



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 11, 2025 – 04:13 PM EDT

PDB ID : 9PDH / pdb_00009pdh
Title : X-ray structure of WT Drosophila Ahcy bound to NAD and adenosine at 2.50 Å
Authors : Singhal, K.; Mesecar, A.D.
Deposited on : 2025-06-30
Resolution : 2.50 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 : 4-5-2 with Phenix2.0rc1
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.45.1

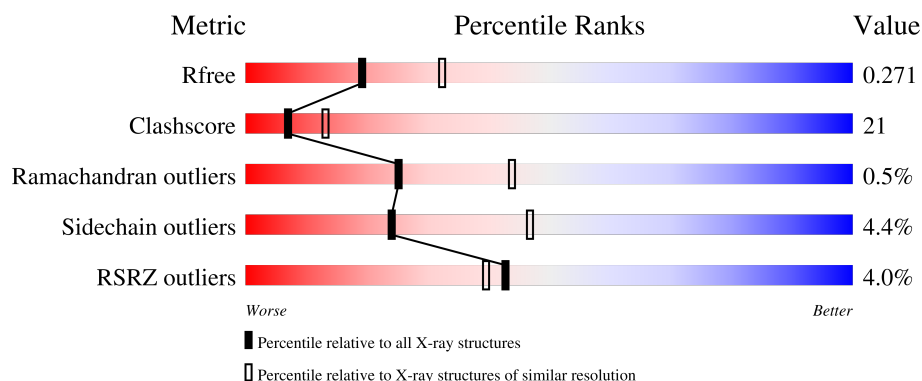
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5504 (2.50-2.50)
Clashscore	180529	6282 (2.50-2.50)
Ramachandran outliers	177936	6191 (2.50-2.50)
Sidechain outliers	177891	6193 (2.50-2.50)
RSRZ outliers	164620	5504 (2.50-2.50)

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	A	434	
1	B	434	
1	C	434	
1	D	434	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14506 atoms, of which 143 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.

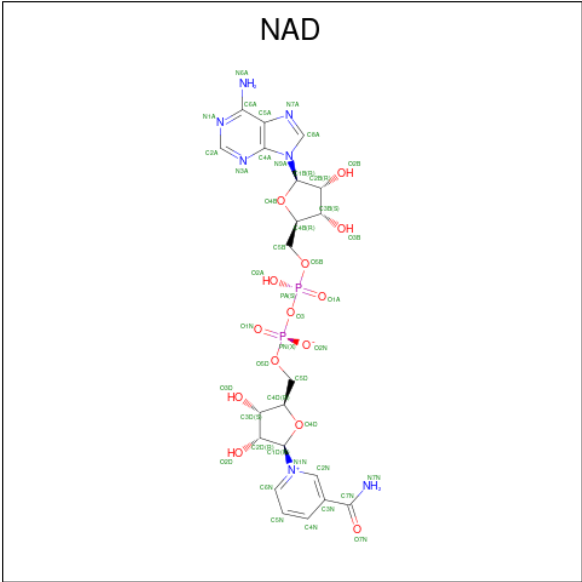
- Molecule 1 is a protein called Adenosylhomocysteinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	431	Total	C	N	O	S	0	1	0
			3320	2102	560	635	23			
1	B	432	Total	C	N	O	S	0	0	0
			3320	2103	560	633	24			
1	C	433	Total	C	N	O	S	0	0	0
			3330	2109	563	634	24			
1	D	430	Total	C	N	O	S	0	1	0
			3316	2101	562	630	23			

There are 8 discrepancies between the modelled and reference sequences:

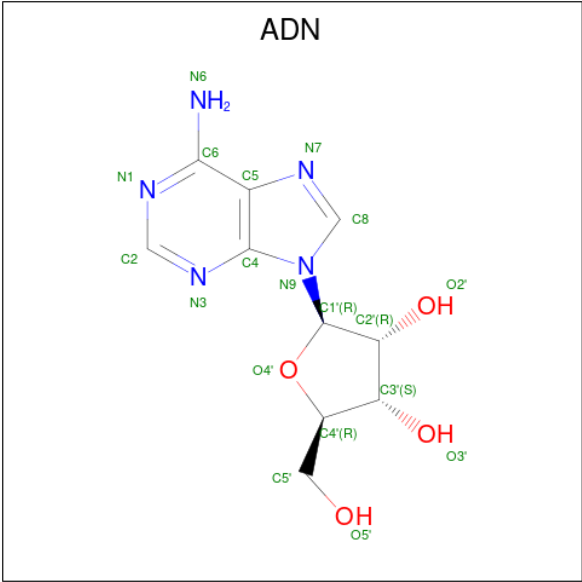
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q27580
A	0	HIS	-	expression tag	UNP Q27580
B	-1	GLY	-	expression tag	UNP Q27580
B	0	HIS	-	expression tag	UNP Q27580
C	-1	GLY	-	expression tag	UNP Q27580
C	0	HIS	-	expression tag	UNP Q27580
D	-1	GLY	-	expression tag	UNP Q27580
D	0	HIS	-	expression tag	UNP Q27580

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	P	0	0
			70	21	26	7	14	2		
2	B	1	Total	C	H	N	O	P	0	0
			70	21	26	7	14	2		
2	C	1	Total	C	H	N	O	P	0	0
			70	21	26	7	14	2		
2	D	1	Total	C	H	N	O	P	0	0
			70	21	26	7	14	2		

- Molecule 3 is ADENOSINE (CCD ID: ADN) (formula: C₁₀H₁₃N₅O₄).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	0	0
			32	10	13	5	4		
3	B	1	Total	C	H	N	O	0	0
			32	10	13	5	4		
3	C	1	Total	C	H	N	O	0	0
			32	10	13	5	4		

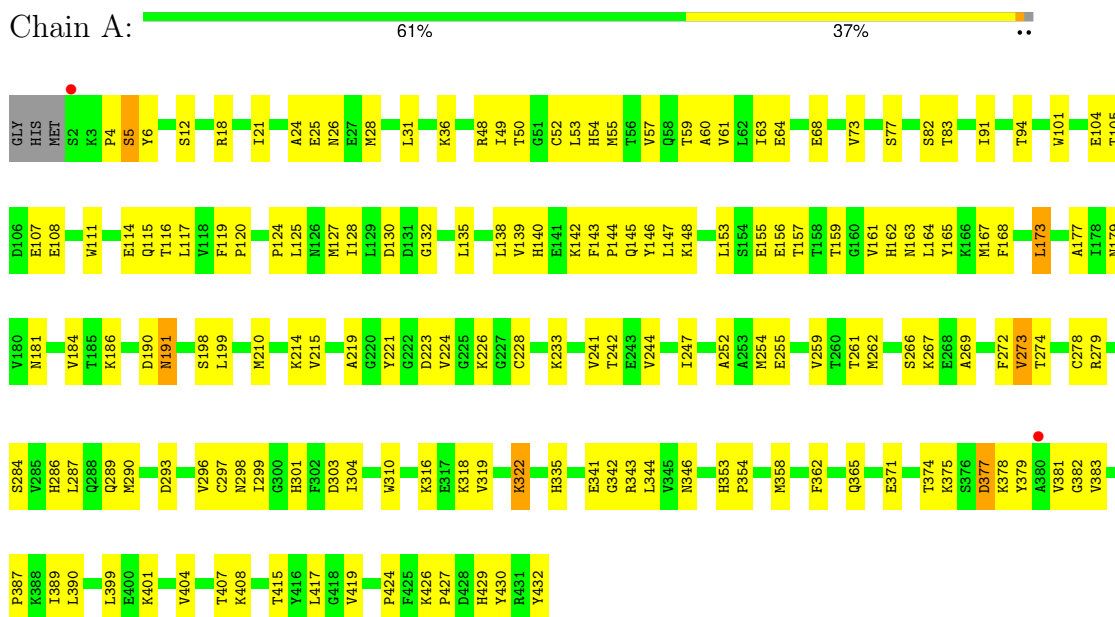
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	228	Total	O	0	0
			228	228		
4	B	263	Total	O	0	0
			263	263		
4	C	198	Total	O	0	0
			198	198		
4	D	155	Total	O	0	0
			155	155		

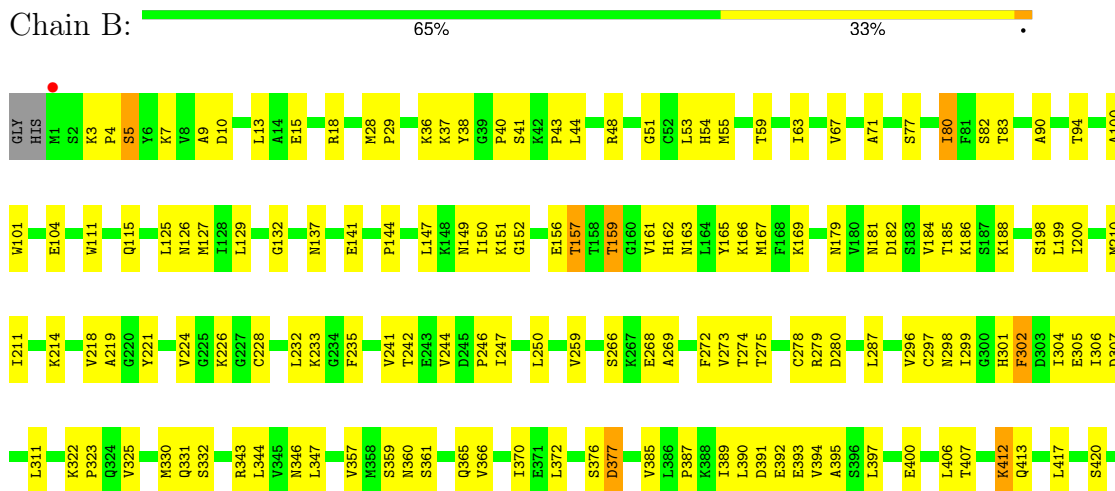
3 Residue-property plots

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: Adenosylhomocysteinase



• Molecule 1: Adenosylhomocysteinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	89.31Å 169.41Å 127.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.45 – 2.50 20.45 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.2 (20.45-2.50) 99.0 (20.45-2.50)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.50Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.230 , 0.270 0.230 , 0.271	Depositor DCC
R_{free} test set	1998 reflections (2.96%)	wwPDB-VP
Wilson B-factor (Å ²)	24.6	Xtriage
Anisotropy	0.193	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 62.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	14506	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	A	0.23	0/3383	0.40	0/4583
1	B	0.21	0/3383	0.40	0/4582
1	C	0.21	0/3394	0.40	0/4597
1	D	0.23	0/3379	0.43	0/4577
All	All	0.22	0/13539	0.41	0/18339

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	A	3320	0	3323	129	0
1	B	3320	0	3332	134	0
1	C	3330	0	3339	138	0
1	D	3316	0	3324	201	0
2	A	44	26	26	1	0
2	B	44	26	26	3	0
2	C	44	26	26	2	0
2	D	44	26	26	3	0
3	A	19	13	13	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	19	13	13	1	0
3	C	19	13	13	1	0
4	A	228	0	0	17	0
4	B	263	0	0	19	0
4	C	198	0	0	16	0
4	D	155	0	0	11	0
All	All	14363	143	13461	558	0

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 21.

