32:CB:184:VAL:HG22	2.02	0.42
32:CB:194:PRO:HB2	32:CB:195:ASP:H	1.54	0.42
41:CK:86:GLY:N	41:CK:112:THR:OG1	2.28	0.42
46:CP:8:ARG:HB3	46:CP:28:ARG:NH1	2.35	0.42
31:DA:123:C:OP1	31:DA:311:C:O2'	2.28	0.42
31:DA:224:C:H2'	31:DA:225:C:H6	1.84	0.42
31:DA:339:C:H2'	31:DA:340:U:H6	1.83	0.42
31:DA:355:C:O4'	31:DA:388:G:O2'	2.33	0.42
31:DA:538:G:H2'	31:DA:539:A:C8	2.55	0.42
31:DA:1086:U:H2'	31:DA:1087:G:H8	1.85	0.42
31:DA:1186:G:O3'	39:DI:113:LYS:NZ	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DD:36:ARG:HG2	34:DD:36:ARG:NH1	2.34	0.42
36:DF:12:PRO:C	36:DF:14:LEU:H	2.22	0.42
38:DH:121:ASP:HB2	38:DH:125:ARG:NH2	2.35	0.42
43:DM:37:THR:O	43:DM:55:ARG:HD3	2.20	0.42
1:AA:88:G:C2	1:AA:89:G:C4	3.07	0.41
1:AA:470:A:H2'	1:AA:471:A:O4'	2.20	0.41
1:AA:775:G:H22	1:AA:794:G:H5'	1.86	0.41
1:AA:815:C:H2'	1:AA:816:C:H6	1.84	0.41
1:AA:1015:G:C6	1:AA:1148:A:C6	3.08	0.41
1:AA:1485:G:C2	1:AA:1486:A:C5	3.08	0.41
1:AA:1973:G:H2'	1:AA:1974:C:C6	2.55	0.41
1:AA:2127:G:N2	1:AA:2162:G:H1'	2.34	0.41
1:AA:2577:A:OP2	27:A5:3:LYS:NZ	2.49	0.41
3:AD:68:LYS:O	3:AD:70:TRP:CD1	2.73	0.41
5:AF:160:ASN:CG	5:AF:163:VAL:HG23	2.40	0.41
14:AS:25:ARG:HG3	14:AS:88:ASP:HB2	2.01	0.41
16:AU:58:ARG:HA	16:AU:61:TRP:CE3	2.54	0.41
19:AX:50:LYS:CG	19:AX:84:ALA:HB2	2.50	0.41
1:BA:139(A):G:O2'	1:BA:140:G:H5'	2.20	0.41
1:BA:245:G:O2'	1:BA:384:U:O2	2.27	0.41
1:BA:1578:U:H2'	1:BA:1579:A:H5'	2.02	0.41
1:BA:2833:G:O2'	1:BA:2834:G:P	2.78	0.41
52:BA:3001:T8B:C12	31:CA:1491:G:N7	2.83	0.41
23:B1:55:GLY:O	23:B1:56:GLN:HG2	2.19	0.41
31:CA:947:G:N2	31:CA:1235:U:H1'	2.35	0.41
31:CA:1095:U:H5''	31:CA:1109:C:O2	2.20	0.41
31:CA:1525:G:OP1	41:CK:120:ARG:NH2	2.53	0.41
34:CD:59:ARG:HH21	34:CD:66:ARG:HH12	1.67	0.41
35:CE:24:ARG:HG3	35:CE:26:PHE:CZ	2.55	0.41
42:CL:90:VAL:O	42:CL:92:ASP:N	2.53	0.41
43:CM:60:VAL:HG13	43:CM:64:TRP:CE3	2.55	0.41
46:CP:17:TYR:HE1	46:CP:41:PRO:HG3	1.85	0.41
31:DA:10:A:H2'	31:DA:11:G:C8	2.55	0.41
31:DA:149:A:HO2'	31:DA:150:C:H6	1.65	0.41
31:DA:474:G:C2	31:DA:475:G:C5	3.08	0.41
31:DA:590:C:OP1	38:DH:29:SER:HA	2.20	0.41
31:DA:602:A:C2	31:DA:603:U:C2	3.08	0.41
31:DA:757:U:H2'	31:DA:758:G:O4'	2.20	0.41
31:DA:904:C:C4	31:DA:905:U:C4	3.08	0.41
31:DA:927:G:N2	31:DA:1391:U:H1'	2.35	0.41
31:DA:1015:A:O2'	31:DA:1219:U:H5'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:1073:U:H2'	31:DA:1074:G:C8	2.43	0.41
31:DA:1202:G:C4	44:DN:42:ILE:HD12	2.54	0.41
31:DA:1262:C:H2'	31:DA:1263:C:H6	1.83	0.41
31:DA:1496:C:H2'	31:DA:1497:G:C1'	2.50	0.41
32:DB:17:PHE:CD2	32:DB:17:PHE:N	2.82	0.41
32:DB:155:LEU:HD21	32:DB:159:PRO:HG3	2.01	0.41
34:DD:108:LEU:HB3	34:DD:110:PHE:CD1	2.55	0.41
1:AA:117:G:C6	1:AA:119:A:C6	3.08	0.41
1:AA:177:G:N3	1:AA:177:G:H2'	2.35	0.41
1:AA:243:U:C2'	1:AA:244:A:H5'	2.50	0.41
1:AA:323:G:H1'	1:AA:1205:U:O2	2.20	0.41
1:AA:489:G:N7	18:AW:49:LYS:NZ	2.67	0.41
1:AA:588:U:H2'	1:AA:589:C:C6	2.55	0.41
1:AA:679:C:O2'	1:AA:680:G:H5'	2.20	0.41
1:AA:776:G:H4'	1:AA:777:A:O5'	2.20	0.41
1:AA:1174:A:N3	1:AA:1175:U:H5''	2.35	0.41
1:AA:1422:G:C4	1:AA:1423:G:C8	3.09	0.41
1:AA:1531:C:N4	1:AA:1532:C:N4	2.68	0.41
1:AA:2263:C:O2'	1:AA:2264:C:H5'	2.19	0.41
1:AA:2319:G:C8	1:AA:2320:A:C2	3.09	0.41
1:AA:2428:G:H5''	1:AA:2429:G:O5'	2.19	0.41
1:AA:2489:G:C6	1:AA:2490:G:C6	3.08	0.41
2:AB:13:A:H2'	2:AB:70:C:O2'	2.19	0.41
7:AH:70:THR:O	7:AH:72:ILE:N	2.53	0.41
16:AU:14:HIS:CD2	16:AU:32:PHE:CE2	3.08	0.41
21:AZ:19:ARG:HE	21:AZ:19:ARG:HB2	1.48	0.41
21:AZ:100:VAL:N	21:AZ:124:ILE:O	2.37	0.41
1:BA:501:A:H8	1:BA:501:A:O5'	2.03	0.41
1:BA:554:U:O2'	1:BA:555:U:H5'	2.21	0.41
1:BA:1223:G:N2	1:BA:1226:A:OP2	2.43	0.41
1:BA:2183:C:O2'	1:BA:2184:G:H5'	2.20	0.41
1:BA:2208:A:N3	1:BA:2219:G:N1	2.68	0.41
2:BB:2:C:H2'	2:BB:3:C:H6	1.85	0.41
4:BE:85:ASN:HA	4:BE:86:PRO:HD2	1.83	0.41
4:BE:111:ARG:HG2	4:BE:160:TYR:O	2.20	0.41
6:BG:103:LEU:HD23	6:BG:103:LEU:HA	1.94	0.41
7:BH:20:ALA:HB3	7:BH:23:ARG:HB3	2.02	0.41
8:BI:111:PRO:O	8:BI:130:TYR:HE1	2.03	0.41
10:BO:112:MET:H	10:BO:112:MET:HG2	1.64	0.41
11:BP:121:LYS:HB3	11:BP:123:LEU:HG	2.01	0.41
12:BQ:37:LEU:HD23	12:BQ:37:LEU:HA	1.71	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:BW:65:LEU:HD12	18:BW:68:ARG:HE	1.85	0.41
21:BZ:68:PRO:O	21:BZ:91:LEU:HB2	2.19	0.41
23:B1:5:CYS:SG	23:B1:62:VAL:HG23	2.60	0.41
31:CA:97:G:O2'	31:CA:98:G:H5''	2.20	0.41
31:CA:242:C:H2'	31:CA:243:A:H5'	2.01	0.41
31:CA:416:G:H2'	31:CA:417:C:C6	2.54	0.41
31:CA:604:G:C2	31:CA:635:G:C5	3.07	0.41
31:CA:634:C:H2'	31:CA:635:G:H8	1.85	0.41
31:CA:1064:G:H5'	31:CA:1066:C:O4'	2.20	0.41
33:CC:129:ALA:HB3	33:CC:132:ARG:HB3	2.01	0.41
38:CH:24:THR:HG22	38:CH:63:LEU:HD21	2.02	0.41
39:CI:4:TYR:CE1	39:CI:88:TYR:HA	2.54	0.41
31:DA:427:U:P	34:DD:13:ARG:HH22	2.43	0.41
31:DA:553:A:H2'	31:DA:554:C:C6	2.55	0.41
31:DA:834:C:C4	31:DA:835:U:C4	3.08	0.41
31:DA:1057:G:C4	31:DA:1204:A:C2	3.08	0.41
31:DA:1364:U:O2'	31:DA:1365:G:H5'	2.20	0.41
1:AA:479:A:H4'	1:AA:480:A:OP1	2.20	0.41
1:AA:1214:A:N7	1:AA:1215:G:N7	2.68	0.41
1:AA:2226:C:H2'	1:AA:2227:A:O4'	2.20	0.41
1:AA:2572:A:OP1	1:AA:2574:G:O2'	2.29	0.41
1:AA:2640:G:OP1	9:AN:97:ARG:NH2	2.53	0.41
1:AA:2788:C:N4	1:AA:2789:C:N4	2.67	0.41
4:AE:201:THR:C	4:AE:202:LYS:HD2	2.41	0.41
7:AH:9:ILE:HA	7:AH:10:PRO:HD2	1.86	0.41
10:AO:70:LYS:HE2	10:AO:70:LYS:HB3	1.62	0.41
11:AP:132:LYS:HE2	11:AP:132:LYS:HB2	1.87	0.41
12:AQ:46:GLN:NE2	12:AQ:126:PRO:HG3	2.35	0.41
15:AT:11:GLU:OE1	15:AT:57:PHE:HB3	2.20	0.41
16:AU:90:VAL:O	16:AU:95:LEU:HD13	2.20	0.41
17:AV:1:MET:HA	17:AV:41:GLY:O	2.20	0.41
1:BA:219:G:H2'	1:BA:220:G:O4'	2.20	0.41
1:BA:468:G:C6	1:BA:469:G:C4	3.08	0.41
1:BA:672:C:O2'	1:BA:673:C:H5'	2.20	0.41
1:BA:768:G:C6	1:BA:769:G:C5	3.08	0.41
1:BA:2133:G:N2	1:BA:2157:G:H2'	2.35	0.41
1:BA:2218:U:H6	1:BA:2218:U:H2'	1.32	0.41
1:BA:2291:U:H2'	1:BA:2292:C:C6	2.55	0.41
3:BD:221:VAL:HG22	3:BD:226:MET:CE	2.50	0.41
3:BD:222:ARG:NH1	3:BD:224:ALA:HB3	2.35	0.41
14:BS:41:ASP:OD1	14:BS:43:GLU:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:BU:59:ARG:HB3	16:BU:59:ARG:HH11	1.84	0.41
21:BZ:111:VAL:O	21:BZ:113:ALA:N	2.44	0.41
23:B1:51:VAL:HG11	23:B1:74:VAL:HG21	2.01	0.41
26:B4:40:HIS:HB3	26:B4:43:TYR:CD1	2.55	0.41
31:CA:119:A:C6	31:CA:240:C:C2	3.08	0.41
31:CA:152:A:N7	31:CA:153:C:C4	2.88	0.41
31:CA:377:G:C2	31:CA:387:U:O2	2.73	0.41
31:CA:836:G:C6	31:CA:851:G:C6	3.08	0.41
31:CA:933:G:C6	31:CA:1385:G:N1	2.88	0.41
31:CA:1036:G:H5'	31:CA:1037:C:OP2	2.19	0.41
31:CA:1055:A:O2'	33:CC:156:ARG:NE	2.52	0.41
31:CA:1147:C:C4	31:CA:1148:U:C4	3.08	0.41
31:CA:1318:A:H62	44:CN:18:VAL:HG11	1.84	0.41
31:CA:1347:G:N1	31:CA:1374:A:OP2	2.45	0.41
48:CR:33:ASP:OD1	48:CR:36:ASN:HB2	2.19	0.41
31:DA:221:C:H2'	31:DA:222:U:H6	1.85	0.41
31:DA:410:G:H5''	31:DA:411:A:OP1	2.21	0.41
31:DA:432:A:H3'	31:DA:433:C:C6	2.55	0.41
33:DC:32:LEU:HB3	33:DC:59:ARG:NH1	2.35	0.41
35:DE:43:LEU:HD21	35:DE:132:ALA:HB1	2.01	0.41
43:DM:43:THR:HB	43:DM:47:ASP:O	2.21	0.41
1:AA:191:A:C2	1:AA:192:C:C2	3.08	0.41
1:AA:329:G:H4'	1:AA:330:A:OP2	2.20	0.41
1:AA:536:A:H2'	1:AA:537:C:C6	2.56	0.41
1:AA:671:C:H2'	1:AA:672:C:C6	2.54	0.41
1:AA:947:G:N3	1:AA:984:A:H2	2.18	0.41
1:AA:1021:A:H61	1:AA:1142(A):A:H61	1.68	0.41
1:AA:1047:G:O2'	1:AA:1048:A:O5'	2.35	0.41
1:AA:1252:G:H8	1:AA:1252:G:OP1	2.04	0.41
1:AA:1374:G:C2	1:AA:1375:C:C2	3.08	0.41
1:AA:1496:A:N3	1:AA:1577:C:O2'	2.38	0.41
1:AA:2744:G:N2	7:AH:143:GLN:OE1	2.52	0.41
1:AA:2820:A:C6	13:AR:4:LEU:HD11	2.55	0.41
8:AI:47:LEU:O	8:AI:51:ILE:HG12	2.21	0.41
9:AN:19:GLU:HA	9:AN:59:LYS:HB2	2.02	0.41
9:AN:73:THR:HA	9:AN:83:LYS:O	2.21	0.41
9:AN:104:LYS:HA	9:AN:107:LEU:HD12	2.01	0.41
15:AT:97:ALA:O	15:AT:98:LYS:HD2	2.19	0.41
19:AX:18:TYR:C	19:AX:20:GLY:N	2.74	0.41
28:A6:44:ARG:HB3	28:A6:44:ARG:NH1	2.34	0.41
1:BA:530:G:C5	1:BA:2022:U:H5''	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:652(C):G:N2	1:BA:653:A:H1'	2.35	0.41
1:BA:734:A:C4	1:BA:735:A:C8	3.09	0.41
1:BA:1186:G:H2'	1:BA:1187:G:O4'	2.20	0.41
1:BA:1337:G:H2'	1:BA:1338:G:O4'	2.20	0.41
1:BA:1425:G:H2'	1:BA:1426:G:O4'	2.20	0.41
1:BA:2708:G:H1'	13:BR:71:GLN:HE22	1.85	0.41
2:BB:108:U:H2'	2:BB:109:C:H5''	2.02	0.41
15:BT:118:ARG:HH11	15:BT:118:ARG:CG	2.26	0.41
16:BU:16:LYS:HE2	16:BU:16:LYS:HB3	1.77	0.41
21:BZ:35:ARG:HA	21:BZ:35:ARG:HD2	1.94	0.41
28:B6:40:CYS:HA	28:B6:41:PRO:HD3	1.80	0.41
31:CA:324:G:N2	31:CA:327:A:OP2	2.53	0.41
31:CA:544:G:H2'	31:CA:545:C:C6	2.55	0.41
31:CA:586:C:C2'	31:CA:587:G:H5'	2.50	0.41
31:CA:620:C:C2	34:CD:135:LEU:HG	2.56	0.41
32:CB:101:MET:HA	32:CB:108:ILE:HG13	2.03	0.41
32:CB:145:LEU:HD13	32:CB:145:LEU:HA	1.82	0.41
38:CH:25:ASP:HA	38:CH:59:LEU:O	2.20	0.41
31:DA:44:G:N3	31:DA:399:G:C2	2.89	0.41
31:DA:44:G:N1	31:DA:45:U:O2	2.53	0.41
31:DA:108:G:N1	50:DT:15:ARG:HG3	2.36	0.41
31:DA:221:C:H2'	31:DA:222:U:C6	2.55	0.41
31:DA:223:U:H2'	31:DA:224:C:H6	1.85	0.41
31:DA:523:A:H61	42:DL:53:ARG:NH1	2.18	0.41
31:DA:543:C:O2'	31:DA:544:G:H5'	2.20	0.41
31:DA:819:A:N7	31:DA:1529:G:N1	2.68	0.41
31:DA:1157:A:H4'	31:DA:1158:C:O5'	2.21	0.41
31:DA:1434:A:H2'	31:DA:1435:G:O4'	2.20	0.41
32:DB:60:ASP:OD2	32:DB:64:ARG:NE	2.37	0.41
32:DB:97:TRP:HH2	32:DB:176:GLU:CD	2.23	0.41
32:DB:170:GLU:O	32:DB:174:VAL:HG23	2.21	0.41
34:DD:79:PHE:HE1	34:DD:204:ILE:HD13	1.86	0.41
37:DG:72:ARG:H	37:DG:72:ARG:HG2	1.61	0.41
37:DG:95:ARG:HE	37:DG:99:LEU:HD11	1.85	0.41
39:DI:44:VAL:CA	39:DI:45:ALA:HB2	2.50	0.41
39:DI:124:GLN:HE21	39:DI:124:GLN:HB3	1.66	0.41
42:DL:54:LYS:O	42:DL:70:ILE:HG13	2.20	0.41
45:DO:24:SER:O	45:DO:28:GLN:HG3	2.21	0.41
1:AA:45:C:H2'	1:AA:47:C:H6	1.86	0.41
1:AA:183:C:H1'	1:AA:433:C:H1'	2.03	0.41
1:AA:195:A:OP1	11:AP:46:LYS:HE2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:271(H):G:O2'	1:AA:271(I):G:H8	2.02	0.41
1:AA:476:G:N2	1:AA:478:A:H3'	2.35	0.41
1:AA:477:A:H2'	1:AA:478:A:C8	2.55	0.41
1:AA:628:G:H2'	1:AA:629:G:H8	1.83	0.41
1:AA:706:A:C2	1:AA:707:G:H1'	2.55	0.41
1:AA:729:G:OP2	3:AD:13:ARG:NH1	2.51	0.41
1:AA:824:A:H1'	1:AA:2358:G:N7	2.36	0.41
1:AA:1160:G:C6	1:AA:1161:C:C4	3.08	0.41
1:AA:1316:U:H2'	1:AA:1317:A:H8	1.85	0.41
1:AA:1575:C:H2'	1:AA:1576:U:H6	1.84	0.41
1:AA:1665:A:C4'	10:AO:67:LYS:HB2	2.49	0.41
1:AA:2023:G:H4'	1:AA:2617:C:O3'	2.21	0.41
1:AA:2078:C:C4	1:AA:2079:U:C4	3.08	0.41
1:AA:2517:C:C6	1:AA:2542:A:N7	2.88	0.41
1:AA:2618:G:H21	4:AE:150:VAL:HG21	1.85	0.41
1:AA:2802:G:H2'	1:AA:2803:C:H6	1.82	0.41
4:AE:60:ASN:OD1	4:AE:63:LEU:HB2	2.20	0.41
4:AE:71:GLY:HA2	4:AE:72:VAL:C	2.41	0.41
12:AQ:35:VAL:HG13	12:AQ:130:LYS:HB3	2.03	0.41
1:BA:30:G:H2'	1:BA:31:C:O4'	2.20	0.41
1:BA:71:A:H5'	1:BA:71:A:C8	2.56	0.41
1:BA:111:A:H4'	24:B2:69:ARG:NH1	2.34	0.41
1:BA:883:G:H1	1:BA:893:C:N4	2.18	0.41
1:BA:1177:A:OP1	1:BA:1177:A:H3'	2.21	0.41
1:BA:2020:A:O2'	1:BA:2021:C:H5'	2.20	0.41
1:BA:2252:G:H2'	1:BA:2253:G:O4'	2.20	0.41
1:BA:2294:C:P	14:BS:89:ARG:HH22	2.43	0.41
7:BH:43:VAL:HG22	7:BH:52:VAL:HG22	2.02	0.41
12:BQ:58:PHE:O	12:BQ:60:ARG:N	2.53	0.41
22:B0:20:ARG:HH11	22:B0:20:ARG:HD3	1.72	0.41
24:B2:3:LEU:HD23	24:B2:3:LEU:HA	1.83	0.41
30:B8:27:THR:O	30:B8:44:LYS:HE2	2.19	0.41
31:CA:166:G:O2'	31:CA:167:G:H5'	2.20	0.41
31:CA:309:G:H2'	31:CA:310:G:H8	1.85	0.41
31:CA:373:A:N1	31:CA:391:G:O2'	2.51	0.41
31:CA:374:A:C4	31:CA:375:U:C5	3.09	0.41
31:CA:461:A:C4	31:CA:471:G:C5	3.09	0.41
31:CA:626:U:H2'	31:CA:627:G:H8	1.84	0.41
31:CA:654:G:C6	31:CA:655:A:C5	3.08	0.41
31:CA:938:A:N6	31:CA:939:G:C6	2.88	0.41
31:CA:1430:C:N3	31:CA:1471:G:C2	2.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:CC:23:TYR:CG	33:CC:24:ALA:N	2.88	0.41
41:CK:33:THR:HA	41:CK:39:PRO:HA	2.03	0.41
31:DA:114:U:H2'	31:DA:115:G:C8	2.56	0.41
31:DA:373:A:N3	31:DA:374:A:C8	2.89	0.41
31:DA:963:G:O2'	31:DA:1199:U:H5''	2.19	0.41
31:DA:1023:G:H2'	31:DA:1023:G:N3	2.34	0.41
31:DA:1028:C:N3	31:DA:1034:G:C2	2.89	0.41
31:DA:1362:C:C2'	31:DA:1363:C:H5''	2.50	0.41
34:DD:117:ALA:O	34:DD:121:VAL:HG23	2.19	0.41
38:DH:86:ILE:CG1	38:DH:133:LEU:HD22	2.49	0.41
45:DO:19:PRO:HB2	45:DO:20:GLY:H	1.72	0.41
47:DQ:67:LYS:HA	47:DQ:70:ARG:NH1	2.35	0.41
1:AA:56:A:H2'	1:AA:57:C:O4'	2.21	0.41
1:AA:455:C:N3	1:AA:473:G:H5'	2.35	0.41
1:AA:565:C:H4'	1:AA:1253:A:N6	2.35	0.41
1:AA:648:G:O2'	1:AA:2351:G:OP1	2.32	0.41
1:AA:1235:G:C2	1:AA:1236:G:N2	2.89	0.41
1:AA:1334:G:H2'	1:AA:1335:U:C6	2.56	0.41
1:AA:1371:G:HO2'	1:AA:1372:U:H5	1.64	0.41
1:AA:1790:C:H2'	1:AA:1791:A:C5	2.56	0.41
1:AA:1992:G:C8	1:AA:1992:G:O5'	2.74	0.41
1:AA:2065:C:H2'	1:AA:2066:C:C6	2.55	0.41
1:AA:2415:G:H2'	1:AA:2416:C:C6	2.56	0.41
1:AA:2506:U:H4'	1:AA:2507:C:OP1	2.21	0.41
4:AE:49:LEU:HD12	4:AE:49:LEU:HA	1.76	0.41
4:AE:63:LEU:HD23	4:AE:63:LEU:HA	1.91	0.41
6:AG:114:ILE:HG12	6:AG:140:ILE:HD13	2.01	0.41
6:AG:132:ASN:HA	6:AG:157:ILE:O	2.21	0.41
12:AQ:57:HIS:NE2	12:AQ:116:GLU:HG2	2.35	0.41
27:A5:36:CYS:SG	27:A5:49:CYS:HB3	2.61	0.41
1:BA:729:G:C6	3:BD:208:LYS:HB2	2.55	0.41
1:BA:829:A:H5''	1:BA:831:G:N7	2.35	0.41
1:BA:864:G:O2'	1:BA:865:C:H5'	2.21	0.41
1:BA:1047:G:H2'	1:BA:1110:G:N1	2.34	0.41
1:BA:1177:A:H8	1:BA:1177:A:O5'	2.04	0.41
1:BA:1754:C:H2'	1:BA:1755:A:O4'	2.20	0.41
1:BA:1790:C:O2'	3:BD:209:ALA:HB2	2.20	0.41
1:BA:1839:G:N7	1:BA:1927:A:H1'	2.35	0.41
1:BA:2035:G:H4'	1:BA:2036:C:OP2	2.20	0.41
1:BA:2104:G:H2'	1:BA:2104:G:N3	2.35	0.41
1:BA:2174:C:H2'	1:BA:2175:C:H6	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:2420:C:H5'	28:B6:54:ILE:CD1	2.50	0.41
1:BA:2479:G:OP1	1:BA:2537:U:H1'	2.20	0.41
1:BA:2638:G:P	4:BE:82:ARG:NH2	2.94	0.41
1:BA:2689:U:C4'	1:BA:2690:C:H5'	2.50	0.41
1:BA:2801(A):A:H1'	1:BA:2895:U:H1'	2.03	0.41
1:BA:2839:G:N2	1:BA:2880:C:C2	2.89	0.41
2:BB:8:U:O2'	14:BS:40:ILE:HD13	2.21	0.41
3:BD:147:LEU:HD13	3:BD:155:LEU:HD21	2.02	0.41
6:BG:126:ASP:HB2	6:BG:130:ASN:O	2.19	0.41
9:BN:23:LEU:HD12	9:BN:99:LEU:HD23	2.03	0.41
11:BP:135:LEU:HD23	11:BP:135:LEU:HA	1.79	0.41
14:BS:101:LEU:O	14:BS:102:ALA:CB	2.69	0.41
19:BX:29:TRP:CE3	19:BX:78:LYS:HB3	2.56	0.41
21:BZ:45:ASP:OD1	21:BZ:49:ARG:HD2	2.20	0.41
29:B7:19:ARG:HG2	29:B7:19:ARG:HH11	1.85	0.41
31:CA:102:G:O2'	31:CA:151:A:N3	2.32	0.41
31:CA:116:A:H2'	31:CA:117:G:H8	1.84	0.41
31:CA:185:A:C6	31:CA:186:C:N4	2.88	0.41
31:CA:195:A:H1'	31:CA:222:U:O2'	2.21	0.41
31:CA:309:G:O2'	31:CA:607:A:N1	2.53	0.41
31:CA:457:C:H2'	31:CA:458:C:C6	2.55	0.41
31:CA:586:C:O2'	31:CA:878:G:H4'	2.21	0.41
31:CA:817:C:C5	31:CA:819:A:H1'	2.56	0.41
31:CA:1016:A:O2'	31:CA:1217:C:O2	2.37	0.41
31:CA:1444:C:H42	31:CA:1458:G:H1	1.68	0.41
32:CB:25:ASN:HA	32:CB:26:PRO:HD2	1.89	0.41
33:CC:23:TYR:HB2	40:CJ:10:GLY:HA2	2.03	0.41
39:CI:46:ALA:HB1	39:CI:78:LYS:CA	2.50	0.41
40:CJ:13:HIS:CG	40:CJ:14:LYS:N	2.88	0.41
41:CK:38:ASN:HA	41:CK:39:PRO:HD3	1.86	0.41
31:DA:45:U:H2'	31:DA:46:G:H8	1.85	0.41
31:DA:560:U:N3	31:DA:566:G:C4	2.88	0.41
31:DA:575:G:O2'	31:DA:821:G:H5'	2.21	0.41
31:DA:673:G:H5''	36:DF:87:ARG:CZ	2.49	0.41
31:DA:685:G:O2'	31:DA:686:U:H5'	2.20	0.41
31:DA:832:C:O2'	31:DA:833:U:OP2	2.34	0.41
31:DA:922:G:H2'	31:DA:923:A:H8	1.85	0.41
31:DA:1206:G:C6	31:DA:1207:G:C5	3.09	0.41
34:DD:189:PRO:CB	34:DD:194:LEU:HD11	2.48	0.41
35:DE:108:ALA:O	35:DE:112:LEU:HG	2.21	0.41
38:DH:88:LYS:HB3	38:DH:89:PRO:HD2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:DQ:59:ILE:HG22	47:DQ:73:VAL:HA	2.03	0.41
1:AA:239:U:H2'	1:AA:240:G:O4'	2.21	0.41
1:AA:301:G:C4	1:AA:302:C:C5	3.09	0.41
1:AA:301:G:HO2'	1:AA:302:C:H6	1.67	0.41
1:AA:341:G:H8	1:AA:341:G:O5'	2.03	0.41
1:AA:623:G:C2	1:AA:624:C:C2	3.09	0.41
1:AA:1327:C:H2'	1:AA:1328:G:O4'	2.21	0.41
1:AA:1439:A:H2'	1:AA:1440:G:O4'	2.20	0.41
1:AA:2236:C:H2'	1:AA:2237:G:O4'	2.20	0.41
1:AA:2813:A:H2'	1:AA:2814:C:O4'	2.21	0.41
7:AH:121:ILE:HD11	7:AH:140:LYS:HG2	2.03	0.41
9:AN:63:THR:HB	9:AN:64:GLY:H	1.69	0.41
11:AP:84:ASN:O	11:AP:87:ASP:N	2.51	0.41
1:BA:479:A:H4'	1:BA:480:A:OP1	2.21	0.41
1:BA:592:G:O2'	30:B8:4:MET:N	2.43	0.41
1:BA:878:A:C2	1:BA:879:G:N7	2.89	0.41
1:BA:1648:C:H2'	1:BA:1649:G:O5'	2.21	0.41
1:BA:1913:A:H61	31:CA:1493:A:C5'	2.31	0.41
1:BA:1914:C:H3'	1:BA:1914:C:OP1	2.21	0.41
1:BA:1925:C:C2'	1:BA:1926:U:H5'	2.50	0.41
1:BA:2791:C:H1'	1:BA:2807:G:N2	2.36	0.41
2:BB:16:G:N2	2:BB:69:G:H1'	2.35	0.41
7:BH:27:LYS:HE2	7:BH:27:LYS:HB3	1.70	0.41
15:BT:93:ARG:NH2	15:BT:95:ARG:HD2	2.36	0.41
31:CA:337:C:H2'	31:CA:338:A:C8	2.55	0.41
31:CA:428:G:H4'	31:CA:429:U:O5'	2.21	0.41
34:CD:33:MET:HA	34:CD:37:PRO:HA	2.02	0.41
37:CG:122:HIS:HA	37:CG:125:MET:SD	2.60	0.41
43:CM:40:ASN:HA	43:CM:41:PRO:HD2	1.75	0.41
45:CO:67:LEU:HB3	45:CO:78:TYR:HE1	1.85	0.41
50:CT:72:LEU:HD11	50:CT:80:ARG:HD2	2.03	0.41
31:DA:445:G:C6	31:DA:446:G:C5	3.08	0.41
31:DA:908:A:O5'	31:DA:908:A:H8	2.03	0.41
31:DA:926:G:C6	31:DA:1505:G:C5	3.08	0.41
31:DA:944:G:C2'	31:DA:1339:A:H61	2.33	0.41
31:DA:1225:A:C5	31:DA:1226:C:N4	2.89	0.41
31:DA:1457:G:H2'	31:DA:1458:G:C8	2.56	0.41
31:DA:1505:G:H4'	31:DA:1506:U:H5''	2.03	0.41
32:DB:106:LYS:O	32:DB:110:GLN:HG3	2.20	0.41
32:DB:153:ARG:HG3	32:DB:154:LEU:N	2.35	0.41
32:DB:204:ASN:HB3	32:DB:210:SER:CB	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:DB:215:LEU:HD23	32:DB:215:LEU:HA	1.88	0.41
34:DD:61:LYS:HD3	34:DD:206:PHE:CE2	2.55	0.41
35:DE:98:THR:HB	35:DE:99:GLY:H	1.55	0.41
37:DG:113:GLU:HG3	37:DG:119:ARG:HG2	2.01	0.41
1:AA:110:G:C2	1:AA:111:A:C8	3.08	0.41
1:AA:284:U:H2'	1:AA:285:C:C6	2.55	0.41
1:AA:393:C:H2'	1:AA:394:A:C8	2.56	0.41
1:AA:414:C:O2'	1:AA:415:A:H5'	2.21	0.41
1:AA:517:C:OP1	27:A5:13:LYS:HD3	2.20	0.41
1:AA:557:U:H2'	1:AA:558:G:C8	2.56	0.41
1:AA:705:A:C8	1:AA:727:A:C2	3.09	0.41
1:AA:1050:A:O2'	1:AA:2752:C:H1'	2.19	0.41
1:AA:1202:C:N4	1:AA:1203:G:C6	2.89	0.41
1:AA:1210:A:H5''	1:AA:1212:G:O4'	2.20	0.41
1:AA:1223:G:C2	1:AA:1227:G:C5	3.08	0.41
1:AA:1243:G:H2'	1:AA:1244:G:O4'	2.20	0.41
1:AA:1355:G:H2'	1:AA:1356:G:O4'	2.20	0.41
1:AA:1478:G:HO2'	1:AA:1558:A:H2	1.66	0.41
1:AA:2199:A:H3'	1:AA:2200:C:C6	2.56	0.41
1:AA:2199:A:H5''	1:AA:2200:C:OP2	2.21	0.41
1:AA:2295:C:C2'	1:AA:2296:U:H5'	2.50	0.41
1:AA:2342:C:H2'	1:AA:2343:C:O4'	2.20	0.41
1:AA:2741:A:H2'	1:AA:2742:C:O4'	2.20	0.41
1:AA:2803:C:H2'	1:AA:2804:C:H6	1.85	0.41
2:AB:52:A:O2'	2:AB:53:A:H5'	2.20	0.41
2:AB:66:A:H61	2:AB:109:C:C5'	2.33	0.41
7:AH:149:ARG:HD2	7:AH:164:TYR:CE1	2.56	0.41
9:AN:29:LYS:NZ	9:AN:140:VAL:HG13	2.35	0.41
13:AR:28:LEU:HD21	13:AR:114:VAL:O	2.20	0.41
14:AS:28:VAL:HG11	14:AS:98:VAL:HG13	2.03	0.41
15:AT:11:GLU:O	15:AT:15:VAL:HG23	2.20	0.41
16:AU:95:LEU:HD12	16:AU:95:LEU:HA	1.83	0.41
24:A2:32:LEU:HD21	24:A2:50:ILE:CG2	2.51	0.41
25:A3:28:LEU:HD23	25:A3:28:LEU:HA	1.82	0.41
1:BA:1364:G:P	23:B1:3:LYS:HG2	2.60	0.41
1:BA:1889:A:H2'	1:BA:1890:A:C8	2.56	0.41
1:BA:2070:G:H2'	1:BA:2071:A:O4'	2.20	0.41
1:BA:2291:U:O3'	1:BA:2379:G:N2	2.53	0.41
1:BA:2691:C:O3'	1:BA:2871:C:H4'	2.20	0.41
2:BB:33:G:N2	2:BB:50:G:C4	2.89	0.41
3:BD:9:TYR:CZ	3:BD:13:ARG:HG2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BD:152:GLY:O	3:BD:154:LYS:HG2	2.20	0.41
3:BD:172:TYR:CD1	3:BD:186:HIS:HA	2.55	0.41
8:BI:85:GLU:O	8:BI:86:THR:OG1	2.38	0.41
9:BN:125:GLY:HA2	9:BN:126:PRO:O	2.20	0.41
12:BQ:37:LEU:HD21	12:BQ:130:LYS:CE	2.51	0.41
31:CA:55:A:C5	31:CA:56:U:C5	3.09	0.41
31:CA:377:G:P	46:CP:3:LYS:HZ2	2.43	0.41
31:CA:874:G:H2'	31:CA:875:C:H6	1.85	0.41
31:CA:923:A:H2'	31:CA:924:C:O4'	2.21	0.41
31:CA:1007:C:N3	31:CA:1022:G:O6	2.54	0.41
31:CA:1052:U:H3	31:CA:1206:G:H1	1.68	0.41
31:CA:1169:A:C6	31:CA:1170:A:C6	3.09	0.41
31:CA:1220:G:N2	49:CS:54:GLY:O	2.53	0.41
31:CA:1255:G:N2	31:CA:1259:C:O2	2.45	0.41
31:CA:1288:A:H2	31:CA:1352:C:O2	2.04	0.41
32:CB:144:ARG:O	32:CB:147:LYS:N	2.53	0.41
34:CD:13:ARG:HB2	34:CD:40:PRO:HD3	2.02	0.41
34:CD:61:LYS:O	34:CD:65:ARG:HB2	2.20	0.41
37:CG:15:ASP:OD1	37:CG:44:TYR:OH	2.37	0.41
37:CG:90:GLU:OE1	37:CG:90:GLU:N	2.52	0.41
41:CK:48:ILE:CG2	41:CK:63:LEU:HD22	2.51	0.41
42:CL:31:PRO:HB2	42:CL:32:PHE:CD2	2.55	0.41
47:CQ:84:LEU:O	47:CQ:87:LYS:HB2	2.21	0.41
31:DA:189(F):U:O4	47:DQ:63:ARG:N	2.39	0.41
31:DA:338:A:H2'	31:DA:339:C:O4'	2.21	0.41
31:DA:577:G:C8	31:DA:816:A:C6	3.09	0.41
31:DA:586:C:HO2'	31:DA:878:G:H4'	1.86	0.41
31:DA:786:G:C2	31:DA:797:C:C2	3.09	0.41
31:DA:1229:A:OP2	43:DM:114:ARG:HD2	2.19	0.41
31:DA:1284:C:H2'	31:DA:1285:A:N7	2.36	0.41
31:DA:1333:A:C2	31:DA:1334:G:H1'	2.56	0.41
34:DD:73:ARG:HA	34:DD:73:ARG:HD2	1.89	0.41
37:DG:111:ARG:HD3	37:DG:113:GLU:OE2	2.20	0.41
39:DI:120:ARG:O	39:DI:122:ALA:N	2.54	0.41
40:DJ:40:LEU:HD12	40:DJ:69:ASN:HB3	2.02	0.41
48:DR:66:LEU:O	48:DR:70:ILE:HG13	2.21	0.41
1:AA:264:C:O2'	1:AA:265:A:H2'	2.21	0.41
1:AA:312:G:H4'	1:AA:331:A:N3	2.36	0.41
1:AA:393:C:H2'	1:AA:394:A:H8	1.86	0.41
1:AA:613:G:O2'	1:AA:614(C):A:N1	2.45	0.41
1:AA:705:A:H1'	3:AD:9:TYR:CE1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:910:A:H2'	1:AA:911:A:C8	2.56	0.41
1:AA:911:A:H2'	12:AQ:9:TYR:OH	2.21	0.41
1:AA:942:G:C5	1:AA:943:U:C5	3.09	0.41
1:AA:945:A:C4	1:AA:2448:A:C2	3.08	0.41
1:AA:977:G:C6	1:AA:987:G:C6	3.08	0.41
1:AA:1221:C:H2'	1:AA:1221(A):C:C6	2.55	0.41
1:AA:1651:G:N2	1:AA:2007:C:C2	2.89	0.41
1:AA:1653:G:H4'	1:AA:1654:A:O5'	2.20	0.41
1:AA:1758:G:C2	1:AA:2696:U:H5'	2.55	0.41
1:AA:1889:A:H2'	1:AA:1890:A:H8	1.83	0.41
1:AA:2131:G:H5'	1:AA:2133:G:O5'	2.21	0.41
1:AA:2139:C:N4	1:AA:2152:G:H1	2.19	0.41
1:AA:2201:C:H2'	1:AA:2202:C:H6	1.86	0.41
1:AA:2282:G:H5'	1:AA:2389:G:H1'	2.03	0.41
1:AA:2298:A:C8	1:AA:2299:G:C8	3.09	0.41
1:AA:2335:A:C8	1:AA:2337:G:N7	2.89	0.41
1:AA:2405:G:O2'	1:AA:2406:U:OP2	2.35	0.41
1:AA:2427:C:H5''	1:AA:2428:G:OP1	2.21	0.41
1:AA:2468:G:H22	1:AA:2481:G:H1'	1.85	0.41
1:AA:2469:A:C2	1:AA:2482:G:C8	3.09	0.41
1:AA:2680:C:H5'	4:AE:189:PRO:HA	2.03	0.41
1:AA:2738:A:N1	1:AA:2739:U:C2	2.89	0.41
1:AA:2811:G:N2	1:AA:2891:G:H1'	2.36	0.41
1:AA:2865:U:C4	1:AA:2866:U:C4	3.09	0.41
2:AB:114:C:O2'	14:AS:46:VAL:HG13	2.21	0.41
3:AD:68:LYS:HD2	3:AD:70:TRP:CZ2	2.56	0.41
3:AD:233:HIS:NE2	3:AD:246:PRO:HA	2.35	0.41
4:AE:21:VAL:HA	4:AE:22:PRO:HD2	1.82	0.41
6:AG:64:THR:HG22	6:AG:94:LEU:HD11	2.03	0.41
6:AG:125:PHE:CZ	6:AG:170:ARG:HA	2.56	0.41
6:AG:141:PHE:HA	6:AG:142:PRO:HD2	1.81	0.41
7:AH:24:VAL:HG21	7:AH:72:ILE:HG12	2.02	0.41
7:AH:69:ARG:HG3	7:AH:70:THR:N	2.36	0.41
7:AH:99:VAL:O	7:AH:102:ALA:HB3	2.20	0.41
11:AP:113:LYS:HA	11:AP:129:ALA:O	2.21	0.41
11:AP:120:ALA:HB1	11:AP:138:LEU:HA	2.02	0.41
14:AS:104:GLY:O	14:AS:107:GLU:N	2.54	0.41
18:AW:9:TYR:HA	18:AW:100:THR:HG23	2.03	0.41
20:AY:68:HIS:ND1	20:AY:70:SER:HB3	2.36	0.41
21:AZ:5:LEU:HD23	21:AZ:47:VAL:CG2	2.42	0.41
23:A1:8:SER:OG	23:A1:10:LYS:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:222:A:N1	1:BA:233:A:H5''	2.36	0.41
1:BA:1173:G:H2'	1:BA:1173:G:OP2	2.21	0.41
1:BA:1212:G:H1'	1:BA:1236:G:N2	2.36	0.41
1:BA:1359:A:N3	1:BA:1359:A:H5'	2.35	0.41
1:BA:1400:G:H2'	1:BA:1401:G:C8	2.55	0.41
1:BA:1491:G:O2'	1:BA:1492:G:H5'	2.20	0.41
1:BA:1539:G:H2'	1:BA:1540:U:O4'	2.21	0.41
1:BA:2126:A:N3	1:BA:2127:G:H1'	2.36	0.41
1:BA:2134:A:H62	1:BA:2157:G:C4'	2.34	0.41
1:BA:2186:G:H2'	1:BA:2187:G:H5''	2.02	0.41
1:BA:2319:G:H22	14:BS:3:ARG:HE	1.67	0.41
2:BB:78:A:C2	2:BB:100:A:C4	3.09	0.41
3:BD:79:VAL:HG12	3:BD:113:VAL:HA	2.02	0.41
3:BD:107:ALA:HA	3:BD:108:PRO:HD2	1.99	0.41
3:BD:246:PRO:HG2	3:BD:255:LYS:HG3	2.02	0.41
5:BF:32:LEU:CD1	5:BF:105:VAL:HG13	2.51	0.41
6:BG:50:ALA:O	6:BG:52:ILE:N	2.54	0.41
13:BR:29:LEU:HD12	13:BR:29:LEU:HA	1.82	0.41
13:BR:56:LYS:HE3	13:BR:87:TYR:O	2.21	0.41
18:BW:18:ARG:NH1	18:BW:76:VAL:O	2.53	0.41
23:B1:3:LYS:HB3	23:B1:4:VAL:H	1.28	0.41
31:CA:45:U:H2'	31:CA:46:G:C8	2.56	0.41
31:CA:344:A:O2'	31:CA:346:G:O6	2.27	0.41
31:CA:452:A:N6	31:CA:480:U:H3	2.19	0.41
31:CA:472:A:C5	31:CA:473:G:C8	3.09	0.41
31:CA:509:A:C8	31:CA:509:A:H3'	2.55	0.41
31:CA:512:U:H2'	31:CA:513:C:C6	2.55	0.41
31:CA:592:G:C2	31:CA:593:G:C8	3.08	0.41
31:CA:991:U:C4	31:CA:1212:U:H1'	2.56	0.41
31:CA:992:U:H1'	31:CA:993:G:C5	2.55	0.41
31:CA:1085:U:H3'	31:CA:1086:U:H5	1.86	0.41
31:CA:1103:C:H5''	32:CB:98:LEU:HD13	2.01	0.41
31:CA:1104:G:C6	31:CA:1105:A:C5	3.09	0.41
31:CA:1128:C:H1'	31:CA:1146:A:N6	2.34	0.41
31:CA:1158:C:H5	31:CA:1181:G:H1	1.67	0.41
31:CA:1492:A:C2'	31:CA:1493:A:OP1	2.69	0.41
32:CB:15:VAL:CG2	32:CB:209:ARG:HG2	2.50	0.41
32:CB:189:ASP:N	32:CB:189:ASP:OD1	2.54	0.41
33:CC:114:PRO:HA	33:CC:185:GLY:HA3	2.02	0.41
34:CD:50:ARG:HA	34:CD:51:PRO:HD3	1.84	0.41
34:CD:98:GLU:HA	34:CD:103:ASN:ND2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:CE:92:LYS:HA	35:CE:93:PRO:HD2	1.86	0.41
37:CG:149:ARG:HB3	41:CK:59:TYR:CZ	2.56	0.41
39:CI:44:VAL:N	39:CI:45:ALA:CA	2.82	0.41
39:CI:89:ASN:HA	39:CI:90:PRO:HD3	1.81	0.41
39:CI:115:GLY:HA2	40:CJ:58:ASP:OD1	2.21	0.41
40:CJ:6:ILE:HA	40:CJ:97:GLU:O	2.21	0.41
45:CO:18:PHE:CE1	45:CO:21:ASP:HB2	2.56	0.41
45:CO:24:SER:O	45:CO:28:GLN:HG3	2.21	0.41
47:CQ:86:GLU:O	47:CQ:90:ILE:HG12	2.21	0.41
49:CS:73:GLU:HB2	49:CS:74:PHE:CD2	2.56	0.41
50:CT:13:LEU:O	50:CT:17:ARG:HG3	2.20	0.41
31:DA:44:G:H2'	31:DA:45:U:O4'	2.21	0.41
31:DA:160:A:N7	31:DA:161:A:N7	2.69	0.41
31:DA:414:A:H2'	31:DA:415:A:C8	2.55	0.41
31:DA:457:C:H2'	31:DA:458:C:C6	2.56	0.41
31:DA:457:C:H42	31:DA:474:G:H1	1.69	0.41
31:DA:657:G:C2	31:DA:750:G:C5	3.09	0.41
31:DA:673:G:C6	31:DA:734:G:C6	3.09	0.41
31:DA:708:C:H2'	31:DA:709:G:H8	1.86	0.41
31:DA:724:G:O2'	31:DA:725:G:H5'	2.20	0.41
31:DA:724:G:N3	31:DA:725:G:C8	2.88	0.41
31:DA:814:A:H2'	31:DA:816:A:H5''	2.02	0.41
31:DA:1069:C:C4	31:DA:1070:U:C5	3.09	0.41
31:DA:1079:G:C6	31:DA:1080:A:N6	2.88	0.41
31:DA:1127:G:H1'	31:DA:1147:C:H42	1.86	0.41
31:DA:1187:G:C6	31:DA:1188:A:C5	3.09	0.41
31:DA:1337:G:H5''	31:DA:1338:G:OP1	2.21	0.41
31:DA:1361:G:H2'	31:DA:1362:C:O4'	2.21	0.41
32:DB:98:LEU:HD23	32:DB:98:LEU:HA	1.92	0.41
32:DB:178:ARG:HH22	38:DH:68:ARG:HH12	1.67	0.41
34:DD:106:TYR:CD2	34:DD:106:TYR:C	2.93	0.41
34:DD:108:LEU:HB3	34:DD:110:PHE:HE1	1.84	0.41
34:DD:148:VAL:HG12	34:DD:149:ALA:N	2.36	0.41
35:DE:33:VAL:HG21	35:DE:109:ILE:HG12	2.02	0.41
37:DG:87:VAL:HA	37:DG:88:PRO:HD2	1.90	0.41
37:DG:122:HIS:HA	37:DG:125:MET:SD	2.61	0.41
38:DH:109:ILE:HD13	38:DH:109:ILE:H	1.85	0.41
39:DI:20:ARG:HA	39:DI:21:PRO:HD2	1.91	0.41
41:DK:38:ASN:HA	41:DK:39:PRO:HD3	1.79	0.41
48:DR:76:LEU:HD13	48:DR:76:LEU:HA	1.74	0.41
50:DT:12:ALA:O	50:DT:15:ARG:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:233:A:H8	1:AA:233:A:OP2	2.04	0.41
1:AA:948:G:H2'	1:AA:949:C:C6	2.56	0.41
1:AA:1257:C:H4'	5:AF:83:PHE:CD2	2.56	0.41
1:AA:1297:C:H2'	1:AA:1298:C:H6	1.86	0.41
1:AA:1308:A:H5''	1:AA:1309:G:OP2	2.21	0.41
1:AA:1385:G:H1'	1:AA:1386:C:C6	2.56	0.41
1:AA:2041:U:H2'	1:AA:2042:A:H8	1.85	0.41
1:AA:2472:G:H2'	1:AA:2475:C:N4	2.36	0.41
1:AA:2543:G:H21	1:AA:2646:C:H5''	1.86	0.41
1:AA:2748:A:N6	1:AA:2749:A:C6	2.89	0.41
2:AB:113:G:H2'	2:AB:114:C:C6	2.56	0.41
4:AE:29:GLY:HA2	4:AE:30:PRO:HA	1.80	0.41
4:AE:68:ALA:O	4:AE:70:ALA:N	2.44	0.41
4:AE:111:ARG:HA	13:AR:1:MET:SD	2.61	0.41
5:AF:129:PHE:HB3	5:AF:132:VAL:HG22	2.03	0.41
8:AI:98:ALA:O	8:AI:102:SER:HB3	2.21	0.41
12:AQ:44:ALA:HB2	12:AQ:70:PRO:HG3	2.02	0.41
21:AZ:144:LEU:HD12	21:AZ:144:LEU:HA	1.84	0.41
24:A2:35:LEU:HD12	24:A2:53:LEU:HD12	2.03	0.41
1:BA:26:G:H1'	1:BA:514:A:N6	2.35	0.41
1:BA:95:G:O2'	24:B2:46:GLN:HA	2.21	0.41
1:BA:229:A:OP1	1:BA:229:A:C8	2.74	0.41
1:BA:244:A:C2	1:BA:255:A:C4	3.09	0.41
1:BA:370:G:H4'	1:BA:371:A:OP2	2.20	0.41
1:BA:755:C:H2'	1:BA:756:C:H6	1.86	0.41
1:BA:1044:G:H5''	1:BA:1045:A:OP2	2.21	0.41
1:BA:1209:G:H21	1:BA:1210:A:H62	1.69	0.41
1:BA:1311:G:O6	29:B7:9:ARG:NH2	2.53	0.41
1:BA:2137:C:O2	1:BA:2137:C:C2'	2.69	0.41
1:BA:2203:U:C2'	1:BA:2205:C:H5'	2.51	0.41
1:BA:2290:G:H4'	1:BA:2381:C:O2'	2.21	0.41
1:BA:2572:A:OP1	1:BA:2574:G:O2'	2.35	0.41
1:BA:2646:C:O5'	1:BA:2646:C:H6	2.04	0.41
5:BF:114:VAL:HG21	5:BF:202:PHE:CZ	2.56	0.41
6:BG:126:ASP:CB	6:BG:130:ASN:H	2.32	0.41
9:BN:15:LEU:HB2	9:BN:135:PRO:HB2	2.03	0.41
9:BN:75:TYR:CE2	9:BN:77:GLY:HA2	2.56	0.41
11:BP:19:VAL:CG2	11:BP:31:ALA:HB1	2.51	0.41
21:BZ:93:ASP:HB2	21:BZ:131:ARG:NH1	2.35	0.41
31:CA:221:C:H2'	31:CA:222:U:H6	1.85	0.41
31:CA:667:G:H4'	45:CO:51:HIS:ND1	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:678:U:H2'	31:CA:679:C:C6	2.55	0.41
31:CA:684:A:N6	31:CA:685:G:C6	2.89	0.41
31:CA:781:A:H4'	31:CA:1522:U:O2'	2.21	0.41
31:CA:1040:U:H2'	31:CA:1041:A:C8	2.55	0.41
31:CA:1041:A:C6	31:CA:1042:G:C5	3.09	0.41
31:CA:1479:C:H2'	31:CA:1480:G:C8	2.56	0.41
32:CB:163:PHE:HD2	32:CB:185:ILE:HG13	1.80	0.41
37:CG:26:PHE:CE2	37:CG:30:ILE:HD11	2.56	0.41
43:CM:40:ASN:ND2	43:CM:43:THR:HG23	2.36	0.41
43:CM:85:GLY:HA3	43:CM:86:CYS:HB3	2.02	0.41
51:CU:9:ARG:NH1	51:CU:13:ILE:HD11	2.36	0.41
31:DA:373:A:C4	31:DA:374:A:C8	3.09	0.41
31:DA:660:G:C4	31:DA:661:G:C8	3.09	0.41
31:DA:681:C:N3	31:DA:710:G:C2	2.88	0.41
31:DA:689:C:H2'	31:DA:690:G:O4'	2.21	0.41
31:DA:1072:G:C4	31:DA:1104:G:N2	2.89	0.41
32:DB:107:THR:HA	32:DB:110:GLN:OE1	2.20	0.41
34:DD:36:ARG:HB3	34:DD:38:TYR:CZ	2.56	0.41
34:DD:135:LEU:C	34:DD:137:SER:H	2.24	0.41
37:DG:140:ASP:O	37:DG:144:MET:HB2	2.21	0.41
40:DJ:45:ARG:HG2	40:DJ:47:PHE:CZ	2.56	0.41
45:DO:61:GLY:O	45:DO:64:ARG:HB3	2.21	0.41
49:DS:20:LEU:HD21	49:DS:43:GLU:HG2	2.03	0.41
1:AA:354:G:H2'	1:AA:355:G:O4'	2.21	0.40
1:AA:401:A:C6	1:AA:402:A:C6	3.09	0.40
1:AA:556:G:C6	1:AA:557:U:C4	3.09	0.40
1:AA:1486:A:H2'	1:AA:1487:G:C8	2.55	0.40
1:AA:1514:U:H2'	1:AA:1515:G:H8	1.86	0.40
1:AA:1572:A:H8	1:AA:1572:A:O5'	2.04	0.40
1:AA:2050:C:H2'	1:AA:2051:A:C8	2.54	0.40
1:AA:2176:A:C6	1:AA:2177:C:C4	3.09	0.40
1:AA:2292:C:H2'	1:AA:2293:C:C6	2.57	0.40
1:AA:2364:C:H4'	22:A0:56:ASP:OD2	2.21	0.40
1:AA:2409:G:H2'	1:AA:2410:G:O4'	2.21	0.40
1:AA:2776:A:H3'	1:AA:2776:A:OP1	2.22	0.40
1:AA:2821:A:OP2	1:AA:2822:G:OP2	2.40	0.40
9:AN:3:THR:CG2	9:AN:4:TYR:H	2.34	0.40
9:AN:30:ILE:HG22	9:AN:34:LEU:HD22	2.03	0.40
30:A8:6:THR:HG23	30:A8:64:TYR:HD2	1.86	0.40
30:A8:23:VAL:HG11	30:A8:47:LYS:HD3	2.03	0.40
1:BA:38:A:H2'	1:BA:39:C:C6	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:71:A:H2	19:BX:31:HIS:CE1	2.37	0.40
1:BA:322:A:H2	1:BA:339:U:O4	2.04	0.40
1:BA:652(Q):G:C6	1:BA:652(R):C:C4	3.09	0.40
1:BA:690:G:H2'	1:BA:691:C:C6	2.57	0.40
1:BA:847:U:OP2	1:BA:928:G:O6	2.39	0.40
1:BA:988:A:H8	1:BA:988:A:O5'	2.03	0.40
1:BA:1149:G:H2'	1:BA:1150:C:C6	2.56	0.40
1:BA:1176:G:N2	1:BA:1178:C:OP2	2.54	0.40
1:BA:1358:G:N2	1:BA:1372:U:OP2	2.47	0.40
1:BA:1885:A:H2'	1:BA:1886:C:O4'	2.21	0.40
5:BF:129:PHE:CD2	5:BF:163:VAL:HG21	2.57	0.40
6:BG:39:ILE:HD11	6:BG:102:PHE:CE1	2.56	0.40
18:BW:12:ILE:HD13	18:BW:12:ILE:HG21	1.76	0.40
21:BZ:156:LYS:O	21:BZ:157:LEU:HB2	2.21	0.40
30:B8:29:LYS:HB2	30:B8:33:ASN:HD21	1.85	0.40
31:CA:425:G:N2	31:CA:426:G:H1'	2.36	0.40
31:CA:613:C:C2	31:CA:628:G:N2	2.89	0.40
31:CA:731:G:H5'	31:CA:766:A:H4'	2.02	0.40
31:CA:741:G:H2'	31:CA:742:G:H8	1.86	0.40
31:CA:826:C:H6	31:CA:826:C:O5'	2.04	0.40
31:CA:1031:G:H2'	31:CA:1032:G:H8	1.80	0.40
31:CA:1079:G:C6	31:CA:1080:A:N6	2.90	0.40
31:CA:1263:C:N3	31:CA:1272:G:O6	2.54	0.40
35:CE:127:ASN:HA	35:CE:128:PRO:HD3	1.92	0.40
37:CG:16:LEU:HD13	39:CI:42:ARG:HA	2.02	0.40
31:DA:105:G:H2'	31:DA:106:C:H6	1.80	0.40
31:DA:253:U:H2'	31:DA:254:G:H8	1.86	0.40
31:DA:665:A:H2'	31:DA:725:G:N2	2.36	0.40
31:DA:709:G:H2'	31:DA:710:G:C8	2.55	0.40
31:DA:885:G:O2'	31:DA:914:A:N1	2.49	0.40
31:DA:1030:C:N4	31:DA:1031:G:C6	2.80	0.40
31:DA:1030:C:C5	31:DA:1030(A):G:H1'	2.56	0.40
31:DA:1356:G:N2	31:DA:1367:C:C2	2.89	0.40
32:DB:87:ARG:NE	32:DB:233:SER:HB2	2.36	0.40
35:DE:11:ILE:HG22	35:DE:12:LEU:HB2	2.02	0.40
38:DH:10:LEU:O	38:DH:14:ARG:HB2	2.22	0.40
44:DN:48:ALA:HB2	44:DN:53:LEU:HD12	2.02	0.40
47:DQ:6:LEU:O	47:DQ:58:GLU:HA	2.21	0.40
1:AA:239:U:O4	1:AA:240:G:C6	2.75	0.40
1:AA:428:A:N6	1:AA:429:A:N1	2.69	0.40
1:AA:638:G:H2'	1:AA:639:U:O4'	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:675:A:C6	1:AA:676:A:C6	3.09	0.40
1:AA:841:A:H2'	1:AA:842:G:C8	2.57	0.40
1:AA:974:G:O3'	1:AA:975(A):G:C8	2.74	0.40
1:AA:1121:C:H2'	1:AA:1122:G:O4'	2.21	0.40
1:AA:1540:U:O2'	1:AA:1541:G:H5'	2.20	0.40
1:AA:1568:G:H5''	3:AD:61:LEU:HD22	2.03	0.40
1:AA:2301:C:H2'	1:AA:2302:G:C8	2.54	0.40
1:AA:2744:G:N3	1:AA:2744:G:H2'	2.36	0.40
1:AA:2768:C:H2'	1:AA:2769:C:O4'	2.22	0.40
2:AB:66:A:C5	2:AB:109:C:C5	3.09	0.40
4:AE:104:VAL:HG11	4:AE:188:VAL:HG22	2.02	0.40
4:AE:111:ARG:HB3	13:AR:1:MET:HE1	2.01	0.40
4:AE:117:MET:O	4:AE:118:LYS:HB3	2.21	0.40
5:AF:34:TRP:CZ3	5:AF:35:GLU:HG2	2.56	0.40
5:AF:118:ALA:C	5:AF:120:GLU:H	2.24	0.40
6:AG:19:LEU:HD23	6:AG:19:LEU:HA	1.92	0.40
7:AH:139:GLN:O	7:AH:143:GLN:N	2.47	0.40
9:AN:3:THR:HG22	9:AN:4:TYR:N	2.36	0.40
9:AN:66:LYS:HB3	9:AN:70:LYS:HB2	2.03	0.40
21:AZ:3:TYR:HB2	21:AZ:56:VAL:O	2.22	0.40
22:A0:68:GLU:N	22:A0:80:HIS:O	2.55	0.40
1:BA:176:G:O2'	1:BA:177:G:H5'	2.21	0.40
1:BA:271(I):G:O2'	1:BA:271(J):C:H5'	2.21	0.40
1:BA:528:A:H3'	1:BA:528:A:H8	1.85	0.40
1:BA:646:A:H2'	1:BA:647:G:O4'	2.22	0.40
1:BA:692:C:C2	1:BA:771:G:C2	3.09	0.40
1:BA:732:C:H2'	1:BA:733:G:O4'	2.21	0.40
1:BA:848:G:C4	1:BA:933:A:H8	2.39	0.40
1:BA:855:G:H2'	1:BA:856:C:C6	2.57	0.40
1:BA:940:G:H2'	1:BA:941:A:O4'	2.22	0.40
1:BA:1109:C:H5'	1:BA:1110:G:P	2.61	0.40
1:BA:1955:U:O4	1:BA:2554:U:H5	2.03	0.40
1:BA:2122:U:O2'	1:BA:2123:G:OP1	2.34	0.40
3:BD:222:ARG:HH11	3:BD:224:ALA:HB3	1.87	0.40
7:BH:144:VAL:O	7:BH:148:ILE:HG13	2.21	0.40
9:BN:28:THR:HG22	9:BN:29:LYS:N	2.36	0.40
29:B7:8:ASN:OD1	29:B7:11:LYS:HB3	2.20	0.40
31:CA:49:U:O4	31:CA:365:U:H5	2.05	0.40
31:CA:68:G:C2	31:CA:69:G:C8	3.10	0.40
31:CA:189(D):C:C4	31:CA:189(E):U:C4	3.09	0.40
31:CA:193:C:C2	31:CA:194:C:C5	3.09	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:CA:547:A:OP2	34:CD:2:GLY:HA2	2.21	0.40
31:CA:584:G:H1	31:CA:757:U:H3	1.69	0.40
31:CA:899:C:H6	31:CA:899:C:O5'	2.03	0.40
31:CA:938:A:C6	31:CA:939:G:C5	3.09	0.40
31:CA:1202:G:C4	44:CN:42:ILE:HD13	2.56	0.40
31:CA:1278:U:H5'	31:CA:1279:A:O4'	2.21	0.40
32:CB:91:PRO:CD	32:CB:154:LEU:HD22	2.50	0.40
34:CD:106:TYR:HE1	34:CD:112:VAL:O	2.03	0.40
34:CD:134:ASP:O	34:CD:136:PRO:HD3	2.21	0.40
34:CD:158:ILE:O	34:CD:162:LEU:HB2	2.21	0.40
35:CE:69:VAL:HA	35:CE:70:PRO:HD3	1.74	0.40
48:CR:21:LYS:HA	48:CR:21:LYS:HD3	1.74	0.40
31:DA:23:C:OP2	31:DA:561:U:N3	2.53	0.40
31:DA:708:C:P	41:DK:85:ARG:HH22	2.45	0.40
31:DA:736:C:H4'	36:DF:89:MET:HE1	2.02	0.40
31:DA:972:C:H2'	40:DJ:55:LYS:HB2	2.03	0.40
31:DA:1003:G:H21	31:DA:1038:C:H42	1.68	0.40
31:DA:1032:G:C2'	31:DA:1033:G:H8	2.30	0.40
31:DA:1248:A:H2'	31:DA:1249:C:C6	2.56	0.40
31:DA:1380:U:O4'	31:DA:1381:U:H5	2.04	0.40
38:DH:81:HIS:ND1	38:DH:138:TRP:OXT	2.41	0.40
41:DK:29:ILE:HG23	41:DK:44:SER:HB3	2.02	0.40
41:DK:85:ARG:HG2	41:DK:111:ASP:O	2.21	0.40
49:DS:40:ILE:HA	49:DS:44:MET:SD	2.61	0.40
1:AA:199:A:N6	1:AA:2434:A:C5	2.89	0.40
1:AA:658:C:H2'	1:AA:659:C:C6	2.56	0.40
1:AA:745:G:H2'	1:AA:746:A:H5'	2.03	0.40
1:AA:754:C:H2'	1:AA:755:C:C6	2.57	0.40
1:AA:815:C:H2'	1:AA:816:C:C6	2.56	0.40
1:AA:996:A:C6	1:AA:1160:G:N1	2.89	0.40
1:AA:1515:G:H2'	1:AA:1516:C:C6	2.56	0.40
1:AA:1672:C:H5''	1:AA:2554:U:OP1	2.21	0.40
1:AA:1815:A:C5	1:AA:1817:G:C6	3.09	0.40
1:AA:2050:C:H2'	1:AA:2051:A:O4'	2.20	0.40
1:AA:2218:U:H4'	1:AA:2219:G:OP2	2.21	0.40
1:AA:2313:C:H2'	1:AA:2314:C:C6	2.56	0.40
1:AA:2630:G:O5'	1:AA:2630:G:H8	2.04	0.40
2:AB:28:C:N3	2:AB:29:A:C5	2.89	0.40
8:AI:128:LEU:HD23	8:AI:128:LEU:HA	1.96	0.40
12:AQ:1:MET:O	12:AQ:2:LEU:HB2	2.21	0.40
13:AR:96:ARG:HG3	13:AR:98:LEU:HG	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:AU:30:LYS:HA	16:AU:30:LYS:HD3	1.93	0.40
20:AY:23:ARG:HG3	20:AY:24:VAL:N	2.35	0.40
23:A1:40:ARG:HE	23:A1:40:ARG:HB2	1.74	0.40
1:BA:9:U:HO2'	1:BA:10:G:P	2.42	0.40
1:BA:413:C:H6	1:BA:413:C:O5'	2.04	0.40
1:BA:459:U:H5''	29:B7:40:TRP:CD2	2.56	0.40
1:BA:602:G:O2'	1:BA:655:A:N6	2.52	0.40
1:BA:699:A:H4'	1:BA:1634:A:N7	2.36	0.40
1:BA:764:A:H5'	3:BD:210:GLY:CA	2.51	0.40
1:BA:861:A:C2	1:BA:917:A:C4	3.09	0.40
1:BA:1487:G:H5''	1:BA:1488:G:OP2	2.21	0.40
1:BA:1658:C:OP1	4:BE:135:HIS:NE2	2.54	0.40
1:BA:1935:G:H1'	1:BA:1964:G:N2	2.37	0.40
1:BA:2126:A:C2	1:BA:2162:G:N3	2.90	0.40
1:BA:2228:G:C5	1:BA:2229:C:C4	3.10	0.40
1:BA:2302:G:C2	1:BA:2315:G:C2	3.10	0.40
6:BG:53:LEU:C	6:BG:55:LYS:H	2.25	0.40
8:BI:62:LYS:O	8:BI:66:GLU:HG2	2.21	0.40
15:BT:56:GLY:O	15:BT:59:THR:HG22	2.22	0.40
21:BZ:70:LEU:HD23	21:BZ:70:LEU:HA	1.74	0.40
24:B2:10:LEU:O	24:B2:14:ARG:HD2	2.22	0.40
26:B4:40:HIS:HB3	26:B4:43:TYR:CG	2.57	0.40
31:CA:73:G:H2'	31:CA:76:C:O4'	2.21	0.40
31:CA:108:G:N2	50:CT:12:ALA:HB1	2.36	0.40
31:CA:134:A:N6	46:CP:25:ARG:NH1	2.70	0.40
31:CA:302:G:N3	31:CA:556:C:H4'	2.36	0.40
31:CA:501:C:H2'	31:CA:502:G:H8	1.86	0.40
31:CA:626:U:H5''	46:CP:38:TYR:CG	2.56	0.40
31:CA:1225:A:OP2	31:CA:1226:C:H5	2.04	0.40
31:CA:1369:C:H2'	31:CA:1370:G:C8	2.57	0.40
31:CA:1446:U:O2'	31:CA:1447:A:C8	2.74	0.40
32:CB:174:VAL:O	32:CB:178:ARG:HB2	2.22	0.40
32:CB:178:ARG:HH22	38:CH:68:ARG:NH1	2.17	0.40
34:CD:28:SER:OG	34:CD:30:LYS:N	2.54	0.40
34:CD:95:GLY:O	34:CD:98:GLU:N	2.54	0.40
34:CD:149:ALA:HB3	34:CD:152:SER:OG	2.21	0.40
36:CF:4:TYR:HD1	36:CF:92:LYS:HA	1.86	0.40
43:CM:4:ILE:HG12	43:CM:5:ALA:N	2.36	0.40
46:CP:17:TYR:CD1	46:CP:17:TYR:N	2.89	0.40
50:CT:24:LEU:HA	50:CT:24:LEU:HD13	1.80	0.40
31:DA:192:U:H2'	31:DA:193:C:C6	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:DA:230:G:H2'	31:DA:231:G:O4'	2.21	0.40
31:DA:355:C:H1'	31:DA:388:G:N3	2.36	0.40
31:DA:552:U:H4'	42:DL:86:ARG:HG2	2.03	0.40
31:DA:688:G:H2'	31:DA:689:C:C6	2.56	0.40
31:DA:696:A:O5'	31:DA:696:A:H8	2.05	0.40
31:DA:707:C:H5''	41:DK:85:ARG:NH1	2.36	0.40
31:DA:825:G:H21	38:DH:11:THR:HG21	1.84	0.40
31:DA:1116:C:O2'	39:DI:108:VAL:HG21	2.21	0.40
35:DE:33:VAL:HG13	35:DE:112:LEU:HD12	2.03	0.40
37:DG:26:PHE:O	37:DG:30:ILE:HG13	2.22	0.40
37:DG:27:ILE:HD13	37:DG:40:ALA:HA	2.03	0.40
37:DG:150:ALA:HA	41:DK:59:TYR:CB	2.49	0.40
39:DI:46:ALA:C	39:DI:49:PRO:HD2	2.41	0.40
40:DJ:13:HIS:CD2	40:DJ:14:LYS:N	2.89	0.40
45:DO:69:TYR:O	45:DO:73:GLU:HB2	2.21	0.40
1:AA:266:G:N2	1:AA:427:U:H1'	2.35	0.40
1:AA:465:G:H2'	1:AA:466:A:C8	2.56	0.40
1:AA:528:A:C8	1:AA:528:A:C3'	3.05	0.40
1:AA:588:U:H1'	5:AF:90:PHE:HB3	2.04	0.40
1:AA:817:C:H2'	1:AA:818:G:H8	1.86	0.40
1:AA:827:U:H4'	1:AA:828:U:C5	2.57	0.40
1:AA:1354:A:C8	1:AA:1355:G:C8	3.09	0.40
1:AA:1355:G:N2	1:AA:1376:C:O2	2.32	0.40
1:AA:1410:G:C5	1:AA:1411:C:C5	3.09	0.40
1:AA:1429:G:H2'	1:AA:1430:C:H6	1.83	0.40
1:AA:1592:C:H2'	1:AA:1593:G:C8	2.56	0.40
1:AA:1700:A:H2'	1:AA:1701:A:O5'	2.22	0.40
1:AA:2031:A:C6	1:AA:2498:C:H1'	2.56	0.40
1:AA:2208:A:H1'	1:AA:2219:G:C4	2.57	0.40
1:AA:2537:U:H2'	1:AA:2538:C:C6	2.56	0.40
4:AE:170:LEU:HD23	4:AE:184:VAL:HG11	2.03	0.40
7:AH:17:VAL:HG21	7:AH:50:VAL:HG21	2.02	0.40
7:AH:35:VAL:O	7:AH:37:VAL:HG23	2.21	0.40
9:AN:20:GLY:HA2	9:AN:61:ARG:HH11	1.87	0.40
9:AN:38:HIS:NE2	9:AN:50:ASP:OD2	2.53	0.40
14:AS:58:LEU:HB2	14:AS:59:LYS:HG3	2.03	0.40
15:AT:22:PHE:HA	15:AT:91:ARG:NH1	2.37	0.40
28:A6:20:ASN:ND2	28:A6:40:CYS:HB2	2.37	0.40
1:BA:724:U:H2'	1:BA:725:G:O4'	2.21	0.40
1:BA:804:A:H2'	1:BA:806:C:C4	2.57	0.40
1:BA:892:G:H2'	1:BA:893:C:H6	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:1109:C:C5	1:BA:1110:G:C6	3.00	0.40
1:BA:1186:G:H8	1:BA:1186:G:O5'	2.04	0.40
1:BA:1482:G:O6	1:BA:1507:A:N6	2.54	0.40
1:BA:1833:U:O2'	1:BA:1969:A:N1	2.39	0.40
1:BA:1840:G:N2	1:BA:1902:C:O2	2.54	0.40
1:BA:2152:G:H2'	1:BA:2153:G:H8	1.86	0.40
1:BA:2313:C:H4'	6:BG:91:ARG:HG3	2.03	0.40
1:BA:2351:G:O6	30:B8:39:LYS:HG3	2.21	0.40
1:BA:2404:C:N3	1:BA:2414:G:C2	2.89	0.40
1:BA:2766:G:H2'	1:BA:2766:G:N3	2.35	0.40
3:BD:70:TRP:HB3	3:BD:190:TYR:CZ	2.56	0.40
6:BG:34:LEU:HD23	6:BG:34:LEU:HA	1.85	0.40
9:BN:36:GLY:O	9:BN:39:ARG:HB2	2.21	0.40
14:BS:92:TYR:HB3	14:BS:98:VAL:HG21	2.02	0.40
15:BT:11:GLU:OE1	15:BT:57:PHE:HB3	2.21	0.40
31:CA:59:A:N3	31:CA:59:A:H2'	2.36	0.40
31:CA:179:A:H2'	31:CA:180:U:C6	2.57	0.40
31:CA:189(D):C:C5	31:CA:189(E):U:C4	3.09	0.40
31:CA:246:A:C2	31:CA:282:A:C5	3.09	0.40
31:CA:575:G:H4'	31:CA:576:G:O5'	2.22	0.40
31:CA:890:G:O2'	31:CA:906:G:O6	2.25	0.40
31:CA:1011:G:H1	31:CA:1018:C:H42	1.69	0.40
31:CA:1030(D):A:N6	31:CA:1031:G:H21	2.11	0.40
31:CA:1097:C:O2'	31:CA:1169:A:N3	2.48	0.40
31:CA:1237:C:H3'	31:CA:1336:C:H41	1.86	0.40
31:CA:1363(A):A:H4'	31:CA:1364:U:H2'	2.04	0.40
32:CB:112:VAL:C	32:CB:114:ARG:H	2.24	0.40
32:CB:187:LEU:HD13	32:CB:205:ASP:HA	2.04	0.40
33:CC:177:THR:HG22	33:CC:179:ARG:H	1.86	0.40
36:CF:11:ASN:HA	36:CF:12:PRO:HD2	1.90	0.40
45:CO:57:LEU:HD23	45:CO:57:LEU:HA	1.96	0.40
31:DA:66:G:N2	31:DA:172:A:N3	2.70	0.40
31:DA:91:C:C2'	31:DA:92:C:H5'	2.50	0.40
31:DA:191:G:C6	31:DA:192:U:C4	3.09	0.40
31:DA:284:G:H2'	31:DA:285:G:H8	1.86	0.40
31:DA:323:U:H2'	31:DA:324:G:O4'	2.21	0.40
31:DA:409:G:N2	31:DA:434:U:C2	2.90	0.40
31:DA:540:G:H2'	31:DA:541:G:C8	2.56	0.40
31:DA:695:A:H2	31:DA:787:A:HO2'	1.68	0.40
32:DB:94:ASN:HB3	32:DB:95:GLN:NE2	2.37	0.40
33:DC:6:HIS:CE1	33:DC:184:TYR:CE2	3.09	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DD:65:ARG:HD3	34:DD:72:GLU:HA	2.03	0.40
39:DI:37:PHE:HB3	39:DI:43:ALA:CB	2.51	0.40
40:DJ:38:ILE:HA	40:DJ:39:PRO:HD3	1.73	0.40
45:DO:26:GLU:H	45:DO:26:GLU:HG2	1.64	0.40
47:DQ:31:LEU:HD23	47:DQ:32:TYR:CZ	2.57	0.40
50:DT:59:ALA:O	50:DT:63:ILE:HG13	2.22	0.40
1:AA:528:A:O2'	1:AA:529:A:H5'	2.21	0.40
1:AA:777:A:H2'	1:AA:778:G:C8	2.56	0.40
1:AA:974:G:N2	1:AA:989:G:H1'	2.36	0.40
1:AA:1637:A:C6	1:AA:1638:C:C4	3.09	0.40
1:AA:1799:G:N2	1:AA:1819:A:OP2	2.47	0.40
1:AA:2103:C:O2	1:AA:2187:G:C2	2.74	0.40
1:AA:2240:C:H2'	1:AA:2241:A:H8	1.85	0.40
1:AA:2489:G:C5	1:AA:2490:G:C6	3.09	0.40
2:AB:8:U:H5'	2:AB:9:G:OP2	2.21	0.40
3:AD:130:ALA:C	3:AD:131:LEU:HD12	2.42	0.40
5:AF:123:LEU:HD12	5:AF:124:LEU:N	2.36	0.40
6:AG:131:TYR:HE2	6:AG:133:LEU:HD23	1.86	0.40
7:AH:16:SER:OG	7:AH:27:LYS:HB2	2.21	0.40
8:AI:77:LEU:HB2	8:AI:142:VAL:HG12	2.04	0.40
9:AN:46:VAL:O	9:AN:47:ALA:HB3	2.21	0.40
9:AN:58:ASP:OD2	9:AN:59:LYS:HE2	2.22	0.40
10:AO:29:ASN:N	10:AO:29:ASN:OD1	2.55	0.40
19:AX:11:PRO:HD3	24:A2:37:PHE:CE2	2.56	0.40
20:AY:28:LYS:CG	20:AY:40:GLU:HG2	2.48	0.40
21:AZ:70:LEU:HD23	21:AZ:70:LEU:HA	1.91	0.40
21:AZ:133:ILE:HA	21:AZ:134:PRO:HD2	1.98	0.40
25:A3:58:VAL:O	1:BA:652(Q):G:H5''	2.22	0.40
1:BA:271(I):G:H2'	1:BA:271(J):C:C6	2.57	0.40
1:BA:1268:A:C2	1:BA:2013:A:C4	3.10	0.40
1:BA:1385:G:H1'	1:BA:1386:C:C6	2.57	0.40
1:BA:1897:G:H2'	1:BA:1898:U:O4'	2.22	0.40
1:BA:2834:G:H5''	1:BA:2834:G:C8	2.55	0.40
3:BD:77:ALA:HB2	3:BD:97:TYR:CG	2.56	0.40
4:BE:73:GLU:HA	4:BE:74:PRO:HD3	1.82	0.40
9:BN:15:LEU:HB3	9:BN:137:LYS:HA	2.03	0.40
9:BN:46:VAL:O	9:BN:47:ALA:HB3	2.21	0.40
12:BQ:30:GLY:HA2	12:BQ:107:ALA:HB2	2.04	0.40
16:BU:108:GLU:OE1	16:BU:112:ARG:NH1	2.54	0.40
18:BW:13:SER:HA	18:BW:14:PRO:HD3	1.88	0.40
20:BY:5:MET:HE2	20:BY:5:MET:HB2	1.71	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:BY:28:LYS:HG3	20:BY:40:GLU:HG2	2.03	0.40
23:B1:3:LYS:HE3	23:B1:4:VAL:HG13	2.03	0.40
31:CA:56:U:H2'	31:CA:57:G:H8	1.86	0.40
31:CA:189:G:H2'	31:CA:189(A):C:O4'	2.21	0.40
31:CA:361:G:H2'	31:CA:362:G:O4'	2.21	0.40
31:CA:546:G:OP1	34:CD:73:ARG:HB2	2.21	0.40
31:CA:592:G:N1	31:CA:648:A:C6	2.90	0.40
31:CA:689:C:H2'	31:CA:690:G:O4'	2.21	0.40
31:CA:1108:G:OP2	33:CC:174:PRO:HA	2.20	0.40
31:CA:1389:C:H2'	31:CA:1390:U:O4'	2.21	0.40
31:CA:1410:G:H2'	31:CA:1411:C:C6	2.57	0.40
33:CC:58:GLU:O	33:CC:59:ARG:HD2	2.22	0.40
42:CL:30:ALA:HA	42:CL:31:PRO:HD3	1.92	0.40
48:CR:56:THR:CB	48:CR:58:LEU:HD13	2.47	0.40
31:DA:163:C:H2'	31:DA:164:U:C6	2.57	0.40
31:DA:336:C:H2'	31:DA:337:C:C6	2.56	0.40
31:DA:1054:C:H2'	31:DA:1055:A:H5''	2.03	0.40
31:DA:1127:G:H2'	31:DA:1128:C:O4'	2.21	0.40
31:DA:1157:A:C2	31:DA:1181:G:H1'	2.56	0.40
31:DA:1202:G:N9	44:DN:42:ILE:HD12	2.37	0.40
31:DA:1213:A:C4	31:DA:1215:G:C8	3.10	0.40
32:DB:112:VAL:C	32:DB:114:ARG:H	2.25	0.40
33:DC:19:GLU:HB2	33:DC:40:ARG:NH2	2.36	0.40
34:DD:50:ARG:HA	34:DD:51:PRO:HD3	1.88	0.40
34:DD:52:SER:O	34:DD:56:VAL:HG23	2.21	0.40
36:DF:12:PRO:O	36:DF:14:LEU:N	2.48	0.40
36:DF:15:ASP:O	36:DF:19:LEU:HB2	2.21	0.40
39:DI:8:GLY:HA3	39:DI:76:ALA:O	2.21	0.40
39:DI:99:LEU:HB3	39:DI:101:PHE:CE1	2.57	0.40
47:DQ:96:GLU:O	47:DQ:97:SER:HB2	2.21	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AI:89:TYR:O	31:CA:357:G:O2'[2_654]	2.11	0.09
8:AI:91:SER:OG	31:CA:368:U:OP2[2_654]	2.12	0.08

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	AD	273/276 (99%)	240 (88%)	27 (10%)	6 (2%)	5	30
3	BD	273/276 (99%)	246 (90%)	23 (8%)	4 (2%)	8	38
4	AE	202/206 (98%)	177 (88%)	18 (9%)	7 (4%)	3	20
4	BE	202/206 (98%)	174 (86%)	21 (10%)	7 (4%)	3	20
5	AF	198/205 (97%)	168 (85%)	25 (13%)	5 (2%)	4	28
5	BF	198/205 (97%)	170 (86%)	21 (11%)	7 (4%)	3	20
6	AG	179/182 (98%)	136 (76%)	33 (18%)	10 (6%)	1	11
6	BG	179/182 (98%)	135 (75%)	30 (17%)	14 (8%)	1	5
7	AH	172/180 (96%)	143 (83%)	21 (12%)	8 (5%)	2	14
7	BH	172/180 (96%)	144 (84%)	22 (13%)	6 (4%)	3	20
8	AI	143/148 (97%)	103 (72%)	28 (20%)	12 (8%)	0	4
8	BI	143/148 (97%)	109 (76%)	24 (17%)	10 (7%)	1	6
9	AN	138/140 (99%)	113 (82%)	16 (12%)	9 (6%)	1	8
9	BN	138/140 (99%)	119 (86%)	13 (9%)	6 (4%)	2	16
10	AO	120/122 (98%)	108 (90%)	8 (7%)	4 (3%)	3	21
10	BO	120/122 (98%)	109 (91%)	7 (6%)	4 (3%)	3	21
11	AP	143/150 (95%)	117 (82%)	18 (13%)	8 (6%)	1	11
11	BP	143/150 (95%)	126 (88%)	12 (8%)	5 (4%)	3	20
12	AQ	139/141 (99%)	126 (91%)	9 (6%)	4 (3%)	3	24
12	BQ	139/141 (99%)	126 (91%)	9 (6%)	4 (3%)	3	24
13	AR	116/118 (98%)	95 (82%)	16 (14%)	5 (4%)	2	16
13	BR	116/118 (98%)	108 (93%)	7 (6%)	1 (1%)	14	49
14	AS	108/112 (96%)	84 (78%)	21 (19%)	3 (3%)	4	25
14	BS	108/112 (96%)	93 (86%)	12 (11%)	3 (3%)	4	25
15	AT	129/146 (88%)	109 (84%)	16 (12%)	4 (3%)	3	22