All (558) 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)
1:B:259:VAL:HB	1:D:404:VAL:HG13	1.41	1.01
1:D:172:ARG:HD3	4:D:643:HOH:O	1.63	0.98
1:D:50:THR:HG22	1:D:135:LEU:HD11	1.45	0.96
1:A:53:LEU:HD22	1:A:59:THR:OG1	1.65	0.96
1:C:1:MET:HE3	1:C:1:MET:HA	1.48	0.94
1:A:279:ARG:HD3	1:C:412:LYS:HG2	1.51	0.91
1:D:79:ASN:HB3	1:D:82:SER:HB2	1.52	0.91
1:A:156:GLU:HG2	1:A:365:GLN:HE21	1.36	0.90
1:D:62:LEU:HD13	1:D:362:PHE:HB3	1.54	0.90
1:D:167:MET:HE2	1:D:172:ARG:HG2	1.55	0.89
1:D:148:LYS:NZ	4:D:601:HOH:O	2.06	0.88
1:C:18:ARG:O	1:C:22:ILE:HG13	1.74	0.87
1:D:159:THR:HG23	2:D:501:NAD:O3D	1.74	0.87
1:C:54:HIS:HB3	1:C:82:SER:HB2	1.57	0.87
1:D:163:ASN:O	1:D:167:MET:HG3	1.74	0.87
1:A:181:ASN:HA	1:A:186:LYS:HD2	1.55	0.86
1:D:156:GLU:O	1:D:181:ASN:HB2	1.76	0.85
1:C:119:PHE:HB3	1:C:120:PRO:HD2	1.57	0.85
1:B:137:ASN:O	1:B:141:GLU:HG2	1.78	0.84
1:B:392:GLU:HA	1:D:247:ILE:HD11	1.58	0.83
1:C:121:ASP:OD2	1:C:123:GLN:HB2	1.78	0.83
1:D:74:GLN:HE22	1:D:97:PRO:HG2	1.43	0.82
1:D:150:ILE:HD11	1:D:175:VAL:HG11	1.60	0.82
1:B:15:GLU:OE1	4:B:601:HOH:O	1.97	0.82
1:D:367:LEU:O	1:D:371:GLU:HG2	1.81	0.81
1:D:11:ILE:H	1:D:11:ILE:HD12	1.47	0.80
1:D:179:ASN:ND2	1:D:182:ASP:OD2	2.14	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:272:PHE:CD2	1:D:290:MET:HE2	2.18	0.78
1:A:53:LEU:CD2	1:A:59:THR:OG1	2.31	0.78
1:D:153:LEU:HD11	1:D:164:LEU:HD21	1.66	0.78
1:D:128:ILE:CD1	1:D:135:LEU:HD22	2.15	0.77
1:C:389:ILE:O	4:C:601:HOH:O	2.03	0.77
1:B:188:LYS:NZ	4:B:602:HOH:O	1.98	0.77
1:C:321:VAL:HA	1:D:19:LYS:NZ	2.00	0.77
1:C:235:PHE:HE1	1:D:197:GLU:HA	1.49	0.77
1:C:1:MET:HA	1:C:1:MET:CE	2.11	0.76
1:C:275:THR:HG22	1:C:299:ILE:HG22	1.66	0.76
1:D:181:ASN:HA	1:D:186:LYS:HG2	1.66	0.76
1:C:148:LYS:NZ	4:C:604:HOH:O	2.18	0.76
1:C:273:VAL:HG22	1:C:297:CYS:SG	2.25	0.76
1:A:60:ALA:HB1	1:A:91:ILE:HD11	1.67	0.76
1:B:268:GLU:O	4:B:603:HOH:O	2.03	0.76
1:A:4:PRO:O	4:A:601:HOH:O	2.02	0.75
1:C:316:LYS:HE2	4:C:760:HOH:O	1.87	0.75
1:C:163:ASN:O	1:C:167:MET:HG3	1.87	0.74
1:C:224:VAL:HG23	1:C:275:THR:HB	1.68	0.74
1:C:317:GLU:HG3	4:C:728:HOH:O	1.86	0.73
1:D:128:ILE:HD12	1:D:135:LEU:HD22	1.68	0.73
1:C:186:LYS:HG3	1:C:361:SER:HB3	1.71	0.73
1:B:417:LEU:HD11	1:D:278:CYS:HB2	1.70	0.73
1:B:301:HIS:O	1:B:302:PHE:HB2	1.89	0.72
1:B:111:TRP:O	1:B:115:GLN:HG2	1.89	0.72
1:A:287:LEU:HD23	1:A:290:MET:HE1	1.72	0.72
1:A:119:PHE:HB3	1:A:120:PRO:HD2	1.70	0.72
1:B:275:THR:HG22	1:B:299:ILE:HG22	1.71	0.72
1:B:275:THR:HG22	1:B:299:ILE:CG2	2.20	0.72
1:C:1:MET:HE3	1:C:1:MET:CA	2.19	0.72
1:C:266:SER:HB2	1:C:289:GLN:HB2	1.72	0.72
1:D:52:CYS:HB2	1:D:135:LEU:HD12	1.72	0.72
1:D:410:THR:HB	1:D:413:GLN:HG3	1.69	0.72
1:A:156:GLU:HG2	1:A:365:GLN:NE2	2.06	0.71
1:A:274:THR:OG1	1:A:298:ASN:ND2	2.24	0.71
1:C:117:LEU:HD13	1:C:125:LEU:HD22	1.73	0.70
1:B:346:ASN:CG	1:B:347:LEU:HD22	2.17	0.70
1:A:378:LYS:HG3	4:A:651:HOH:O	1.91	0.70
1:D:74:GLN:HE22	1:D:97:PRO:CG	2.04	0.70
1:A:26:ASN:O	1:A:401:LYS:NZ	2.25	0.70
1:D:53:LEU:HD23	1:D:59:THR:HG23	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:MET:HE2	1:D:254:MET:HE1	1.72	0.69
1:A:135:LEU:O	1:A:139:VAL:HG23	1.92	0.69
1:D:52:CYS:HA	1:D:76:SER:O	1.92	0.69
1:D:215:VAL:HB	1:D:269:ALA:HA	1.73	0.69
1:A:272:PHE:HB2	1:A:296:VAL:HG12	1.75	0.68
1:C:224:VAL:CG2	1:C:275:THR:HB	2.24	0.68
1:D:167:MET:CE	1:D:172:ARG:HG2	2.23	0.68
1:D:167:MET:HE2	1:D:172:ARG:CG	2.21	0.68
1:D:208:ASP:O	4:D:602:HOH:O	2.11	0.68
1:A:165:TYR:CD1	1:A:383:VAL:HG11	2.28	0.67
1:C:354:PRO:O	1:C:358:MET:HG2	1.94	0.67
1:B:346:ASN:ND2	1:B:347:LEU:HD22	2.09	0.67
1:B:43:PRO:HD3	4:B:604:HOH:O	1.95	0.67
1:D:143:PHE:HB2	1:D:145:GLN:OE1	1.95	0.67
4:A:669:HOH:O	1:C:166:LYS:HE3	1.94	0.67
1:D:45:LYS:HA	1:D:70:GLY:O	1.95	0.66
1:B:186:LYS:HG2	1:B:361:SER:OG	1.95	0.66
1:C:321:VAL:HA	1:D:19:LYS:HZ1	1.60	0.66
1:B:38:TYR:O	4:B:604:HOH:O	2.12	0.66
1:A:407:THR:HG22	1:C:244:VAL:HG12	1.78	0.66
1:C:354:PRO:HA	4:D:602:HOH:O	1.94	0.66
1:A:215:VAL:HB	1:A:269:ALA:HA	1.78	0.66
1:B:246:PRO:HD3	1:D:425:PHE:CE2	2.31	0.65
1:D:58:GLN:HG2	1:D:355:SER:HA	1.77	0.65
1:A:272:PHE:CE2	1:A:290:MET:HG2	2.31	0.65
1:A:404:VAL:HG13	1:C:259:VAL:HB	1.76	0.65
1:D:127:MET:SD	1:D:151:LYS:HB2	2.36	0.65
1:D:431:ARG:O	4:D:603:HOH:O	2.14	0.65
1:D:48:ARG:HG3	1:D:125:LEU:HA	1.76	0.65
1:C:105:THR:HG22	1:C:107:GLU:H	1.61	0.64
1:B:125:LEU:HG	1:B:150:ILE:HD11	1.80	0.64
1:A:157:THR:HG23	2:A:501:NAD:O2D	1.97	0.64
1:D:125:LEU:O	1:D:150:ILE:HG22	1.98	0.64
1:B:4:PRO:O	1:B:5:SER:HB3	1.96	0.64
1:C:117:LEU:CD1	1:C:125:LEU:HD22	2.28	0.63
1:C:156:GLU:HG3	1:C:156:GLU:O	1.99	0.63
1:D:218:VAL:HG22	1:D:273:VAL:CG1	2.28	0.63
1:A:198:SER:HB2	1:A:346:ASN:HB2	1.81	0.63
1:B:41:SER:HB2	4:B:604:HOH:O	1.98	0.63
1:D:128:ILE:HG23	1:D:153:LEU:CB	2.29	0.63
1:A:164:LEU:HD21	1:A:177:ALA:HB1	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:156:GLU:O	1:B:156:GLU:HG3	1.99	0.62
1:A:261:THR:HG22	1:C:407:THR:HG22	1.79	0.62
1:B:412:LYS:HE3	1:D:279:ARG:HD2	1.82	0.62
1:D:189:PHE:HA	1:D:193:TYR:HD2	1.64	0.62
1:D:164:LEU:HD12	1:D:167:MET:SD	2.39	0.62
1:D:401:LYS:NZ	4:D:605:HOH:O	2.23	0.62
1:A:266:SER:HA	1:A:272:PHE:CZ	2.34	0.62
1:D:125:LEU:HD13	1:D:127:MET:O	2.00	0.62
1:D:39:GLY:O	1:D:42:LYS:HD2	1.99	0.62
1:A:101:TRP:O	1:A:104:GLU:HG3	1.99	0.61
1:C:127:MET:HE2	1:C:151:LYS:HB2	1.81	0.61
4:A:658:HOH:O	1:C:407:THR:HG23	1.99	0.61
1:D:380:ALA:O	1:D:384:HIS:NE2	2.32	0.61
1:C:301:HIS:O	1:C:302:PHE:HB2	2.01	0.61
1:D:126:ASN:HA	1:D:150:ILE:HA	1.83	0.61
1:D:128:ILE:HG23	1:D:153:LEU:HB2	1.83	0.60
1:B:246:PRO:HD3	1:D:425:PHE:HE2	1.66	0.60
1:A:244:VAL:HG11	1:C:409:LEU:HG	1.83	0.60
1:D:50:THR:CG2	1:D:135:LEU:HD11	2.26	0.60
1:B:40:PRO:HA	4:B:777:HOH:O	2.00	0.60
1:B:9:ALA:O	4:B:605:HOH:O	2.17	0.60
1:D:354:PRO:O	1:D:358:MET:HG2	2.02	0.60
1:D:304:ILE:HG13	1:D:304:ILE:O	2.00	0.59
1:C:45:LYS:NZ	4:C:602:HOH:O	2.16	0.59
1:C:210:MET:HB2	1:D:354:PRO:HB2	1.85	0.59
1:D:158:THR:HG22	1:D:181:ASN:ND2	2.17	0.59
1:D:50:THR:OG1	1:D:116:THR:O	2.21	0.59
1:A:157:THR:O	1:A:161:VAL:HG13	2.02	0.59
1:A:156:GLU:O	1:A:186:LYS:NZ	2.30	0.59
1:A:278:CYS:SG	1:C:417:LEU:HD11	2.43	0.59
1:A:303:ASP:C	1:A:304:ILE:HD13	2.28	0.59
1:D:426:LYS:HE2	1:D:430:TYR:CZ	2.38	0.58
1:A:219:ALA:HA	1:A:242:THR:OG1	2.03	0.58
1:B:157:THR:O	1:B:161:VAL:HG23	2.03	0.58
1:A:68:GLU:HA	4:A:731:HOH:O	2.04	0.58
1:A:301:HIS:HA	1:A:344:LEU:HD21	1.84	0.58
1:C:90:ALA:O	1:C:94:THR:HG23	2.02	0.58
1:C:166:LYS:HA	1:C:169:LYS:HB3	1.86	0.58
1:C:301:HIS:HA	1:C:344:LEU:HD21	1.86	0.58
1:D:200:ILE:O	1:D:204:LYS:HG2	2.04	0.58
1:D:167:MET:HB3	1:D:172:ARG:HB3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:GLY:HA3	1:A:301:HIS:NE2	2.19	0.58
1:C:373:TRP:O	4:C:603:HOH:O	2.17	0.57
4:A:761:HOH:O	1:C:407:THR:HG21	2.03	0.57
1:A:408:LYS:NZ	4:A:602:HOH:O	2.06	0.57
1:C:181:ASN:HA	1:C:186:LYS:HD2	1.86	0.57
1:A:184:VAL:HG11	1:C:247:ILE:HD13	1.86	0.57
1:B:126:ASN:C	1:B:150:ILE:HD13	2.29	0.57
1:D:171:GLY:HA2	4:D:725:HOH:O	2.04	0.57
1:D:279:ARG:HD3	1:D:280:ASP:OD2	2.04	0.57
1:D:53:LEU:HD21	1:D:362:PHE:CZ	2.40	0.57
1:A:210:MET:O	1:A:214:LYS:HG3	2.05	0.56
1:B:127:MET:HE1	1:B:372:LEU:HB2	1.86	0.56
1:C:105:THR:HG22	1:C:107:GLU:N	2.19	0.56
1:C:113:ILE:HG23	1:C:135:LEU:HD13	1.87	0.56
1:C:224:VAL:HG22	2:C:501:NAD:O2N	2.06	0.56
1:D:272:PHE:CE2	1:D:290:MET:HG2	2.40	0.56
1:A:407:THR:HG23	4:A:649:HOH:O	2.05	0.56
1:D:356:PHE:HE2	1:D:401:LYS:HG2	1.70	0.56
1:B:101:TRP:O	1:B:104:GLU:HG3	2.05	0.56
1:B:169:LYS:HG2	4:B:862:HOH:O	2.06	0.56
1:C:42:LYS:HG3	1:C:45:LYS:HE3	1.87	0.56
1:C:111:TRP:NE1	4:C:610:HOH:O	2.32	0.56
1:A:287:LEU:HA	1:A:290:MET:SD	2.46	0.56
1:C:412:LYS:HA	4:C:748:HOH:O	2.06	0.55
1:B:218:VAL:HG22	1:B:273:VAL:HB	1.88	0.55
1:B:431:ARG:CZ	1:D:184:VAL:HG22	2.36	0.55
1:C:35:ARG:HG3	1:C:65:THR:HG23	1.87	0.55
1:C:327:ARG:NH1	1:D:22:ILE:HD12	2.21	0.55
1:C:274:THR:OG1	1:C:298:ASN:HA	2.06	0.55
1:B:389:ILE:HG23	1:B:390:LEU:HG	1.88	0.55
1:C:54:HIS:HA	1:C:77:SER:OG	2.07	0.55
1:C:354:PRO:HG2	1:D:210:MET:HE3	1.89	0.55
1:A:377:ASP:N	1:A:377:ASP:OD1	2.40	0.55
1:B:287:LEU:HD13	1:B:311:LEU:HD13	1.89	0.55
1:C:133:GLY:HA2	4:C:674:HOH:O	2.07	0.55