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
15	BT	129/146 (88%)	116 (90%)	12 (9%)	1 (1%)	16	51
16	AU	114/118 (97%)	104 (91%)	10 (9%)	0	100	100
16	BU	114/118 (97%)	111 (97%)	3 (3%)	0	100	100
17	AV	99/101 (98%)	89 (90%)	8 (8%)	2 (2%)	6	32
17	BV	99/101 (98%)	89 (90%)	9 (9%)	1 (1%)	13	47
18	AW	110/113 (97%)	101 (92%)	8 (7%)	1 (1%)	14	49
18	BW	110/113 (97%)	104 (94%)	6 (6%)	0	100	100
19	AX	93/96 (97%)	82 (88%)	9 (10%)	2 (2%)	5	30
19	BX	93/96 (97%)	82 (88%)	9 (10%)	2 (2%)	5	30
20	AY	105/110 (96%)	90 (86%)	10 (10%)	5 (5%)	2	14
20	BY	105/110 (96%)	88 (84%)	13 (12%)	4 (4%)	2	18
21	AZ	196/206 (95%)	153 (78%)	32 (16%)	11 (6%)	1	11
21	BZ	196/206 (95%)	158 (81%)	31 (16%)	7 (4%)	3	20
22	A0	74/85 (87%)	67 (90%)	6 (8%)	1 (1%)	9	40
22	B0	74/85 (87%)	67 (90%)	6 (8%)	1 (1%)	9	40
23	A1	95/98 (97%)	88 (93%)	5 (5%)	2 (2%)	5	31
23	B1	95/98 (97%)	86 (90%)	6 (6%)	3 (3%)	3	22
24	A2	68/72 (94%)	59 (87%)	8 (12%)	1 (2%)	8	38
24	B2	68/72 (94%)	63 (93%)	5 (7%)	0	100	100
25	A3	57/60 (95%)	54 (95%)	3 (5%)	0	100	100
25	B3	57/60 (95%)	54 (95%)	2 (4%)	1 (2%)	7	35
26	A4	44/71 (62%)	30 (68%)	10 (23%)	4 (9%)	0	3
26	B4	44/71 (62%)	30 (68%)	10 (23%)	4 (9%)	0	3
27	A5	57/60 (95%)	50 (88%)	6 (10%)	1 (2%)	7	35
27	B5	57/60 (95%)	49 (86%)	6 (10%)	2 (4%)	3	20
28	A6	51/54 (94%)	46 (90%)	4 (8%)	1 (2%)	6	32
28	B6	51/54 (94%)	47 (92%)	4 (8%)	0	100	100
29	A7	46/49 (94%)	41 (89%)	4 (9%)	1 (2%)	5	30
29	B7	46/49 (94%)	42 (91%)	3 (6%)	1 (2%)	5	30
30	A8	62/65 (95%)	48 (77%)	13 (21%)	1 (2%)	8	37
30	B8	62/65 (95%)	57 (92%)	5 (8%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
32	CB	227/256 (89%)	171 (75%)	36 (16%)	20 (9%)	0	3
32	DB	227/256 (89%)	173 (76%)	39 (17%)	15 (7%)	1	7
33	CC	204/239 (85%)	163 (80%)	30 (15%)	11 (5%)	1	12
33	DC	204/239 (85%)	144 (71%)	42 (21%)	18 (9%)	0	3
34	CD	206/209 (99%)	154 (75%)	40 (19%)	12 (6%)	1	10
34	DD	206/209 (99%)	152 (74%)	47 (23%)	7 (3%)	3	21
35	CE	146/162 (90%)	112 (77%)	23 (16%)	11 (8%)	1	6
35	DE	146/162 (90%)	116 (80%)	25 (17%)	5 (3%)	3	21
36	CF	98/101 (97%)	87 (89%)	8 (8%)	3 (3%)	3	22
36	DF	98/101 (97%)	85 (87%)	9 (9%)	4 (4%)	2	17
37	CG	153/156 (98%)	127 (83%)	16 (10%)	10 (6%)	1	8
37	DG	153/156 (98%)	129 (84%)	18 (12%)	6 (4%)	2	18
38	CH	136/138 (99%)	116 (85%)	18 (13%)	2 (2%)	8	38
38	DH	136/138 (99%)	118 (87%)	13 (10%)	5 (4%)	2	19
39	CI	123/128 (96%)	93 (76%)	20 (16%)	10 (8%)	1	4
39	DI	123/128 (96%)	94 (76%)	21 (17%)	8 (6%)	1	8
40	CJ	94/105 (90%)	66 (70%)	19 (20%)	9 (10%)	0	3
40	DJ	94/105 (90%)	74 (79%)	14 (15%)	6 (6%)	1	8
41	CK	112/129 (87%)	96 (86%)	12 (11%)	4 (4%)	3	20
41	DK	112/129 (87%)	90 (80%)	19 (17%)	3 (3%)	4	26
42	CL	120/132 (91%)	98 (82%)	17 (14%)	5 (4%)	2	17
42	DL	120/132 (91%)	100 (83%)	15 (12%)	5 (4%)	2	17
43	CM	112/126 (89%)	84 (75%)	18 (16%)	10 (9%)	0	3
43	DM	112/126 (89%)	80 (71%)	20 (18%)	12 (11%)	0	2
44	CN	58/61 (95%)	43 (74%)	10 (17%)	5 (9%)	0	3
44	DN	58/61 (95%)	48 (83%)	7 (12%)	3 (5%)	1	12
45	CO	86/89 (97%)	71 (83%)	12 (14%)	3 (4%)	3	20
45	DO	86/89 (97%)	68 (79%)	16 (19%)	2 (2%)	5	29
46	CP	80/88 (91%)	50 (62%)	23 (29%)	7 (9%)	0	3
46	DP	80/88 (91%)	52 (65%)	25 (31%)	3 (4%)	2	18
47	CQ	97/105 (92%)	84 (87%)	8 (8%)	5 (5%)	1	12

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
47	DQ	97/105 (92%)	81 (84%)	13 (13%)	3 (3%)	3	22
48	CR	66/88 (75%)	55 (83%)	10 (15%)	1 (2%)	8	38
48	DR	66/88 (75%)	60 (91%)	6 (9%)	0	100	100
49	CS	76/93 (82%)	48 (63%)	17 (22%)	11 (14%)	0	1
49	DS	76/93 (82%)	54 (71%)	17 (22%)	5 (7%)	1	7
50	CT	94/106 (89%)	73 (78%)	13 (14%)	8 (8%)	0	3
50	DT	94/106 (89%)	72 (77%)	15 (16%)	7 (7%)	1	6
51	CU	21/27 (78%)	18 (86%)	3 (14%)	0	100	100
51	DU	21/27 (78%)	16 (76%)	3 (14%)	2 (10%)	0	3
All	All	11280/12044 (94%)	9338 (83%)	1460 (13%)	482 (4%)	2	16

All (482) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	AD	239	ARG
3	AD	275	LYS
5	AF	60	SER
6	AG	14	GLU
6	AG	78	SER
7	AH	47	GLU
7	AH	126	PRO
8	AI	10	GLU
8	AI	104	GLN
8	AI	106	GLY
8	AI	111	PRO
9	AN	18	ALA
10	AO	48	PRO
11	AP	102	ARG
11	AP	103	ALA
13	AR	14	SER
14	AS	62	LYS
14	AS	83	LYS
17	AV	43	GLU
19	AX	94	GLY
20	AY	2	ARG
20	AY	91	GLU
21	AZ	161	VAL
23	A1	3	LYS
29	A7	46	VAL

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Mol	Chain	Res	Type
3	BD	156	ALA
3	BD	239	ARG
3	BD	275	LYS
6	BG	49	ASP
7	BH	126	PRO
7	BH	170	ARG
8	BI	10	GLU
8	BI	111	PRO
9	BN	76	SER
10	BO	48	PRO
11	BP	103	ALA
12	BQ	2	LEU
13	BR	2	ARG
14	BS	83	LYS
20	BY	2	ARG
21	BZ	158	PRO
21	BZ	159	PRO
21	BZ	161	VAL
23	B1	3	LYS
26	B4	28	LYS
29	B7	46	VAL
32	CB	15	VAL
32	CB	26	PRO
32	CB	83	MET
32	CB	132	LYS
32	CB	194	PRO
32	CB	195	ASP
33	CC	156	ARG
33	CC	190	ARG
34	CD	14	ARG
34	CD	32	ALA
34	CD	42	GLN
35	CE	72	GLN
35	CE	73	ASN
35	CE	98	THR
36	CF	40	VAL
37	CG	56	GLN
39	CI	43	ALA
39	CI	103	THR
40	CJ	56	HIS
40	CJ	90	LEU
40	CJ	91	PRO

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Mol	Chain	Res	Type
41	CK	49	GLY
42	CL	27	LEU
43	CM	49	THR
44	CN	59	ALA
45	CO	19	PRO
46	CP	53	VAL
46	CP	73	LEU
49	CS	30	LEU
49	CS	40	ILE
49	CS	45	VAL
49	CS	73	GLU
50	CT	9	ASN
50	CT	71	THR
32	DB	15	VAL
32	DB	24	TRP
32	DB	26	PRO
32	DB	75	LYS
32	DB	120	ALA
32	DB	169	LYS
32	DB	194	PRO
32	DB	195	ASP
33	DC	3	ASN
33	DC	49	SER
33	DC	156	ARG
33	DC	181	ASN
34	DD	14	ARG
34	DD	42	GLN
35	DE	73	ASN
35	DE	98	THR
36	DF	40	VAL
37	DG	16	LEU
37	DG	35	LYS
39	DI	45	ALA
40	DJ	34	VAL
40	DJ	56	HIS
40	DJ	90	LEU
41	DK	49	GLY
42	DL	27	LEU
42	DL	91	LYS
43	DM	7	VAL
43	DM	10	PRO
43	DM	45	VAL

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Mol	Chain	Res	Type
43	DM	49	THR
43	DM	66	LEU
43	DM	84	ILE
44	DN	41	ARG
44	DN	59	ALA
45	DO	19	PRO
46	DP	53	VAL
47	DQ	97	SER
50	DT	9	ASN
50	DT	71	THR
51	DU	7	ARG
4	AE	69	LYS
5	AF	17	ARG
6	AG	49	ASP
6	AG	51	ARG
6	AG	117	PHE
6	AG	150	ASP
7	AH	65	HIS
7	AH	92	ILE
7	AH	108	GLY
7	AH	170	ARG
8	AI	68	LEU
8	AI	98	ALA
8	AI	107	VAL
8	AI	112	LYS
9	AN	2	LYS
9	AN	5	VAL
9	AN	19	GLU
10	AO	5	GLN
11	AP	27	HIS
11	AP	39	LYS
12	AQ	2	LEU
12	AQ	62	GLY
13	AR	2	ARG
13	AR	11	ASN
13	AR	86	ARG
17	AV	79	VAL
20	AY	92	ASN
21	AZ	158	PRO
21	AZ	159	PRO
21	AZ	191	VAL
26	A4	39	CYS

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Mol	Chain	Res	Type
27	A5	35	GLU
3	BD	3	VAL
4	BE	71	GLY
5	BF	7	TYR
5	BF	15	SER
5	BF	89	VAL
5	BF	168	ARG
6	BG	14	GLU
6	BG	51	ARG
6	BG	115	ARG
8	BI	42	SER
8	BI	85	GLU
8	BI	107	VAL
8	BI	112	LYS
9	BN	5	VAL
9	BN	18	ALA
9	BN	127	ASP
11	BP	27	HIS
11	BP	29	LYS
11	BP	39	LYS
11	BP	102	ARG
12	BQ	62	GLY
15	BT	126	ALA
17	BV	79	VAL
20	BY	91	GLU
26	B4	39	CYS
32	CB	10	LEU
32	CB	76	GLN
32	CB	165	VAL
33	CC	50	ALA
33	CC	52	LEU
33	CC	75	VAL
35	CE	71	LEU
35	CE	146	ALA
37	CG	6	ARG
37	CG	69	VAL
38	CH	51	VAL
38	CH	133	LEU
39	CI	54	ASP
39	CI	118	LYS
40	CJ	55	LYS
40	CJ	57	LYS

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Mol	Chain	Res	Type
40	CJ	89	ASP
42	CL	28	LYS
42	CL	91	LYS
43	CM	12	ASN
43	CM	84	ILE
45	CO	21	ASP
47	CQ	14	LYS
47	CQ	97	SER
49	CS	26	GLY
49	CS	81	ARG
50	CT	10	LEU
50	CT	94	ALA
32	DB	83	MET
32	DB	134	GLU
33	DC	100	ALA
33	DC	175	LEU
33	DC	206	GLU
34	DD	32	ALA
34	DD	129	ASN
35	DE	71	LEU
35	DE	146	ALA
37	DG	81	GLY
38	DH	50	ARG
38	DH	51	VAL
38	DH	133	LEU
39	DI	121	ARG
40	DJ	77	PRO
41	DK	45	GLY
41	DK	87	THR
42	DL	28	LYS
43	DM	3	ARG
43	DM	69	GLU
46	DP	24	ALA
46	DP	78	GLY
47	DQ	68	ARG
49	DS	24	ALA
49	DS	26	GLY
50	DT	10	LEU
50	DT	100	ILE
5	AF	7	TYR
6	AG	80	PHE
6	AG	115	ARG

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Mol	Chain	Res	Type
8	AI	85	GLU
9	AN	63	THR
10	AO	26	LYS
10	AO	104	ARG
11	AP	101	VAL
11	AP	122	PRO
12	AQ	59	ARG
12	AQ	135	ASP
13	AR	107	ASP
15	AT	127	ALA
20	AY	54	LYS
21	AZ	13	GLU
21	AZ	120	ILE
4	BE	52	LEU
5	BF	130	ALA
6	BG	78	SER
6	BG	117	PHE
6	BG	138	GLN
8	BI	86	THR
8	BI	104	GLN
9	BN	19	GLU
10	BO	5	GLN
12	BQ	59	ARG
14	BS	62	LYS
20	BY	92	ASN
21	BZ	151	HIS
21	BZ	152	ALA
23	B1	83	GLU
26	B4	36	CYS
26	B4	43	TYR
27	B5	37	LYS
32	CB	11	LEU
32	CB	78	GLN
32	CB	169	LYS
33	CC	24	ALA
33	CC	53	ALA
33	CC	60	ALA
34	CD	124	GLY
34	CD	186	LEU
34	CD	189	PRO
35	CE	8	GLU
35	CE	68	GLU

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Mol	Chain	Res	Type
37	CG	7	ALA
37	CG	33	ASP
37	CG	146	GLU
39	CI	11	LYS
39	CI	23	ASN
39	CI	24	GLY
39	CI	102	LEU
39	CI	120	ARG
43	CM	10	PRO
43	CM	40	ASN
43	CM	66	LEU
43	CM	67	GLU
44	CN	26	ARG
47	CQ	49	GLU
47	CQ	68	ARG
48	CR	32	ARG
49	CS	12	ASP
33	DC	12	LEU
33	DC	26	LYS
34	DD	136	PRO
36	DF	13	ASN
37	DG	7	ALA
38	DH	41	ARG
39	DI	23	ASN
39	DI	90	PRO
40	DJ	39	PRO
43	DM	12	ASN
49	DS	25	LYS
50	DT	102	GLY
3	AD	3	VAL
3	AD	30	GLU
3	AD	156	ALA
3	AD	233	HIS
4	AE	52	LEU
4	AE	70	ALA
9	AN	64	GLY
9	AN	77	GLY
9	AN	127	ASP
11	AP	29	LYS
11	AP	140	ALA
15	AT	36	GLU
15	AT	84	GLN

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Mol	Chain	Res	Type
20	AY	90	LEU
21	AZ	152	ALA
22	A0	44	ARG
23	A1	83	GLU
26	A4	41	PRO
30	A8	11	LYS
4	BE	72	VAL
5	BF	18	ARG
6	BG	43	LEU
6	BG	54	GLU
6	BG	86	MET
6	BG	124	SER
6	BG	150	ASP
6	BG	151	ALA
7	BH	70	THR
8	BI	109	ILE
9	BN	77	GLY
10	BO	29	ASN
14	BS	57	LYS
19	BX	15	GLU
20	BY	54	LYS
22	B0	55	ARG
23	B1	45	ASN
32	CB	34	ALA
32	CB	52	GLU
32	CB	120	ALA
32	CB	131	PRO
32	CB	202	PRO
33	CC	108	ASN
34	CD	159	ARG
34	CD	166	LYS
35	CE	140	ARG
35	CE	148	VAL
37	CG	104	LEU
37	CG	145	ALA
37	CG	147	ALA
40	CJ	80	LYS
41	CK	106	LYS
43	CM	7	VAL
44	CN	14	PRO
46	CP	67	THR
46	CP	72	ARG

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Mol	Chain	Res	Type
47	CQ	34	LYS
49	CS	14	HIS
32	DB	76	GLN
32	DB	131	PRO
32	DB	165	VAL
33	DC	45	LYS
33	DC	61	ALA
33	DC	101	LEU
34	DD	186	LEU
36	DF	14	LEU
39	DI	102	LEU
43	DM	40	ASN
44	DN	14	PRO
47	DQ	96	GLU
49	DS	40	ILE
49	DS	47	HIS
4	AE	73	GLU
4	AE	94	GLU
5	AF	10	PRO
5	AF	141	ALA
6	AG	43	LEU
7	AH	145	ALA
8	AI	87	LYS
8	AI	114	LEU
15	AT	100	TYR
18	AW	58	ALA
21	AZ	101	PRO
21	AZ	157	LEU
24	A2	6	VAL
28	A6	44	ARG
4	BE	69	LYS
4	BE	118	LYS
6	BG	30	GLU
7	BH	159	GLU
8	BI	117	GLU
10	BO	26	LYS
21	BZ	157	LEU
21	BZ	191	VAL
32	CB	227	GLY
33	CC	127	ARG
34	CD	107	ARG
34	CD	136	PRO

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Mol	Chain	Res	Type
39	CI	45	ALA
40	CJ	77	PRO
41	CK	62	GLN
42	CL	9	GLN
42	CL	26	ALA
43	CM	70	LEU
43	CM	87	TYR
44	CN	15	LYS
45	CO	88	ARG
46	CP	76	GLN
49	CS	72	GLY
49	CS	74	PHE
33	DC	24	ALA
33	DC	99	VAL
33	DC	107	GLN
33	DC	117	ALA
35	DE	85	GLY
37	DG	56	GLN
38	DH	86	ILE
40	DJ	57	LYS
42	DL	19	ARG
45	DO	88	ARG
4	AE	72	VAL
6	AG	87	PRO
9	AN	76	SER
19	AX	40	LYS
21	AZ	93	ASP
26	A4	28	LYS
4	BE	73	GLU
7	BH	108	GLY
27	B5	36	CYS
33	CC	98	ASN
35	CE	49	PRO
35	CE	85	GLY
36	CF	13	ASN
37	CG	52	GLU
40	CJ	78	ASN
44	CN	60	SER
50	CT	102	GLY
39	DI	54	ASP
39	DI	103	THR
43	DM	85	GLY

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Mol	Chain	Res	Type
43	DM	99	ARG
50	DT	94	ALA
50	DT	98	PRO
8	AI	109	ILE
14	AS	104	GLY
5	BF	66	PRO
32	CB	80	ILE
34	CD	28	SER
34	CD	142	PRO
39	DI	24	GLY
46	CP	36	ILE
46	CP	78	GLY
50	CT	96	GLY
33	DC	81	GLY
4	AE	71	GLY
7	AH	76	VAL
26	A4	29	PRO
19	BX	94	GLY
25	B3	13	ILE
32	CB	167	PRO
32	DB	202	PRO
32	DB	227	GLY
33	DC	108	ASN
34	DD	167	GLY
42	DL	90	VAL
21	AZ	193	GLU
36	CF	12	PRO
41	CK	45	GLY
49	CS	9	VAL
50	CT	98	PRO
50	CT	100	ILE
36	DF	12	PRO
37	DG	55	GLY
7	BH	92	ILE
12	BQ	99	PRO
51	DU	23	PRO
4	BE	30	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar

resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	AD	215/218 (99%)	185 (86%)	30 (14%)	3	14
3	BD	215/218 (99%)	189 (88%)	26 (12%)	4	19
4	AE	163/166 (98%)	139 (85%)	24 (15%)	2	12
4	BE	163/166 (98%)	134 (82%)	29 (18%)	1	7
5	AF	158/162 (98%)	133 (84%)	25 (16%)	2	10
5	BF	158/162 (98%)	138 (87%)	20 (13%)	3	17
6	AG	128/156 (82%)	110 (86%)	18 (14%)	3	14
6	BG	128/156 (82%)	114 (89%)	14 (11%)	5	23
7	AH	141/148 (95%)	125 (89%)	16 (11%)	4	22
7	BH	141/148 (95%)	124 (88%)	17 (12%)	4	19
8	AI	102/124 (82%)	76 (74%)	26 (26%)	0	2
8	BI	102/124 (82%)	76 (74%)	26 (26%)	0	2
9	AN	117/119 (98%)	94 (80%)	23 (20%)	1	6
9	BN	117/119 (98%)	91 (78%)	26 (22%)	1	4
10	AO	98/100 (98%)	89 (91%)	9 (9%)	7	29
10	BO	98/100 (98%)	89 (91%)	9 (9%)	7	29
11	AP	113/116 (97%)	96 (85%)	17 (15%)	2	12
11	BP	113/116 (97%)	97 (86%)	16 (14%)	2	13
12	AQ	111/111 (100%)	94 (85%)	17 (15%)	2	11
12	BQ	111/111 (100%)	94 (85%)	17 (15%)	2	11
13	AR	101/101 (100%)	80 (79%)	21 (21%)	1	5
13	BR	101/101 (100%)	79 (78%)	22 (22%)	1	4
14	AS	84/88 (96%)	67 (80%)	17 (20%)	1	5
14	BS	84/88 (96%)	72 (86%)	12 (14%)	2	13
15	AT	110/127 (87%)	99 (90%)	11 (10%)	6	26
15	BT	110/127 (87%)	100 (91%)	10 (9%)	7	29
16	AU	93/94 (99%)	83 (89%)	10 (11%)	5	23
16	BU	93/94 (99%)	77 (83%)	16 (17%)	1	8
17	AV	80/82 (98%)	63 (79%)	17 (21%)	1	5

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles		
17	BV	80/82 (98%)	63 (79%)	17 (21%)	1	5	
18	AW	89/92 (97%)	75 (84%)	14 (16%)	2	10	
18	BW	89/92 (97%)	78 (88%)	11 (12%)	4	18	
19	AX	75/78 (96%)	67 (89%)	8 (11%)	5	24	
19	BX	75/78 (96%)	66 (88%)	9 (12%)	4	19	
20	AY	80/91 (88%)	70 (88%)	10 (12%)	3	18	
20	BY	80/91 (88%)	72 (90%)	8 (10%)	6	26	
21	AZ	159/179 (89%)	139 (87%)	20 (13%)	3	18	
21	BZ	159/179 (89%)	137 (86%)	22 (14%)	3	14	
22	A0	59/67 (88%)	52 (88%)	7 (12%)	4	19	
22	B0	59/67 (88%)	52 (88%)	7 (12%)	4	19	
23	A1	78/83 (94%)	68 (87%)	10 (13%)	3	17	
23	B1	78/83 (94%)	70 (90%)	8 (10%)	6	25	
24	A2	65/67 (97%)	56 (86%)	9 (14%)	3	14	
24	B2	65/67 (97%)	57 (88%)	8 (12%)	4	19	
25	A3	49/52 (94%)	43 (88%)	6 (12%)	4	19	
25	B3	49/52 (94%)	43 (88%)	6 (12%)	4	19	
26	A4	39/63 (62%)	32 (82%)	7 (18%)	1	7	
26	B4	39/63 (62%)	30 (77%)	9 (23%)	0	3	
27	A5	50/52 (96%)	43 (86%)	7 (14%)	3	14	
27	B5	50/52 (96%)	43 (86%)	7 (14%)	3	14	
28	A6	50/52 (96%)	40 (80%)	10 (20%)	1	5	
28	B6	50/52 (96%)	41 (82%)	9 (18%)	1	7	
29	A7	41/42 (98%)	35 (85%)	6 (15%)	2	13	
29	B7	41/42 (98%)	32 (78%)	9 (22%)	1	4	
30	A8	52/55 (94%)	45 (86%)	7 (14%)	3	15	
30	B8	52/55 (94%)	47 (90%)	5 (10%)	7	28	
32	CB	177/220 (80%)	141 (80%)	36 (20%)	1	5	
32	DB	177/220 (80%)	142 (80%)	35 (20%)	1	6	
33	CC	114/188 (61%)	95 (83%)	19 (17%)	2	9	
33	DC	114/188 (61%)	96 (84%)	18 (16%)	2	10	

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
34	CD	139/181 (77%)	117 (84%)	22 (16%)	2	10
34	DD	139/181 (77%)	120 (86%)	19 (14%)	3	14
35	CE	108/123 (88%)	86 (80%)	22 (20%)	1	5
35	DE	108/123 (88%)	88 (82%)	20 (18%)	1	7
36	CF	77/90 (86%)	65 (84%)	12 (16%)	2	10
36	DF	77/90 (86%)	65 (84%)	12 (16%)	2	10
37	CG	104/127 (82%)	86 (83%)	18 (17%)	1	8
37	DG	104/127 (82%)	88 (85%)	16 (15%)	2	11
38	CH	103/119 (87%)	85 (82%)	18 (18%)	1	8
38	DH	103/119 (87%)	84 (82%)	19 (18%)	1	7
39	CI	62/99 (63%)	53 (86%)	9 (14%)	2	13
39	DI	62/99 (63%)	54 (87%)	8 (13%)	3	17
40	CJ	52/92 (56%)	42 (81%)	10 (19%)	1	6
40	DJ	52/92 (56%)	41 (79%)	11 (21%)	1	5
41	CK	81/99 (82%)	69 (85%)	12 (15%)	2	12
41	DK	81/99 (82%)	71 (88%)	10 (12%)	4	19
42	CL	92/109 (84%)	81 (88%)	11 (12%)	4	19
42	DL	92/109 (84%)	83 (90%)	9 (10%)	6	27
43	CM	63/101 (62%)	49 (78%)	14 (22%)	1	4
43	DM	63/101 (62%)	49 (78%)	14 (22%)	1	4
44	CN	46/50 (92%)	35 (76%)	11 (24%)	0	2
44	DN	46/50 (92%)	33 (72%)	13 (28%)	0	1
45	CO	77/80 (96%)	65 (84%)	12 (16%)	2	10
45	DO	77/80 (96%)	66 (86%)	11 (14%)	2	13
46	CP	63/74 (85%)	49 (78%)	14 (22%)	1	4
46	DP	63/74 (85%)	53 (84%)	10 (16%)	2	10
47	CQ	94/97 (97%)	90 (96%)	4 (4%)	25	57
47	DQ	94/97 (97%)	84 (89%)	10 (11%)	5	24
48	CR	49/77 (64%)	44 (90%)	5 (10%)	6	26
48	DR	49/77 (64%)	46 (94%)	3 (6%)	15	47
49	CS	42/80 (52%)	25 (60%)	17 (40%)	0	0

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
49	DS	42/80 (52%)	36 (86%)	6 (14%)	2	13
50	CT	65/82 (79%)	53 (82%)	12 (18%)	1	7
50	DT	65/82 (79%)	57 (88%)	8 (12%)	4	19
51	CU	18/22 (82%)	16 (89%)	2 (11%)	5	22
51	DU	18/22 (82%)	14 (78%)	4 (22%)	1	4
All	All	8652/9990 (87%)	7319 (85%)	1333 (15%)	2	11

All (1333) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	AD	4	LYS
3	AD	10	THR
3	AD	12	SER
3	AD	13	ARG
3	AD	54	ARG
3	AD	61	LEU
3	AD	89	SER
3	AD	94	LEU
3	AD	103	ARG
3	AD	106	ILE
3	AD	111	LEU
3	AD	141	VAL
3	AD	142	VAL
3	AD	150	LYS
3	AD	154	LYS
3	AD	155	LEU
3	AD	192	THR
3	AD	200	ASP
3	AD	208	LYS
3	AD	211	ARG
3	AD	212	SER
3	AD	217	ARG
3	AD	221	VAL
3	AD	229	VAL
3	AD	242	ARG
3	AD	253	GLN
3	AD	257	LEU
3	AD	259	THR
3	AD	260	ARG
3	AD	267	SER

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Mol	Chain	Res	Type
4	AE	1	MET
4	AE	7	VAL
4	AE	12	THR
4	AE	21	VAL
4	AE	33	VAL
4	AE	34	VAL
4	AE	45	THR
4	AE	49	LEU
4	AE	52	LEU
4	AE	75	VAL
4	AE	77	ILE
4	AE	82	ARG
4	AE	87	GLU
4	AE	93	VAL
4	AE	113	PHE
4	AE	116	VAL
4	AE	119	ARG
4	AE	144	ARG
4	AE	154	LYS
4	AE	163	GLU
4	AE	170	LEU
4	AE	175	VAL
4	AE	181	LEU
4	AE	184	VAL
5	AF	13	SER
5	AF	18	ARG
5	AF	24	LEU
5	AF	33	LEU
5	AF	38	ARG
5	AF	50	SER
5	AF	53	THR
5	AF	74	ARG
5	AF	78	ILE
5	AF	82	ILE
5	AF	98	SER
5	AF	106	ARG
5	AF	108	LYS
5	AF	110	LEU
5	AF	125	LEU
5	AF	140	LEU
5	AF	158	THR
5	AF	162	LEU

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Mol	Chain	Res	Type
5	AF	170	LEU
5	AF	175	THR
5	AF	183	VAL
5	AF	192	LEU
5	AF	197	ASP
5	AF	201	VAL
5	AF	205	ARG
6	AG	5	VAL
6	AG	7	LEU
6	AG	20	ILE
6	AG	21	ARG
6	AG	31	VAL
6	AG	35	GLU
6	AG	60	LEU
6	AG	71	THR
6	AG	128	ARG
6	AG	138	GLN
6	AG	139	LEU
6	AG	143	GLU
6	AG	146	TYR
6	AG	148	MET
6	AG	152	LEU
6	AG	159	VAL
6	AG	161	THR
6	AG	170	ARG
7	AH	4	ILE
7	AH	6	ARG
7	AH	41	MET
7	AH	44	VAL
7	AH	45	VAL
7	AH	68	THR
7	AH	69	ARG
7	AH	88	LEU
7	AH	98	LEU
7	AH	105	LEU
7	AH	106	THR
7	AH	116	GLU
7	AH	127	GLU
7	AH	139	GLN
7	AH	169	VAL
7	AH	171	LEU
8	AI	1	MET

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Mol	Chain	Res	Type
8	AI	3	VAL
8	AI	9	LEU
8	AI	15	VAL
8	AI	38	LEU
8	AI	44	LEU
8	AI	47	LEU
8	AI	54	GLN
8	AI	57	ARG
8	AI	61	ARG
8	AI	72	LEU
8	AI	73	GLU
8	AI	75	LEU
8	AI	82	ARG
8	AI	101	LEU
8	AI	102	SER
8	AI	105	HIS
8	AI	112	LYS
8	AI	114	LEU
8	AI	116	LEU
8	AI	123	LEU
8	AI	125	GLU
8	AI	129	THR
8	AI	133	HIS
8	AI	136	VAL
8	AI	144	VAL
9	AN	1	MET
9	AN	2	LYS
9	AN	9	VAL
9	AN	12	ARG
9	AN	15	LEU
9	AN	28	THR
9	AN	33	LEU
9	AN	34	LEU
9	AN	43	THR
9	AN	48	MET
9	AN	55	VAL
9	AN	60	ILE
9	AN	61	ARG
9	AN	62	VAL
9	AN	63	THR
9	AN	73	THR
9	AN	76	SER

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Mol	Chain	Res	Type
9	AN	85	ILE
9	AN	87	LEU
9	AN	93	THR
9	AN	99	LEU
9	AN	120	LEU
9	AN	133	GLN
10	AO	3	GLN
10	AO	10	VAL
10	AO	17	ARG
10	AO	24	VAL
10	AO	53	LYS
10	AO	70	LYS
10	AO	94	ARG
10	AO	98	VAL
10	AO	113	LYS
11	AP	1	MET
11	AP	3	LEU
11	AP	21	ARG
11	AP	42	SER
11	AP	55	ARG
11	AP	56	SER
11	AP	58	THR
11	AP	59	LEU
11	AP	65	ARG
11	AP	76	LYS
11	AP	83	VAL
11	AP	106	LEU
11	AP	112	LEU
11	AP	119	GLU
11	AP	144	GLU
11	AP	146	VAL
11	AP	148	LEU
12	AQ	1	MET
12	AQ	16	ARG
12	AQ	18	LYS
12	AQ	21	THR
12	AQ	25	ASP
12	AQ	31	ASP
12	AQ	45	GLN
12	AQ	55	VAL
12	AQ	56	ARG
12	AQ	59	ARG

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Mol	Chain	Res	Type
12	AQ	63	LYS
12	AQ	89	ASN
12	AQ	106	VAL
12	AQ	109	VAL
12	AQ	110	THR
12	AQ	134	ARG
12	AQ	135	ASP
13	AR	1	MET
13	AR	6	SER
13	AR	15	SER
13	AR	18	LEU
13	AR	23	ASN
13	AR	27	SER
13	AR	28	LEU
13	AR	29	LEU
13	AR	33	ARG
13	AR	36	THR
13	AR	44	LEU
13	AR	49	ASP
13	AR	60	LEU
13	AR	65	LEU
13	AR	67	LEU
13	AR	75	LEU
13	AR	79	LEU
13	AR	86	ARG
13	AR	96	ARG
13	AR	100	LEU
13	AR	111	LEU
14	AS	3	ARG
14	AS	5	THR
14	AS	8	GLU
14	AS	12	PHE
14	AS	15	ARG
14	AS	20	ARG
14	AS	21	THR
14	AS	25	ARG
14	AS	32	LEU
14	AS	38	GLN
14	AS	50	SER
14	AS	52	SER
14	AS	59	LYS
14	AS	78	LEU

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Mol	Chain	Res	Type
14	AS	84	GLN
14	AS	95	HIS
14	AS	101	LEU
15	AT	6	LEU
15	AT	8	LYS
15	AT	34	VAL
15	AT	40	THR
15	AT	49	VAL
15	AT	59	THR
15	AT	64	ARG
15	AT	74	ARG
15	AT	78	LEU
15	AT	96	ARG
15	AT	118	ARG
16	AU	8	VAL
16	AU	16	LYS
16	AU	36	ARG
16	AU	59	ARG
16	AU	60	LEU
16	AU	74	LEU
16	AU	83	LEU
16	AU	95	LEU
16	AU	104	GLN
16	AU	108	GLU
17	AV	5	VAL
17	AV	12	TYR
17	AV	18	LEU
17	AV	21	ARG
17	AV	28	GLU
17	AV	35	LEU
17	AV	38	LEU
17	AV	46	VAL
17	AV	52	VAL
17	AV	61	VAL
17	AV	62	LEU
17	AV	72	VAL
17	AV	79	VAL
17	AV	85	LYS
17	AV	89	GLN
17	AV	95	LEU
17	AV	100	ARG
18	AW	1	MET

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Mol	Chain	Res	Type
18	AW	11	ARG
18	AW	17	VAL
18	AW	23	LEU
18	AW	51	LEU
18	AW	52	GLU
18	AW	60	ASN
18	AW	63	ASP
18	AW	67	ASP
18	AW	83	LYS
18	AW	92	ARG
18	AW	100	THR
18	AW	103	ILE
18	AW	107	LEU
19	AX	25	LYS
19	AX	52	VAL
19	AX	54	VAL
19	AX	57	LEU
19	AX	65	ARG
19	AX	66	LEU
19	AX	68	ARG
19	AX	76	ARG
20	AY	9	LYS
20	AY	23	ARG
20	AY	43	ASN
20	AY	47	LYS
20	AY	61	ILE
20	AY	76	CYS
20	AY	90	LEU
20	AY	97	ARG
20	AY	101	LYS
20	AY	107	ASP
21	AZ	6	LYS
21	AZ	11	GLU
21	AZ	13	GLU
21	AZ	19	ARG
21	AZ	61	LEU
21	AZ	72	ARG
21	AZ	73	GLN
21	AZ	78	LYS
21	AZ	82	ARG
21	AZ	86	VAL
21	AZ	87	ASP

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Mol	Chain	Res	Type
21	AZ	91	LEU
21	AZ	121	HIS
21	AZ	126	VAL
21	AZ	144	LEU
21	AZ	161	VAL
21	AZ	162	GLU
21	AZ	165	VAL
21	AZ	170	THR
21	AZ	185	GLU
22	A0	20	ARG
22	A0	41	ARG
22	A0	46	LYS
22	A0	55	ARG
22	A0	74	ARG
22	A0	77	ARG
22	A0	81	VAL
23	A1	4	VAL
23	A1	5	CYS
23	A1	17	SER
23	A1	21	ARG
23	A1	40	ARG
23	A1	58	ILE
23	A1	59	THR
23	A1	62	VAL
23	A1	81	LYS
23	A1	95	LEU
24	A2	17	SER
24	A2	30	ARG
24	A2	32	LEU
24	A2	34	GLU
24	A2	44	LEU
24	A2	50	ILE
24	A2	53	LEU
24	A2	55	ARG
24	A2	64	LEU
25	A3	8	LEU
25	A3	23	LEU
25	A3	31	LEU
25	A3	40	THR
25	A3	44	ARG
25	A3	58	VAL
26	A4	3	GLU

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Mol	Chain	Res	Type
26	A4	14	ILE
26	A4	18	CYS
26	A4	22	ILE
26	A4	33	VAL
26	A4	35	VAL
26	A4	43	TYR
27	A5	15	ARG
27	A5	23	HIS
27	A5	29	THR
27	A5	33	CYS
27	A5	37	LYS
27	A5	40	LYS
27	A5	46	CYS
28	A6	4	GLU
28	A6	9	LEU
28	A6	13	CYS
28	A6	14	THR
28	A6	28	ARG
28	A6	30	THR
28	A6	34	LEU
28	A6	35	GLU
28	A6	40	CYS
28	A6	44	ARG
29	A7	4	THR
29	A7	8	ASN
29	A7	23	ARG
29	A7	24	THR
29	A7	32	LYS
29	A7	41	ARG
30	A8	4	MET
30	A8	14	VAL
30	A8	31	HIS
30	A8	32	LEU
30	A8	34	TRP
30	A8	35	GLN
30	A8	41	ILE
3	BD	4	LYS
3	BD	12	SER
3	BD	14	ARG
3	BD	54	ARG
3	BD	61	LEU
3	BD	94	LEU

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Mol	Chain	Res	Type
3	BD	103	ARG
3	BD	106	ILE
3	BD	138	VAL
3	BD	142	VAL
3	BD	150	LYS
3	BD	155	LEU
3	BD	165	ILE
3	BD	173	VAL
3	BD	192	THR
3	BD	208	LYS
3	BD	211	ARG
3	BD	212	SER
3	BD	221	VAL
3	BD	229	VAL
3	BD	242	ARG
3	BD	253	GLN
3	BD	257	LEU
3	BD	259	THR
3	BD	260	ARG
3	BD	267	SER
4	BE	1	MET
4	BE	2	LYS
4	BE	9	VAL
4	BE	12	THR
4	BE	21	VAL
4	BE	24	THR
4	BE	33	VAL
4	BE	45	THR
4	BE	49	LEU
4	BE	52	LEU
4	BE	73	GLU
4	BE	75	VAL
4	BE	77	ILE
4	BE	82	ARG
4	BE	87	GLU
4	BE	93	VAL
4	BE	111	ARG
4	BE	116	VAL
4	BE	119	ARG
4	BE	128	SER
4	BE	144	ARG
4	BE	152	LYS

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Mol	Chain	Res	Type
4	BE	154	LYS
4	BE	163	GLU
4	BE	170	LEU
4	BE	175	VAL
4	BE	181	LEU
4	BE	184	VAL
4	BE	195	LEU
5	BF	18	ARG
5	BF	24	LEU
5	BF	32	LEU
5	BF	33	LEU
5	BF	38	ARG
5	BF	50	SER
5	BF	53	THR
5	BF	74	ARG
5	BF	82	ILE
5	BF	108	LYS
5	BF	110	LEU
5	BF	117	ARG
5	BF	158	THR
5	BF	162	LEU
5	BF	168	ARG
5	BF	175	THR
5	BF	183	VAL
5	BF	192	LEU
5	BF	197	ASP
5	BF	201	VAL
6	BG	5	VAL
6	BG	7	LEU
6	BG	19	LEU
6	BG	21	ARG
6	BG	43	LEU
6	BG	60	LEU
6	BG	71	THR
6	BG	139	LEU
6	BG	143	GLU
6	BG	146	TYR
6	BG	148	MET
6	BG	152	LEU
6	BG	161	THR
6	BG	170	ARG
7	BH	3	ARG

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Mol	Chain	Res	Type
7	BH	6	ARG
7	BH	24	VAL
7	BH	41	MET
7	BH	44	VAL
7	BH	45	VAL
7	BH	69	ARG
7	BH	88	LEU
7	BH	98	LEU
7	BH	105	LEU
7	BH	106	THR
7	BH	122	THR
7	BH	127	GLU
7	BH	139	GLN
7	BH	149	ARG
7	BH	169	VAL
7	BH	171	LEU
8	BI	1	MET
8	BI	3	VAL
8	BI	9	LEU
8	BI	15	VAL
8	BI	38	LEU
8	BI	47	LEU
8	BI	57	ARG
8	BI	61	ARG
8	BI	75	LEU
8	BI	77	LEU
8	BI	78	THR
8	BI	85	GLU
8	BI	92	VAL
8	BI	97	ILE
8	BI	101	LEU
8	BI	105	HIS
8	BI	110	GLU
8	BI	112	LYS
8	BI	114	LEU
8	BI	116	LEU
8	BI	117	GLU
8	BI	121	LYS
8	BI	123	LEU
8	BI	133	HIS
8	BI	140	LEU
8	BI	142	VAL

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Mol	Chain	Res	Type
9	BN	1	MET
9	BN	2	LYS
9	BN	9	VAL
9	BN	12	ARG
9	BN	28	THR
9	BN	33	LEU
9	BN	34	LEU
9	BN	48	MET
9	BN	55	VAL
9	BN	60	ILE
9	BN	61	ARG
9	BN	62	VAL
9	BN	63	THR
9	BN	67	LEU
9	BN	73	THR
9	BN	76	SER
9	BN	87	LEU
9	BN	93	THR
9	BN	96	GLU
9	BN	97	ARG
9	BN	99	LEU
9	BN	115	ARG
9	BN	120	LEU
9	BN	121	LYS
9	BN	133	GLN
9	BN	137	LYS
10	BO	10	VAL
10	BO	17	ARG
10	BO	24	VAL
10	BO	42	SER
10	BO	48	PRO
10	BO	53	LYS
10	BO	94	ARG
10	BO	98	VAL
10	BO	113	LYS
11	BP	19	VAL
11	BP	21	ARG
11	BP	27	HIS
11	BP	40	SER
11	BP	42	SER
11	BP	59	LEU
11	BP	65	ARG

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Mol	Chain	Res	Type
11	BP	75	ILE
11	BP	76	LYS
11	BP	83	VAL
11	BP	95	VAL
11	BP	106	LEU
11	BP	112	LEU
11	BP	119	GLU
11	BP	144	GLU
11	BP	148	LEU
12	BQ	1	MET
12	BQ	7	MET
12	BQ	8	LYS
12	BQ	16	ARG
12	BQ	21	THR
12	BQ	42	ILE
12	BQ	45	GLN
12	BQ	56	ARG
12	BQ	59	ARG
12	BQ	63	LYS
12	BQ	75	THR
12	BQ	106	VAL
12	BQ	109	VAL
12	BQ	110	THR
12	BQ	133	ARG
12	BQ	134	ARG
12	BQ	135	ASP
13	BR	1	MET
13	BR	6	SER
13	BR	15	SER
13	BR	17	ARG
13	BR	18	LEU
13	BR	27	SER
13	BR	28	LEU
13	BR	29	LEU
13	BR	33	ARG
13	BR	36	THR
13	BR	44	LEU
13	BR	54	LEU
13	BR	60	LEU
13	BR	65	LEU
13	BR	67	LEU
13	BR	78	LYS

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Mol	Chain	Res	Type
13	BR	79	LEU
13	BR	86	ARG
13	BR	96	ARG
13	BR	100	LEU
13	BR	105	ARG
13	BR	111	LEU
14	BS	3	ARG
14	BS	5	THR
14	BS	13	ARG
14	BS	20	ARG
14	BS	25	ARG
14	BS	36	TYR
14	BS	38	GLN
14	BS	50	SER
14	BS	52	SER
14	BS	69	VAL
14	BS	78	LEU
14	BS	84	GLN
15	BT	8	LYS
15	BT	49	VAL
15	BT	53	ARG
15	BT	59	THR
15	BT	64	ARG
15	BT	74	ARG
15	BT	78	LEU
15	BT	89	VAL
15	BT	96	ARG
15	BT	118	ARG
16	BU	8	VAL
16	BU	16	LYS
16	BU	27	LEU
16	BU	31	SER
16	BU	36	ARG
16	BU	51	LYS
16	BU	55	ARG
16	BU	59	ARG
16	BU	60	LEU
16	BU	69	CYS
16	BU	74	LEU
16	BU	83	LEU
16	BU	95	LEU
16	BU	104	GLN

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Mol	Chain	Res	Type
16	BU	108	GLU
16	BU	112	ARG
17	BV	7	THR
17	BV	18	LEU
17	BV	21	ARG
17	BV	28	GLU
17	BV	35	LEU
17	BV	43	GLU
17	BV	46	VAL
17	BV	49	THR
17	BV	52	VAL
17	BV	61	VAL
17	BV	62	LEU
17	BV	70	ILE
17	BV	72	VAL
17	BV	79	VAL
17	BV	85	LYS
17	BV	95	LEU
17	BV	100	ARG
18	BW	11	ARG
18	BW	15	ARG
18	BW	17	VAL
18	BW	18	ARG
18	BW	23	LEU
18	BW	51	LEU
18	BW	59	VAL
18	BW	63	ASP
18	BW	83	LYS
18	BW	97	LYS
18	BW	107	LEU
19	BX	45	THR
19	BX	50	LYS
19	BX	52	VAL
19	BX	57	LEU
19	BX	60	ARG
19	BX	65	ARG
19	BX	66	LEU
19	BX	68	ARG
19	BX	92	LEU
20	BY	23	ARG
20	BY	43	ASN
20	BY	47	LYS

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Mol	Chain	Res	Type
20	BY	61	ILE
20	BY	72	VAL
20	BY	90	LEU
20	BY	97	ARG
20	BY	101	LYS
21	BZ	5	LEU
21	BZ	6	LYS
21	BZ	11	GLU
21	BZ	18	LEU
21	BZ	19	ARG
21	BZ	61	LEU
21	BZ	72	ARG
21	BZ	73	GLN
21	BZ	81	ARG
21	BZ	82	ARG
21	BZ	86	VAL
21	BZ	87	ASP
21	BZ	91	LEU
21	BZ	121	HIS
21	BZ	126	VAL
21	BZ	128	VAL
21	BZ	144	LEU
21	BZ	161	VAL
21	BZ	162	GLU
21	BZ	165	VAL
21	BZ	170	THR
21	BZ	185	GLU
22	B0	10	THR
22	B0	19	LYS
22	B0	20	ARG
22	B0	46	LYS
22	B0	55	ARG
22	B0	74	ARG
22	B0	77	ARG
23	B1	4	VAL
23	B1	21	ARG
23	B1	40	ARG
23	B1	41	ARG
23	B1	51	VAL
23	B1	58	ILE
23	B1	59	THR
23	B1	95	LEU

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Mol	Chain	Res	Type
24	B2	28	LYS
24	B2	30	ARG
24	B2	34	GLU
24	B2	45	SER
24	B2	50	ILE
24	B2	52	ASP
24	B2	53	LEU
24	B2	64	LEU
25	B3	8	LEU
25	B3	23	LEU
25	B3	35	ARG
25	B3	40	THR
25	B3	44	ARG
25	B3	58	VAL
26	B4	8	LYS
26	B4	18	CYS
26	B4	22	ILE
26	B4	24	THR
26	B4	27	THR
26	B4	33	VAL
26	B4	34	GLU
26	B4	35	VAL
26	B4	43	TYR
27	B5	6	VAL
27	B5	15	ARG
27	B5	16	ARG
27	B5	23	HIS
27	B5	29	THR
27	B5	37	LYS
27	B5	40	LYS
28	B6	4	GLU
28	B6	6	ARG
28	B6	9	LEU
28	B6	13	CYS
28	B6	14	THR
28	B6	30	THR
28	B6	35	GLU
28	B6	40	CYS
28	B6	44	ARG
29	B7	4	THR
29	B7	8	ASN
29	B7	9	ARG

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Mol	Chain	Res	Type
29	B7	10	ARG
29	B7	24	THR
29	B7	32	LYS
29	B7	42	LEU
29	B7	43	THR
29	B7	47	ARG
30	B8	14	VAL
30	B8	31	HIS
30	B8	34	TRP
30	B8	35	GLN
30	B8	41	ILE
32	CB	15	VAL
32	CB	16	HIS
32	CB	17	PHE
32	CB	21	ARG
32	CB	24	TRP
32	CB	67	THR
32	CB	73	THR
32	CB	74	LYS
32	CB	80	ILE
32	CB	87	ARG
32	CB	93	VAL
32	CB	94	ASN
32	CB	105	PHE
32	CB	111	ARG
32	CB	121	LEU
32	CB	126	GLU
32	CB	133	LYS
32	CB	137	ARG
32	CB	139	LYS
32	CB	142	LEU
32	CB	145	LEU
32	CB	155	LEU
32	CB	160	ASP
32	CB	170	GLU
32	CB	185	ILE
32	CB	187	LEU
32	CB	190	THR
32	CB	191	ASP
32	CB	193	ASP
32	CB	198	ASP
32	CB	200	ILE

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Mol	Chain	Res	Type
32	CB	208	ILE
32	CB	210	SER
32	CB	215	LEU
32	CB	221	LEU
32	CB	224	GLN
33	CC	8	ILE
33	CC	19	GLU
33	CC	30	ARG
33	CC	32	LEU
33	CC	35	GLU
33	CC	47	LEU
33	CC	59	ARG
33	CC	70	VAL
33	CC	102	ASN
33	CC	104	GLN
33	CC	112	SER
33	CC	118	GLN
33	CC	130	VAL
33	CC	131	ARG
33	CC	150	LYS
33	CC	154	SER
33	CC	179	ARG
33	CC	184	TYR
33	CC	196	LEU
34	CD	11	LEU
34	CD	12	CYS
34	CD	13	ARG
34	CD	28	SER
34	CD	33	MET
34	CD	36	ARG
34	CD	53	ASP
34	CD	58	LEU
34	CD	65	ARG
34	CD	70	ILE
34	CD	76	ARG
34	CD	83	SER
34	CD	97	LEU
34	CD	106	TYR
34	CD	107	ARG
34	CD	110	PHE
34	CD	127	THR
34	CD	135	LEU

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Mol	Chain	Res	Type
34	CD	137	SER
34	CD	158	ILE
34	CD	188	LEU
34	CD	196	LEU
35	CE	12	LEU
35	CE	16	THR
35	CE	18	ARG
35	CE	31	LEU
35	CE	34	VAL
35	CE	41	VAL
35	CE	47	LYS
35	CE	51	VAL
35	CE	57	LYS
35	CE	75	THR
35	CE	76	ILE
35	CE	78	HIS
35	CE	79	GLU
35	CE	89	ILE
35	CE	91	LEU
35	CE	116	THR
35	CE	120	THR
35	CE	135	THR
35	CE	137	GLU
35	CE	144	THR
35	CE	147	ASP
35	CE	152	ARG
36	CF	14	LEU
36	CF	15	ASP
36	CF	16	GLN
36	CF	25	ILE
36	CF	36	ARG
36	CF	55	ASP
36	CF	64	GLN
36	CF	69	GLU
36	CF	70	ASP
36	CF	75	LEU
36	CF	89	MET
36	CF	98	LEU
37	CG	6	ARG
37	CG	15	ASP
37	CG	35	LYS
37	CG	37	ASN

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Mol	Chain	Res	Type
37	CG	51	GLN
37	CG	57	GLU
37	CG	59	LEU
37	CG	69	VAL
37	CG	77	SER
37	CG	80	VAL
37	CG	90	GLU
37	CG	104	LEU
37	CG	113	GLU
37	CG	114	ARG
37	CG	138	LYS
37	CG	142	GLU
37	CG	144	MET
37	CG	146	GLU
38	CH	19	VAL
38	CH	21	LYS
38	CH	25	ASP
38	CH	26	VAL
38	CH	52	ASP
38	CH	54	ASP
38	CH	77	GLU
38	CH	78	GLN
38	CH	84	ARG
38	CH	91	ARG
38	CH	95	VAL
38	CH	99	GLU
38	CH	107	LEU
38	CH	109	ILE
38	CH	112	LEU
38	CH	127	LEU
38	CH	133	LEU
38	CH	137	VAL
39	CI	3	GLN
39	CI	31	GLN
39	CI	34	ASN
39	CI	40	LEU
39	CI	60	ASP
39	CI	104	ARG
39	CI	107	ARG
39	CI	108	VAL
39	CI	117	HIS
40	CJ	16	LEU

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Mol	Chain	Res	Type
40	CJ	34	VAL
40	CJ	43	ARG
40	CJ	55	LYS
40	CJ	66	ARG
40	CJ	67	THR
40	CJ	68	HIS
40	CJ	96	ILE
40	CJ	97	GLU
40	CJ	100	THR
41	CK	14	VAL
41	CK	33	THR
41	CK	48	ILE
41	CK	63	LEU
41	CK	70	LYS
41	CK	91	ARG
41	CK	96	ARG
41	CK	104	GLN
41	CK	109	VAL
41	CK	114	VAL
41	CK	117	ASN
41	CK	121	PRO
42	CL	33	ARG
42	CL	43	VAL
42	CL	47	LYS
42	CL	53	ARG
42	CL	55	VAL
42	CL	66	VAL
42	CL	70	ILE
42	CL	80	HIS
42	CL	84	LEU
42	CL	97	ARG
42	CL	117	ARG
43	CM	3	ARG
43	CM	4	ILE
43	CM	14	ARG
43	CM	15	VAL
43	CM	16	ASP
43	CM	17	VAL
43	CM	19	LEU
43	CM	47	ASP
43	CM	55	ARG
43	CM	78	ILE

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Mol	Chain	Res	Type
43	CM	91	ARG
43	CM	101	GLN
43	CM	105	THR
43	CM	110	ARG
44	CN	3	ARG
44	CN	6	LEU
44	CN	7	ILE
44	CN	8	GLU
44	CN	17	LYS
44	CN	18	VAL
44	CN	22	THR
44	CN	23	ARG
44	CN	24	CYS
44	CN	25	VAL
44	CN	33	VAL
45	CO	3	ILE
45	CO	8	LYS
45	CO	24	SER
45	CO	26	GLU
45	CO	35	ARG
45	CO	39	LEU
45	CO	40	SER
45	CO	41	GLU
45	CO	73	GLU
45	CO	76	GLU
45	CO	83	GLU
45	CO	87	ILE
46	CP	1	MET
46	CP	2	VAL
46	CP	6	LEU
46	CP	8	ARG
46	CP	11	SER
46	CP	27	LYS
46	CP	29	ASP
46	CP	45	THR
46	CP	47	ASP
46	CP	62	VAL
46	CP	67	THR
46	CP	69	THR
46	CP	76	GLN
46	CP	79	VAL
47	CQ	13	ASP

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Mol	Chain	Res	Type
47	CQ	45	HIS
47	CQ	57	VAL
47	CQ	60	ILE
48	CR	32	ARG
48	CR	53	ARG
48	CR	74	ARG
48	CR	76	LEU
48	CR	82	THR
49	CS	7	LYS
49	CS	11	VAL
49	CS	31	ILE
49	CS	32	LYS
49	CS	33	THR
49	CS	36	ARG
49	CS	37	ARG
49	CS	43	GLU
49	CS	44	MET
49	CS	47	HIS
49	CS	52	TYR
49	CS	57	HIS
49	CS	66	MET
49	CS	73	GLU
49	CS	77	THR
49	CS	79	THR
49	CS	83	HIS
50	CT	10	LEU
50	CT	13	LEU
50	CT	24	LEU
50	CT	30	LYS
50	CT	31	SER
50	CT	37	SER
50	CT	38	LYS
50	CT	39	LYS
50	CT	62	LEU
50	CT	71	THR
50	CT	75	ASN
50	CT	80	ARG
51	CU	9	ARG
51	CU	10	ARG
32	DB	15	VAL
32	DB	16	HIS
32	DB	17	PHE

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Mol	Chain	Res	Type
32	DB	21	ARG
32	DB	24	TRP
32	DB	39	ILE
32	DB	67	THR
32	DB	80	ILE
32	DB	87	ARG
32	DB	90	MET
32	DB	93	VAL
32	DB	94	ASN
32	DB	105	PHE
32	DB	126	GLU
32	DB	133	LYS
32	DB	137	ARG
32	DB	139	LYS
32	DB	142	LEU
32	DB	145	LEU
32	DB	155	LEU
32	DB	158	LEU
32	DB	160	ASP
32	DB	170	GLU
32	DB	185	ILE
32	DB	187	LEU
32	DB	190	THR
32	DB	191	ASP
32	DB	193	ASP
32	DB	198	ASP
32	DB	200	ILE
32	DB	208	ILE
32	DB	214	ILE
32	DB	215	LEU
32	DB	221	LEU
32	DB	224	GLN
33	DC	4	LYS
33	DC	8	ILE
33	DC	10	PHE
33	DC	17	ASP
33	DC	30	ARG
33	DC	36	ASP
33	DC	52	LEU
33	DC	54	ARG
33	DC	59	ARG
33	DC	104	GLN

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Mol	Chain	Res	Type
33	DC	131	ARG
33	DC	135	LYS
33	DC	144	SER
33	DC	165	THR
33	DC	178	LEU
33	DC	188	LEU
33	DC	202	ILE
33	DC	207	VAL
34	DD	11	LEU
34	DD	13	ARG
34	DD	27	TYR
34	DD	28	SER
34	DD	33	MET
34	DD	36	ARG
34	DD	58	LEU
34	DD	65	ARG
34	DD	70	ILE
34	DD	76	ARG
34	DD	83	SER
34	DD	106	TYR
34	DD	107	ARG
34	DD	126	ILE
34	DD	127	THR
34	DD	135	LEU
34	DD	158	ILE
34	DD	188	LEU
34	DD	193	ASP
35	DE	12	LEU
35	DE	16	THR
35	DE	31	LEU
35	DE	34	VAL
35	DE	41	VAL
35	DE	47	LYS
35	DE	51	VAL
35	DE	57	LYS
35	DE	75	THR
35	DE	76	ILE
35	DE	78	HIS
35	DE	79	GLU
35	DE	80	ILE
35	DE	89	ILE
35	DE	91	LEU