1:A:143:PHE:N	1:A:144:PRO:HD3	2.22	0.55
1:D:86:ASN:OD1	1:D:86:ASN:N	2.37	0.55
1:C:230:GLN:NE2	4:C:615:HOH:O	2.40	0.54
1:B:144:PRO:HA	1:B:147:LEU:HD12	1.89	0.54
1:A:254:MET:HE3	1:C:402:LEU:HD11	1.90	0.54
1:C:119:PHE:HB3	1:C:120:PRO:CD	2.34	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:CYS:HB3	1:A:130:ASP:OD1	2.08	0.54
1:D:55:MET:HB3	1:D:83:THR:HG23	1.89	0.54
1:D:414:ALA:HB1	1:D:419:VAL:O	2.08	0.53
1:D:176:PRO:HA	1:D:381:VAL:HA	1.89	0.53
1:D:206:ALA:HB1	1:D:325:VAL:HG22	1.90	0.53
1:A:31:LEU:HD13	1:A:61:VAL:HG12	1.90	0.53
1:D:48:ARG:HB2	1:D:119:PHE:CE2	2.44	0.53
1:A:21:ILE:O	1:A:25:GLU:HG3	2.07	0.53
1:D:112:CYS:HB2	4:D:622:HOH:O	2.08	0.53
1:D:282:ILE:HB	1:D:306:ILE:HG23	1.89	0.53
1:D:287:LEU:HA	1:D:290:MET:SD	2.47	0.53
1:C:354:PRO:CB	1:D:210:MET:HB2	2.39	0.53
1:A:224:VAL:HG11	1:A:299:ILE:HG23	1.89	0.53
1:B:198:SER:HB2	1:B:346:ASN:HB2	1.91	0.53
1:B:387:PRO:HD2	1:B:390:LEU:HD12	1.90	0.53
1:C:156:GLU:HG3	1:C:186:LYS:HE3	1.89	0.53
1:A:146:TYR:N	4:A:603:HOH:O	2.15	0.53
1:D:241:VAL:O	1:D:259:VAL:HA	2.08	0.53
1:A:130:ASP:OD1	1:A:135:LEU:HB3	2.08	0.52
1:A:247:ILE:HD13	1:C:184:VAL:HG11	1.90	0.52
1:B:413:GLN:HB3	1:D:278:CYS:SG	2.49	0.52
1:D:53:LEU:O	1:D:77:SER:HA	2.08	0.52
1:B:247:ILE:HD12	1:D:392:GLU:HG2	1.91	0.52
1:D:48:ARG:HB2	1:D:119:PHE:CZ	2.44	0.52
1:D:127:MET:HE1	1:D:372:LEU:HB2	1.91	0.52
1:B:179:ASN:ND2	1:B:182:ASP:HB2	2.23	0.52
1:D:151:LYS:O	1:D:176:PRO:HD2	2.10	0.52
1:B:346:ASN:ND2	1:B:347:LEU:CD2	2.72	0.52
1:A:371:GLU:HG2	1:A:379:TYR:OH	2.10	0.52
1:B:179:ASN:HD21	1:B:182:ASP:HB2	1.74	0.52
1:B:296:VAL:HG12	1:B:306:ILE:HD13	1.92	0.52
1:D:159:THR:CG2	2:D:501:NAD:O3D	2.55	0.52
1:D:292:ASP:OD1	1:D:335:HIS:N	2.42	0.52
1:B:100:ALA:HA	4:B:609:HOH:O	2.09	0.52
1:D:18[B]:ARG:O	1:D:22:ILE:HG23	2.10	0.52
1:B:181:ASN:HA	1:B:186:LYS:HD2	1.92	0.52
1:C:167:MET:HA	1:C:172:ARG:HG2	1.91	0.51
1:D:384:HIS:C	1:D:385:VAL:HG23	2.34	0.51
1:A:125:LEU:H	1:A:125:LEU:HD23	1.75	0.51
1:C:278:CYS:C	1:C:304:ILE:HD12	2.35	0.51
1:D:158:THR:HG22	1:D:181:ASN:HD21	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:224:VAL:HG11	1:B:299:ILE:HG23	1.93	0.51
1:C:235:PHE:CE1	1:D:197:GLU:HA	2.37	0.51
1:A:55:MET:HE1	1:A:63:ILE:HD12	1.91	0.51
1:C:8:VAL:HG12	4:C:635:HOH:O	2.11	0.51
1:D:11:ILE:HD12	1:D:11:ILE:N	2.21	0.51
1:D:58:GLN:HG3	1:D:355:SER:HB3	1.92	0.51
1:A:255:GLU:HB2	4:A:630:HOH:O	2.11	0.51
1:B:279:ARG:HA	1:B:304:ILE:CD1	2.41	0.51
1:D:37:LYS:O	4:D:604:HOH:O	2.18	0.51
1:A:148:LYS:HB2	4:A:776:HOH:O	2.11	0.51
1:A:399:LEU:HB3	1:A:404:VAL:HB	1.92	0.51
1:D:124:PRO:HB2	1:D:149:ASN:HD21	1.75	0.51
1:A:374:THR:HG21	4:A:781:HOH:O	2.10	0.51
1:D:180:VAL:HG22	1:D:364:ASN:HB3	1.93	0.51
1:A:153:LEU:C	1:A:153:LEU:HD12	2.36	0.50
1:A:358:MET:HE2	1:A:362:PHE:CZ	2.47	0.50
1:C:211:ILE:HB	1:C:235:PHE:HB3	1.93	0.50
1:D:11:ILE:HA	1:D:89:ALA:CB	2.42	0.50
1:D:57:VAL:HG13	1:D:87:ALA:HB2	1.92	0.50
1:B:54:HIS:HA	1:B:77:SER:OG	2.11	0.50
1:D:172:ARG:O	1:D:173:LEU:HB2	2.10	0.50
1:A:318:LYS:HB3	4:A:605:HOH:O	2.10	0.50
1:B:266:SER:HA	1:B:272:PHE:CZ	2.46	0.50
1:C:161:VAL:HG11	1:C:179:ASN:OD1	2.11	0.50
1:C:322:LYS:HB2	1:C:323:PRO:CD	2.42	0.50
1:D:303:ASP:HB3	1:D:343:ARG:HG2	1.93	0.50
1:B:431:ARG:HG2	1:D:431:ARG:HA	1.93	0.50
1:C:321:VAL:HA	1:D:19:LYS:HZ3	1.72	0.50
1:A:273:VAL:HB	1:A:297:CYS:SG	2.52	0.50
1:B:278:CYS:O	1:B:305:GLU:HG2	2.12	0.50
1:D:48:ARG:NH1	1:D:123:GLN:OE1	2.44	0.50
1:B:431:ARG:HA	1:D:431:ARG:HG2	1.93	0.50
1:D:363:THR:HG22	1:D:394:VAL:HG22	1.94	0.50
1:C:354:PRO:HB2	1:D:210:MET:HB2	1.93	0.50
1:D:275:THR:HG22	1:D:299:ILE:CG2	2.42	0.50
1:D:311:LEU:HD12	1:D:330:MET:HE1	1.92	0.50
1:A:387:PRO:HD2	1:A:390:LEU:HD12	1.94	0.49
1:B:307:ASP:HB2	4:B:681:HOH:O	2.12	0.49
1:C:94:THR:O	4:C:605:HOH:O	2.20	0.49
1:C:181:ASN:O	1:C:187:SER:HB3	2.12	0.49
1:D:111:TRP:NE1	1:D:115:GLN:OE1	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:ASN:HB2	1:A:223:ASP:HB3	1.95	0.49
1:B:241:VAL:O	1:B:259:VAL:HA	2.13	0.49
1:C:353:HIS:HB2	1:C:358:MET:SD	2.52	0.49
1:C:266:SER:O	1:C:290:MET:HA	2.11	0.49
1:D:125:LEU:HD12	1:D:126:ASN:N	2.27	0.49
1:C:219:ALA:HA	1:C:242:THR:OG1	2.13	0.49
1:A:145:GLN:HG2	4:A:603:HOH:O	2.12	0.49
1:B:101:TRP:CE2	1:B:104:GLU:HG2	2.46	0.49
1:D:11:ILE:HA	1:D:89:ALA:HB1	1.94	0.49
1:B:63:ILE:O	1:B:67:VAL:HG23	2.12	0.49
1:A:298:ASN:HB3	1:A:342:GLY:O	2.13	0.49
1:A:407:THR:CG2	1:C:244:VAL:HG12	2.41	0.49
1:A:54:HIS:HB3	1:A:82:SER:HB2	1.95	0.48
1:B:48:ARG:H	1:B:126:ASN:CG	2.21	0.48
1:B:199:LEU:HD22	1:B:228:CYS:HB3	1.95	0.48
1:B:279:ARG:HA	1:B:304:ILE:HD12	1.94	0.48
1:B:420:SER:HA	4:B:623:HOH:O	2.13	0.48
1:B:156:GLU:HB2	1:B:365:GLN:HE21	1.78	0.48
1:B:331:GLN:HA	1:B:331:GLN:HE21	1.77	0.48
1:C:197:GLU:HG2	1:D:235:PHE:CD2	2.48	0.48
1:C:430:TYR:CE2	1:C:432:TYR:HA	2.48	0.48
1:A:107:GLU:HG2	1:A:108:GLU:H	1.78	0.48
1:A:224:VAL:HG11	1:A:299:ILE:CG2	2.43	0.48
1:A:273:VAL:HA	1:A:297:CYS:O	2.14	0.48
1:B:322:LYS:HB2	1:B:323:PRO:CD	2.44	0.48
1:C:53:LEU:O	1:C:54:HIS:C	2.56	0.48
1:D:179:ASN:HD21	1:D:182:ASP:CG	2.21	0.48
1:A:266:SER:HA	1:A:272:PHE:HZ	1.78	0.48
1:B:210:MET:O	1:B:214:LYS:HG3	2.14	0.48
1:C:107:GLU:O	1:C:108:GLU:HB2	2.13	0.48
1:C:191:ASN:OD1	1:C:191:ASN:N	2.44	0.48
1:C:263:GLU:O	1:C:267:LYS:HE3	2.13	0.48
1:D:68:GLU:HG2	4:D:668:HOH:O	2.12	0.48
1:A:284:SER:HB3	1:A:310:TRP:CD2	2.49	0.48
1:A:341:GLU:HB2	1:A:343:ARG:NH1	2.27	0.48
1:B:53:LEU:HD22	1:B:59:THR:OG1	2.13	0.48
1:B:125:LEU:HD23	1:B:125:LEU:N	2.28	0.48
1:B:184:VAL:HG21	1:B:432:TYR:CD1	2.48	0.48
1:D:18[A]:ARG:O	1:D:22:ILE:HG23	2.13	0.48
1:D:198:SER:HB2	1:D:346:ASN:HB2	1.96	0.48
1:B:55:MET:HB3	1:B:83:THR:HG23	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:279:ARG:HD3	1:B:280:ASP:OD1	2.13	0.48
1:B:250:LEU:HD13	1:D:399:LEU:HD23	1.95	0.47
1:C:233:LYS:HE2	1:C:257:TYR:CE1	2.48	0.47
1:D:149:ASN:OD1	1:D:149:ASN:N	2.47	0.47
1:A:140:HIS:HA	1:A:147:LEU:HD11	1.96	0.47
1:D:74:GLN:HG2	1:D:119:PHE:CE2	2.49	0.47
1:D:288:GLN:O	1:D:334:LYS:NZ	2.46	0.47
1:B:412:LYS:HD2	1:B:412:LYS:HA	1.48	0.47
1:D:53:LEU:HD11	1:D:54:HIS:NE2	2.29	0.47
1:A:191:ASN:OD1	1:A:191:ASN:N	2.47	0.47
1:B:51:GLY:HA2	1:B:129:LEU:O	2.14	0.47
1:B:54:HIS:HB3	1:B:82:SER:HB2	1.96	0.47
1:C:275:THR:HG22	1:C:299:ILE:CG2	2.41	0.47
1:A:55:MET:HB3	1:A:83:THR:HG23	1.95	0.47
1:A:389:ILE:HD13	1:A:424:PRO:HD3	1.95	0.47
1:B:278:CYS:HB2	1:D:417:LEU:HD11	1.96	0.47
1:C:38:TYR:HB3	1:C:43:PRO:HG2	1.96	0.47
1:D:272:PHE:CG	1:D:290:MET:HE2	2.49	0.47
1:D:322:LYS:HB2	1:D:323:PRO:HD2	1.96	0.47
1:A:50:THR:HG23	1:A:116:THR:HB	1.96	0.47
1:A:267:LYS:HG2	1:A:289:GLN:HB3	1.97	0.47
1:B:152:GLY:HA3	1:B:372:LEU:HD13	1.96	0.47
1:B:218:VAL:HA	1:B:273:VAL:O	2.14	0.47
1:C:311:LEU:HG	1:C:328:TYR:CE2	2.50	0.47
1:D:301:HIS:O	1:D:302:PHE:HB2	2.14	0.47
1:A:117:LEU:HA	1:A:125:LEU:HD22	1.97	0.47
1:B:80:ILE:HD11	1:B:343:ARG:NH1	2.29	0.47
1:A:161:VAL:HG11	1:A:179:ASN:OD1	2.15	0.47
1:C:213:GLY:O	1:D:402:LEU:HD22	2.14	0.47
1:C:409:LEU:CD2	1:C:413:GLN:HB3	2.45	0.47
1:D:82:SER:OG	1:D:347:LEU:HG	2.14	0.47
1:D:92:ALA:HB2	1:D:98:VAL:CG1	2.45	0.47
1:A:105:THR:OG1	1:A:107:GLU:HG2	2.14	0.47
1:C:111:TRP:O	1:C:115:GLN:HG2	2.15	0.47
1:C:243:GLU:OE2	1:C:248:ASN:HB2	2.15	0.46
1:A:54:HIS:HA	1:A:77:SER:OG	2.14	0.46
1:D:101:TRP:CZ2	1:D:108:GLU:HB3	2.50	0.46
1:A:111:TRP:O	1:A:115:GLN:HG2	2.15	0.46
1:C:39:GLY:N	1:C:40:PRO:HD2	2.30	0.46
1:D:153:LEU:HD23	1:D:175:VAL:HG23	1.98	0.46
1:A:430:TYR:CE2	1:A:432:TYR:HA	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:346:ASN:CG	1:B:347:LEU:CD2	2.88	0.46
1:C:56:THR:HG21	3:C:502:ADN:N1	2.31	0.46
1:D:157:THR:HG21	1:D:301:HIS:NE2	2.30	0.46
1:B:346:ASN:OD1	1:B:347:LEU:CD2	2.64	0.46
1:D:296:VAL:HG12	1:D:306:ILE:HD13	1.98	0.46
1:B:149:ASN:OD1	1:B:151:LYS:NZ	2.49	0.46
1:D:27:GLU:O	1:D:356:PHE:HA	2.16	0.46
1:B:311:LEU:HD12	1:B:330:MET:HE1	1.96	0.46
1:B:376:SER:O	1:B:377:ASP:CB	2.64	0.46
1:B:370:ILE:HG21	4:B:825:HOH:O	2.16	0.46
1:B:161:VAL:HG12	1:B:165:TYR:CE2	2.50	0.46
1:C:408:LYS:HB3	1:C:408:LYS:HE2	1.73	0.46
1:D:15:GLU:HA	1:D:18[A]:ARG:CZ	2.46	0.46
1:A:148:LYS:HA	1:A:148:LYS:HD3	1.72	0.45
1:A:262:MET:HG3	1:A:286:HIS:CD2	2.50	0.45