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Mol	Chain	Res	Type
35	DE	98	THR
35	DE	120	THR
35	DE	137	GLU
35	DE	144	THR
35	DE	147	ASP
36	DF	14	LEU
36	DF	15	ASP
36	DF	16	GLN
36	DF	25	ILE
36	DF	36	ARG
36	DF	55	ASP
36	DF	64	GLN
36	DF	69	GLU
36	DF	70	ASP
36	DF	75	LEU
36	DF	89	MET
36	DF	98	LEU
37	DG	3	ARG
37	DG	9	VAL
37	DG	16	LEU
37	DG	41	ARG
37	DG	45	ASP
37	DG	51	GLN
37	DG	59	LEU
37	DG	72	ARG
37	DG	75	VAL
37	DG	90	GLU
37	DG	98	SER
37	DG	102	ARG
37	DG	104	LEU
37	DG	113	GLU
37	DG	119	ARG
37	DG	126	ASP
38	DH	21	LYS
38	DH	25	ASP
38	DH	29	SER
38	DH	52	ASP
38	DH	54	ASP
38	DH	63	LEU
38	DH	77	GLU
38	DH	78	GLN
38	DH	84	ARG

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Mol	Chain	Res	Type
38	DH	91	ARG
38	DH	95	VAL
38	DH	97	VAL
38	DH	99	GLU
38	DH	107	LEU
38	DH	109	ILE
38	DH	112	LEU
38	DH	120	THR
38	DH	127	LEU
38	DH	133	LEU
39	DI	27	THR
39	DI	31	GLN
39	DI	64	THR
39	DI	66	ARG
39	DI	104	ARG
39	DI	108	VAL
39	DI	113	LYS
39	DI	124	GLN
40	DJ	16	LEU
40	DJ	21	GLN
40	DJ	38	ILE
40	DJ	55	LYS
40	DJ	59	SER
40	DJ	67	THR
40	DJ	68	HIS
40	DJ	72	VAL
40	DJ	96	ILE
40	DJ	97	GLU
40	DJ	100	THR
41	DK	14	VAL
41	DK	33	THR
41	DK	48	ILE
41	DK	63	LEU
41	DK	70	LYS
41	DK	96	ARG
41	DK	104	GLN
41	DK	107	SER
41	DK	109	VAL
41	DK	114	VAL
42	DL	33	ARG
42	DL	53	ARG
42	DL	55	VAL

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Mol	Chain	Res	Type
42	DL	66	VAL
42	DL	70	ILE
42	DL	80	HIS
42	DL	84	LEU
42	DL	97	ARG
42	DL	104	VAL
43	DM	4	ILE
43	DM	19	LEU
43	DM	47	ASP
43	DM	49	THR
43	DM	52	GLU
43	DM	53	VAL
43	DM	56	LEU
43	DM	64	TRP
43	DM	65	LYS
43	DM	70	LEU
43	DM	96	LEU
43	DM	108	ARG
43	DM	110	ARG
43	DM	114	ARG
44	DN	3	ARG
44	DN	7	ILE
44	DN	8	GLU
44	DN	18	VAL
44	DN	22	THR
44	DN	24	CYS
44	DN	25	VAL
44	DN	27	CYS
44	DN	33	VAL
44	DN	35	ARG
44	DN	40	CYS
44	DN	41	ARG
44	DN	42	ILE
45	DO	3	ILE
45	DO	24	SER
45	DO	26	GLU
45	DO	35	ARG
45	DO	39	LEU
45	DO	41	GLU
45	DO	65	ARG
45	DO	66	LEU
45	DO	76	GLU

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Mol	Chain	Res	Type
45	DO	83	GLU
45	DO	87	ILE
46	DP	1	MET
46	DP	2	VAL
46	DP	6	LEU
46	DP	8	ARG
46	DP	27	LYS
46	DP	62	VAL
46	DP	67	THR
46	DP	69	THR
46	DP	76	GLN
46	DP	79	VAL
47	DQ	6	LEU
47	DQ	9	VAL
47	DQ	12	SER
47	DQ	13	ASP
47	DQ	34	LYS
47	DQ	45	HIS
47	DQ	57	VAL
47	DQ	60	ILE
47	DQ	72	ARG
47	DQ	74	LEU
48	DR	21	LYS
48	DR	76	LEU
48	DR	82	THR
49	DS	31	ILE
49	DS	37	ARG
49	DS	66	MET
49	DS	73	GLU
49	DS	79	THR
49	DS	83	HIS
50	DT	13	LEU
50	DT	30	LYS
50	DT	31	SER
50	DT	38	LYS
50	DT	39	LYS
50	DT	71	THR
50	DT	75	ASN
50	DT	80	ARG
51	DU	9	ARG
51	DU	10	ARG
51	DU	12	LYS

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Mol	Chain	Res	Type
51	DU	15	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (63) such sidechains are listed below:

Mol	Chain	Res	Type
3	AD	143	HIS
3	AD	253	GLN
4	AE	143	ASN
5	AF	75	HIS
5	AF	169	ASN
6	AG	26	GLN
8	AI	105	HIS
9	AN	133	GLN
10	AO	5	GLN
11	AP	35	HIS
12	AQ	123	HIS
13	AR	3	HIS
13	AR	13	HIS
14	AS	68	GLN
19	AX	31	HIS
20	AY	43	ASN
24	A2	46	GLN
28	A6	20	ASN
28	A6	49	HIS
29	A7	6	GLN
3	BD	143	HIS
3	BD	253	GLN
5	BF	75	HIS
5	BF	169	ASN
6	BG	138	GLN
8	BI	105	HIS
10	BO	5	GLN
12	BQ	123	HIS
13	BR	13	HIS
19	BX	31	HIS
21	BZ	34	ASN
30	B8	35	GLN
32	CB	25	ASN
34	CD	129	ASN
35	CE	20	GLN
35	CE	78	HIS
40	CJ	68	HIS

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Mol	Chain	Res	Type
43	CM	106	ASN
45	CO	28	GLN
45	CO	62	GLN
49	CS	57	HIS
32	DB	94	ASN
32	DB	95	GLN
32	DB	113	HIS
33	DC	37	GLN
33	DC	181	ASN
34	DD	129	ASN
35	DE	78	HIS
36	DF	73	ASN
36	DF	100	ASN
37	DG	28	ASN
37	DG	109	ASN
38	DH	82	HIS
39	DI	31	GLN
39	DI	124	GLN
40	DJ	13	HIS
40	DJ	56	HIS
43	DM	40	ASN
43	DM	106	ASN
45	DO	28	GLN
46	DP	13	HIS
49	DS	47	HIS
49	DS	83	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	AA	2819/2915 (96%)	581 (20%)	59 (2%)
1	BA	2819/2915 (96%)	586 (20%)	60 (2%)
2	AB	119/122 (97%)	25 (21%)	0
2	BB	119/122 (97%)	21 (17%)	0
31	CA	1496/1521 (98%)	339 (22%)	31 (2%)
31	DA	1496/1521 (98%)	341 (22%)	27 (1%)
All	All	8868/9116 (97%)	1893 (21%)	177 (1%)

All (1893) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	AA	10	G
1	AA	14	A
1	AA	15	G
1	AA	34	C
1	AA	45	C
1	AA	61	G
1	AA	69	C
1	AA	71	A
1	AA	72	U
1	AA	74	A
1	AA	75	G
1	AA	78	A
1	AA	84	A
1	AA	90	U
1	AA	94	C
1	AA	94(A)	G
1	AA	102	G
1	AA	103	A
1	AA	118	A
1	AA	120	U
1	AA	121	G
1	AA	131	G
1	AA	139(A)	G
1	AA	141	A
1	AA	154	G
1	AA	175	G
1	AA	181	A
1	AA	182	A
1	AA	196	A
1	AA	197	A
1	AA	199	A
1	AA	201	C
1	AA	204	A
1	AA	205	G
1	AA	214	G
1	AA	215	G
1	AA	216	A
1	AA	221	A
1	AA	222	A
1	AA	225	A
1	AA	229	A
1	AA	233	A
1	AA	248	G

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Mol	Chain	Res	Type
1	AA	249	C
1	AA	250	G
1	AA	252	G
1	AA	271(I)	G
1	AA	271(K)	U
1	AA	271(L)	U
1	AA	271(M)	G
1	AA	271(N)	U
1	AA	271(O)	C
1	AA	271(P)	C
1	AA	271(R)	G
1	AA	272(B)	G
1	AA	272(H)	C
1	AA	275	G
1	AA	279	C
1	AA	287	C
1	AA	310	A
1	AA	311	A
1	AA	324	A
1	AA	329	G
1	AA	330	A
1	AA	332	A
1	AA	333	G
1	AA	342	G
1	AA	352	G
1	AA	363	G
1	AA	363(C)	G
1	AA	363(F)	A
1	AA	366	C
1	AA	372	G
1	AA	386	G
1	AA	404	C
1	AA	405	U
1	AA	406	G
1	AA	411	G
1	AA	412	A
1	AA	415	A
1	AA	428	A
1	AA	443	A
1	AA	444	C
1	AA	448	U
1	AA	454	A

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Mol	Chain	Res	Type
1	AA	455	C
1	AA	457	A
1	AA	470	A
1	AA	474	G
1	AA	475	U
1	AA	479	A
1	AA	481	G
1	AA	482	A
1	AA	505	A
1	AA	508	G
1	AA	509	C
1	AA	530	G
1	AA	531	C
1	AA	532	A
1	AA	533	G
1	AA	545	G
1	AA	546	C
1	AA	549	G
1	AA	563	G
1	AA	573	G
1	AA	575	A
1	AA	586	A
1	AA	587	C
1	AA	588	U
1	AA	602	G
1	AA	603	A
1	AA	604	G
1	AA	605	C
1	AA	606	U
1	AA	607	U
1	AA	614(B)	G
1	AA	615	G
1	AA	616	G
1	AA	619	G
1	AA	627	A
1	AA	634	C
1	AA	637	A
1	AA	639	U
1	AA	645	C
1	AA	646	A
1	AA	652(B)	A
1	AA	652(C)	G

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Mol	Chain	Res	Type
1	AA	652(D)	C
1	AA	652(F)	G
1	AA	652(I)	C
1	AA	652(P)	G
1	AA	652(Q)	G
1	AA	652(R)	C
1	AA	652(T)	C
1	AA	652(U)	G
1	AA	657	U
1	AA	669	G
1	AA	670	A
1	AA	686	G
1	AA	708	C
1	AA	717	G
1	AA	726	G
1	AA	730	C
1	AA	752	A
1	AA	753	C
1	AA	764	A
1	AA	765	G
1	AA	775	G
1	AA	776	G
1	AA	782	A
1	AA	784	A
1	AA	785	G
1	AA	787	U
1	AA	792	G
1	AA	794	G
1	AA	796	C
1	AA	805	G
1	AA	812	C
1	AA	819	A
1	AA	822	U
1	AA	827	U
1	AA	828	U
1	AA	829	A
1	AA	830	G
1	AA	857	C
1	AA	859	G
1	AA	879	G
1	AA	884	C
1	AA	885	C

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Mol	Chain	Res	Type
1	AA	886	C
1	AA	888	C
1	AA	889	C
1	AA	890	A
1	AA	893	C
1	AA	896	A
1	AA	900	A
1	AA	901	A
1	AA	910	A
1	AA	916	G
1	AA	917	A
1	AA	932	G
1	AA	938	G
1	AA	941	A
1	AA	945	A
1	AA	946	G
1	AA	953	A
1	AA	958	U
1	AA	959	A
1	AA	961	C
1	AA	974	G
1	AA	975	C
1	AA	975(A)	G
1	AA	983	A
1	AA	991	C
1	AA	996	A
1	AA	1005	C
1	AA	1012	U
1	AA	1013	C
1	AA	1016	G
1	AA	1021	A
1	AA	1022	G
1	AA	1025	G
1	AA	1026	U
1	AA	1033	U
1	AA	1038	C
1	AA	1039	G
1	AA	1042	G
1	AA	1043	C
1	AA	1044	G
1	AA	1045	A
1	AA	1046	A

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Mol	Chain	Res	Type
1	AA	1047	G
1	AA	1048	A
1	AA	1049	C
1	AA	1050	A
1	AA	1051	G
1	AA	1052	C
1	AA	1107	G
1	AA	1108	U
1	AA	1109	C
1	AA	1110	G
1	AA	1111	A
1	AA	1112	G
1	AA	1115	G
1	AA	1130	U
1	AA	1135	C
1	AA	1136	G
1	AA	1139	G
1	AA	1141	U
1	AA	1144	G
1	AA	1155	A
1	AA	1171	G
1	AA	1173	G
1	AA	1174	A
1	AA	1175	U
1	AA	1176	G
1	AA	1177	A
1	AA	1178	C
1	AA	1205	U
1	AA	1210	A
1	AA	1211	U
1	AA	1213	A
1	AA	1220	A
1	AA	1230	C
1	AA	1240	U
1	AA	1241	A
1	AA	1249	U
1	AA	1250	G
1	AA	1253	A
1	AA	1256	G
1	AA	1271	G
1	AA	1272	A
1	AA	1273	U

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Mol	Chain	Res	Type
1	AA	1286	A
1	AA	1300	U
1	AA	1301	A
1	AA	1308	A
1	AA	1313	U
1	AA	1314	C
1	AA	1329	U
1	AA	1332	G
1	AA	1359	A
1	AA	1360	A
1	AA	1365	A
1	AA	1368	G
1	AA	1370	C
1	AA	1379	A
1	AA	1384	A
1	AA	1385	G
1	AA	1386	C
1	AA	1395	A
1	AA	1405	U
1	AA	1416	G
1	AA	1417	C
1	AA	1419	A
1	AA	1420	U
1	AA	1421	G
1	AA	1428	C
1	AA	1437	C
1	AA	1445	A
1	AA	1449	A
1	AA	1450	G
1	AA	1455	G
1	AA	1459	G
1	AA	1461	G
1	AA	1467	C
1	AA	1471	A
1	AA	1478	G
1	AA	1482	G
1	AA	1487	G
1	AA	1488	G
1	AA	1490	A
1	AA	1493	C
1	AA	1495	A
1	AA	1496	A

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Mol	Chain	Res	Type
1	AA	1497	U
1	AA	1507	A
1	AA	1508	A
1	AA	1509	C
1	AA	1509(A)	A
1	AA	1531	C
1	AA	1541	G
1	AA	1542	A
1	AA	1558	A
1	AA	1559	G
1	AA	1569	A
1	AA	1577	C
1	AA	1578	U
1	AA	1579	A
1	AA	1580	A
1	AA	1581	G
1	AA	1584	C
1	AA	1586	A
1	AA	1598	C
1	AA	1607	C
1	AA	1608	A
1	AA	1609	A
1	AA	1610	A
1	AA	1617	C
1	AA	1618	A
1	AA	1640	C
1	AA	1647	G
1	AA	1648	C
1	AA	1654	A
1	AA	1674	G
1	AA	1696	G
1	AA	1700	A
1	AA	1701	A
1	AA	1703	G
1	AA	1721	G
1	AA	1722	A
1	AA	1739	U
1	AA	1740	G
1	AA	1742	G
1	AA	1746	G
1	AA	1756	G
1	AA	1762	A

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Mol	Chain	Res	Type
1	AA	1763	G
1	AA	1764	G
1	AA	1773	A
1	AA	1780	A
1	AA	1781	C
1	AA	1782	C
1	AA	1786	A
1	AA	1791	A
1	AA	1799	G
1	AA	1800	C
1	AA	1801	G
1	AA	1816	G
1	AA	1820	U
1	AA	1829	A
1	AA	1834	U
1	AA	1835	G
1	AA	1836	C
1	AA	1840	G
1	AA	1847	A
1	AA	1858	G
1	AA	1877	A
1	AA	1878	G
1	AA	1882	C
1	AA	1900	A
1	AA	1906	G
1	AA	1913	A
1	AA	1914	C
1	AA	1915	U
1	AA	1929	G
1	AA	1930	G
1	AA	1931	U
1	AA	1936	A
1	AA	1937	A
1	AA	1938	A
1	AA	1942	C
1	AA	1955	U
1	AA	1963	U
1	AA	1964	G
1	AA	1965	C
1	AA	1967	C
1	AA	1969	A
1	AA	1970	A

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Mol	Chain	Res	Type
1	AA	1971	A
1	AA	1972	A
1	AA	1982	C
1	AA	1991	U
1	AA	1992	G
1	AA	1993	U
1	AA	1997	G
1	AA	2018	G
1	AA	2020	A
1	AA	2021	C
1	AA	2023	G
1	AA	2028	U
1	AA	2030	A
1	AA	2031	A
1	AA	2032	G
1	AA	2033	A
1	AA	2036	C
1	AA	2043	C
1	AA	2049	G
1	AA	2052	G
1	AA	2055	C
1	AA	2056	G
1	AA	2060	A
1	AA	2061	G
1	AA	2062	A
1	AA	2063	C
1	AA	2069	G
1	AA	2086	U
1	AA	2098	U
1	AA	2102	U
1	AA	2103	C
1	AA	2104	G
1	AA	2105	C
1	AA	2107	C
1	AA	2108	C
1	AA	2116	G
1	AA	2117	A
1	AA	2118	U
1	AA	2119	A
1	AA	2121	G
1	AA	2123	G
1	AA	2126	A

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Mol	Chain	Res	Type
1	AA	2127	G
1	AA	2130	U
1	AA	2131	G
1	AA	2133	G
1	AA	2134	A
1	AA	2142	C
1	AA	2144	U
1	AA	2145	C
1	AA	2146	C
1	AA	2147	G
1	AA	2148	G
1	AA	2157	G
1	AA	2159	G
1	AA	2160	G
1	AA	2161	C
1	AA	2162	G
1	AA	2163	C
1	AA	2165	G
1	AA	2166	G
1	AA	2170	A
1	AA	2171	A
1	AA	2172	U
1	AA	2173	A
1	AA	2174	C
1	AA	2184	G
1	AA	2186	G
1	AA	2187	G
1	AA	2189	U
1	AA	2191	G
1	AA	2192	G
1	AA	2198	A
1	AA	2199	A
1	AA	2200	C
1	AA	2201	C
1	AA	2206	G
1	AA	2207	G
1	AA	2208	A
1	AA	2218	U
1	AA	2219	G
1	AA	2225	A
1	AA	2235	G
1	AA	2238	G

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Mol	Chain	Res	Type
1	AA	2239	G
1	AA	2261	C
1	AA	2268	A
1	AA	2275	C
1	AA	2279	G
1	AA	2282	G
1	AA	2283	C
1	AA	2287	A
1	AA	2288	A
1	AA	2289	G
1	AA	2294	C
1	AA	2296	U
1	AA	2305	A
1	AA	2309	A
1	AA	2311	A
1	AA	2319	G
1	AA	2320	A
1	AA	2325	G
1	AA	2334	G
1	AA	2336	A
1	AA	2347	C
1	AA	2348	U
1	AA	2350	C
1	AA	2354	G
1	AA	2355	C
1	AA	2379	G
1	AA	2383	G
1	AA	2385	C
1	AA	2388	A
1	AA	2406	U
1	AA	2410	G
1	AA	2414	G
1	AA	2422	A
1	AA	2423	U
1	AA	2425	A
1	AA	2429	G
1	AA	2430	A
1	AA	2431	U
1	AA	2434	A
1	AA	2435	A
1	AA	2439	A
1	AA	2440	C

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Mol	Chain	Res	Type
1	AA	2441	C
1	AA	2446	G
1	AA	2448	A
1	AA	2449	U
1	AA	2468	G
1	AA	2469	A
1	AA	2474	C
1	AA	2476	A
1	AA	2477	C
1	AA	2487	G
1	AA	2489	G
1	AA	2490	G
1	AA	2502	G
1	AA	2505	G
1	AA	2506	U
1	AA	2518	A
1	AA	2520	C
1	AA	2525	G
1	AA	2529	G
1	AA	2535	G
1	AA	2554	U
1	AA	2566	A
1	AA	2567	G
1	AA	2573	C
1	AA	2585	U
1	AA	2608	G
1	AA	2610	C
1	AA	2611	U
1	AA	2612	C
1	AA	2615	U
1	AA	2620	C
1	AA	2629	A
1	AA	2630	G
1	AA	2631	G
1	AA	2638	G
1	AA	2663	G
1	AA	2673	G
1	AA	2685	G
1	AA	2689	U
1	AA	2690	C
1	AA	2707	G
1	AA	2712(A)	A

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Mol	Chain	Res	Type
1	AA	2713	A
1	AA	2714	G
1	AA	2726	U
1	AA	2732	G
1	AA	2733	A
1	AA	2739	U
1	AA	2757	A
1	AA	2758	A
1	AA	2761	G
1	AA	2764	A
1	AA	2765	A
1	AA	2766	G
1	AA	2778	A
1	AA	2790	A
1	AA	2791	C
1	AA	2802	G
1	AA	2803	C
1	AA	2808	U
1	AA	2818	G
1	AA	2820	A
1	AA	2821	A
1	AA	2833	G
1	AA	2834	G
1	AA	2835	A
1	AA	2839	G
1	AA	2846	G
1	AA	2847	U
1	AA	2872	G
1	AA	2873	A
1	AA	2875	C
1	AA	2880	C
1	AA	2892	A
1	AA	2895	U
1	AA	2897	U
2	AB	2	C
2	AB	5	C
2	AB	7	G
2	AB	8	U
2	AB	9	G
2	AB	12	C
2	AB	13	A
2	AB	20	C

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Mol	Chain	Res	Type
2	AB	24	G
2	AB	25	A
2	AB	28	C
2	AB	29	A
2	AB	33	G
2	AB	40	U
2	AB	53	A
2	AB	54	G
2	AB	56	G
2	AB	57	A
2	AB	64	C
2	AB	73	A
2	AB	75	G
2	AB	99	G
2	AB	106	G
2	AB	108	U
2	AB	110	G
1	BA	10	G
1	BA	15	G
1	BA	34	C
1	BA	36	G
1	BA	45	C
1	BA	61	G
1	BA	63	U
1	BA	69	C
1	BA	70	G
1	BA	71	A
1	BA	72	U
1	BA	74	A
1	BA	75	G
1	BA	84	A
1	BA	90	U
1	BA	94	C
1	BA	94(A)	G
1	BA	95	G
1	BA	102	G
1	BA	103	A
1	BA	118	A
1	BA	120	U
1	BA	122	G
1	BA	123	G
1	BA	125	G

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Mol	Chain	Res	Type
1	BA	128	C
1	BA	131	G
1	BA	141	A
1	BA	154	G
1	BA	181	A
1	BA	182	A
1	BA	196	A
1	BA	199	A
1	BA	201	C
1	BA	204	A
1	BA	205	G
1	BA	214	G
1	BA	215	G
1	BA	216	A
1	BA	221	A
1	BA	222	A
1	BA	225	A
1	BA	229	A
1	BA	233	A
1	BA	248	G
1	BA	249	C
1	BA	250	G
1	BA	252	G
1	BA	271(I)	G
1	BA	271(K)	U
1	BA	271(L)	U
1	BA	271(M)	G
1	BA	271(N)	U
1	BA	271(O)	C
1	BA	271(R)	G
1	BA	272(B)	G
1	BA	272(H)	C
1	BA	275	G
1	BA	279	C
1	BA	280	C
1	BA	287	C
1	BA	311	A
1	BA	324	A
1	BA	329	G
1	BA	330	A
1	BA	332	A
1	BA	333	G

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Mol	Chain	Res	Type
1	BA	342	G
1	BA	352	G
1	BA	363	G
1	BA	363(C)	G
1	BA	363(F)	A
1	BA	366	C
1	BA	371	A
1	BA	372	G
1	BA	386	G
1	BA	396	G
1	BA	405	U
1	BA	406	G
1	BA	411	G
1	BA	412	A
1	BA	415	A
1	BA	416	C
1	BA	428	A
1	BA	443	A
1	BA	444	C
1	BA	448	U
1	BA	455	C
1	BA	457	A
1	BA	463	G
1	BA	470	A
1	BA	475	U
1	BA	481	G
1	BA	482	A
1	BA	491	G
1	BA	494	G
1	BA	505	A
1	BA	508	G
1	BA	509	C
1	BA	530	G
1	BA	531	C
1	BA	532	A
1	BA	533	G
1	BA	545	G
1	BA	546	C
1	BA	549	G
1	BA	563	G
1	BA	571	A
1	BA	573	G

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Mol	Chain	Res	Type
1	BA	575	A
1	BA	585	G
1	BA	586	A
1	BA	587	C
1	BA	588	U
1	BA	602	G
1	BA	603	A
1	BA	604	G
1	BA	606	U
1	BA	607	U
1	BA	614(A)	U
1	BA	614(B)	G
1	BA	615	G
1	BA	616	G
1	BA	619	G
1	BA	627	A
1	BA	634	C
1	BA	637	A
1	BA	645	C
1	BA	646	A
1	BA	652(B)	A
1	BA	652(C)	G
1	BA	652(D)	C
1	BA	652(F)	G
1	BA	652(I)	C
1	BA	652(P)	G
1	BA	652(Q)	G
1	BA	652(R)	C
1	BA	652(T)	C
1	BA	652(U)	G
1	BA	669	G
1	BA	673	C
1	BA	686	G
1	BA	694	U
1	BA	708	C
1	BA	717	G
1	BA	722	A
1	BA	730	C
1	BA	739	G
1	BA	752	A
1	BA	753	C
1	BA	764	A

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Mol	Chain	Res	Type
1	BA	765	G
1	BA	775	G
1	BA	776	G
1	BA	777	A
1	BA	782	A
1	BA	784	A
1	BA	785	G
1	BA	787	U
1	BA	792	G
1	BA	805	G
1	BA	812	C
1	BA	819	A
1	BA	827	U
1	BA	828	U
1	BA	829	A
1	BA	830	G
1	BA	857	C
1	BA	859	G
1	BA	863	A
1	BA	879	G
1	BA	884	C
1	BA	885	C
1	BA	886	C
1	BA	888	C
1	BA	889	C
1	BA	890	A
1	BA	893	C
1	BA	896	A
1	BA	899	A
1	BA	900	A
1	BA	901	A
1	BA	910	A
1	BA	915	C
1	BA	916	G
1	BA	917	A
1	BA	932	G
1	BA	934	G
1	BA	938	G
1	BA	941	A
1	BA	945	A
1	BA	946	G
1	BA	953	A

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Mol	Chain	Res	Type
1	BA	958	U
1	BA	959	A
1	BA	961	C
1	BA	974	G
1	BA	975	C
1	BA	975(A)	G
1	BA	983	A
1	BA	990	A
1	BA	991	C
1	BA	996	A
1	BA	1005	C
1	BA	1012	U
1	BA	1013	C
1	BA	1020	A
1	BA	1022	G
1	BA	1024	G
1	BA	1025	G
1	BA	1026	U
1	BA	1033	U
1	BA	1038	C
1	BA	1039	G
1	BA	1042	G
1	BA	1043	C
1	BA	1044	G
1	BA	1045	A
1	BA	1046	A
1	BA	1047	G
1	BA	1048	A
1	BA	1049	C
1	BA	1050	A
1	BA	1051	G
1	BA	1052	C
1	BA	1107	G
1	BA	1108	U
1	BA	1109	C
1	BA	1110	G
1	BA	1111	A
1	BA	1112	G
1	BA	1115	G
1	BA	1129	A
1	BA	1130	U
1	BA	1135	C

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Mol	Chain	Res	Type
1	BA	1136	G
1	BA	1139	G
1	BA	1141	U
1	BA	1149	G
1	BA	1155	A
1	BA	1171	G
1	BA	1173	G
1	BA	1174	A
1	BA	1175	U
1	BA	1176	G
1	BA	1177	A
1	BA	1178	C
1	BA	1210	A
1	BA	1211	U
1	BA	1218	C
1	BA	1220	A
1	BA	1240	U
1	BA	1252	G
1	BA	1253	A
1	BA	1256	G
1	BA	1271	G
1	BA	1272	A
1	BA	1273	U
1	BA	1286	A
1	BA	1298	C
1	BA	1300	U
1	BA	1301	A
1	BA	1307	A
1	BA	1313	U
1	BA	1314	C
1	BA	1329	U
1	BA	1332	G
1	BA	1349	A
1	BA	1351	C
1	BA	1359	A
1	BA	1360	A
1	BA	1365	A
1	BA	1368	G
1	BA	1370	C
1	BA	1373	A
1	BA	1380	G
1	BA	1384	A

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Mol	Chain	Res	Type
1	BA	1385	G
1	BA	1386	C
1	BA	1395	A
1	BA	1416	G
1	BA	1417	C
1	BA	1419	A
1	BA	1420	U
1	BA	1421	G
1	BA	1422	G
1	BA	1428	C
1	BA	1437	C
1	BA	1445	A
1	BA	1449	A
1	BA	1450	G
1	BA	1455	G
1	BA	1459	G
1	BA	1461	G
1	BA	1467	C
1	BA	1471	A
1	BA	1478	G
1	BA	1482	G
1	BA	1487	G
1	BA	1488	G
1	BA	1490	A
1	BA	1493	C
1	BA	1496	A
1	BA	1497	U
1	BA	1507	A
1	BA	1508	A
1	BA	1509	C
1	BA	1509(A)	A
1	BA	1531	C
1	BA	1541	G
1	BA	1542	A
1	BA	1545	A
1	BA	1558	A
1	BA	1559	G
1	BA	1566	A
1	BA	1569	A
1	BA	1577	C
1	BA	1578	U
1	BA	1579	A

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Mol	Chain	Res	Type
1	BA	1580	A
1	BA	1581	G
1	BA	1582	C
1	BA	1584	C
1	BA	1586	A
1	BA	1598	C
1	BA	1608	A
1	BA	1609	A
1	BA	1610	A
1	BA	1617	C
1	BA	1639	U
1	BA	1640	C
1	BA	1647	G
1	BA	1648	C
1	BA	1654	A
1	BA	1674	G
1	BA	1688	U
1	BA	1696	G
1	BA	1700	A
1	BA	1701	A
1	BA	1721	G
1	BA	1722	A
1	BA	1739	U
1	BA	1742	G
1	BA	1746	G
1	BA	1756	G
1	BA	1762	A
1	BA	1763	G
1	BA	1764	G
1	BA	1773	A
1	BA	1780	A
1	BA	1782	C
1	BA	1786	A
1	BA	1791	A
1	BA	1799	G
1	BA	1800	C
1	BA	1801	G
1	BA	1816	G
1	BA	1820	U
1	BA	1827	C
1	BA	1829	A
1	BA	1834	U

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Mol	Chain	Res	Type
1	BA	1835	G
1	BA	1840	G
1	BA	1847	A
1	BA	1858	G
1	BA	1877	A
1	BA	1878	G
1	BA	1882	C
1	BA	1889	A
1	BA	1896	G
1	BA	1900	A
1	BA	1906	G
1	BA	1913	A
1	BA	1914	C
1	BA	1915	U
1	BA	1929	G
1	BA	1930	G
1	BA	1931	U
1	BA	1936	A
1	BA	1937	A
1	BA	1938	A
1	BA	1955	U
1	BA	1963	U
1	BA	1964	G
1	BA	1967	C
1	BA	1969	A
1	BA	1970	A
1	BA	1971	A
1	BA	1972	A
1	BA	1982	C
1	BA	1983	C
1	BA	1991	U
1	BA	1992	G
1	BA	1993	U
1	BA	1997	G
1	BA	2020	A
1	BA	2023	G
1	BA	2024	G
1	BA	2030	A
1	BA	2031	A
1	BA	2032	G
1	BA	2033	A
1	BA	2036	C

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Mol	Chain	Res	Type
1	BA	2043	C
1	BA	2055	C
1	BA	2056	G
1	BA	2060	A
1	BA	2061	G
1	BA	2062	A
1	BA	2063	C
1	BA	2069	G
1	BA	2098	U
1	BA	2102	U
1	BA	2103	C
1	BA	2104	G
1	BA	2105	C
1	BA	2107	C
1	BA	2108	C
1	BA	2116	G
1	BA	2117	A
1	BA	2118	U
1	BA	2119	A
1	BA	2121	G
1	BA	2123	G
1	BA	2124	G
1	BA	2126	A
1	BA	2127	G
1	BA	2130	U
1	BA	2131	G
1	BA	2133	G
1	BA	2134	A
1	BA	2142	C
1	BA	2144	U
1	BA	2145	C
1	BA	2146	C
1	BA	2147	G
1	BA	2148	G
1	BA	2157	G
1	BA	2159	G
1	BA	2160	G
1	BA	2161	C
1	BA	2163	C
1	BA	2165	G
1	BA	2166	G
1	BA	2167	U

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Mol	Chain	Res	Type
1	BA	2170	A
1	BA	2172	U
1	BA	2173	A
1	BA	2174	C
1	BA	2180	U
1	BA	2184	G
1	BA	2186	G
1	BA	2187	G
1	BA	2191	G
1	BA	2192	G
1	BA	2198	A
1	BA	2199	A
1	BA	2200	C
1	BA	2201	C
1	BA	2205	C
1	BA	2206	G
1	BA	2207	G
1	BA	2208	A
1	BA	2218	U
1	BA	2219	G
1	BA	2225	A
1	BA	2226	C
1	BA	2235	G
1	BA	2238	G
1	BA	2239	G
1	BA	2261	C
1	BA	2268	A
1	BA	2273	A
1	BA	2275	C
1	BA	2282	G
1	BA	2283	C
1	BA	2287	A
1	BA	2288	A
1	BA	2289	G
1	BA	2296	U
1	BA	2305	A
1	BA	2309	A
1	BA	2311	A
1	BA	2319	G
1	BA	2320	A
1	BA	2325	G
1	BA	2334	G

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Mol	Chain	Res	Type
1	BA	2336	A
1	BA	2347	C
1	BA	2350	C
1	BA	2354	G
1	BA	2355	C
1	BA	2377	A
1	BA	2379	G
1	BA	2383	G
1	BA	2385	C
1	BA	2406	U
1	BA	2410	G
1	BA	2414	G
1	BA	2422	A
1	BA	2424	C
1	BA	2425	A
1	BA	2429	G
1	BA	2430	A
1	BA	2434	A
1	BA	2435	A
1	BA	2439	A
1	BA	2440	C
1	BA	2441	C
1	BA	2446	G
1	BA	2448	A
1	BA	2449	U
1	BA	2468	G
1	BA	2469	A
1	BA	2471	C
1	BA	2474	C
1	BA	2476	A
1	BA	2477	C
1	BA	2478	A
1	BA	2490	G
1	BA	2498	C
1	BA	2502	G
1	BA	2505	G
1	BA	2506	U
1	BA	2518	A
1	BA	2520	C
1	BA	2525	G
1	BA	2529	G
1	BA	2535	G

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Mol	Chain	Res	Type
1	BA	2554	U
1	BA	2564	A
1	BA	2566	A
1	BA	2567	G
1	BA	2569	G
1	BA	2573	C
1	BA	2582	G
1	BA	2585	U
1	BA	2608	G
1	BA	2609	U
1	BA	2610	C
1	BA	2611	U
1	BA	2612	C
1	BA	2615	U
1	BA	2629	A
1	BA	2630	G
1	BA	2654	A
1	BA	2663	G
1	BA	2673	G
1	BA	2679	A
1	BA	2682	U
1	BA	2690	C
1	BA	2702	U
1	BA	2703	C
1	BA	2712(A)	A
1	BA	2713	A
1	BA	2714	G
1	BA	2718	G
1	BA	2726	U
1	BA	2733	A
1	BA	2758	A
1	BA	2761	G
1	BA	2765	A
1	BA	2766	G
1	BA	2778	A
1	BA	2780	G
1	BA	2789	C
1	BA	2790	A
1	BA	2791	C
1	BA	2802	G
1	BA	2803	C
1	BA	2808	U

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Mol	Chain	Res	Type
1	BA	2818	G
1	BA	2820	A
1	BA	2821	A
1	BA	2824	C
1	BA	2832	U
1	BA	2833	G
1	BA	2834	G
1	BA	2835	A
1	BA	2847	U
1	BA	2872	G
1	BA	2873	A
1	BA	2880	C
1	BA	2892	A
1	BA	2895	U
1	BA	2897	U
2	BB	2	C
2	BB	8	U
2	BB	9	G
2	BB	12	C
2	BB	13	A
2	BB	20	C
2	BB	25	A
2	BB	29	A
2	BB	33	G
2	BB	40	U
2	BB	47	C
2	BB	53	A
2	BB	54	G
2	BB	56	G
2	BB	73	A
2	BB	75	G
2	BB	88	C
2	BB	99	G
2	BB	106	G
2	BB	108	U
2	BB	110	G
31	CA	4	U
31	CA	7	G
31	CA	8	A
31	CA	9	G
31	CA	10	A
31	CA	14	U

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Mol	Chain	Res	Type
31	CA	32	A
31	CA	39	G
31	CA	44	G
31	CA	47	C
31	CA	48	C
31	CA	51	A
31	CA	60	A
31	CA	61	G
31	CA	65	U
31	CA	67	C
31	CA	70	G
31	CA	72	C
31	CA	76	C
31	CA	77	G
31	CA	78	G
31	CA	79	G
31	CA	92	C
31	CA	96	U
31	CA	97	G
31	CA	98	G
31	CA	115	G
31	CA	116	A
31	CA	119	A
31	CA	120	A
31	CA	121	C
31	CA	131	C
31	CA	138	G
31	CA	142	G
31	CA	144	G
31	CA	146	G
31	CA	150	C
31	CA	162	A
31	CA	163	C
31	CA	167	G
31	CA	169	C
31	CA	170	U
31	CA	173	U
31	CA	182	U
31	CA	189(G)	G
31	CA	189(H)	G
31	CA	195	A
31	CA	197	A

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Mol	Chain	Res	Type
31	CA	199	G
31	CA	201	C
31	CA	203	U
31	CA	204	U
31	CA	216	G
31	CA	231	G
31	CA	243	A
31	CA	244	U
31	CA	247	G
31	CA	251	G
31	CA	266	G
31	CA	267	C
31	CA	281	G
31	CA	289	G
31	CA	321	A
31	CA	328	C
31	CA	330	C
31	CA	332	G
31	CA	341	C
31	CA	345	C
31	CA	347	G
31	CA	348	G
31	CA	350	G
31	CA	352	C
31	CA	353	A
31	CA	354	G
31	CA	355	C
31	CA	363	A
31	CA	366	C
31	CA	367	U
31	CA	372	C
31	CA	373	A
31	CA	384	G
31	CA	388	G
31	CA	392	G
31	CA	397	A
31	CA	398	C
31	CA	406	G
31	CA	411	A
31	CA	412	A
31	CA	413	G
31	CA	414	A

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Mol	Chain	Res	Type
31	CA	422	C
31	CA	423	G
31	CA	424	G
31	CA	428	G
31	CA	429	U
31	CA	433	C
31	CA	435	C
31	CA	436	C
31	CA	437	U
31	CA	439	A
31	CA	442	C
31	CA	451	A
31	CA	452	A
31	CA	461	A
31	CA	473	G
31	CA	475	G
31	CA	476	G
31	CA	483	C
31	CA	484	G
31	CA	485	G
31	CA	496	A
31	CA	498	U
31	CA	505	G
31	CA	509	A
31	CA	510	A
31	CA	511	C
31	CA	518	C
31	CA	520	A
31	CA	527	G
31	CA	531	U
31	CA	532	A
31	CA	533	A
31	CA	534	U
31	CA	547	A
31	CA	559	A
31	CA	560	U
31	CA	561	U
31	CA	562	C
31	CA	567	G
31	CA	572	A
31	CA	573	A
31	CA	576	G

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Mol	Chain	Res	Type
31	CA	577	G
31	CA	592	G
31	CA	596	C
31	CA	630	G
31	CA	631	G
31	CA	632	A
31	CA	653	A
31	CA	665	A
31	CA	687	A
31	CA	688	G
31	CA	723	U
31	CA	724	G
31	CA	731	G
31	CA	749	C
31	CA	752	G
31	CA	753	A
31	CA	754	C
31	CA	755	G
31	CA	777	A
31	CA	786	G
31	CA	792	A
31	CA	793	U
31	CA	794	A
31	CA	796	C
31	CA	817	C
31	CA	818	G
31	CA	821	G
31	CA	827	U
31	CA	828	A
31	CA	829	G
31	CA	833	U
31	CA	836	G
31	CA	839	U
31	CA	840	C
31	CA	841	U
31	CA	848	C
31	CA	851	G
31	CA	859	A
31	CA	870	U
31	CA	872	A
31	CA	902	G
31	CA	913	A

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Mol	Chain	Res	Type
31	CA	914	A
31	CA	926	G
31	CA	927	G
31	CA	934	C
31	CA	935	A
31	CA	961	U
31	CA	962	C
31	CA	969	A
31	CA	971	G
31	CA	972	C
31	CA	974	A
31	CA	975	A
31	CA	976	G
31	CA	977	A
31	CA	989	C
31	CA	992	U
31	CA	993	G
31	CA	1000	U
31	CA	1001(A)	G
31	CA	1004	A
31	CA	1005	A
31	CA	1009	G
31	CA	1010	G
31	CA	1011	G
31	CA	1012	U
31	CA	1016	A
31	CA	1019	C
31	CA	1023	G
31	CA	1025	U
31	CA	1026	G
31	CA	1027	C
31	CA	1028	C
31	CA	1029	C
31	CA	1030(A)	G
31	CA	1030(B)	C
31	CA	1030(C)	G
31	CA	1032	G
31	CA	1034	G
31	CA	1035	A
31	CA	1036	G
31	CA	1044	A
31	CA	1053	G

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Mol	Chain	Res	Type
31	CA	1054	C
31	CA	1064	G
31	CA	1065	U
31	CA	1066	C
31	CA	1068	G
31	CA	1077	G
31	CA	1081	G
31	CA	1086	U
31	CA	1091	U
31	CA	1094	G
31	CA	1095	U
31	CA	1101	A
31	CA	1103	C
31	CA	1110	A
31	CA	1124	G
31	CA	1125	U
31	CA	1126	U
31	CA	1127	G
31	CA	1129	C
31	CA	1136	U
31	CA	1137	C
31	CA	1138	G
31	CA	1139	G
31	CA	1140	C
31	CA	1145	C
31	CA	1146	A
31	CA	1147	C
31	CA	1151	A
31	CA	1152	A
31	CA	1157	A
31	CA	1159	U
31	CA	1160	G
31	CA	1166	G
31	CA	1174	G
31	CA	1180	A
31	CA	1181	G
31	CA	1182	G
31	CA	1193	G
31	CA	1196	U
31	CA	1197	G
31	CA	1201	A
31	CA	1202	G

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Mol	Chain	Res	Type
31	CA	1203	C
31	CA	1209	C
31	CA	1212	U
31	CA	1213	A
31	CA	1214	C
31	CA	1215	G
31	CA	1225	A
31	CA	1227	A
31	CA	1238	A
31	CA	1240	U
31	CA	1248	A
31	CA	1250	A
31	CA	1256	A
31	CA	1257	U
31	CA	1258	G
31	CA	1263	C
31	CA	1270	C
31	CA	1273	G
31	CA	1278	U
31	CA	1280	A
31	CA	1282	C
31	CA	1286	A
31	CA	1287	A
31	CA	1292	U
31	CA	1294	G
31	CA	1297	C
31	CA	1300	G
31	CA	1301	U
31	CA	1302	U
31	CA	1305	G
31	CA	1317	C
31	CA	1320	C
31	CA	1322	C
31	CA	1323	G
31	CA	1326	C
31	CA	1331	G
31	CA	1336	C
31	CA	1338	G
31	CA	1346	A
31	CA	1347	G
31	CA	1351	U
31	CA	1353	G

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Mol	Chain	Res	Type
31	CA	1358	U
31	CA	1360	A
31	CA	1363	C
31	CA	1363(A)	A
31	CA	1364	U
31	CA	1365	G
31	CA	1370	G
31	CA	1379	G
31	CA	1381	U
31	CA	1397	C
31	CA	1398	A
31	CA	1399	C
31	CA	1400	C
31	CA	1402	C
31	CA	1419	G
31	CA	1423	G
31	CA	1442	G
31	CA	1442(A)	G
31	CA	1442(B)	A
31	CA	1447	A
31	CA	1452	C
31	CA	1456	G
31	CA	1457	G
31	CA	1487	G
31	CA	1492	A
31	CA	1493	A
31	CA	1494	G
31	CA	1497	G
31	CA	1502	A
31	CA	1503	A
31	CA	1504	G
31	CA	1505	G
31	CA	1506	U
31	CA	1507	A
31	CA	1517	G
31	CA	1519	A
31	CA	1520	G
31	CA	1529	G
31	CA	1530	G
31	DA	4	U
31	DA	7	G
31	DA	8	A

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Mol	Chain	Res	Type
31	DA	9	G
31	DA	10	A
31	DA	14	U
31	DA	22	G
31	DA	26	A
31	DA	32	A
31	DA	39	G
31	DA	47	C
31	DA	48	C
31	DA	51	A
31	DA	61	G
31	DA	70	G
31	DA	72	C
31	DA	76	C
31	DA	77	G
31	DA	78	G
31	DA	79	G
31	DA	92	C
31	DA	96	U
31	DA	97	G
31	DA	98	G
31	DA	101	A
31	DA	115	G
31	DA	116	A
31	DA	119	A
31	DA	120	A
31	DA	121	C
31	DA	131	C
31	DA	144	G
31	DA	146	G
31	DA	150	C
31	DA	162	A
31	DA	163	C
31	DA	167	G
31	DA	169	C
31	DA	173	U
31	DA	182	U
31	DA	189(F)	U
31	DA	189(G)	G
31	DA	189(H)	G
31	DA	195	A
31	DA	197	A

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Mol	Chain	Res	Type
31	DA	201	C
31	DA	203	U
31	DA	204	U
31	DA	216	G
31	DA	217	C
31	DA	231	G
31	DA	243	A
31	DA	244	U
31	DA	247	G
31	DA	251	G
31	DA	266	G
31	DA	267	C
31	DA	280	C
31	DA	281	G
31	DA	289	G
31	DA	301	G
31	DA	321	A
31	DA	328	C
31	DA	330	C
31	DA	332	G
31	DA	341	C
31	DA	344	A
31	DA	345	C
31	DA	347	G
31	DA	348	G
31	DA	350	G
31	DA	351	G
31	DA	352	C
31	DA	353	A
31	DA	354	G
31	DA	355	C
31	DA	363	A
31	DA	366	C
31	DA	367	U
31	DA	372	C
31	DA	373	A
31	DA	392	G
31	DA	397	A
31	DA	398	C
31	DA	406	G
31	DA	411	A
31	DA	412	A

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Mol	Chain	Res	Type
31	DA	413	G
31	DA	414	A
31	DA	421	U
31	DA	422	C
31	DA	423	G
31	DA	424	G
31	DA	428	G
31	DA	429	U
31	DA	433	C
31	DA	435	C
31	DA	437	U
31	DA	439	A
31	DA	442	C
31	DA	452	A
31	DA	461	A
31	DA	473	G
31	DA	475	G
31	DA	476	G
31	DA	483	C
31	DA	484	G
31	DA	485	G
31	DA	496	A
31	DA	498	U
31	DA	505	G
31	DA	509	A
31	DA	510	A
31	DA	511	C
31	DA	518	C
31	DA	527	G
31	DA	531	U
31	DA	532	A
31	DA	533	A
31	DA	534	U
31	DA	547	A
31	DA	559	A
31	DA	560	U
31	DA	561	U
31	DA	562	C
31	DA	567	G
31	DA	572	A
31	DA	573	A
31	DA	575	G

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Mol	Chain	Res	Type
31	DA	576	G
31	DA	596	C
31	DA	607	A
31	DA	623	C
31	DA	630	G
31	DA	631	G
31	DA	632	A
31	DA	653	A
31	DA	665	A
31	DA	687	A
31	DA	688	G
31	DA	723	U
31	DA	724	G
31	DA	731	G
31	DA	749	C
31	DA	753	A
31	DA	755	G
31	DA	777	A
31	DA	786	G
31	DA	792	A
31	DA	793	U
31	DA	794	A
31	DA	801	U
31	DA	810	C
31	DA	817	C
31	DA	821	G
31	DA	827	U
31	DA	828	A
31	DA	829	G
31	DA	833	U
31	DA	836	G
31	DA	839	U
31	DA	840	C
31	DA	841	U
31	DA	848	C
31	DA	851	G
31	DA	859	A
31	DA	870	U
31	DA	872	A
31	DA	902	G
31	DA	914	A
31	DA	926	G

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Mol	Chain	Res	Type
31	DA	927	G
31	DA	934	C
31	DA	935	A
31	DA	940	C
31	DA	942	G
31	DA	958	A
31	DA	960	U
31	DA	961	U
31	DA	966	G
31	DA	969	A
31	DA	971	G
31	DA	974	A
31	DA	975	A
31	DA	976	G
31	DA	977	A
31	DA	978	A
31	DA	991	U
31	DA	992	U
31	DA	993	G
31	DA	997	U
31	DA	1000	U
31	DA	1001	A
31	DA	1003	G
31	DA	1004	A
31	DA	1005	A
31	DA	1010	G
31	DA	1011	G
31	DA	1012	U
31	DA	1023	G
31	DA	1025	U
31	DA	1026	G
31	DA	1027	C
31	DA	1028	C
31	DA	1029	C
31	DA	1030	C
31	DA	1030(A)	G
31	DA	1030(B)	C
31	DA	1030(C)	G
31	DA	1031	G
31	DA	1032	G
31	DA	1033	G
31	DA	1034	G

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Mol	Chain	Res	Type
31	DA	1036	G
31	DA	1037	C
31	DA	1050	G
31	DA	1051	C
31	DA	1054	C
31	DA	1064	G
31	DA	1065	U
31	DA	1066	C
31	DA	1068	G
31	DA	1081	G
31	DA	1086	U
31	DA	1087	G
31	DA	1091	U
31	DA	1092	A
31	DA	1094	G
31	DA	1095	U
31	DA	1100	C
31	DA	1101	A
31	DA	1112	C
31	DA	1117	G
31	DA	1121	U
31	DA	1124	G
31	DA	1125	U
31	DA	1126	U
31	DA	1127	G
31	DA	1129	C
31	DA	1136	U
31	DA	1137	C
31	DA	1138	G
31	DA	1139	G
31	DA	1140	C
31	DA	1143	G
31	DA	1146	A
31	DA	1147	C
31	DA	1152	A
31	DA	1154	G
31	DA	1159	U
31	DA	1160	G
31	DA	1161	C
31	DA	1166	G
31	DA	1170	A
31	DA	1171	G

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Mol	Chain	Res	Type
31	DA	1180	A
31	DA	1181	G
31	DA	1182	G
31	DA	1184	G
31	DA	1185	G
31	DA	1186	G
31	DA	1188	A
31	DA	1196	U
31	DA	1197	G
31	DA	1200	C
31	DA	1202	G
31	DA	1209	C
31	DA	1211	U
31	DA	1212	U
31	DA	1213	A
31	DA	1225	A
31	DA	1226	C
31	DA	1236	A
31	DA	1238	A
31	DA	1240	U
31	DA	1256	A
31	DA	1257	U
31	DA	1258	G
31	DA	1260	C
31	DA	1268	A
31	DA	1272	G
31	DA	1273	G
31	DA	1278	U
31	DA	1280	A
31	DA	1281	U
31	DA	1282	C
31	DA	1284	C
31	DA	1286	A
31	DA	1287	A
31	DA	1290	G
31	DA	1293	G
31	DA	1294	G
31	DA	1297	C
31	DA	1299	A
31	DA	1300	G
31	DA	1301	U
31	DA	1302	U

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Mol	Chain	Res	Type
31	DA	1312	G
31	DA	1317	C
31	DA	1320	C
31	DA	1322	C
31	DA	1323	G
31	DA	1331	G
31	DA	1336	C
31	DA	1338	G
31	DA	1346	A
31	DA	1347	G
31	DA	1358	U
31	DA	1363	C
31	DA	1364	U
31	DA	1368	G
31	DA	1379	G
31	DA	1386	G
31	DA	1389	C
31	DA	1397	C
31	DA	1399	C
31	DA	1419	G
31	DA	1442	G
31	DA	1442(A)	G
31	DA	1442(B)	A
31	DA	1447	A
31	DA	1452	C
31	DA	1456	G
31	DA	1457	G
31	DA	1487	G
31	DA	1492	A
31	DA	1493	A
31	DA	1494	G
31	DA	1497	G
31	DA	1502	A
31	DA	1503	A
31	DA	1504	G
31	DA	1505	G
31	DA	1506	U
31	DA	1507	A
31	DA	1508	G
31	DA	1517	G
31	DA	1519	A
31	DA	1520	G

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Mol	Chain	Res	Type
31	DA	1529	G
31	DA	1530	G

All (177) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	AA	9	U
1	AA	71	A
1	AA	102	G
1	AA	196	A
1	AA	221	A
1	AA	271(K)	U
1	AA	271(M)	G
1	AA	278	A
1	AA	310	A
1	AA	363(E)	U
1	AA	474	G
1	AA	481	G
1	AA	587	C
1	AA	669	G
1	AA	752	A
1	AA	827	U
1	AA	856	C
1	AA	900	A
1	AA	1047	G
1	AA	1049	C
1	AA	1106	A
1	AA	1108	U
1	AA	1174	A
1	AA	1175	U
1	AA	1176	G
1	AA	1210	A
1	AA	1240	U
1	AA	1300	U
1	AA	1301	A
1	AA	1378	A
1	AA	1420	U
1	AA	1427	A
1	AA	1507	A
1	AA	1530	C
1	AA	1558	A
1	AA	1559	G

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Mol	Chain	Res	Type
1	AA	1608	A
1	AA	1653	G
1	AA	1786	A
1	AA	1799	G
1	AA	1819	A
1	AA	1992	G
1	AA	2122	U
1	AA	2126	A
1	AA	2171	A
1	AA	2172	U
1	AA	2207	G
1	AA	2218	U
1	AA	2282	G
1	AA	2318	G
1	AA	2405	G
1	AA	2439	A
1	AA	2610	C
1	AA	2689	U
1	AA	2726	U
1	AA	2756	U
1	AA	2778	A
1	AA	2789	C
1	AA	2802	G
1	BA	9	U
1	BA	71	A
1	BA	102	G
1	BA	196	A
1	BA	221	A
1	BA	249	C
1	BA	271(K)	U
1	BA	271(M)	G
1	BA	278	A
1	BA	363(E)	U
1	BA	405	U
1	BA	474	G
1	BA	481	G
1	BA	587	C
1	BA	746	A
1	BA	752	A
1	BA	827	U
1	BA	856	C
1	BA	900	A

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Mol	Chain	Res	Type
1	BA	974	G
1	BA	1026	U
1	BA	1047	G
1	BA	1049	C
1	BA	1106	A
1	BA	1108	U
1	BA	1174	A
1	BA	1175	U
1	BA	1176	G
1	BA	1210	A
1	BA	1300	U
1	BA	1301	A
1	BA	1378	A
1	BA	1420	U
1	BA	1427	A
1	BA	1507	A
1	BA	1530	C
1	BA	1558	A
1	BA	1608	A
1	BA	1653	G
1	BA	1799	G
1	BA	1819	A
1	BA	1914	C
1	BA	1992	G
1	BA	2104	G
1	BA	2122	U
1	BA	2126	A
1	BA	2171	A
1	BA	2172	U
1	BA	2207	G
1	BA	2218	U
1	BA	2282	G
1	BA	2318	G
1	BA	2319	G
1	BA	2405	G
1	BA	2439	A
1	BA	2610	C
1	BA	2689	U
1	BA	2778	A
1	BA	2789	C
1	BA	2802	G
31	CA	60	A

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Mol	Chain	Res	Type
31	CA	76	C
31	CA	93	G
31	CA	96	U
31	CA	115	G
31	CA	119	A
31	CA	243	A
31	CA	250	A
31	CA	266	G
31	CA	344	A
31	CA	366	C
31	CA	428	G
31	CA	438	G
31	CA	509	A
31	CA	533	A
31	CA	560	U
31	CA	687	A
31	CA	748	C
31	CA	793	U
31	CA	913	A
31	CA	991	U
31	CA	1064	G
31	CA	1065	U
31	CA	1067	A
31	CA	1165	C
31	CA	1201	A
31	CA	1279	A
31	CA	1285	A
31	CA	1300	G
31	CA	1492	A
31	CA	1504	G
31	DA	60	A
31	DA	76	C
31	DA	93	G
31	DA	96	U
31	DA	115	G
31	DA	119	A
31	DA	243	A
31	DA	266	G
31	DA	344	A
31	DA	366	C
31	DA	428	G
31	DA	509	A

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Mol	Chain	Res	Type
31	DA	533	A
31	DA	560	U
31	DA	687	A
31	DA	748	C
31	DA	913	A
31	DA	991	U
31	DA	1065	U
31	DA	1067	A
31	DA	1165	C
31	DA	1201	A
31	DA	1211	U
31	DA	1285	A
31	DA	1300	G
31	DA	1492	A
31	DA	1504	G

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
52	T8B	BA	3001	53	48,48,48	1.00	2 (4%)	63,71,71	1.28	9 (14%)
52	T8B	AA	3001	53	48,48,48	1.01	2 (4%)	63,71,71	1.28	9 (14%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
52	T8B	BA	3001	53	-	0/26/26/26	0/5/5/5
52	T8B	AA	3001	53	-	0/26/26/26	0/5/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
52	AA	3001	T8B	O11-C26	3.13	1.38	1.31
52	BA	3001	T8B	O11-C26	3.12	1.38	1.31
52	BA	3001	T8B	C25-C24	2.27	1.50	1.43
52	AA	3001	T8B	C25-C24	2.25	1.50	1.43

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
52	BA	3001	T8B	C21-C23-C24	-4.27	120.19	125.09
52	AA	3001	T8B	C21-C23-C24	-4.25	120.21	125.09
52	BA	3001	T8B	O5-C14-O6	3.22	120.29	116.45
52	AA	3001	T8B	O5-C14-O6	3.22	120.29	116.45
52	BA	3001	T8B	O11-C26-C25	2.94	124.96	121.09
52	AA	3001	T8B	O11-C26-C25	2.92	124.92	121.09
52	AA	3001	T8B	O5-C11-C12	2.65	117.33	112.19
52	BA	3001	T8B	O5-C11-C12	2.64	117.32	112.19
52	BA	3001	T8B	C28-C27-C32	2.62	121.45	118.15
52	AA	3001	T8B	C28-C27-C32	2.62	121.44	118.15
52	AA	3001	T8B	C26-C25-C24	2.31	124.06	120.81
52	BA	3001	T8B	C26-C25-C24	2.30	124.05	120.81
52	BA	3001	T8B	O4-C12-C11	-2.17	108.73	111.77
52	AA	3001	T8B	O4-C12-C11	-2.14	108.77	111.77
52	BA	3001	T8B	O5-C11-C10	2.13	124.04	122.25
52	BA	3001	T8B	O4-C12-O3	2.12	127.55	123.50
52	AA	3001	T8B	O4-C12-O3	2.10	127.51	123.50
52	AA	3001	T8B	O5-C11-C10	2.09	124.01	122.25

There are no chirality outliers.

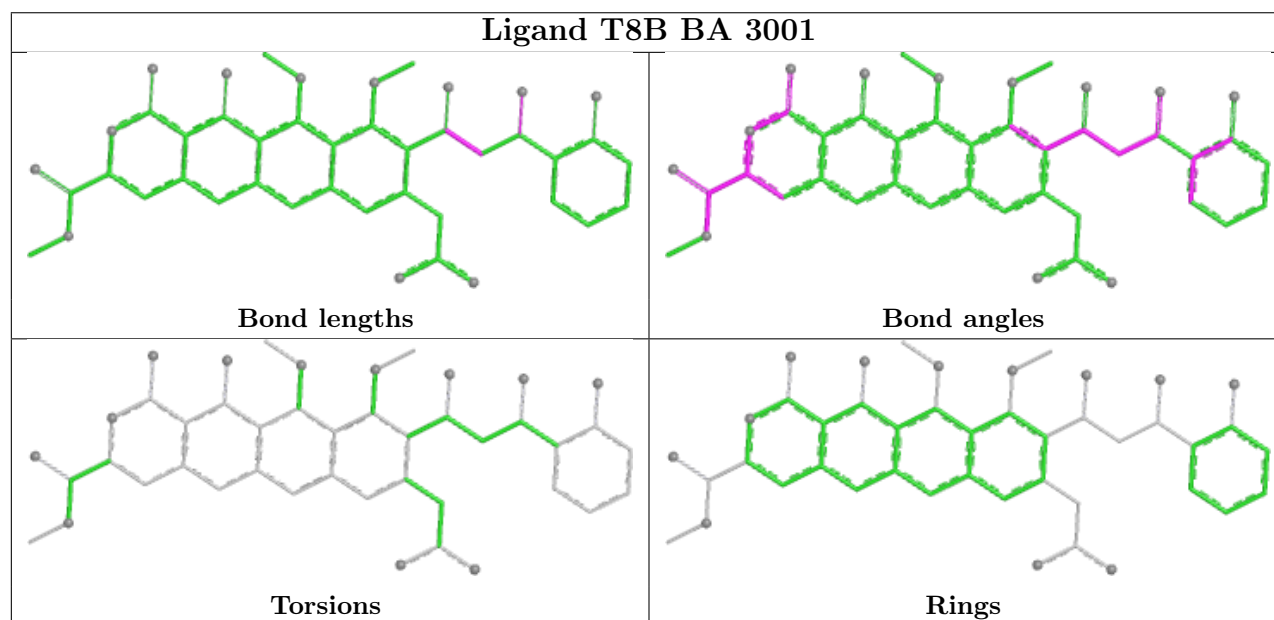
There are no torsion outliers.

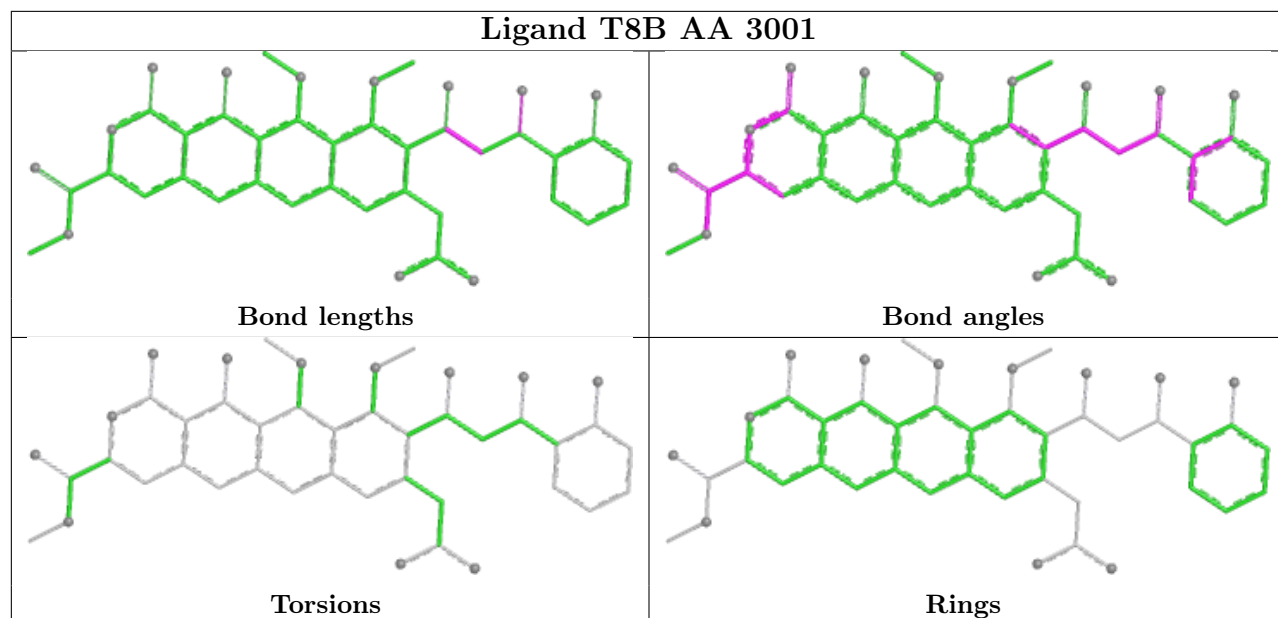
There are no ring outliers.

2 monomers are involved in 54 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
52	BA	3001	T8B	23	0
52	AA	3001	T8B	31	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	AA	2827/2915 (96%)	0.03	112 (3%) 43 29	46, 70, 113, 128	0
1	BA	2827/2915 (96%)	-0.06	93 (3%) 49 34	27, 56, 110, 126	0
2	AB	120/122 (98%)	0.77	11 (9%) 16 11	68, 95, 108, 111	0
2	BB	120/122 (98%)	0.14	3 (2%) 58 42	46, 80, 95, 103	0
3	AD	275/276 (99%)	0.20	8 (2%) 54 38	45, 67, 82, 103	0
3	BD	275/276 (99%)	0.21	5 (1%) 67 52	36, 59, 78, 103	0
4	AE	204/206 (99%)	0.08	4 (1%) 64 49	47, 72, 87, 98	0
4	BE	204/206 (99%)	0.21	5 (2%) 58 42	33, 60, 81, 96	0
5	AF	203/205 (99%)	0.12	3 (1%) 71 56	44, 77, 95, 112	0
5	BF	203/205 (99%)	0.19	3 (1%) 71 56	26, 64, 90, 110	0
6	AG	181/182 (99%)	1.51	50 (27%) 2 1	89, 106, 114, 116	0
6	BG	181/182 (99%)	0.71	22 (12%) 10 7	81, 101, 110, 118	0
7	AH	174/180 (96%)	0.84	14 (8%) 20 14	80, 93, 101, 108	0
7	BH	174/180 (96%)	0.65	14 (8%) 20 14	63, 78, 91, 98	0
8	AI	145/148 (97%)	1.31	33 (22%) 2 2	72, 104, 116, 123	0
8	BI	145/148 (97%)	0.30	3 (2%) 63 47	68, 90, 98, 100	0
9	AN	140/140 (100%)	0.19	3 (2%) 63 47	58, 73, 90, 94	0
9	BN	140/140 (100%)	0.32	4 (2%) 54 38	38, 57, 82, 85	0
10	AO	122/122 (100%)	0.15	2 (1%) 70 55	56, 73, 85, 91	0
10	BO	122/122 (100%)	0.26	4 (3%) 49 34	43, 63, 82, 89	0
11	AP	147/150 (98%)	0.51	10 (6%) 25 17	47, 81, 96, 105	0
11	BP	147/150 (98%)	0.47	9 (6%) 28 19	28, 68, 91, 100	0
12	AQ	141/141 (100%)	0.39	5 (3%) 47 33	60, 78, 91, 97	0
12	BQ	141/141 (100%)	0.20	2 (1%) 73 58	44, 65, 78, 88	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	AR	118/118 (100%)	0.23	3 (2%) 58 42	47, 66, 78, 88	0
13	BR	118/118 (100%)	0.17	3 (2%) 58 42	35, 53, 69, 86	0
14	AS	110/112 (98%)	1.19	20 (18%) 4 3	74, 91, 102, 109	0
14	BS	110/112 (98%)	0.54	6 (5%) 32 21	58, 79, 93, 100	0
15	AT	131/146 (89%)	0.23	3 (2%) 61 44	66, 76, 99, 109	0
15	BT	131/146 (89%)	0.39	6 (4%) 38 26	55, 68, 91, 102	0
16	AU	116/118 (98%)	0.11	1 (0%) 81 68	53, 70, 84, 89	0
16	BU	116/118 (98%)	0.05	1 (0%) 81 68	34, 50, 70, 84	0
17	AV	101/101 (100%)	0.09	1 (0%) 79 66	48, 79, 93, 103	0
17	BV	101/101 (100%)	0.05	1 (0%) 79 66	31, 61, 81, 95	0
18	AW	112/113 (99%)	0.01	2 (1%) 67 52	50, 59, 80, 104	0
18	BW	112/113 (99%)	0.07	1 (0%) 81 68	36, 45, 75, 106	0
19	AX	95/96 (98%)	0.52	5 (5%) 33 22	54, 71, 89, 93	0
19	BX	95/96 (98%)	0.33	2 (2%) 63 47	33, 59, 83, 91	0
20	AY	107/110 (97%)	0.64	8 (7%) 22 15	71, 81, 94, 105	0
20	BY	107/110 (97%)	0.61	8 (7%) 22 15	56, 71, 89, 103	0
21	AZ	198/206 (96%)	0.55	7 (3%) 47 33	80, 92, 103, 108	0
21	BZ	198/206 (96%)	0.35	7 (3%) 47 33	64, 82, 97, 103	0
22	A0	76/85 (89%)	0.32	2 (2%) 57 41	59, 74, 84, 89	0
22	B0	76/85 (89%)	0.20	1 (1%) 74 60	46, 60, 74, 82	0
23	A1	97/98 (98%)	0.55	6 (6%) 28 18	54, 71, 94, 99	0
23	B1	97/98 (98%)	0.31	3 (3%) 51 36	42, 65, 91, 95	0
24	A2	70/72 (97%)	0.48	5 (7%) 23 16	65, 81, 92, 101	0
24	B2	70/72 (97%)	0.48	6 (8%) 18 12	52, 70, 85, 102	0
25	A3	59/60 (98%)	0.62	6 (10%) 13 9	63, 74, 92, 101	0
25	B3	59/60 (98%)	0.36	1 (1%) 69 53	43, 56, 84, 94	0
26	A4	46/71 (64%)	1.02	5 (10%) 12 9	101, 109, 113, 119	0
26	B4	46/71 (64%)	0.78	4 (8%) 17 12	98, 107, 113, 118	0
27	A5	59/60 (98%)	0.10	0 100 100	46, 64, 78, 91	0
27	B5	59/60 (98%)	0.19	0 100 100	28, 52, 71, 89	0
28	A6	53/54 (98%)	0.37	4 (7%) 22 15	62, 76, 87, 89	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
28	B6	53/54 (98%)	0.22	0	100 100	52, 66, 77, 85	0
29	A7	48/49 (97%)	0.27	4 (8%)	19 13	43, 54, 79, 99	0
29	B7	48/49 (97%)	0.17	2 (4%)	41 28	29, 44, 73, 89	0
30	A8	64/65 (98%)	0.67	6 (9%)	15 11	60, 68, 76, 85	0
30	B8	64/65 (98%)	0.44	2 (3%)	51 36	45, 56, 65, 79	0
31	CA	1498/1521 (98%)	0.76	160 (10%)	12 9	59, 99, 120, 126	0
31	DA	1498/1521 (98%)	0.72	136 (9%)	16 11	65, 99, 120, 128	0
32	CB	229/256 (89%)	0.67	13 (5%)	30 20	93, 103, 111, 117	0
32	DB	229/256 (89%)	0.96	26 (11%)	11 8	95, 105, 112, 115	0
33	CC	206/239 (86%)	1.15	33 (16%)	6 4	92, 106, 112, 115	0
33	DC	206/239 (86%)	1.20	38 (18%)	4 3	97, 108, 115, 120	0
34	CD	208/209 (99%)	1.24	42 (20%)	3 3	85, 98, 108, 114	0
34	DD	208/209 (99%)	1.15	39 (18%)	4 3	87, 97, 107, 123	0
35	CE	148/162 (91%)	0.37	3 (2%)	64 49	74, 93, 103, 110	0
35	DE	148/162 (91%)	0.63	12 (8%)	19 13	84, 96, 104, 110	0
36	CF	100/101 (99%)	0.46	4 (4%)	43 29	80, 92, 101, 107	0
36	DF	100/101 (99%)	0.58	6 (6%)	29 19	84, 93, 103, 110	0
37	CG	155/156 (99%)	1.93	60 (38%)	1 1	99, 109, 115, 120	0
37	DG	155/156 (99%)	1.58	45 (29%)	1 1	97, 109, 115, 119	0
38	CH	138/138 (100%)	0.53	6 (4%)	40 27	82, 94, 101, 104	0
38	DH	138/138 (100%)	0.76	8 (5%)	30 20	86, 96, 102, 106	0
39	CI	125/128 (97%)	2.10	59 (47%)	0 1	97, 111, 117, 120	0
39	DI	125/128 (97%)	2.58	78 (62%)	0 1	100, 112, 118, 120	0
40	CJ	96/105 (91%)	1.89	39 (40%)	1 1	100, 109, 115, 117	0
40	DJ	96/105 (91%)	2.03	45 (46%)	0 1	97, 110, 115, 119	0
41	CK	114/129 (88%)	0.64	6 (5%)	33 22	72, 93, 101, 105	0
41	DK	114/129 (88%)	0.49	7 (6%)	28 19	75, 95, 103, 107	0
42	CL	122/132 (92%)	0.80	12 (9%)	14 10	72, 85, 96, 100	0
42	DL	122/132 (92%)	0.81	12 (9%)	14 10	77, 87, 96, 105	0
43	CM	114/126 (90%)	2.21	56 (49%)	0 1	102, 111, 118, 126	0
43	DM	114/126 (90%)	1.89	43 (37%)	1 1	101, 109, 115, 117	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
44	CN	60/61 (98%)	2.08	28 (46%) 0 1	99, 108, 116, 118	0
44	DN	60/61 (98%)	2.65	37 (61%) 0 1	103, 111, 116, 122	0
45	CO	88/89 (98%)	0.48	5 (5%) 30 20	72, 91, 101, 107	0
45	DO	88/89 (98%)	0.65	7 (7%) 20 14	81, 93, 103, 107	0
46	CP	82/88 (93%)	1.06	11 (13%) 8 6	88, 96, 106, 112	0
46	DP	82/88 (93%)	1.26	18 (21%) 3 2	86, 94, 103, 111	0
47	CQ	99/105 (94%)	0.57	4 (4%) 43 29	78, 90, 98, 102	0
47	DQ	99/105 (94%)	0.67	6 (6%) 28 19	79, 92, 100, 103	0
48	CR	68/88 (77%)	0.65	6 (8%) 17 12	80, 91, 102, 103	0
48	DR	68/88 (77%)	0.42	1 (1%) 71 56	85, 93, 104, 106	0
49	CS	78/93 (83%)	2.39	47 (60%) 0 1	107, 111, 117, 123	0
49	DS	78/93 (83%)	1.92	33 (42%) 1 1	91, 112, 117, 119	0
50	CT	96/106 (90%)	0.91	16 (16%) 5 4	84, 93, 99, 102	0
50	DT	96/106 (90%)	0.85	13 (13%) 8 6	82, 92, 100, 101	0
51	CU	23/27 (85%)	4.80	22 (95%) 0 0	106, 112, 117, 119	0
51	DU	23/27 (85%)	3.68	21 (91%) 0 0	104, 109, 112, 113	0
All	All	20372/21160 (96%)	0.51	1827 (8%) 17 11	26, 84, 115, 128	0