1:B:211:ILE:HB	1:B:235:PHE:HB3	1.97	0.45
1:D:53:LEU:HD23	1:D:59:THR:CG2	2.45	0.45
1:D:221:TYR:HB2	1:D:241:VAL:HG11	1.98	0.45
1:C:26:ASN:ND2	1:D:292:ASP:OD2	2.50	0.45
1:D:292:ASP:HB2	1:D:334:LYS:HG2	1.98	0.45
1:A:210:MET:HE2	1:D:254:MET:CE	2.44	0.45
1:B:266:SER:HA	1:B:272:PHE:HZ	1.80	0.45
1:C:166:LYS:HG2	1:C:169:LYS:HD3	1.98	0.45
1:B:157:THR:HG23	1:B:159:THR:H	1.80	0.45
1:C:74:GLN:HG2	1:C:97:PRO:HB2	1.97	0.45
1:C:357:VAL:HB	1:D:210:MET:HE1	1.99	0.45
1:D:127:MET:HE1	1:D:372:LEU:CB	2.47	0.45
1:A:101:TRP:CE2	1:A:104:GLU:HG2	2.52	0.45
1:C:269:ALA:HB3	1:C:272:PHE:CZ	2.52	0.45
1:C:358:MET:HA	1:C:358:MET:HE3	1.99	0.45
1:D:137:ASN:HB2	4:D:647:HOH:O	2.17	0.45
1:D:322:LYS:HB2	1:D:323:PRO:CD	2.46	0.45
1:B:273:VAL:HG22	1:B:297:CYS:SG	2.57	0.45
1:D:176:PRO:HD3	1:D:381:VAL:HG12	1.97	0.45
1:D:331:GLN:HA	1:D:331:GLN:NE2	2.31	0.45
1:D:353:HIS:HB2	1:D:358:MET:SD	2.57	0.45
1:B:376:SER:O	1:B:377:ASP:HB2	2.17	0.45
1:D:143:PHE:HB2	1:D:145:GLN:CD	2.41	0.45
1:A:48:ARG:NH1	1:A:124:PRO:O	2.50	0.45
1:A:241:VAL:O	1:A:259:VAL:HA	2.16	0.45
1:B:125:LEU:HD23	1:B:125:LEU:H	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:233:LYS:NZ	1:D:255:GLU:O	2.50	0.44
1:D:430:TYR:CE2	1:D:432:TYR:HA	2.52	0.44
1:A:221:TYR:CD2	1:A:252:ALA:HB2	2.52	0.44
1:C:38:TYR:HB3	1:C:43:PRO:CG	2.47	0.44
1:B:200:ILE:HD11	1:B:232:LEU:HD23	1.98	0.44
1:C:231:ALA:HA	4:C:680:HOH:O	2.17	0.44
1:A:24:ALA:O	1:A:28:MET:HG3	2.17	0.44
1:C:17:GLY:O	1:C:21:ILE:HD12	2.18	0.44
1:C:274:THR:HB	1:C:305:GLU:OE1	2.18	0.44
1:D:51:GLY:HA3	1:D:75:TRP:CE3	2.52	0.44
1:D:53:LEU:CD1	1:D:54:HIS:CD2	3.00	0.44
1:B:7:LYS:HD3	1:B:101:TRP:CH2	2.53	0.44
1:D:218:VAL:HG22	1:D:273:VAL:HG13	1.97	0.44
1:D:344:LEU:HD23	1:D:344:LEU:HA	1.84	0.44
1:D:399:LEU:HB3	1:D:404:VAL:HB	1.99	0.44
1:B:37:LYS:HD2	1:B:38:TYR:CZ	2.52	0.44
1:D:262:MET:HE2	1:D:262:MET:HB2	1.89	0.44
1:B:279:ARG:CA	1:B:304:ILE:HD13	2.48	0.44
1:C:38:TYR:O	1:C:43:PRO:HD3	2.18	0.44
1:D:54:HIS:HB3	1:D:82:SER:OG	2.18	0.44
1:A:287:LEU:HD23	1:A:290:MET:CE	2.45	0.43
1:B:322:LYS:HE3	1:B:325:VAL:HG21	2.00	0.43
1:D:20:ALA:HB3	1:D:57:VAL:HG11	2.00	0.43
1:D:49:ILE:HG23	1:D:127:MET:HB2	2.00	0.43
1:D:81:PHE:HE1	1:D:343:ARG:HD2	1.83	0.43
1:A:55:MET:HE1	1:A:63:ILE:CD1	2.47	0.43
1:B:162:HIS:CE1	1:D:417:LEU:O	2.71	0.43
1:C:387:PRO:HG2	1:C:390:LEU:HG	1.99	0.43
1:A:378:LYS:N	4:A:624:HOH:O	2.52	0.43
1:B:244:VAL:HG11	1:D:409:LEU:HG	1.99	0.43
1:A:163:ASN:HD21	1:A:167:MET:CE	2.30	0.43
1:B:360:ASN:HD22	1:B:394:VAL:HG12	1.83	0.43
1:C:107:GLU:O	1:C:108:GLU:CB	2.66	0.43
1:D:48:ARG:CZ	1:D:123:GLN:OE1	2.66	0.43
1:A:371:GLU:CD	1:A:375:LYS:HD2	2.43	0.43
1:B:274:THR:OG1	1:B:298:ASN:HA	2.18	0.43
1:B:343:ARG:HA	4:B:746:HOH:O	2.19	0.43
1:C:219:ALA:HB1	1:C:281:ILE:HG21	2.01	0.43
1:C:299:ILE:HG23	1:C:299:ILE:O	2.19	0.43
1:B:385:VAL:HG11	4:B:676:HOH:O	2.19	0.43
1:A:233:LYS:NZ	4:A:614:HOH:O	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:28:MET:HA	1:B:359:SER:HB2	1.99	0.43
1:C:173:LEU:HD11	1:C:175:VAL:O	2.18	0.43
1:D:125:LEU:HD12	1:D:127:MET:N	2.33	0.43
1:D:170:GLU:CG	1:D:171:GLY:N	2.81	0.43
1:A:4:PRO:C	1:A:6:TYR:H	2.26	0.43
1:A:53:LEU:O	1:A:54:HIS:C	2.61	0.43
1:A:426:LYS:HD2	1:A:430:TYR:CG	2.53	0.43
1:B:301:HIS:O	1:B:302:PHE:CB	2.61	0.43
1:C:200:ILE:HD12	1:C:231:ALA:HB1	2.01	0.43
1:B:221:TYR:CZ	1:B:226:LYS:HG2	2.53	0.43
1:A:168:PHE:CE2	1:A:382:GLY:HA2	2.53	0.43
1:B:132:GLY:HA3	1:B:301:HIS:NE2	2.34	0.43
1:C:409:LEU:HD23	1:C:413:GLN:HB3	2.00	0.43
1:D:170:GLU:HG3	1:D:171:GLY:N	2.34	0.43
1:D:426:LYS:HD2	1:D:430:TYR:CD1	2.54	0.43
1:B:185:THR:HG21	1:B:391:ASP:O	2.19	0.42
1:B:269:ALA:HB3	1:B:272:PHE:CZ	2.54	0.42
1:C:99:TYR:O	1:C:112:CYS:HB3	2.19	0.42
1:D:48:ARG:HB3	1:D:72:GLN:OE1	2.19	0.42
1:A:18:ARG:NH2	1:A:64:GLU:OE1	2.51	0.42
1:B:219:ALA:HA	1:B:242:THR:OG1	2.19	0.42
2:B:501:NAD:H6N	2:B:501:NAD:H2D	1.87	0.42
1:C:279:ARG:HE	1:C:279:ARG:HB2	1.37	0.42
1:C:301:HIS:O	1:C:302:PHE:CB	2.67	0.42
1:D:53:LEU:HD12	1:D:54:HIS:CD2	2.53	0.42
1:D:204:LYS:O	1:D:208:ASP:N	2.47	0.42
1:D:387:PRO:HD2	1:D:390:LEU:HD12	2.02	0.42
1:A:162:HIS:NE2	1:C:417:LEU:O	2.46	0.42
1:B:233:LYS:HD2	4:C:740:HOH:O	2.19	0.42
1:C:126:ASN:O	1:C:150:ILE:HA	2.19	0.42
1:A:173:LEU:HD23	1:A:173:LEU:HA	1.91	0.42
1:A:293:ASP:N	1:A:335:HIS:O	2.50	0.42
1:B:10:ASP:HB3	1:B:13:LEU:HG	2.01	0.42
1:B:259:VAL:CB	1:D:404:VAL:HG13	2.31	0.42
1:D:221:TYR:CZ	1:D:226:LYS:HG2	2.53	0.42
1:A:371:GLU:CG	1:A:379:TYR:HE2	2.33	0.42
1:B:344:LEU:HD12	4:B:746:HOH:O	2.19	0.42
1:B:366:VAL:O	1:B:370:ILE:HG13	2.19	0.42
1:A:4:PRO:O	1:A:5:SER:HB3	2.19	0.42
1:A:266:SER:O	1:A:290:MET:HA	2.19	0.42
1:C:318:LYS:NZ	1:C:328:TYR:OH	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:201:ASP:O	1:D:205:ARG:HG3	2.19	0.42
1:D:334:LYS:HB2	1:D:334:LYS:HE3	1.66	0.42
1:A:53:LEU:HD21	3:A:502:ADN:N3	2.35	0.42
1:A:138:LEU:O	1:A:142:LYS:O	2.38	0.42
1:A:417:LEU:HB2	1:A:419:VAL:HG22	2.02	0.42
1:B:4:PRO:O	1:B:5:SER:CB	2.64	0.42
1:B:36:LYS:HE2	4:B:791:HOH:O	2.20	0.42
1:B:393:GLU:O	1:B:397:LEU:HG	2.20	0.42
1:C:293:ASP:OD1	1:C:327:ARG:HD2	2.20	0.42
1:D:367:LEU:O	1:D:368:ALA:C	2.62	0.42
1:C:322:LYS:HB2	1:C:323:PRO:HD2	2.00	0.42
1:D:54:HIS:CD2	1:D:54:HIS:H	2.36	0.42
1:A:190:ASP:OD2	1:A:353:HIS:HD2	2.03	0.42
1:B:125:LEU:H	1:B:125:LEU:CD2	2.33	0.42
1:C:44:LEU:HB3	1:C:71:ALA:HB2	2.02	0.42
1:C:158:THR:O	1:C:162:HIS:CD2	2.73	0.42
1:C:360:ASN:O	1:C:364:ASN:ND2	2.53	0.42
1:C:375:LYS:HA	4:C:621:HOH:O	2.20	0.42
1:A:146:TYR:HB3	4:A:616:HOH:O	2.20	0.41
4:C:686:HOH:O	1:D:204:LYS:HE2	2.19	0.41
1:D:11:ILE:H	1:D:11:ILE:CD1	2.22	0.41
1:D:131:ASP:OD1	1:D:156:GLU:HB3	2.20	0.41
1:D:402:LEU:HB2	1:D:404:VAL:HG23	2.02	0.41
1:A:128:ILE:O	1:A:153:LEU:HA	2.19	0.41
1:C:217:CYS:HA	1:C:240:ILE:O	2.19	0.41
1:B:101:TRP:NE1	1:B:104:GLU:HG2	2.36	0.41
1:B:346:ASN:HD21	1:B:347:LEU:CD2	2.33	0.41
1:D:82:SER:HA	1:D:347:LEU:HB3	2.03	0.41
1:A:49:ILE:HG12	1:A:127:MET:HB2	2.02	0.41
1:A:354:PRO:HB2	1:B:210:MET:HB2	2.03	0.41
1:C:184:VAL:HG12	1:C:391:ASP:CG	2.45	0.41
1:D:35:ARG:HG3	1:D:65:THR:HG23	2.01	0.41
1:A:114:GLU:HG3	1:A:138:LEU:HD21	2.03	0.41
1:B:29:PRO:HA	4:B:645:HOH:O	2.20	0.41
1:C:188:LYS:HE3	1:C:188:LYS:HB3	1.92	0.41
1:D:8:VAL:O	1:D:9:ALA:HB3	2.19	0.41
1:A:145:GLN:H	1:A:145:GLN:CD	2.29	0.41
1:A:274:THR:OG1	1:A:298:ASN:HA	2.21	0.41
1:D:287:LEU:HD23	1:D:290:MET:SD	2.60	0.41
1:A:322:LYS:HE2	1:A:322:LYS:HB3	1.16	0.41
1:D:53:LEU:HB3	1:D:75:TRP:HZ2	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:148:LYS:HE3	1:D:148:LYS:HB3	1.73	0.41
1:A:199:LEU:HD22	1:A:228:CYS:SG	2.61	0.41
1:B:278:CYS:HB2	1:D:417:LEU:HD21	2.03	0.41
2:B:501:NAD:C4N	3:B:502:ADN:H3'	2.51	0.41
1:C:185:THR:HG21	1:C:394:VAL:HB	2.03	0.41
1:C:279:ARG:HB3	1:C:304:ILE:HB	2.03	0.41
1:D:143:PHE:HB2	1:D:145:GLN:NE2	2.35	0.41
1:A:107:GLU:HG2	1:A:108:GLU:N	2.35	0.41
1:A:221:TYR:CE1	1:A:226:LYS:HA	2.56	0.41
1:A:371:GLU:HG3	1:A:379:TYR:HE2	1.86	0.41
1:B:163:ASN:OD1	1:B:167:MET:HE3	2.21	0.41
1:B:299:ILE:HG13	2:B:501:NAD:N7N	2.36	0.41
1:B:332:SER:HA	4:B:770:HOH:O	2.21	0.41
1:B:344:LEU:HD12	1:B:344:LEU:H	1.86	0.41
1:C:191:ASN:O	1:C:224:VAL:HA	2.21	0.41
1:D:3:LYS:N	1:D:4:PRO:CD	2.84	0.41
1:D:18[A]:ARG:HE	1:D:18[A]:ARG:HB3	1.27	0.41
1:D:53:LEU:HB3	1:D:75:TRP:CZ2	2.55	0.41
1:D:266:SER:O	1:D:290:MET:HA	2.21	0.41
1:D:328:TYR:O	1:D:335:HIS:HA	2.21	0.41
1:B:406:LEU:HG	1:B:407:THR:H	1.86	0.41
1:C:240:ILE:HG21	1:C:260:THR:HG22	2.03	0.41
1:D:158:THR:HG23	2:D:501:NAD:H2D	2.03	0.41
1:B:18:ARG:HE	1:B:18:ARG:HA	1.86	0.40
1:B:90:ALA:O	1:B:94:THR:HG23	2.21	0.40
1:B:395:ALA:HB3	1:D:247:ILE:HD13	2.02	0.40
1:B:188:LYS:HA	1:B:188:LYS:HE2	2.03	0.40
1:C:363:THR:O	1:C:367:LEU:HG	2.22	0.40
1:A:155:GLU:OE2	1:A:157:THR:HB	2.21	0.40
1:B:44:LEU:HB3	1:B:71:ALA:HB2	2.03	0.40
1:C:54:HIS:O	1:C:56:THR:HG23	2.21	0.40
1:D:180:VAL:HG11	1:D:365:GLN:HG2	2.02	0.40
1:B:3:LYS:HA	1:B:4:PRO:HD3	1.89	0.40
1:B:346:ASN:OD1	1:B:347:LEU:HD23	2.21	0.40
1:C:47:ALA:HB2	1:C:373:TRP:CD2	2.57	0.40
1:C:244:VAL:HG13	2:C:501:NAD:C2A	2.52	0.40
1:A:427:PRO:HB2	1:A:429:HIS:CE1	2.57	0.40
1:C:58:GLN:H	1:C:58:GLN:CD	2.30	0.40
1:D:126:ASN:O	1:D:150:ILE:HB	2.22	0.40