All (1827) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
31	CA	1286	A	9.2
51	CU	14	TRP	8.8
49	CS	40	ILE	8.7
39	DI	126	SER	8.2
31	CA	1353	G	8.1
51	CU	23	PRO	8.0
39	DI	119	ALA	7.7
39	CI	12	GLU	7.6
31	DA	1149	C	7.5
6	AG	155	MET	7.5
39	CI	126	SER	7.5
37	CG	8	GLU	7.4
31	CA	1288	A	7.0
39	DI	105	ASP	6.9
39	CI	106	ALA	6.9

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Mol	Chain	Res	Type	RSRZ
39	CI	109	VAL	6.8
51	CU	24	ARG	6.8
1	BA	2116	G	6.7
44	DN	30	ALA	6.7
20	BY	1	MET	6.6
31	CA	1290	G	6.4
51	CU	20	LYS	6.4
31	DA	1287	A	6.4
8	AI	121	LYS	6.4
23	A1	2	SER	6.3
34	DD	19	LEU	6.3
10	BO	5	GLN	6.3
43	DM	7	VAL	6.2
31	DA	1286	A	6.2
1	AA	2125	G	6.1
44	DN	2	ALA	6.1
31	CA	1287	A	6.1
1	BA	2172	U	6.1
34	CD	31	CYS	6.1
51	CU	17	THR	6.1
51	CU	13	ILE	6.0
39	DI	109	VAL	6.0
31	DA	1148	U	6.0
1	AA	2147	G	5.9
43	CM	96	LEU	5.9
8	AI	100	ALA	5.9
51	DU	6	ARG	5.9
39	DI	108	VAL	5.9
51	CU	3	LYS	5.9
34	CD	12	CYS	5.8
43	CM	87	TYR	5.8
34	CD	4	TYR	5.8
31	DA	1353	G	5.7
34	CD	26	CYS	5.7
11	AP	66	GLY	5.6
1	BA	2168	G	5.6
34	DD	26	CYS	5.6
26	B4	17	GLY	5.6
37	CG	120	ILE	5.6
51	DU	13	ILE	5.6
43	CM	102	ARG	5.6
31	DA	1202	G	5.6

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Mol	Chain	Res	Type	RSRZ
51	DU	11	GLY	5.6
43	CM	107	ALA	5.5
8	AI	119	PRO	5.5
37	DG	5	ARG	5.5
51	DU	14	TRP	5.5
8	AI	79	ILE	5.4
3	AD	276	LYS	5.4
39	DI	17	VAL	5.4
37	CG	115	ARG	5.4
39	DI	12	GLU	5.4
37	CG	9	VAL	5.3
37	DG	8	GLU	5.3
49	CS	13	ASP	5.3
31	DA	79	G	5.3
39	CI	72	GLY	5.3
42	CL	91	LYS	5.3
39	DI	111	ARG	5.3
37	CG	118	VAL	5.2
34	DD	12	CYS	5.2
40	CJ	44	VAL	5.2
49	CS	76	PRO	5.2
51	DU	8	THR	5.2
1	BA	2125	G	5.2
6	AG	41	GLN	5.2
1	AA	2174	C	5.1
44	CN	16	PHE	5.1
49	CS	49	ILE	5.1
1	BA	652(G)	G	5.1
51	CU	8	THR	5.1
34	DD	69	GLY	5.1
39	DI	70	LYS	5.1
31	CA	1352	C	5.1
37	DG	27	ILE	5.0
39	CI	119	ALA	5.0
37	CG	13	GLN	5.0
40	DJ	65	LEU	5.0
51	CU	9	ARG	5.0
39	CI	105	ASP	5.0
31	DA	1036	G	5.0
1	BA	2153	G	4.9
39	DI	11	LYS	4.9
1	BA	2117	A	4.9

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Mol	Chain	Res	Type	RSRZ
51	CU	18	TYR	4.9
34	DD	206	PHE	4.9
1	BA	2167	U	4.8
34	CD	9	CYS	4.8
14	AS	33	LYS	4.8
6	AG	28	VAL	4.8
33	CC	90	GLU	4.8
37	CG	101	LEU	4.8
31	CA	1371	G	4.8
31	CA	1354	C	4.8
43	DM	5	ALA	4.8
39	CI	116	LYS	4.8
31	CA	1349	A	4.7
44	DN	14	PRO	4.7
39	DI	106	ALA	4.7
44	DN	34	TYR	4.7
34	CD	181	MET	4.7
1	AA	2124	G	4.7
1	BA	652(I)	C	4.7
39	DI	13	ALA	4.7
44	DN	28	GLY	4.6
37	DG	38	LEU	4.6
31	CA	1030(B)	C	4.6
40	DJ	20	ALA	4.6
40	CJ	49	VAL	4.6
39	DI	7	THR	4.6
8	AI	146	ALA	4.6
39	DI	41	VAL	4.6
51	DU	3	LYS	4.6
38	CH	1	MET	4.6
39	DI	10	ARG	4.6
51	CU	15	ARG	4.6
32	DB	133	LYS	4.6
31	CA	1149	C	4.6
1	BA	2173	A	4.6
34	DD	181	MET	4.6
44	CN	61	TRP	4.6
40	DJ	58	ASP	4.6
1	BA	2108	C	4.6
31	CA	1289	A	4.6
37	CG	42	ILE	4.5
42	CL	28	LYS	4.5

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Mol	Chain	Res	Type	RSRZ
1	AA	2110	G	4.5
1	BA	652(J)	G	4.5
39	CI	123	PRO	4.5
43	CM	27	LYS	4.5
51	CU	12	LYS	4.5
37	CG	117	ALA	4.5
43	CM	30	ALA	4.5
37	CG	25	ALA	4.4
49	CS	75	ALA	4.4
44	CN	13	THR	4.4
49	CS	74	PHE	4.4
1	BA	2147	G	4.4
49	CS	41	VAL	4.4
39	DI	37	PHE	4.4
3	BD	276	LYS	4.4
8	AI	84	GLY	4.4
39	DI	115	GLY	4.4
24	A2	52	ASP	4.4
40	DJ	49	VAL	4.4
31	DA	91	C	4.4
44	DN	8	GLU	4.4
4	BE	194	GLY	4.4
49	DS	13	ASP	4.4
29	B7	48	LYS	4.4
1	BA	888	C	4.4
14	AS	3	ARG	4.3
39	DI	66	ARG	4.3
51	CU	21	TYR	4.3
44	CN	15	LYS	4.3
34	CD	2	GLY	4.3
49	DS	65	ASN	4.3
39	CI	108	VAL	4.3
44	DN	10	ALA	4.3
40	CJ	50	ILE	4.3
51	CU	11	GLY	4.3
8	AI	93	THR	4.3
49	DS	9	VAL	4.3
51	CU	10	ARG	4.3
1	AA	2162	G	4.3
33	CC	2	GLY	4.3
40	DJ	30	SER	4.3
49	DS	11	VAL	4.3

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Mol	Chain	Res	Type	RSRZ
39	DI	125	TYR	4.3
1	AA	2170	A	4.3
31	DA	97	G	4.3
39	CI	30	GLY	4.3
37	CG	12	LEU	4.3
40	DJ	71	LEU	4.3
31	DA	1028	C	4.3
33	CC	39	ILE	4.3
39	DI	118	LYS	4.2
44	DN	11	LYS	4.2
1	AA	2148	G	4.2
32	CB	163	PHE	4.2
51	CU	16	GLY	4.2
51	DU	9	ARG	4.2
9	AN	44	PRO	4.2
37	CG	116	ALA	4.2
43	CM	101	GLN	4.2
31	DA	1311	G	4.2
49	CS	58	VAL	4.2
51	CU	6	ARG	4.2
49	DS	68	GLY	4.2
6	AG	2	PRO	4.2
31	CA	985	C	4.2
43	CM	45	VAL	4.2
11	AP	65	ARG	4.2
1	BA	2171	A	4.2
43	CM	10	PRO	4.2
51	DU	23	PRO	4.2
33	DC	152	ILE	4.2
6	AG	134	GLY	4.1
31	DA	78	G	4.1
37	CG	37	ASN	4.1
39	DI	40	LEU	4.1
40	DJ	57	LYS	4.1
44	DN	15	LYS	4.1
24	B2	70	GLN	4.1
49	CS	42	PRO	4.1
39	DI	112	LYS	4.1
31	DA	344	A	4.1
37	CG	66	VAL	4.1
43	CM	98	VAL	4.1
1	AA	2108	C	4.1

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Mol	Chain	Res	Type	RSRZ
31	DA	1128	C	4.1
39	CI	11	LYS	4.1
31	DA	1034	G	4.1
31	CA	1248	A	4.1
44	CN	41	ARG	4.1
1	AA	652(I)	C	4.1
31	DA	1354	C	4.1
46	DP	82	GLN	4.1
31	DA	1310	G	4.1
50	CT	10	LEU	4.1
1	AA	2126	A	4.1
31	DA	1352	C	4.0
37	CG	10	ARG	4.0
30	A8	53	PRO	4.0
1	AA	652(G)	G	4.0
37	CG	36	LYS	4.0
39	DI	107	ARG	4.0
35	DE	20	GLN	4.0
6	AG	76	SER	4.0
44	DN	25	VAL	4.0
39	DI	116	LYS	4.0
39	CI	40	LEU	4.0
37	DG	14	PRO	4.0
37	CG	21	VAL	4.0
39	DI	113	LYS	4.0
51	DU	17	THR	4.0
43	CM	99	ARG	4.0
15	BT	37	GLY	4.0
1	AA	2107	C	4.0
31	DA	1030(B)	C	4.0
37	DG	40	ALA	4.0
40	CJ	65	LEU	4.0
49	CS	77	THR	4.0
43	DM	101	GLN	3.9
31	DA	1150	U	3.9
31	CA	1183	A	3.9
42	CL	29	GLY	3.9
43	DM	100	GLY	3.9
6	AG	133	LEU	3.9
1	AA	2118	U	3.9
39	CI	13	ALA	3.9
1	AA	2176	A	3.9

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Mol	Chain	Res	Type	RSRZ
34	DD	209	ARG	3.9
43	CM	31	LYS	3.9
1	AA	652(H)	C	3.9
31	CA	1030	C	3.9
51	CU	5	ASP	3.9
41	DK	126	ARG	3.9
1	BA	548	A	3.9
40	DJ	33	GLN	3.9
47	DQ	59	ILE	3.9
34	DD	177	ASP	3.9
34	DD	190	ASP	3.9
31	CA	1363	C	3.9
1	BA	2155	G	3.9
32	DB	163	PHE	3.9
44	DN	16	PHE	3.9
46	CP	69	THR	3.9
1	AA	2506	U	3.9
34	DD	21	LEU	3.9
1	AA	2320	A	3.9
39	CI	111	ARG	3.9
42	DL	71	PRO	3.8
31	DA	76	C	3.8
46	DP	13	HIS	3.8
25	A3	27	GLY	3.8
39	CI	70	LYS	3.8
40	DJ	37	PRO	3.8
43	CM	53	VAL	3.8
34	CD	152	SER	3.8
1	AA	652(S)	C	3.8
31	DA	1027	C	3.8
31	DA	1363	C	3.8
39	CI	107	ARG	3.8
41	CK	60	ALA	3.8
44	DN	31	ARG	3.8
39	DI	14	VAL	3.8
49	DS	16	LEU	3.8
29	A7	48	LYS	3.8
39	DI	9	ARG	3.8
1	BA	2174	C	3.8
39	CI	117	HIS	3.8
51	CU	4	GLY	3.8
31	CA	1350	A	3.8

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Mol	Chain	Res	Type	RSRZ
31	DA	1248	A	3.8
37	DG	109	ASN	3.8
44	CN	14	PRO	3.8
34	CD	177	ASP	3.8
31	DA	1235	U	3.8
43	CM	100	GLY	3.8
49	CS	35	SER	3.7
33	DC	3	ASN	3.7
43	DM	10	PRO	3.7
43	DM	102	ARG	3.7
51	DU	22	ARG	3.7
31	CA	1129	C	3.7
44	DN	21	TYR	3.7
37	DG	13	GLN	3.7
33	CC	26	LYS	3.7
3	AD	28	GLU	3.7
43	CM	8	GLU	3.7
14	AS	55	ALA	3.7
37	CG	99	LEU	3.7
37	DG	42	ILE	3.7
39	DI	110	GLU	3.7
1	AA	2119	A	3.7
31	CA	1317	C	3.7
42	DL	72	GLY	3.7
34	CD	21	LEU	3.7
49	CS	51	VAL	3.7
51	DU	10	ARG	3.7
40	CJ	35	SER	3.7
1	BA	2124	G	3.7
1	BA	2152	G	3.7
31	CA	947	G	3.7
33	CC	158	GLY	3.6
43	DM	60	VAL	3.6
45	CO	88	ARG	3.6
37	DG	106	GLN	3.6
37	DG	84	ASN	3.6
7	BH	166	GLY	3.6
51	DU	20	LYS	3.6
39	DI	102	LEU	3.6
43	CM	56	LEU	3.6
1	BA	2112	G	3.6
32	CB	96	ARG	3.6

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Mol	Chain	Res	Type	RSRZ
40	CJ	56	HIS	3.6
49	CS	69	HIS	3.6
37	CG	11	GLN	3.6
43	CM	12	ASN	3.6
43	CM	85	GLY	3.6
1	BA	2119	A	3.6
8	AI	27	ARG	3.6
37	CG	22	LEU	3.6
39	CI	110	GLU	3.6
39	DI	5	TYR	3.6
1	BA	2133	G	3.6
31	CA	971	G	3.6
43	DM	105	THR	3.6
1	BA	2130	U	3.6
31	CA	1247	U	3.6
37	CG	35	LYS	3.6
37	CG	7	ALA	3.6
37	DG	124	LEU	3.6
39	DI	80	GLY	3.6
41	CK	90	GLY	3.6
49	CS	84	GLY	3.6
1	AA	275	G	3.6
1	AA	2139	C	3.6
31	CA	1002	G	3.6
31	CA	1224	G	3.6
43	DM	98	VAL	3.6
1	BA	2135	A	3.5
37	DG	2	ALA	3.5
34	CD	155	LEU	3.5
37	DG	16	LEU	3.5
1	BA	652(H)	C	3.5
31	CA	1030(A)	G	3.5
31	CA	1138	G	3.5
31	DA	98	G	3.5
14	BS	3	ARG	3.5
20	BY	16	ALA	3.5
49	DS	15	LEU	3.5
50	CT	12	ALA	3.5
1	BA	887	A	3.5
40	DJ	17	ASP	3.5
1	BA	2113	U	3.5
31	CA	1148	U	3.5

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Mol	Chain	Res	Type	RSRZ
39	DI	71	SER	3.5
50	DT	11	SER	3.5
32	CB	133	LYS	3.5
41	DK	117	ASN	3.5
37	CG	6	ARG	3.5
39	DI	123	PRO	3.5
37	DG	11	GLN	3.5
40	CJ	63	PHE	3.5
32	CB	229	VAL	3.5
31	DA	1030(C)	G	3.5
34	DD	134	ASP	3.5
6	AG	88	ILE	3.5
7	AH	136	ILE	3.5
44	DN	9	LYS	3.5
34	DD	150	GLU	3.5
43	DM	73	GLU	3.5
51	DU	24	ARG	3.5
34	CD	19	LEU	3.5
50	DT	12	ALA	3.5
1	AA	2111	C	3.5
31	DA	1037	C	3.5
50	CT	74	LYS	3.5
4	BE	195	LEU	3.5
31	CA	3	G	3.5
40	CJ	37	PRO	3.5
46	DP	67	THR	3.5
34	CD	69	GLY	3.5
42	CL	126	LYS	3.5
37	CG	49	ILE	3.5
1	AA	2136	C	3.4
5	AF	56	GLU	3.4
31	DA	90	C	3.4
31	DA	1249	C	3.4
33	CC	52	LEU	3.4
33	DC	12	LEU	3.4
43	DM	97	PRO	3.4
8	AI	81	VAL	3.4
40	CJ	36	GLY	3.4
1	AA	548	A	3.4
46	DP	34	GLU	3.4
31	DA	1141	C	3.4
39	DI	63	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
40	CJ	33	GLN	3.4
1	BA	2118	U	3.4
1	AA	278	A	3.4
31	CA	1003	G	3.4
33	CC	55	VAL	3.4
34	CD	207	TYR	3.4
39	DI	62	TYR	3.4
14	AS	52	SER	3.4
40	DJ	59	SER	3.4
37	DG	6	ARG	3.4
37	DG	59	LEU	3.4
6	AG	17	PRO	3.4
37	CG	112	PRO	3.4
40	DJ	64	GLU	3.4
44	DN	20	ALA	3.4
37	DG	9	VAL	3.4
44	DN	38	GLY	3.4
1	AA	2121	G	3.4
3	BD	32	SER	3.4
20	AY	79	CYS	3.4
39	CI	113	LYS	3.4
39	DI	15	ALA	3.4
43	CM	33	ALA	3.4
40	DJ	44	VAL	3.4
49	CS	17	GLU	3.4
8	AI	97	ILE	3.4
12	AQ	22	LYS	3.3
31	CA	1030(C)	G	3.3
31	CA	1233	G	3.3
43	DM	8	GLU	3.3
39	DI	67	GLY	3.3
31	CA	1209	C	3.3
46	DP	19	ILE	3.3
1	AA	2173	A	3.3
1	BA	2169	A	3.3
39	CI	112	LYS	3.3
39	CI	14	VAL	3.3
39	CI	41	VAL	3.3
1	BA	2120	G	3.3
44	CN	34	TYR	3.3
39	DI	47	LEU	3.3
43	CM	19	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
6	BG	75	LYS	3.3
31	DA	1129	C	3.3
40	CJ	30	SER	3.3
28	A6	20	ASN	3.3
38	DH	2	LEU	3.3
40	CJ	71	LEU	3.3
37	DG	118	VAL	3.3
1	AA	2146	C	3.3
1	AA	2175	C	3.3
40	CJ	53	PRO	3.3
5	BF	20	LEU	3.3
39	CI	35	GLU	3.3
6	AG	69	ALA	3.3
49	CS	9	VAL	3.3
49	DS	51	VAL	3.3
1	AA	2116	G	3.3
1	AA	2165	G	3.3
1	BA	2148	G	3.3
31	CA	1276	G	3.3
49	DS	69	HIS	3.3
1	AA	1049	C	3.2
31	CA	1244	C	3.2
31	DA	972	C	3.2
8	AI	78	THR	3.2
31	CA	958	A	3.2
38	DH	1	MET	3.2
8	AI	88	ILE	3.2
40	CJ	59	SER	3.2
43	DM	70	LEU	3.2
38	DH	99	GLU	3.2
43	CM	32	GLU	3.2
11	BP	15	ARG	3.2
31	CA	1137	C	3.2
31	CA	1243	C	3.2
44	CN	22	THR	3.2
49	CS	39	THR	3.2
36	CF	94	GLN	3.2
10	BO	22	ILE	3.2
30	A8	29	LYS	3.2
37	DG	99	LEU	3.2
11	AP	43	GLY	3.2
35	DE	107	ARG	3.2

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Mol	Chain	Res	Type	RSRZ
46	DP	1	MET	3.2
43	DM	18	ALA	3.2
31	DA	72	C	3.2
34	CD	70	ILE	3.2
50	DT	74	LYS	3.2
48	CR	76	LEU	3.2
46	DP	29	ASP	3.2
1	AA	883	G	3.2
4	BE	193	GLY	3.2
31	CA	1331	G	3.2
51	CU	19	GLY	3.2
51	DU	2	GLY	3.2
34	CD	153	ARG	3.2
50	CT	80	ARG	3.2
8	AI	85	GLU	3.2
8	AI	120	ILE	3.2
33	DC	39	ILE	3.2
37	CG	14	PRO	3.2
39	CI	21	PRO	3.2
43	DM	113	PRO	3.2
39	DI	72	GLY	3.2
23	B1	26	ARG	3.2
29	B7	23	ARG	3.2
40	CJ	89	ASP	3.2
23	A1	65	SER	3.2
34	CD	154	ASN	3.2
49	CS	44	MET	3.2
31	CA	949	A	3.2
31	CA	1291	G	3.2
31	DA	347	G	3.2
31	DA	1026	G	3.2
31	DA	1124	G	3.2
40	DJ	47	PHE	3.2
44	DN	58	LYS	3.2
32	DB	221	LEU	3.2
11	AP	15	ARG	3.1
43	DM	24	GLY	3.1
31	CA	1065	U	3.1
39	DI	117	HIS	3.1
43	DM	82	MET	3.1
33	CC	3	ASN	3.1
42	DL	91	LYS	3.1

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Mol	Chain	Res	Type	RSRZ
1	BA	2111	C	3.1
31	DA	1140	C	3.1
32	CB	196	LEU	3.1
39	DI	77	ILE	3.1
40	DJ	40	LEU	3.1
1	BA	2131	G	3.1
31	CA	1220	G	3.1
31	CA	1370	G	3.1
31	DA	1030(A)	G	3.1
45	CO	89	GLY	3.1
15	AT	1	MET	3.1
40	CJ	54	PHE	3.1
7	BH	105	LEU	3.1
50	CT	79	ARG	3.1
43	CM	103	THR	3.1
1	AA	229	A	3.1
31	CA	946	A	3.1
31	CA	1046	A	3.1
31	DA	1130	A	3.1
34	CD	140	VAL	3.1
2	AB	54	G	3.1
6	BG	49	ASP	3.1
31	DA	73	G	3.1
31	DA	1058	G	3.1
31	DA	1139	G	3.1
6	BG	46	ALA	3.1
36	CF	50	TYR	3.1
39	DI	81	ILE	3.1
50	CT	8	ARG	3.1
40	DJ	67	THR	3.1
33	CC	28	GLN	3.1
1	BA	2136	C	3.1
7	AH	24	VAL	3.1
9	BN	46	VAL	3.1
31	CA	1112	C	3.1
31	CA	1362	C	3.1
36	CF	6	VAL	3.1
1	AA	652(B)	A	3.1
1	BA	529	A	3.1
14	AS	56	LEU	3.1
19	BX	92	LEU	3.1
31	DA	1492	A	3.1

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Mol	Chain	Res	Type	RSRZ
44	DN	6	LEU	3.1
49	DS	10	PHE	3.1
37	CG	83	ALA	3.1
39	CI	82	ALA	3.1
32	CB	172	ILE	3.1
44	CN	26	ARG	3.1
44	CN	32	SER	3.1
7	BH	153	LYS	3.1
23	A1	48	LYS	3.1
37	CG	29	LYS	3.1
44	CN	25	VAL	3.1
40	CJ	40	LEU	3.1
44	DN	5	ALA	3.1
43	CM	25	ILE	3.1
46	DP	8	ARG	3.1
49	DS	31	ILE	3.1
39	DI	2	GLU	3.0
31	CA	1035	A	3.0
43	CM	97	PRO	3.0
44	DN	13	THR	3.0
6	AG	82	LEU	3.0
20	BY	79	CYS	3.0
44	CN	39	LEU	3.0
49	CS	15	LEU	3.0
1	AA	1176	G	3.0
1	AA	2154	G	3.0
32	DB	96	ARG	3.0
40	CJ	43	ARG	3.0
40	CJ	46	ARG	3.0
10	AO	33	ALA	3.0
38	CH	4	ASP	3.0
33	CC	4	LYS	3.0
40	DJ	55	LYS	3.0
3	AD	266	SER	3.0
31	DA	1119	C	3.0
31	DA	1288	A	3.0
32	DB	67	THR	3.0
33	CC	66	VAL	3.0
37	DG	80	VAL	3.0
49	CS	19	VAL	3.0
34	CD	73	ARG	3.0
43	CM	34	LEU	3.0

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Mol	Chain	Res	Type	RSRZ
44	DN	36	PHE	3.0
45	DO	64	ARG	3.0
51	DU	7	ARG	3.0
1	AA	2150	U	3.0
31	CA	1364	U	3.0
42	CL	120	TYR	3.0
33	CC	167	TRP	3.0
1	BA	2166	G	3.0
15	BT	1	MET	3.0
43	CM	7	VAL	3.0
49	CS	33	THR	3.0
6	AG	152	LEU	3.0
37	CG	32	ARG	3.0
44	DN	29	ARG	3.0
1	AA	652(R)	C	3.0
1	AA	1043	C	3.0
1	AA	2416	C	3.0
11	BP	70	GLN	3.0
40	CJ	47	PHE	3.0
15	BT	130	ALA	3.0
15	BT	131	ALA	3.0
3	AD	38	LYS	3.0
39	DI	36	TYR	3.0
33	CC	89	GLU	3.0
8	BI	20	ASP	3.0
37	CG	38	LEU	3.0
1	BA	2115	G	3.0
1	BA	2165	G	3.0
31	DA	346	G	3.0
31	DA	1033	G	3.0
43	DM	107	ALA	3.0
44	CN	5	ALA	3.0
1	AA	1052	C	3.0
1	BA	2139	C	3.0
31	CA	330	C	3.0
31	CA	1246	C	3.0
31	DA	1289	A	3.0
11	AP	26	GLY	3.0
39	DI	6	GLY	3.0
51	DU	15	ARG	3.0
34	DD	162	LEU	3.0
43	CM	16	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
21	AZ	166	SER	3.0
32	CB	235	SER	3.0
43	DM	4	ILE	3.0
13	BR	42	LYS	2.9
33	CC	37	GLN	2.9
1	AA	2319	G	2.9
31	CA	1241	G	2.9
39	CI	62	TYR	2.9
49	CS	52	TYR	2.9
33	DC	155	GLY	2.9
1	AA	888	C	2.9
31	DA	1327	C	2.9
32	DB	10	LEU	2.9
32	DB	136	VAL	2.9
37	CG	16	LEU	2.9
51	DU	5	ASP	2.9
37	DG	35	LYS	2.9
40	DJ	26	ALA	2.9
37	CG	86	GLN	2.9
6	AG	142	PRO	2.9
37	DG	21	VAL	2.9
7	AH	159	GLU	2.9
20	AY	34	LYS	2.9
33	DC	202	ILE	2.9
37	DG	50	ILE	2.9
38	CH	116	LYS	2.9
44	CN	11	LYS	2.9
4	BE	204	ALA	2.9
31	DA	1029	C	2.9
31	DA	1117	G	2.9
31	CA	1308	U	2.9
49	DS	35	SER	2.9
40	CJ	76	ASN	2.9
36	DF	90	VAL	2.9
40	DJ	16	LEU	2.9
49	CS	11	VAL	2.9
49	DS	67	VAL	2.9
50	DT	10	LEU	2.9
37	DG	112	PRO	2.9
40	DJ	50	ILE	2.9
37	DG	83	ALA	2.9
40	CJ	26	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
43	DM	42	ALA	2.9
44	CN	30	ALA	2.9
37	CG	78	ARG	2.9
37	CG	94	ARG	2.9
37	CG	114	ARG	2.9
37	DG	32	ARG	2.9
39	CI	66	ARG	2.9
1	AA	2155	G	2.9
1	AA	2171	A	2.9
1	AA	2333	A	2.9
31	CA	1029	C	2.9
31	DA	1116	C	2.9
31	DA	1183	A	2.9
44	CN	44	LEU	2.9
6	AG	160	VAL	2.9
34	CD	178	VAL	2.9
34	DD	4	TYR	2.9
43	CM	41	PRO	2.9
19	AX	34	ALA	2.9
49	CS	48	THR	2.9
25	A3	29	ARG	2.9
10	BO	1	MET	2.9
24	A2	1	MET	2.9
40	DJ	62	HIS	2.9
25	A3	2	PRO	2.9
31	DA	1125	U	2.9
49	DS	76	PRO	2.9
31	CA	1001	A	2.9
43	CM	64	TRP	2.9
1	AA	652(Q)	G	2.8
1	AA	1115	G	2.8
4	AE	204	ALA	2.8
31	DA	3	G	2.8
11	BP	65	ARG	2.8
40	CJ	66	ARG	2.8
40	CJ	92	THR	2.8
43	CM	105	THR	2.8
6	BG	3	LEU	2.8
11	AP	59	LEU	2.8
21	BZ	1	MET	2.8
7	AH	45	VAL	2.8
8	AI	3	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
43	CM	60	VAL	2.8
43	DM	27	LYS	2.8
49	CS	12	ASP	2.8
6	AG	100	TRP	2.8
38	CH	94	TYR	2.8
11	AP	74	GLU	2.8
31	DA	956	U	2.8
1	AA	2164	C	2.8
31	CA	1267	C	2.8
31	CA	1275	A	2.8
31	DA	1118	C	2.8
31	DA	1349	A	2.8
44	CN	40	CYS	2.8
31	CA	380	G	2.8
31	CA	1001(A)	G	2.8
31	DA	1185	G	2.8
32	CB	165	VAL	2.8
34	DD	178	VAL	2.8
43	CM	15	VAL	2.8
1	AA	2109	U	2.8
31	CA	1351	U	2.8
6	AG	7	LEU	2.8
35	DE	53	LEU	2.8
40	CJ	75	LEU	2.8
34	DD	33	MET	2.8
1	BA	2161	C	2.8
31	DA	1191	A	2.8
11	BP	68	GLN	2.8
23	A1	47	GLN	2.8
46	CP	13	HIS	2.8
6	BG	80	PHE	2.8
26	A4	5	ILE	2.8
39	DI	74	ILE	2.8
40	CJ	38	ILE	2.8
21	BZ	166	SER	2.8
1	AA	1042	G	2.8
1	AA	2894	G	2.8
1	BA	2154	G	2.8
8	AI	94	ALA	2.8
31	CA	1036	G	2.8
37	CG	109	ASN	2.8
39	CI	46	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
14	AS	8	GLU	2.8
5	AF	20	LEU	2.8
13	AR	42	LYS	2.8
19	AX	33	LYS	2.8
39	CI	118	LYS	2.8
20	BY	3	VAL	2.8
20	BY	85	VAL	2.8
33	DC	68	VAL	2.8
36	CF	90	VAL	2.8
44	DN	18	VAL	2.8
43	CM	37	THR	2.8
31	CA	1000	U	2.8
50	CT	103	GLY	2.8
11	AP	68	GLN	2.8
31	CA	1333	A	2.8
1	BA	886	C	2.8
51	DU	18	TYR	2.8
22	B0	12	ASN	2.8
39	DI	23	ASN	2.8
6	AG	8	LYS	2.8
6	BG	82	LEU	2.8
41	DK	98	LEU	2.8
1	AA	1044	G	2.8
31	CA	1186	G	2.8
31	CA	1293	G	2.8
31	DA	1064	G	2.8
34	DD	2	GLY	2.8
24	A2	51	ARG	2.7
44	DN	41	ARG	2.7
31	DA	1065	U	2.7
31	DA	1194	U	2.7
33	DC	149	ALA	2.7
7	BH	171	LEU	2.7
8	AI	68	LEU	2.7
31	CA	1111	A	2.7
1	BA	652(S)	C	2.7
1	BA	2164	C	2.7
6	AG	35	GLU	2.7
6	AG	86	MET	2.7
30	A8	25	MET	2.7
31	CA	1260	C	2.7
40	DJ	61	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
44	CN	33	VAL	2.7
30	A8	45	GLY	2.7
34	DD	141	ARG	2.7
37	DG	4	ARG	2.7
44	CN	7	ILE	2.7
49	CS	10	PHE	2.7
23	A1	87	PRO	2.7
1	AA	2120	G	2.7
1	BA	530	G	2.7
1	BA	2149	G	2.7
2	AB	24	G	2.7
31	CA	331	G	2.7
31	CA	1253	G	2.7
39	CI	15	ALA	2.7
49	DS	75	ALA	2.7
1	BA	1026	U	2.7
44	DN	61	TRP	2.7
13	AR	73	VAL	2.7
24	B2	5	GLU	2.7
2	AB	53	A	2.7
31	CA	1318	A	2.7
3	AD	217	ARG	2.7
37	CG	5	ARG	2.7
39	DI	42	ARG	2.7
43	CM	4	ILE	2.7
48	CR	75	ILE	2.7
49	DS	33	THR	2.7
35	DE	21	ALA	2.7
39	DI	76	ALA	2.7
46	CP	27	LYS	2.7
6	AG	11	TYR	2.7
6	AG	12	TYR	2.7
32	CB	197	VAL	2.7
32	DB	165	VAL	2.7
34	CD	112	VAL	2.7
34	DD	140	VAL	2.7
1	BA	2109	U	2.7
31	CA	950	U	2.7
1	BA	2110	G	2.7
6	BG	79	ASN	2.7
31	CA	1117	G	2.7
31	CA	1274	G	2.7