There are no symmetry-related clashes.

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	
1	A	430/434 (99%)	418 (97%)	12 (3%)	0	100	100
1	B	430/434 (99%)	413 (96%)	15 (4%)	2 (0%)	25	44
1	C	431/434 (99%)	414 (96%)	15 (4%)	2 (0%)	25	44
1	D	429/434 (99%)	404 (94%)	21 (5%)	4 (1%)	14	28
All	All	1720/1736 (99%)	1649 (96%)	63 (4%)	8 (0%)	25	44

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	377	ASP
1	C	108	GLU
1	D	143	PHE
1	B	302	PHE
1	D	124	PRO
1	D	385	VAL
1	C	302	PHE
1	D	9	ALA

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	
1	A	361/362 (100%)	345 (96%)	16 (4%)	24	47
1	B	361/362 (100%)	353 (98%)	8 (2%)	47	73

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	362/362 (100%)	346 (96%)	16 (4%)	24	47
1	D	359/362 (99%)	334 (93%)	25 (7%)	12	26
All	All	1443/1448 (100%)	1378 (96%)	65 (4%)	24	46

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	SER
1	A	12	SER
1	A	36	LYS
1	A	57	VAL
1	A	73	VAL
1	A	94	THR
1	A	159	THR
1	A	173	LEU
1	A	191	ASN
1	A	273	VAL
1	A	316	LYS
1	A	319	VAL
1	A	322	LYS
1	A	377	ASP
1	A	381	VAL
1	A	415	THR
1	B	5	SER
1	B	80	ILE
1	B	157	THR
1	B	159	THR
1	B	166	LYS
1	B	357	VAL
1	B	400	GLU
1	B	412	LYS
1	C	8	VAL
1	C	27	GLU
1	C	142	LYS
1	C	159	THR
1	C	163	ASN
1	C	191	ASN
1	C	241	VAL
1	C	266	SER
1	C	267	LYS
1	C	279	ARG

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Mol	Chain	Res	Type
1	C	299	ILE
1	C	322	LYS
1	C	361	SER
1	C	381	VAL
1	C	385	VAL
1	C	397	LEU
1	D	18[A]	ARG
1	D	18[B]	ARG
1	D	22	ILE
1	D	49	ILE
1	D	57	VAL
1	D	82	SER
1	D	86	ASN
1	D	118	VAL
1	D	129	LEU
1	D	144	PRO
1	D	146	TYR
1	D	158	THR
1	D	180	VAL
1	D	187	SER
1	D	199	LEU
1	D	200	ILE
1	D	239	VAL
1	D	282	ILE
1	D	285	VAL
1	D	347	LEU
1	D	355	SER
1	D	357	VAL
1	D	389	ILE
1	D	396	SER
1	D	422	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	163	ASN
1	A	298	ASN
1	A	353	HIS
1	A	365	GLN
1	B	72	GLN
1	B	84	GLN
1	B	115	GLN

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Mol	Chain	Res	Type
1	B	123	GLN
1	B	162	HIS
1	B	320	ASN
1	B	324	GLN
1	B	331	GLN
1	B	335	HIS
1	B	351	HIS
1	B	360	ASN
1	B	365	GLN
1	B	369	GLN
1	C	115	GLN
1	C	126	ASN
1	C	230	GLN
1	C	298	ASN
1	C	324	GLN
1	C	331	GLN
1	C	360	ASN
1	D	54	HIS
1	D	74	GLN
1	D	84	GLN
1	D	181	ASN
1	D	314	ASN
1	D	331	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

7 ligands are modelled in this entry.

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
2	NAD	D	501	-	42,48,48	3.21	18 (42%)	50,73,73	1.63	10 (20%)
3	ADN	C	502	-	17,21,21	0.80	1 (5%)	17,31,31	0.80	1 (5%)
3	ADN	B	502	-	17,21,21	0.74	1 (5%)	17,31,31	1.05	1 (5%)
2	NAD	A	501	-	42,48,48	3.30	17 (40%)	50,73,73	1.53	8 (16%)
2	NAD	C	501	-	42,48,48	2.95	17 (40%)	50,73,73	1.60	10 (20%)
3	ADN	A	502	-	17,21,21	0.84	1 (5%)	17,31,31	0.76	0
2	NAD	B	501	-	42,48,48	3.28	17 (40%)	50,73,73	1.64	11 (22%)

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
2	NAD	D	501	-	-	8/26/62/62	0/5/5/5
3	ADN	C	502	-	-	2/2/22/22	0/3/3/3
3	ADN	B	502	-	-	2/2/22/22	0/3/3/3
2	NAD	A	501	-	-	8/26/62/62	0/5/5/5
2	NAD	C	501	-	-	10/26/62/62	0/5/5/5
3	ADN	A	502	-	-	2/2/22/22	0/3/3/3
2	NAD	B	501	-	-	8/26/62/62	0/5/5/5

All (72) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	NAD	O4B-C1B	10.48	1.54	1.40
2	D	501	NAD	O4B-C1B	10.20	1.54	1.40
2	B	501	NAD	O4B-C1B	9.67	1.53	1.40
2	B	501	NAD	C2N-N1N	7.62	1.43	1.35
2	C	501	NAD	O4D-C1D	7.43	1.50	1.40
2	B	501	NAD	O4D-C1D	7.40	1.50	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	NAD	O4D-C1D	7.32	1.50	1.40
2	C	501	NAD	C2N-N1N	7.19	1.42	1.35
2	C	501	NAD	C7N-N7N	6.99	1.45	1.33
2	D	501	NAD	C7N-N7N	6.91	1.45	1.33
2	B	501	NAD	C7N-N7N	6.84	1.45	1.33
2	A	501	NAD	C2N-N1N	6.84	1.42	1.35
2	A	501	NAD	C7N-N7N	6.81	1.45	1.33
2	D	501	NAD	C2N-N1N	6.63	1.42	1.35
2	D	501	NAD	O4D-C1D	6.07	1.48	1.40
2	D	501	NAD	C8A-N7A	5.69	1.45	1.34
2	A	501	NAD	C8A-N7A	5.59	1.44	1.34
2	C	501	NAD	C8A-N7A	5.52	1.44	1.34
2	B	501	NAD	C8A-N7A	5.50	1.44	1.34
2	A	501	NAD	PA-O3	5.27	1.65	1.59
2	B	501	NAD	PA-O3	5.12	1.65	1.59
2	A	501	NAD	PN-O3	5.11	1.65	1.59
2	D	501	NAD	PA-O3	5.01	1.64	1.59
2	D	501	NAD	PN-O3	4.96	1.64	1.59
2	B	501	NAD	PN-O3	4.93	1.64	1.59
2	C	501	NAD	PN-O3	4.73	1.64	1.59
2	C	501	NAD	PA-O3	4.67	1.64	1.59
2	C	501	NAD	O4B-C1B	4.52	1.46	1.40
2	A	501	NAD	C5A-N7A	4.25	1.55	1.39
2	D	501	NAD	C5A-N7A	4.23	1.55	1.39
2	C	501	NAD	C5A-N7A	4.23	1.55	1.39
2	B	501	NAD	C5A-N7A	4.18	1.54	1.39
2	A	501	NAD	C6A-N6A	3.23	1.45	1.34
2	D	501	NAD	C6A-N6A	3.22	1.45	1.34
2	C	501	NAD	C6A-N6A	3.19	1.45	1.34
2	B	501	NAD	C6A-N6A	3.17	1.45	1.34
2	B	501	NAD	C1B-N9A	-3.17	1.42	1.49
2	C	501	NAD	C1B-N9A	-3.15	1.42	1.49
2	D	501	NAD	C1B-N9A	-3.03	1.42	1.49
2	A	501	NAD	C1B-N9A	-3.00	1.42	1.49
2	D	501	NAD	O7N-C7N	-2.87	1.18	1.24
2	D	501	NAD	C6N-N1N	2.82	1.41	1.35
2	C	501	NAD	O7N-C7N	-2.78	1.18	1.24
2	A	501	NAD	O7N-C7N	-2.77	1.19	1.24
2	C	501	NAD	C6N-N1N	2.75	1.41	1.35
2	A	501	NAD	C6N-N1N	2.73	1.41	1.35
2	B	501	NAD	O7N-C7N	-2.68	1.19	1.24
2	B	501	NAD	C6N-N1N	2.60	1.41	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	ADN	C8-N7	-2.58	1.30	1.34
2	D	501	NAD	O4B-C4B	2.52	1.50	1.45
2	A	501	NAD	C4A-N3A	-2.51	1.32	1.35
2	B	501	NAD	C4A-N3A	-2.48	1.32	1.35
2	D	501	NAD	C4A-N3A	-2.48	1.32	1.35
2	B	501	NAD	O4B-C4B	2.48	1.50	1.45
2	D	501	NAD	C2B-C3B	-2.45	1.46	1.53
2	B	501	NAD	C2B-C3B	-2.42	1.46	1.53
2	A	501	NAD	O4B-C4B	2.40	1.50	1.45
3	C	502	ADN	C8-N7	-2.39	1.30	1.34
2	C	501	NAD	C4A-N3A	-2.39	1.32	1.35
2	A	501	NAD	C2B-C3B	-2.37	1.46	1.53
2	C	501	NAD	C2B-C3B	-2.35	1.47	1.53
2	C	501	NAD	C4N-C3N	2.33	1.42	1.39
2	D	501	NAD	C4N-C3N	2.32	1.42	1.39
2	D	501	NAD	C2D-C3D	-2.28	1.47	1.53
2	A	501	NAD	C4N-C3N	2.26	1.42	1.39
3	B	502	ADN	C8-N7	-2.16	1.30	1.34
2	C	501	NAD	C5N-C4N	2.14	1.42	1.38
2	C	501	NAD	C3N-C7N	2.10	1.53	1.50
2	D	501	NAD	C5N-C4N	2.08	1.42	1.38
2	B	501	NAD	C3N-C7N	2.08	1.53	1.50
2	B	501	NAD	C4N-C3N	2.06	1.42	1.39
2	A	501	NAD	C5N-C4N	2.02	1.42	1.38

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	NAD	C4B-O4B-C1B	-5.54	104.85	109.92
2	D	501	NAD	C4B-O4B-C1B	-5.05	105.30	109.92
2	C	501	NAD	C4B-O4B-C1B	-5.03	105.32	109.92
2	A	501	NAD	C4B-O4B-C1B	-4.69	105.63	109.92
2	C	501	NAD	C1B-N9A-C4A	-3.98	119.65	126.64
2	D	501	NAD	C1B-N9A-C4A	-3.92	119.76	126.64
2	A	501	NAD	C1B-N9A-C4A	-3.70	120.14	126.64
2	B	501	NAD	C6N-N1N-C2N	-3.61	118.80	121.88
2	A	501	NAD	C4A-C5A-N7A	-3.59	105.54	109.34
2	D	501	NAD	C4A-C5A-N7A	-3.57	105.56	109.34
2	B	501	NAD	C1B-N9A-C4A	-3.44	120.61	126.64
2	C	501	NAD	C4A-C5A-N7A	-3.28	105.88	109.34
2	B	501	NAD	O4B-C1B-N9A	3.09	112.84	108.75
2	B	501	NAD	C4A-C5A-N7A	-3.09	106.07	109.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	NAD	C6N-N1N-C2N	-3.07	119.27	121.88
2	C	501	NAD	C6N-N1N-C2N	-3.04	119.29	121.88
2	A	501	NAD	C4D-O4D-C1D	-3.01	107.17	109.92
2	A	501	NAD	C6N-N1N-C2N	-2.80	119.50	121.88
2	D	501	NAD	C4D-O4D-C1D	-2.73	107.42	109.92
2	D	501	NAD	C2B-C3B-C4B	2.69	107.80	102.61
2	C	501	NAD	C4D-O4D-C1D	-2.67	107.48	109.92
2	A	501	NAD	C2B-C3B-C4B	2.65	107.74	102.61
2	B	501	NAD	C2B-C3B-C4B	2.62	107.68	102.61
2	D	501	NAD	C2D-C3D-C4D	2.57	107.58	102.61
2	C	501	NAD	C3N-C7N-N7N	2.56	120.89	117.74
2	C	501	NAD	C2B-C3B-C4B	2.48	107.39	102.61
3	B	502	ADN	C5-C6-N6	2.46	124.06	120.31
2	A	501	NAD	O2A-PA-O1A	-2.45	101.03	112.44
2	B	501	NAD	C3N-C7N-N7N	2.45	120.76	117.74
2	B	501	NAD	O2A-PA-O1A	-2.42	101.17	112.44
2	C	501	NAD	O2N-PN-O1N	-2.36	101.45	112.44
2	C	501	NAD	O7N-C7N-N7N	-2.32	119.26	122.62
2	D	501	NAD	O2A-PA-O1A	-2.31	101.72	112.44
2	A	501	NAD	O2N-PN-O1N	-2.30	101.75	112.44
2	D	501	NAD	O2N-PN-O1N	-2.28	101.83	112.44
2	B	501	NAD	C4D-O4D-C1D	-2.21	107.90	109.92
2	B	501	NAD	O2N-PN-O1N	-2.18	102.29	112.44
2	B	501	NAD	O7N-C7N-N7N	-2.17	119.48	122.62
2	C	501	NAD	O2A-PA-O1A	-2.15	102.45	112.44
2	D	501	NAD	C5D-C4D-C3D	-2.08	107.74	115.21
3	C	502	ADN	C5-C6-N6	2.05	123.44	120.31

There are no chirality outliers.

All (40) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	NAD	O4D-C1D-N1N-C2N
2	A	501	NAD	O4D-C1D-N1N-C6N
2	A	501	NAD	C2D-C1D-N1N-C2N
2	A	501	NAD	C2D-C1D-N1N-C6N
2	B	501	NAD	O4D-C1D-N1N-C2N
2	B	501	NAD	O4D-C1D-N1N-C6N
2	B	501	NAD	C2D-C1D-N1N-C2N
2	B	501	NAD	C2D-C1D-N1N-C6N
2	C	501	NAD	C5D-O5D-PN-O2N
2	C	501	NAD	O4D-C1D-N1N-C2N

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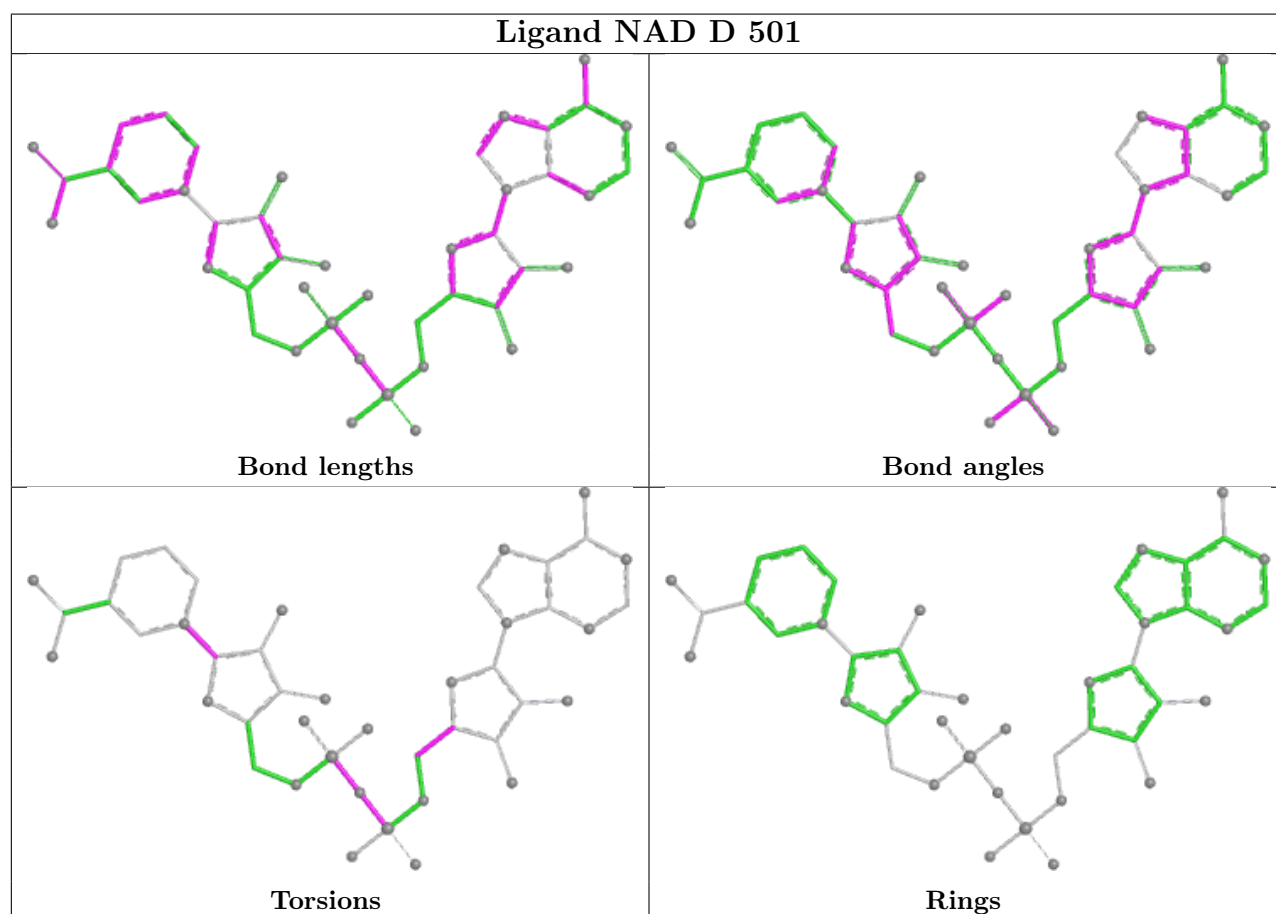
Mol	Chain	Res	Type	Atoms
2	C	501	NAD	O4D-C1D-N1N-C6N
2	C	501	NAD	C2D-C1D-N1N-C2N
2	C	501	NAD	C2D-C1D-N1N-C6N
2	D	501	NAD	O4D-C1D-N1N-C2N
2	D	501	NAD	O4D-C1D-N1N-C6N
2	D	501	NAD	C2D-C1D-N1N-C2N
2	D	501	NAD	C2D-C1D-N1N-C6N
3	B	502	ADN	O4'-C4'-C5'-O5'
3	B	502	ADN	C3'-C4'-C5'-O5'
3	C	502	ADN	C3'-C4'-C5'-O5'
2	C	501	NAD	O4D-C4D-C5D-O5D
2	B	501	NAD	O4B-C4B-C5B-O5B
2	B	501	NAD	C3B-C4B-C5B-O5B
3	C	502	ADN	O4'-C4'-C5'-O5'
2	D	501	NAD	O4B-C4B-C5B-O5B
3	A	502	ADN	C3'-C4'-C5'-O5'
2	A	501	NAD	O4B-C4B-C5B-O5B
2	C	501	NAD	C3D-C4D-C5D-O5D
2	D	501	NAD	C3B-C4B-C5B-O5B
2	A	501	NAD	C3B-C4B-C5B-O5B
2	C	501	NAD	O4B-C4B-C5B-O5B
2	C	501	NAD	C3B-C4B-C5B-O5B
2	D	501	NAD	PN-O3-PA-O5B
2	A	501	NAD	C5B-O5B-PA-O2A
2	B	501	NAD	C5B-O5B-PA-O2A
2	C	501	NAD	C5D-O5D-PN-O3
2	D	501	NAD	PA-O3-PN-O2N
3	A	502	ADN	O4'-C4'-C5'-O5'
2	A	501	NAD	PA-O3-PN-O2N
2	B	501	NAD	PA-O3-PN-O2N