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Mol	Chain	Res	Type	RSRZ
34	CD	3	ARG	2.7
43	CM	110	ARG	2.7
43	DM	12	ASN	2.7
33	DC	2	GLY	2.7
3	BD	275	LYS	2.7
50	DT	14	LYS	2.7
43	CM	28	ALA	2.7
46	CP	82	GLN	2.7
14	AS	17	ARG	2.7
37	DG	10	ARG	2.7
6	BG	35	GLU	2.7
6	AG	157	ILE	2.7
35	DE	76	ILE	2.7
7	AH	120	GLY	2.7
31	DA	1364	U	2.7
20	AY	4	LYS	2.7
37	CG	53	LYS	2.7
6	BG	87	PRO	2.7
10	BO	4	PRO	2.7
1	AA	2151	G	2.7
2	AB	23	G	2.7
19	BX	95	LEU	2.7
31	CA	108	G	2.7
31	CA	630	G	2.7
31	CA	1026	G	2.7
31	CA	1031	G	2.7
31	CA	1222	G	2.7
31	DA	971	G	2.7
31	DA	1002	G	2.7
37	CG	59	LEU	2.7
48	CR	31	LEU	2.7
1	BA	2114	A	2.7
32	CB	7	VAL	2.7
39	CI	17	VAL	2.7
40	CJ	34	VAL	2.7
40	CJ	72	VAL	2.7
1	AA	886	C	2.7
1	BA	2138	C	2.7
31	CA	1114	C	2.7
31	CA	1249	C	2.7
31	CA	1342	C	2.7
18	AW	112	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
40	DJ	63	PHE	2.6
41	DK	36	ASP	2.6
36	DF	7	ASN	2.6
37	CG	84	ASN	2.6
40	CJ	78	ASN	2.6
5	AF	166	ALA	2.6
6	AG	175	LEU	2.6
25	B3	2	PRO	2.6
26	A4	7	PRO	2.6
33	CC	175	LEU	2.6
31	CA	1025	U	2.6
40	DJ	48	THR	2.6
40	DJ	92	THR	2.6
44	DN	49	HIS	2.6
21	BZ	196	VAL	2.6
33	CC	195	VAL	2.6
33	DC	167	TRP	2.6
42	CL	19	ARG	2.6
43	DM	88	ARG	2.6
50	DT	8	ARG	2.6
43	DM	87	TYR	2.6
1	BA	2123	G	2.6
31	CA	347	G	2.6
31	CA	1295	G	2.6
31	DA	1312	G	2.6
31	DA	1371	G	2.6
6	AG	39	ILE	2.6
40	DJ	38	ILE	2.6
43	DM	65	LYS	2.6
10	AO	74	GLY	2.6
1	BA	2142	C	2.6
6	BG	76	SER	2.6
7	BH	2	SER	2.6
31	DA	345	C	2.6
33	DC	33	LEU	2.6
33	CC	169	ALA	2.6
42	CL	26	ALA	2.6
43	CM	93	ARG	2.6
49	CS	37	ARG	2.6
26	B4	35	VAL	2.6
43	DM	54	VAL	2.6
33	CC	33	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
34	DD	31	CYS	2.6
39	DI	79	LEU	2.6
6	BG	142	PRO	2.6
20	AY	53	PRO	2.6
31	DA	1318	A	2.6
31	DA	1363(A)	A	2.6
34	DD	137	SER	2.6
39	DI	82	ALA	2.6
44	CN	2	ALA	2.6
44	DN	32	SER	2.6
31	CA	1185	G	2.6
31	DA	324	G	2.6
31	DA	1013	G	2.6
49	DS	23	ASN	2.6
29	A7	23	ARG	2.6
31	DA	110	C	2.6
34	CD	141	ARG	2.6
43	CM	29	ARG	2.6
49	CS	47	HIS	2.6
37	CG	91	VAL	2.6
44	DN	33	VAL	2.6
24	A2	70	GLN	2.6
37	CG	97	GLN	2.6
12	AQ	98	LYS	2.6
43	CM	13	LYS	2.6
46	DP	27	LYS	2.6
1	AA	2172	U	2.6
31	CA	1330	U	2.6
31	DA	1219	U	2.6
35	CE	118	ILE	2.6
47	CQ	59	ILE	2.6
35	CE	85	GLY	2.6
43	CM	90	LEU	2.6
50	DT	13	LEU	2.6
49	CS	59	PRO	2.6
43	CM	91	ARG	2.6
32	DB	101	MET	2.6
33	CC	86	VAL	2.6
1	AA	2169	A	2.6
31	CA	1332	A	2.6
33	DC	4	LYS	2.6
42	DL	126	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	AA	652(P)	G	2.6
1	AA	652(T)	C	2.6
1	AA	2152	G	2.6
1	BA	2157	G	2.6
7	BH	148	ILE	2.6
28	A6	54	ILE	2.6
31	CA	1300	G	2.6
31	CA	1365	G	2.6
31	DA	1127	G	2.6
34	CD	5	ILE	2.6
37	CG	26	PHE	2.6
32	DB	99	GLY	2.6
43	DM	6	GLY	2.6
43	DM	96	LEU	2.6
49	CS	20	LEU	2.6
1	BA	2132	U	2.6
34	CD	66	ARG	2.6
37	DG	25	ALA	2.6
42	DL	73	GLU	2.6
43	CM	50	GLU	2.6
13	AR	68	ARG	2.6
32	DB	202	PRO	2.6
39	DI	121	ARG	2.6
8	AI	96	ASP	2.6
32	DB	8	LYS	2.6
33	DC	135	LYS	2.6
39	CI	32	ASP	2.6
39	DI	97	LYS	2.6
33	CC	192	THR	2.6
39	CI	64	THR	2.6
43	DM	63	THR	2.6
33	DC	28	GLN	2.5
39	CI	63	ILE	2.5
1	BA	653	A	2.5
2	AB	52	A	2.5
6	AG	19	LEU	2.5
8	AI	77	LEU	2.5
31	CA	1157	A	2.5
31	CA	1285	A	2.5
32	DB	138	LEU	2.5
42	DL	69	TYR	2.5
49	CS	16	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	AA	2105	C	2.5
1	AA	2138	C	2.5
31	CA	163	C	2.5
33	CC	160	ALA	2.5
33	DC	168	ALA	2.5
37	CG	4	ARG	2.5
40	DJ	43	ARG	2.5
40	DJ	66	ARG	2.5
43	CM	72	ALA	2.5
1	AA	2104	G	2.5
1	BA	272(A)	G	2.5
1	BA	2151	G	2.5
31	DA	570	G	2.5
31	DA	1190	G	2.5
31	DA	1193	G	2.5
33	CC	68	VAL	2.5
41	DK	14	VAL	2.5
43	DM	17	VAL	2.5
37	DG	28	ASN	2.5
49	CS	53	ASN	2.5
49	CS	62	ILE	2.5
50	CT	9	ASN	2.5
50	CT	18	GLN	2.5
8	AI	101	LEU	2.5
32	DB	70	PHE	2.5
34	CD	162	LEU	2.5
44	CN	53	LEU	2.5
6	BG	146	TYR	2.5
39	CI	114	TYR	2.5
11	AP	16	ARG	2.5
20	BY	2	ARG	2.5
1	BA	2170	A	2.5
31	DA	1030(D)	A	2.5
34	DD	18	LYS	2.5
40	CJ	61	GLU	2.5
9	BN	53	VAL	2.5
37	CG	17	VAL	2.5
37	CG	135	VAL	2.5
47	DQ	11	VAL	2.5
1	BA	1049	C	2.5
3	AD	220	HIS	2.5
39	CI	81	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
43	CM	92	HIS	2.5
44	DN	42	ILE	2.5
1	AA	558	G	2.5
39	DI	32	ASP	2.5
6	AG	79	ASN	2.5
6	AG	172	LEU	2.5
25	A3	26	LEU	2.5
37	DG	97	GLN	2.5
42	CL	32	PHE	2.5
48	DR	76	LEU	2.5
14	BS	96	GLY	2.5
36	DF	4	TYR	2.5
37	DG	154	TYR	2.5
7	AH	85	LYS	2.5
39	CI	45	ALA	2.5
21	BZ	28	MET	2.5
8	AI	144	VAL	2.5
34	CD	39	PRO	2.5
37	CG	105	VAL	2.5
38	DH	5	PRO	2.5
1	AA	1048	A	2.5
31	DA	1447	A	2.5
14	AS	21	THR	2.5
37	CG	124	LEU	2.5
39	DI	33	PHE	2.5
1	AA	2161	C	2.5
1	BA	2107	C	2.5
1	BA	2150	U	2.5
14	BS	42	ASP	2.5
31	CA	1320	C	2.5
39	CI	10	ARG	2.5
31	CA	1240	U	2.5
20	AY	35	TYR	2.5
51	DU	4	GLY	2.5
20	AY	65	ALA	2.5
33	DC	160	ALA	2.5
35	DE	86	ALA	2.5
51	DU	12	LYS	2.5
1	AA	652(C)	G	2.5
1	AA	2106	G	2.5
29	A7	1	MET	2.5
31	CA	1034	G	2.5

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Mol	Chain	Res	Type	RSRZ
31	DA	1031	G	2.5
31	DA	1184	G	2.5
36	DF	6	VAL	2.5
39	CI	28	VAL	2.5
43	CM	54	VAL	2.5
43	DM	41	PRO	2.5
25	A3	60	GLU	2.5
6	AG	101	ILE	2.5
38	DH	112	LEU	2.5
47	CQ	6	LEU	2.5
37	DG	79	ARG	2.5
39	DI	27	THR	2.5
40	DJ	15	THR	2.5
44	DN	22	THR	2.5
31	CA	1130	A	2.5
31	CA	1493	A	2.5
31	DA	1250	A	2.5
18	BW	112	GLY	2.5
40	CJ	55	LYS	2.5
40	DJ	19	SER	2.5
49	CS	38	SER	2.5
43	DM	16	ASP	2.5
46	CP	17	TYR	2.5
1	AA	2142	C	2.5
1	AA	2296	U	2.5
1	BA	2129	C	2.5
31	CA	984	C	2.5
31	CA	1242	C	2.5
31	DA	1109	C	2.5
39	CI	86	VAL	2.5
34	DD	29	PRO	2.5
33	DC	5	ILE	2.4
43	DM	25	ILE	2.4
45	DO	3	ILE	2.4
1	AA	2655	G	2.4
31	CA	993	G	2.4
31	DA	111	G	2.4
31	DA	662	G	2.4
31	DA	1131	G	2.4
44	DN	26	ARG	2.4
46	DP	22	THR	2.4
49	DS	39	THR	2.4

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Mol	Chain	Res	Type	RSRZ
40	CJ	52	GLY	2.4
43	CM	24	GLY	2.4
49	DS	8	GLY	2.4
6	AG	131	TYR	2.4
39	DI	114	TYR	2.4
33	DC	66	VAL	2.4
31	CA	1257	U	2.4
1	AA	2791	C	2.4
6	AG	18	GLU	2.4
6	BG	144	ILE	2.4
6	BG	133	LEU	2.4
31	CA	1027	C	2.4
31	CA	1147	C	2.4
31	DA	1203	C	2.4
37	DG	146	GLU	2.4
38	DH	22	GLU	2.4
50	CT	55	ILE	2.4
32	CB	70	PHE	2.4
19	AX	31	HIS	2.4
6	BG	73	ALA	2.4
39	DI	61	ALA	2.4
40	DJ	84	GLN	2.4
42	DL	26	ALA	2.4
44	CN	10	ALA	2.4
1	AA	2149	G	2.4
1	AA	2156	G	2.4
7	AH	43	VAL	2.4
8	AI	92	VAL	2.4
31	CA	988	G	2.4
31	DA	1373	G	2.4
34	CD	54	TYR	2.4
39	DI	26	VAL	2.4
6	AG	20	ILE	2.4
6	AG	83	ARG	2.4
7	BH	170	ARG	2.4
37	DG	120	ILE	2.4
46	DP	26	ARG	2.4
47	DQ	70	ARG	2.4
50	CT	17	ARG	2.4
50	DT	84	LEU	2.4
49	CS	64	GLU	2.4
31	DA	1281	U	2.4

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Mol	Chain	Res	Type	RSRZ
46	DP	3	LYS	2.4
50	CT	73	HIS	2.4
1	BA	889	C	2.4
31	CA	980	C	2.4
14	AS	5	THR	2.4
37	DG	7	ALA	2.4
49	CS	34	TRP	2.4
7	BH	133	VAL	2.4
40	DJ	34	VAL	2.4
49	CS	61	TYR	2.4
20	BY	70	SER	2.4
39	DI	21	PRO	2.4
50	CT	11	SER	2.4
34	DD	122	ARG	2.4
39	CI	54	ASP	2.4
39	DI	93	ARG	2.4
44	CN	47	LEU	2.4
49	DS	71	LEU	2.4
9	BN	118	LYS	2.4
1	AA	2802	G	2.4
1	BA	2894	G	2.4
31	CA	587	G	2.4
31	CA	1048	G	2.4
31	CA	1181	G	2.4
31	DA	1224	G	2.4
31	DA	1309	G	2.4
14	AS	34	HIS	2.4
31	DA	974	A	2.4
31	CA	956	U	2.4
31	CA	1150	U	2.4
31	DA	1126	U	2.4
47	DQ	33	GLY	2.4
9	AN	43	THR	2.4
37	CG	128	ALA	2.4
33	CC	153	VAL	2.4
49	DS	60	VAL	2.4
16	BU	117	GLN	2.4
1	AA	898	C	2.4
31	CA	1018	C	2.4
39	CI	92	TYR	2.4
40	DJ	46	ARG	2.4
46	DP	39	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
6	AG	3	LEU	2.4
6	AG	139	LEU	2.4
23	B1	2	SER	2.4
33	CC	144	SER	2.4
33	DC	34	LEU	2.4
43	CM	84	ILE	2.4
14	AS	42	ASP	2.4
32	DB	156	LYS	2.4
34	DD	77	ASN	2.4
8	BI	10	GLU	2.4
15	AT	128	GLU	2.4
23	B1	83	GLU	2.4
49	DS	17	GLU	2.4
24	B2	1	MET	2.4
35	DE	104	ALA	2.4
1	AA	2141	G	2.4
8	AI	86	THR	2.4
31	DA	70	G	2.4
31	DA	1057	G	2.4
31	DA	1175	G	2.4
43	DM	64	TRP	2.4
43	DM	109	THR	2.4
1	BA	2790	A	2.4
31	CA	1204	A	2.4
31	DA	1285	A	2.4
31	DA	1377	A	2.4
34	CD	209	ARG	2.4
37	CG	106	GLN	2.4
44	CN	12	ARG	2.4
40	CJ	80	LYS	2.3
47	DQ	100	LYS	2.3
14	AS	31	SER	2.3
21	BZ	88	PHE	2.3
31	CA	1140	C	2.3
31	DA	1326	C	2.3
46	DP	9	PHE	2.3
37	CG	20	ASP	2.3
7	BH	150	ALA	2.3
17	AV	67	GLY	2.3
20	AY	42	VAL	2.3
29	A7	45	ALA	2.3
44	DN	51	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
49	CS	14	HIS	2.3
32	DB	112	VAL	2.3
50	DT	40	ALA	2.3
6	AG	29	TRP	2.3
11	BP	62	LEU	2.3
42	CL	27	LEU	2.3
32	CB	127	ILE	2.3
43	CM	22	ILE	2.3
44	CN	9	LYS	2.3
4	BE	151	TYR	2.3
32	DB	159	PRO	2.3
1	AA	1175	U	2.3
2	AB	59	A	2.3
14	AS	87	PHE	2.3
1	AA	2127	G	2.3
2	AB	51	G	2.3
31	DA	77	G	2.3
34	DD	144	ASP	2.3
21	BZ	168	GLU	2.3
34	DD	98	GLU	2.3
12	BQ	33	GLY	2.3
21	AZ	74	VAL	2.3
31	CA	106	C	2.3
31	CA	1028	C	2.3
31	CA	1115	C	2.3
31	CA	1234	C	2.3
31	DA	1066	C	2.3
31	DA	1260	C	2.3
37	CG	108	ALA	2.3
39	DI	46	ALA	2.3
42	CL	95	GLY	2.3
39	CI	120	ARG	2.3
45	DO	50	HIS	2.3
46	DP	5	ARG	2.3
48	CR	32	ARG	2.3
8	AI	108	THR	2.3
40	CJ	84	GLN	2.3
39	CI	36	TYR	2.3
49	DS	74	PHE	2.3
7	AH	134	SER	2.3
1	AA	2801(A)	A	2.3
2	BB	53	A	2.3

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Mol	Chain	Res	Type	RSRZ
6	AG	48	GLU	2.3
6	AG	54	GLU	2.3
8	AI	122	GLU	2.3
16	AU	97	ASP	2.3
31	CA	1044	A	2.3
31	CA	1179	A	2.3
31	DA	975	A	2.3
33	DC	138	VAL	2.3
33	DC	156	ARG	2.3
34	CD	156	GLU	2.3
37	DG	115	ARG	2.3
37	DG	150	ALA	2.3
43	DM	74	VAL	2.3
51	CU	22	ARG	2.3
1	BA	2121	G	2.3
14	AS	19	LYS	2.3
31	CA	1139	G	2.3
31	DA	1386	G	2.3
31	DA	1401	G	2.3
33	DC	31	HIS	2.3
47	CQ	87	LYS	2.3
49	DS	7	LYS	2.3
32	DB	44	LEU	2.3
32	DB	98	LEU	2.3
3	BD	271	ILE	2.3
39	CI	74	ILE	2.3
1	BA	2146	C	2.3
2	AB	88	C	2.3
31	CA	435	C	2.3
31	CA	1327	C	2.3
31	DA	1192	C	2.3
30	A8	35	GLN	2.3
49	CS	80	TYR	2.3
50	DT	18	GLN	2.3
6	AG	171	ALA	2.3
13	BR	7	GLY	2.3
33	CC	207	VAL	2.3
41	CK	49	GLY	2.3
42	DL	29	GLY	2.3
43	CM	51	ALA	2.3
4	AE	121	ASN	2.3
43	DM	106	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	BA	2122	U	2.3
1	AA	2117	A	2.3
2	AB	25	A	2.3
6	BG	88	ILE	2.3
31	CA	134	A	2.3
31	CA	1261	A	2.3
31	CA	1360	A	2.3
31	DA	1374	A	2.3
6	AG	145	THR	2.3
7	AH	36	PRO	2.3
39	DI	64	THR	2.3
39	CI	101	PHE	2.3
31	CA	1294	G	2.3
31	DA	1138	G	2.3
1	AA	279	C	2.3
1	AA	2177	C	2.3
1	BA	2477	C	2.3
31	CA	1359	C	2.3
31	DA	824	C	2.3
3	AD	268	ARG	2.3
40	CJ	29	ARG	2.3
45	CO	65	ARG	2.3
33	DC	120	VAL	2.3
34	DD	166	LYS	2.3
37	CG	65	ALA	2.3
37	DG	17	VAL	2.3
42	DL	21	LYS	2.3
42	DL	23	LYS	2.3
49	DS	50	ALA	2.3
44	DN	39	LEU	2.3
34	DD	5	ILE	2.2
40	CJ	58	ASP	2.2
31	CA	1235	U	2.2
31	DA	981	U	2.2
31	DA	1240	U	2.2
7	AH	112	PRO	2.2
39	CI	98	PRO	2.2
40	CJ	41	PRO	2.2
43	CM	113	PRO	2.2
45	DO	4	THR	2.2
49	DS	48	THR	2.2
31	DA	1319	A	2.2

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Mol	Chain	Res	Type	RSRZ
33	CC	29	TYR	2.2
14	BS	30	ARG	2.2
41	DK	120	ARG	2.2
44	CN	3	ARG	2.2
49	CS	7	LYS	2.2
33	DC	116	VAL	2.2
37	CG	80	VAL	2.2
39	DI	65	VAL	2.2
7	BH	7	LEU	2.2
14	AS	58	LEU	2.2
19	AX	95	LEU	2.2
33	CC	47	LEU	2.2
34	CD	11	LEU	2.2
34	CD	124	GLY	2.2
37	CG	104	LEU	2.2
39	DI	43	ALA	2.2
1	AA	1888	G	2.2
1	AA	2145	C	2.2
1	BA	546	C	2.2
1	BA	2162	G	2.2
2	BB	24	G	2.2
31	CA	631	G	2.2
31	CA	1013	G	2.2
31	CA	1033	G	2.2
31	CA	1037	C	2.2
15	BT	128	GLU	2.2
31	CA	1047	G	2.2
31	DA	1253	G	2.2
37	CG	27	ILE	2.2
40	DJ	68	HIS	2.2
49	CS	23	ASN	2.2
43	CM	109	THR	2.2
7	BH	175	LYS	2.2
9	AN	89	LYS	2.2
14	AS	36	TYR	2.2
37	DG	86	GLN	2.2
32	DB	132	LYS	2.2
37	CG	125	MET	2.2
33	DC	151	VAL	2.2
33	DC	153	VAL	2.2
38	CH	118	VAL	2.2
42	CL	18	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
46	DP	2	VAL	2.2
1	AA	2790	A	2.2
1	BA	2126	A	2.2
15	AT	130	ALA	2.2
34	DD	11	LEU	2.2
39	DI	96	LEU	2.2
50	CT	77	ALA	2.2
26	A4	3	GLU	2.2
32	DB	35	GLU	2.2
33	DC	161	GLU	2.2
34	CD	99	SER	2.2
1	AA	34	C	2.2
7	BH	147	ASN	2.2
46	CP	47	ASP	2.2
49	DS	42	PRO	2.2
38	DH	30	ARG	2.2
45	DO	68	ARG	2.2
50	CT	25	ARG	2.2
1	BA	892	G	2.2
1	BA	2319	G	2.2
14	AS	57	LYS	2.2
31	CA	987	G	2.2
31	DA	1001(A)	G	2.2
6	BG	131	TYR	2.2
49	DS	52	TYR	2.2
33	DC	207	VAL	2.2
8	AI	123	LEU	2.2
8	AI	128	LEU	2.2
14	AS	54	LEU	2.2
40	DJ	8	LEU	2.2
31	DA	793	U	2.2
31	DA	1012	U	2.2
31	DA	1372	U	2.2
35	DE	85	GLY	2.2
49	CS	68	GLY	2.2
8	AI	109	ILE	2.2
32	DB	201	ILE	2.2
33	DC	182	ILE	2.2
49	DS	49	ILE	2.2
34	CD	125	HIS	2.2
40	CJ	62	HIS	2.2
43	DM	92	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
34	DD	153	ARG	2.2
39	CI	104	ARG	2.2
40	CJ	45	ARG	2.2
43	DM	79	LYS	2.2
50	DT	9	ASN	2.2
13	BR	49	ASP	2.2
24	B2	52	ASP	2.2
34	CD	190	ASP	2.2
12	AQ	21	THR	2.2
33	CC	95	THR	2.2
43	CM	20	THR	2.2
1	AA	280	C	2.2
1	AA	884	C	2.2
1	BA	2137	C	2.2
6	AG	90	LEU	2.2
12	BQ	13	GLN	2.2
14	BS	36	TYR	2.2
31	DA	1006	C	2.2
33	CC	34	LEU	2.2
37	DG	22	LEU	2.2
49	DS	19	VAL	2.2
8	AI	83	ALA	2.2
28	A6	40	CYS	2.2
33	DC	137	ALA	2.2
43	CM	86	CYS	2.2
6	AG	44	GLY	2.2
1	AA	2123	G	2.2
1	AA	2153	G	2.2
31	CA	348	G	2.2
31	CA	1131	G	2.2
31	CA	1190	G	2.2
31	CA	1305	G	2.2
31	DA	1182	G	2.2
33	DC	124	ILE	2.2
37	DG	30	ILE	2.2
1	AA	2585	U	2.2
6	BG	143	GLU	2.2
11	BP	149	GLU	2.2
33	CC	31	HIS	2.2
40	DJ	56	HIS	2.2
41	CK	126	ARG	2.2
34	CD	22	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
46	DP	12	LYS	2.2
46	DP	31	LYS	2.2
21	AZ	167	PRO	2.2
39	DI	49	PRO	2.2
1	BA	6	A	2.2
31	CA	1245	A	2.2
6	BG	139	LEU	2.2
26	A4	35	VAL	2.2
26	B4	33	VAL	2.2
43	DM	53	VAL	2.2
49	DS	77	THR	2.2
6	AG	146	TYR	2.2
9	BN	72	TYR	2.2
21	BZ	99	TYR	2.2
41	CK	13	GLN	2.2
48	CR	34	TYR	2.2
39	CI	76	ALA	2.1
11	BP	43	GLY	2.1
34	DD	87	GLY	2.1
46	CP	10	GLY	2.1
12	AQ	127	ILE	2.1
28	A6	43	CYS	2.1
36	DF	8	ILE	2.1
49	CS	31	ILE	2.1
31	CA	1214	C	2.1
31	DA	1038	C	2.1
31	DA	1320	C	2.1
6	AG	84	LYS	2.1
18	AW	111	HIS	2.1
21	AZ	193	GLU	2.1
30	A8	54	GLU	2.1
34	CD	145	GLU	2.1
35	CE	81	GLU	2.1
1	AA	362	U	2.1
1	BA	2506	U	2.1
31	DA	1025	U	2.1
1	AA	652(F)	G	2.1
31	CA	945	G	2.1
40	DJ	53	PRO	2.1
6	BG	135	LEU	2.1
11	BP	99	LEU	2.1
38	CH	2	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
39	DI	19	LEU	2.1
46	CP	22	THR	2.1
6	AG	50	ALA	2.1
25	A3	25	ALA	2.1
32	DB	188	ALA	2.1
37	CG	2	ALA	2.1
39	DI	45	ALA	2.1
43	CM	21	TYR	2.1
49	DS	12	ASP	2.1
31	CA	1329	A	2.1
43	DM	78	ILE	2.1
23	A1	26	ARG	2.1
24	B2	51	ARG	2.1
49	CS	36	ARG	2.1
46	CP	12	LYS	2.1
1	BA	2105	C	2.1
11	BP	122	PRO	2.1
31	DA	526	C	2.1
42	DL	25	PRO	2.1
45	CO	56	LEU	2.1
35	DE	51	VAL	2.1
5	BF	175	THR	2.1
7	BH	156	ALA	2.1
21	AZ	152	ALA	2.1
33	DC	92	ALA	2.1
33	DC	184	TYR	2.1
34	CD	68	TYR	2.1
1	AA	272(A)	G	2.1
7	AH	22	GLY	2.1
35	DE	109	ILE	2.1
39	CI	77	ILE	2.1
46	CP	19	ILE	2.1
7	BH	6	ARG	2.1
22	A0	55	ARG	2.1
31	CA	324	G	2.1
35	DE	18	ARG	2.1
40	DJ	7	LYS	2.1
40	DJ	70	ARG	2.1
50	DT	25	ARG	2.1
1	AA	2158	A	2.1
1	BA	2134	A	2.1
31	CA	1180	A	2.1

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Mol	Chain	Res	Type	RSRZ
39	DI	18	PHE	2.1
41	DK	119	CYS	2.1
44	CN	27	CYS	2.1
32	DB	140	HIS	2.1
40	DJ	13	HIS	2.1
4	AE	49	LEU	2.1
8	AI	116	LEU	2.1
19	AX	1	MET	2.1
8	BI	19	VAL	2.1
33	CC	130	VAL	2.1
1	AA	2143	C	2.1
1	AA	2163	C	2.1
1	BA	2145	C	2.1
31	CA	1039	C	2.1
31	CA	1210	C	2.1
39	DI	84	ALA	2.1
6	BG	145	THR	2.1
33	CC	193	TYR	2.1
44	DN	7	ILE	2.1
1	AA	90	U	2.1
6	AG	42	GLY	2.1
34	CD	23	GLY	2.1
39	DI	104	ARG	2.1
43	CM	95	GLY	2.1
51	DU	16	GLY	2.1
39	CI	75	ASP	2.1
1	BA	226	G	2.1
1	BA	652(Q)	G	2.1
1	BA	2141	G	2.1
2	AB	118	G	2.1
31	CA	1215	G	2.1
31	DA	973	G	2.1
31	DA	1094	G	2.1
34	DD	135	LEU	2.1
33	DC	70	VAL	2.1
33	DC	130	VAL	2.1
34	CD	98	GLU	2.1
45	DO	2	PRO	2.1
20	AY	49	VAL	2.1
31	DA	1236	A	2.1
37	CG	61	VAL	2.1
39	CI	53	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
39	DI	86	VAL	2.1
49	DS	64	GLU	2.1
14	AS	37	ALA	2.1
33	CC	189	ALA	2.1
50	CT	67	ALA	2.1
4	AE	79	ARG	2.1
12	AQ	18	LYS	2.1
20	BY	4	LYS	2.1
34	DD	50	ARG	2.1
34	DD	158	ILE	2.1
47	CQ	63	ARG	2.1
37	DG	24	THR	2.1
42	CL	14	GLY	2.1
49	CS	82	GLY	2.1
44	CN	52	GLN	2.1
1	BA	885	C	2.1
31	CA	624	C	2.1
1	AA	2122	U	2.1
1	AA	2167	U	2.1
2	AB	1	U	2.1
39	CI	37	PHE	2.1
39	DI	54	ASP	2.1
8	AI	80	PRO	2.1
21	AZ	56	VAL	2.1
32	DB	197	VAL	2.1
33	DC	109	PRO	2.1
40	DJ	24	VAL	2.1
42	DL	18	VAL	2.1
6	AG	137	GLU	2.1
36	DF	41	GLU	2.1
39	CI	2	GLU	2.1
44	DN	27	CYS	2.1
45	DO	6	GLU	2.1
3	BD	176	ARG	2.0
43	CM	94	ARG	2.0
43	DM	80	ARG	2.0
14	AS	35	ILE	2.0
44	DN	17	LYS	2.0
44	DN	35	ARG	2.0
34	DD	67	ILE	2.0
38	DH	111	ILE	2.0
41	CK	21	ILE	2.0

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Mol	Chain	Res	Type	RSRZ
1	BA	652(B)	A	2.0
1	BA	890	A	2.0
1	BA	2062	A	2.0
17	BV	63	GLY	2.0
22	A0	76	GLY	2.0
31	DA	1001	A	2.0
1	AA	1107	G	2.0
1	AA	2159	G	2.0
1	AA	2166	G	2.0
1	BA	2207	G	2.0
31	CA	1017	G	2.0
31	CA	1266	G	2.0
31	DA	331	G	2.0
8	AI	89	TYR	2.0
43	DM	20	THR	2.0
49	CS	63	THR	2.0
34	CD	175	SER	2.0
39	DI	73	GLN	2.0
37	DG	43	PHE	2.0
15	BT	104	ASN	2.0
30	B8	33	ASN	2.0
3	AD	147	LEU	2.0
6	AG	34	LEU	2.0
6	AG	135	LEU	2.0
33	DC	36	ASP	2.0
34	CD	135	LEU	2.0
40	DJ	75	LEU	2.0
6	AG	70	VAL	2.0
8	AI	145	VAL	2.0
21	AZ	161	VAL	2.0
31	CA	972	C	2.0
31	CA	1141	C	2.0
31	CA	1270	C	2.0
31	DA	1217	C	2.0
5	BF	34	TRP	2.0
24	B2	69	ARG	2.0
34	DD	22	LYS	2.0
34	DD	66	ARG	2.0
39	DI	120	ARG	2.0
7	AH	146	ALA	2.0
7	AH	165	ALA	2.0
14	BS	40	ILE	2.0

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Mol	Chain	Res	Type	RSRZ
45	CO	82	ILE	2.0
39	DI	30	GLY	2.0
46	CP	39	TYR	2.0
33	DC	123	GLN	2.0
31	CA	977	A	2.0
31	DA	968	A	2.0
31	DA	1044	A	2.0
32	CB	98	LEU	2.0
32	DB	121	LEU	2.0
39	CI	79	LEU	2.0
35	DE	10	MET	2.0
39	CI	29	ASN	2.0
39	CI	89	ASN	2.0
1	AA	11	G	2.0
1	AA	1051	G	2.0
1	AA	1559	G	2.0
7	AH	169	VAL	2.0
24	A2	19	VAL	2.0
31	CA	64	G	2.0
31	CA	1385	G	2.0
39	CI	91	ASP	2.0
49	CS	45	VAL	2.0
26	A4	41	PRO	2.0
30	B8	53	PRO	2.0
39	DI	90	PRO	2.0
6	AG	81	LYS	2.0
11	AP	39	LYS	2.0
34	DD	139	ARG	2.0
40	DJ	29	ARG	2.0
47	DQ	91	ARG	2.0
50	DT	16	HIS	2.0
51	CU	7	ARG	2.0
6	BG	48	GLU	2.0
1	BA	958	U	2.0
37	CG	134	ALA	2.0
40	DJ	27	ALA	2.0
48	CR	20	ALA	2.0
1	AA	1118	C	2.0
1	AA	2140	C	2.0
1	BA	34	C	2.0
1	BA	2128	C	2.0
2	BB	5	C	2.0

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Mol	Chain	Res	Type	RSRZ
26	B4	18	CYS	2.0
31	CA	488	C	2.0
31	CA	1118	C	2.0
31	CA	1321	C	2.0
31	DA	1400	C	2.0
40	DJ	10	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

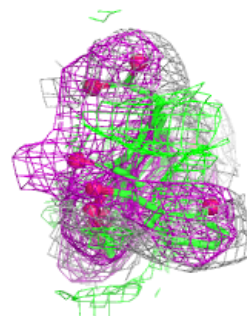
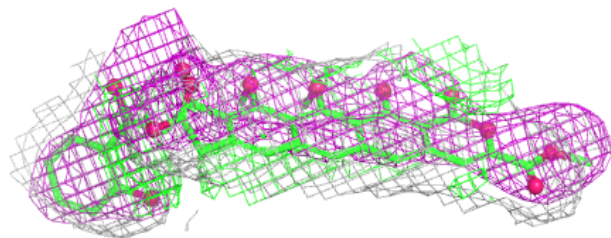
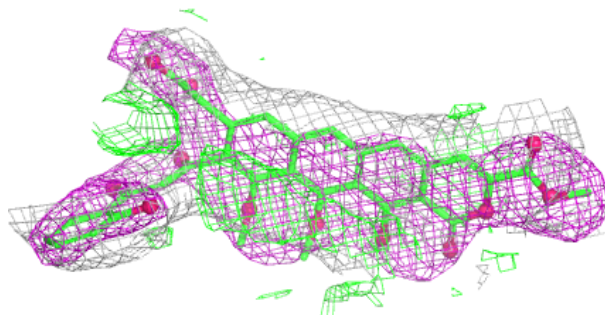
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
52	T8B	AA	3001	44/44	0.74	0.15	20,20,20,20	0
53	MG	BA	3002	1/1	0.77	0.13	30,30,30,30	0
52	T8B	BA	3001	44/44	0.79	0.13	20,20,20,20	0
53	MG	AA	3002	1/1	0.89	0.16	30,30,30,30	0
53	MG	AA	3003	1/1	0.97	0.15	30,30,30,30	0
53	MG	BA	3003	1/1	0.97	0.20	30,30,30,30	0

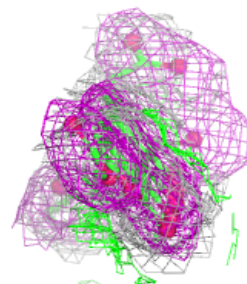
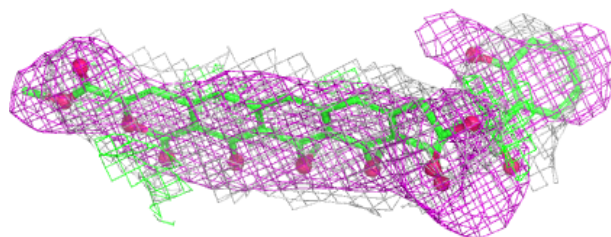
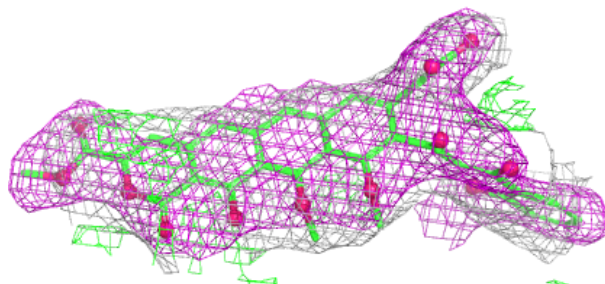
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around T8B AA 3001:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around T8B BA 3001:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.