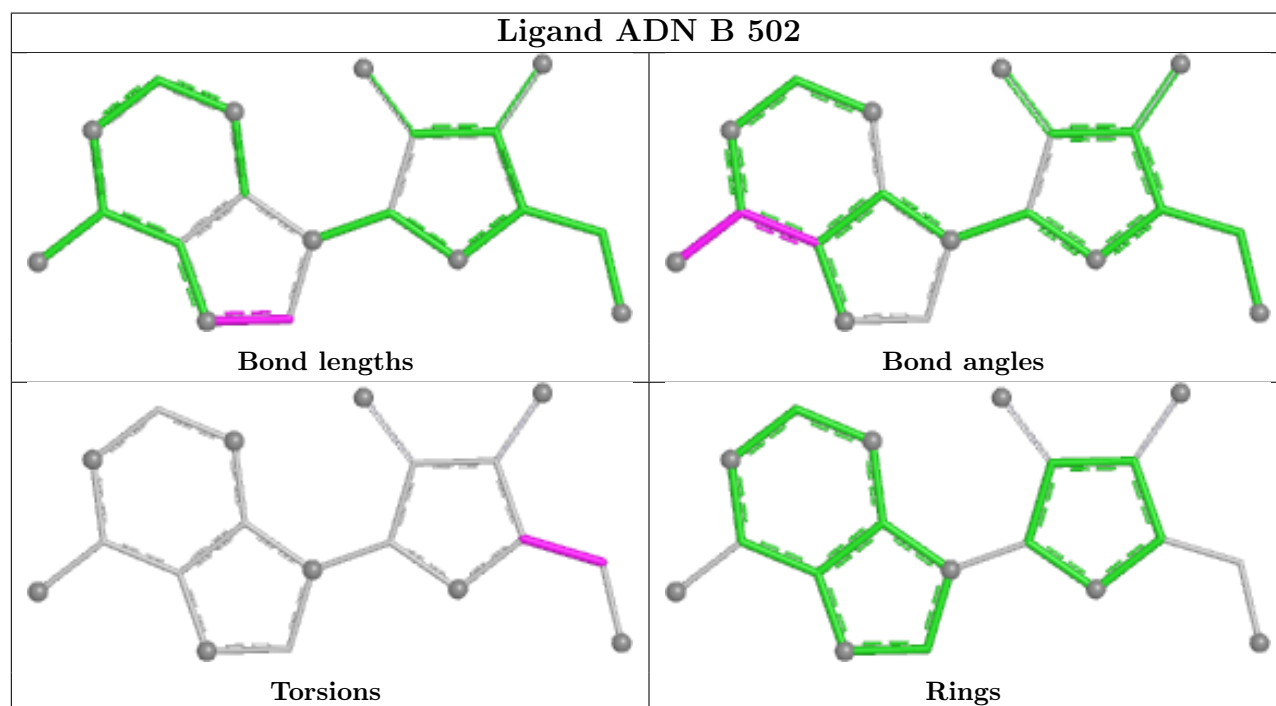
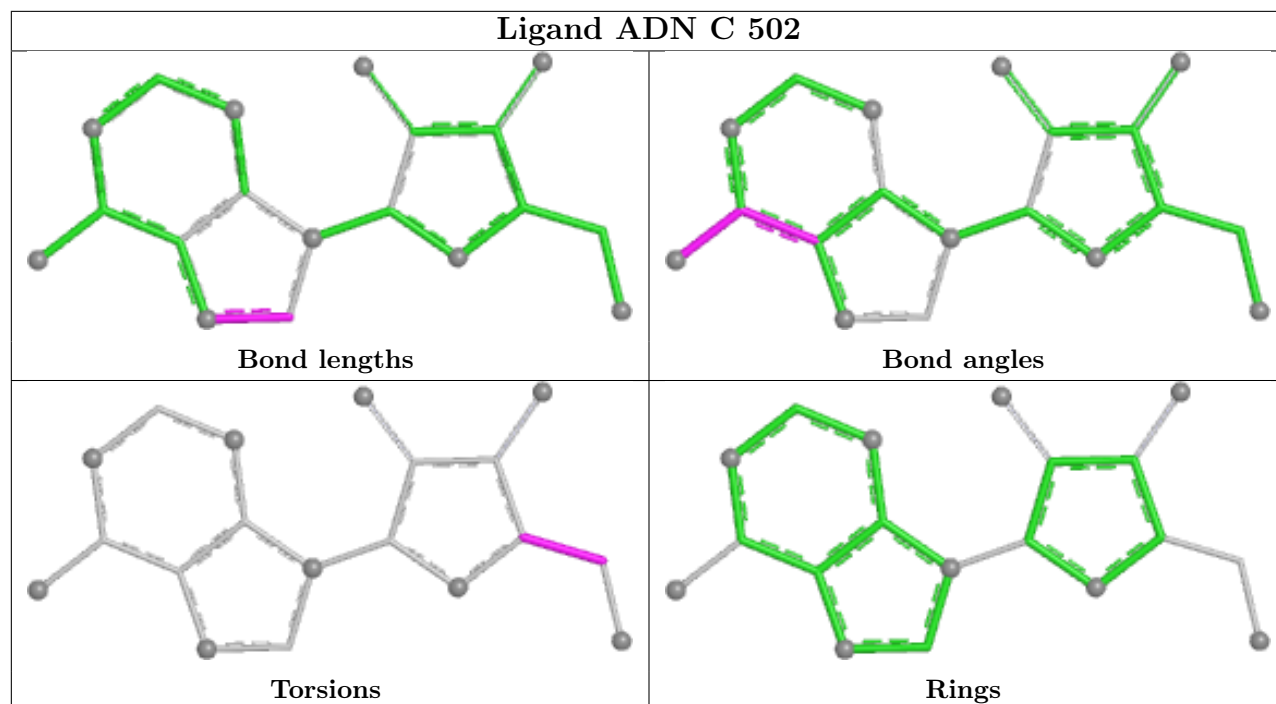
There are no ring outliers.

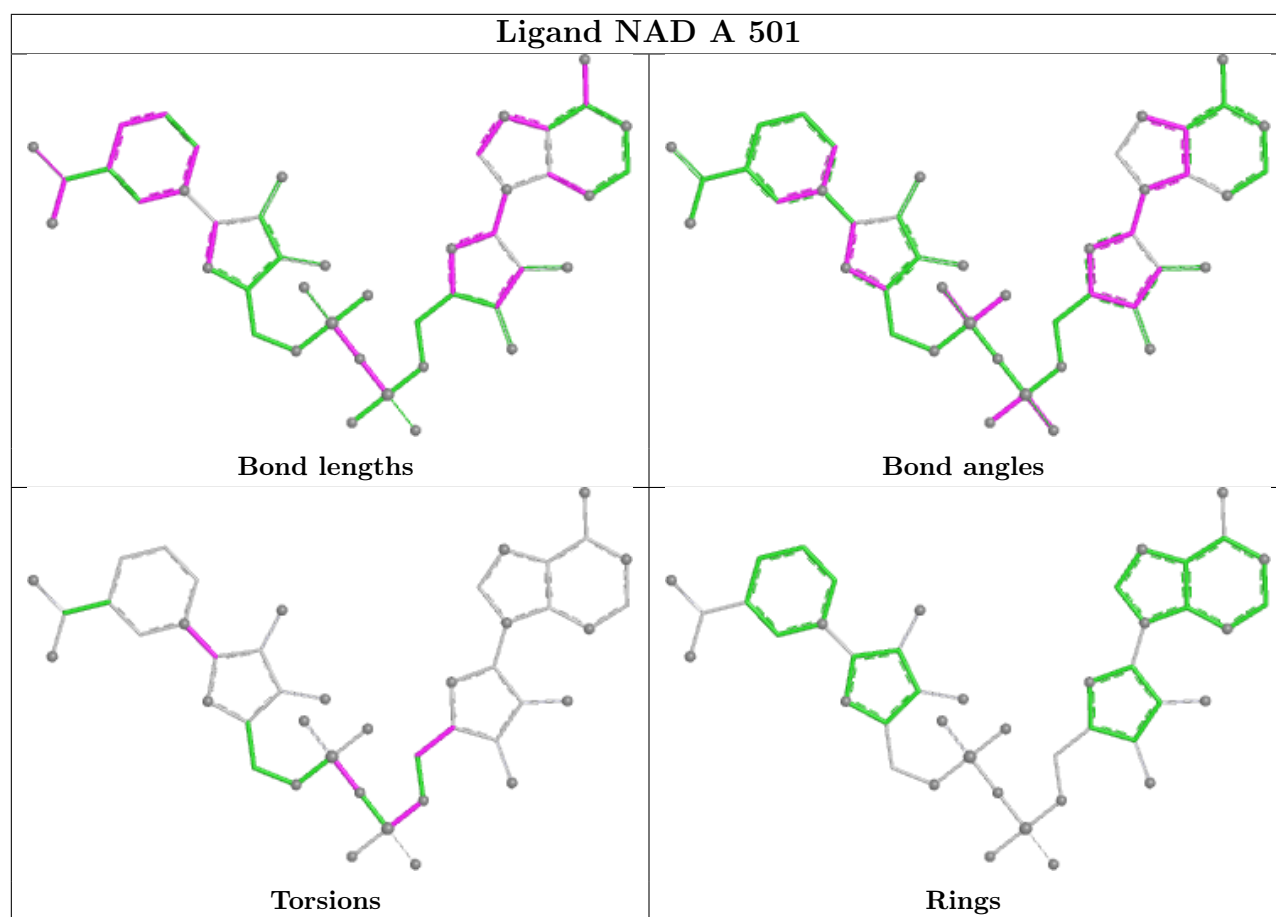
7 monomers are involved in 11 short contacts:

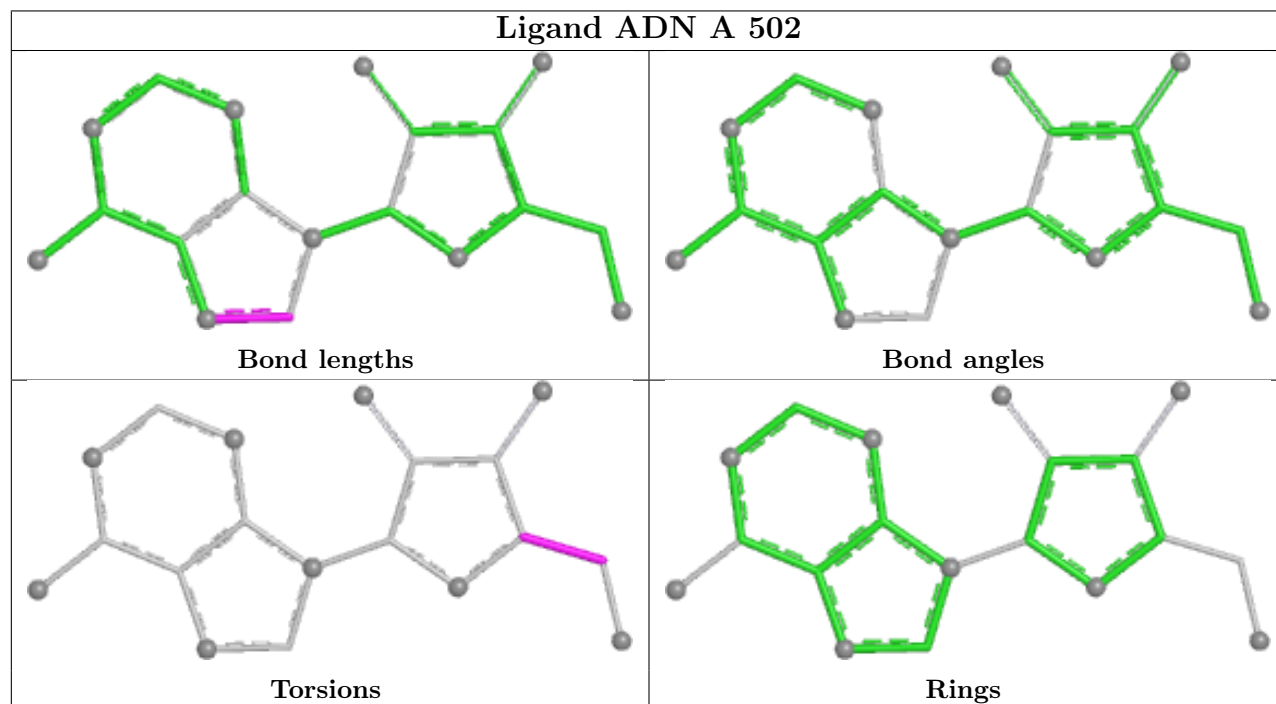
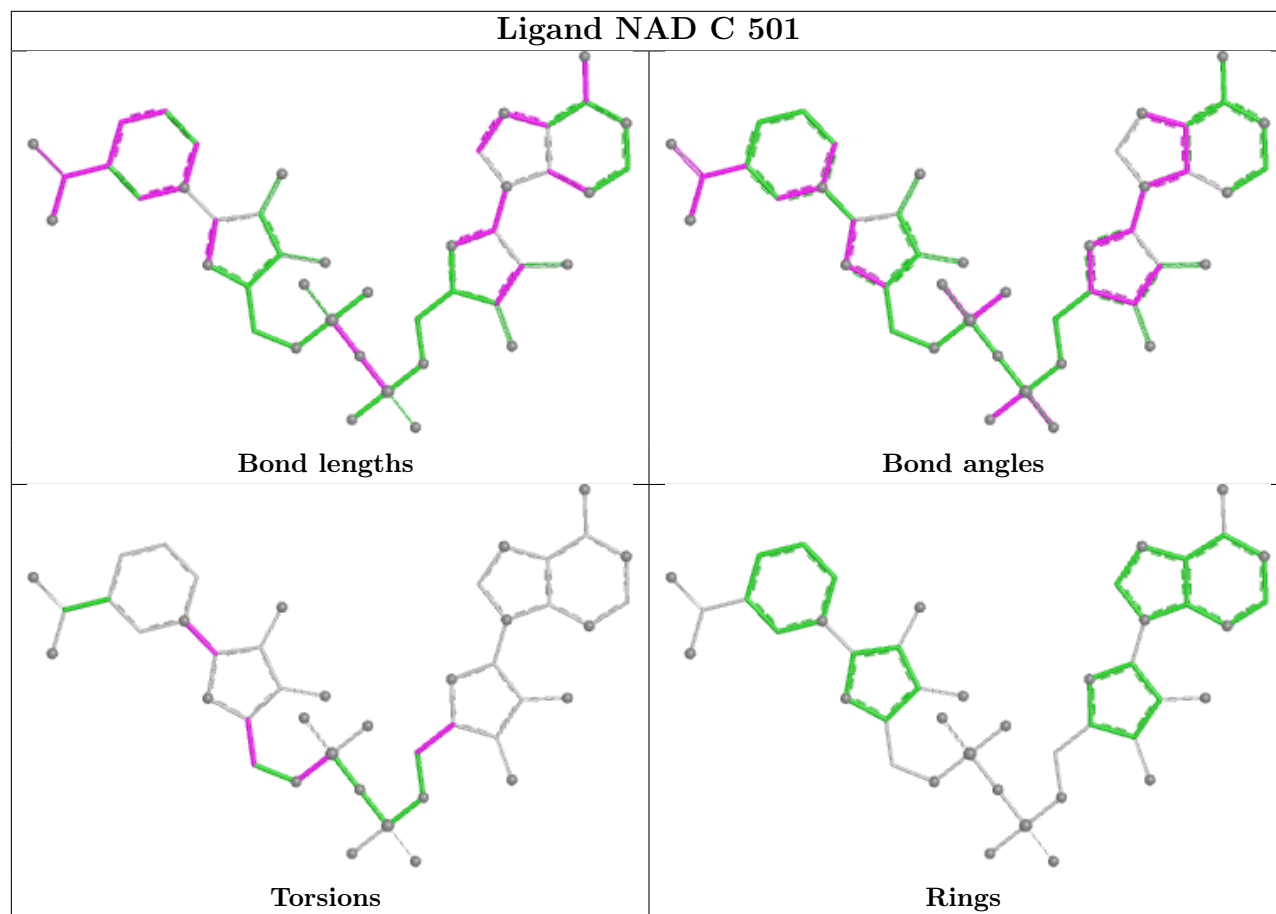
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	NAD	3	0
3	C	502	ADN	1	0
3	B	502	ADN	1	0
2	A	501	NAD	1	0
2	C	501	NAD	2	0
3	A	502	ADN	1	0
2	B	501	NAD	3	0

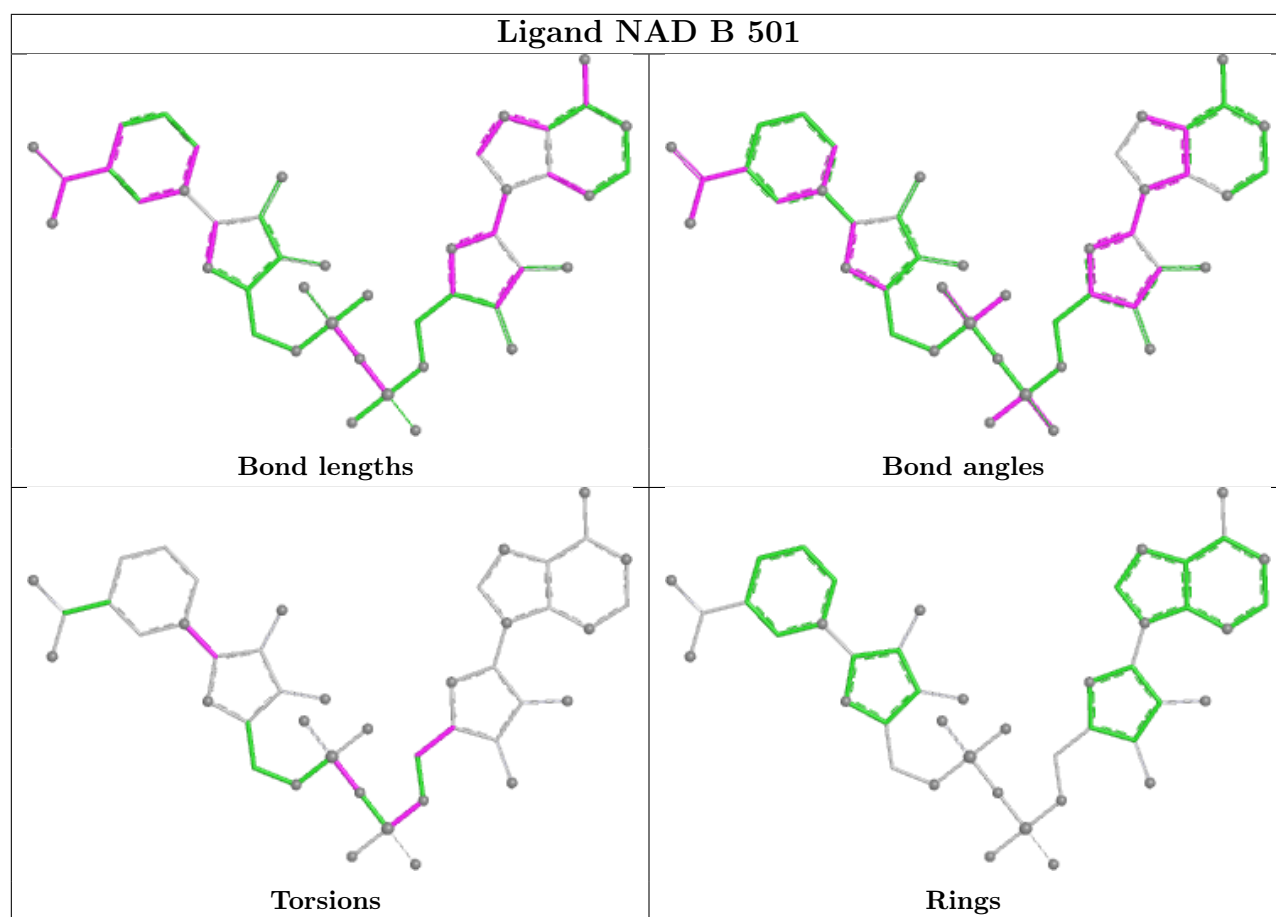
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	A	431/434 (99%)	0.14	2 (0%) 87 85	14, 25, 42, 77	1 (0%)
1	B	432/434 (99%)	0.10	1 (0%) 92 90	15, 25, 39, 80	0
1	C	433/434 (99%)	0.22	5 (1%) 76 73	16, 28, 46, 73	0
1	D	430/434 (99%)	0.84	61 (14%) 7 7	15, 39, 75, 95	1 (0%)
All	All	1726/1736 (99%)	0.32	69 (3%) 43 39	14, 28, 60, 95	2 (0%)

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	118	VAL	4.8
1	D	173	LEU	4.8
1	D	157	THR	4.5
1	D	168	PHE	4.5
1	D	132	GLY	4.5
1	D	143	PHE	4.1
1	D	373	TRP	4.0
1	D	146	TYR	3.8
1	D	18[A]	ARG	3.7
1	D	82	SER	3.6
1	D	175	VAL	3.5
1	D	136	THR	3.5
1	C	380	ALA	3.3
1	D	302	PHE	3.3
1	A	2	SER	3.2
1	D	174	GLY	3.1
1	D	98	VAL	3.1
1	D	189	PHE	3.1
1	D	106	ASP	3.0
1	D	54	HIS	3.0
1	D	301	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	70	GLY	2.9
1	C	107	GLU	2.8
1	D	109	TYR	2.8
1	D	140	HIS	2.7
1	D	122	GLY	2.7
1	D	376	SER	2.7
1	D	383	VAL	2.7
1	C	1	MET	2.6
1	D	164	LEU	2.6
1	D	380	ALA	2.5
1	D	165	TYR	2.5
1	D	171	GLY	2.5
1	C	0	HIS	2.5
1	D	172	ARG	2.5
1	D	75	TRP	2.5
1	D	144	PRO	2.5
1	A	380	ALA	2.5
1	D	158	THR	2.5
1	D	162	HIS	2.4
1	D	124	PRO	2.4
1	D	9	ALA	2.4
1	D	16	TRP	2.4
1	D	384	HIS	2.3
1	D	6	TYR	2.3
1	B	1	MET	2.3
1	D	117	LEU	2.3
1	D	374	THR	2.3
1	D	134	ASP	2.3
1	D	120	PRO	2.3
1	D	371	GLU	2.3
1	D	381	VAL	2.3
1	D	53	LEU	2.2
1	D	149	ASN	2.2
1	D	80	ILE	2.2
1	D	150	ILE	2.2
1	D	4	PRO	2.2
1	D	163	ASN	2.2
1	D	116	THR	2.2
1	D	52	CYS	2.1
1	D	170	GLU	2.1
1	D	379	TYR	2.1
1	D	74	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	151	LYS	2.1
1	D	121	ASP	2.1
1	D	110	MET	2.1
1	D	419	VAL	2.0
1	C	377	ASP	2.0
1	D	161	VAL	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

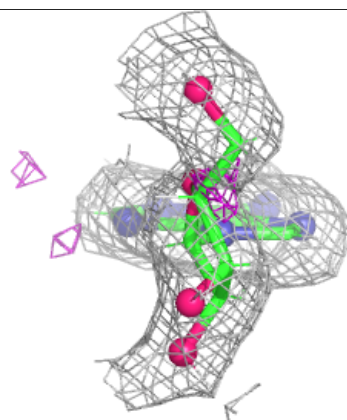
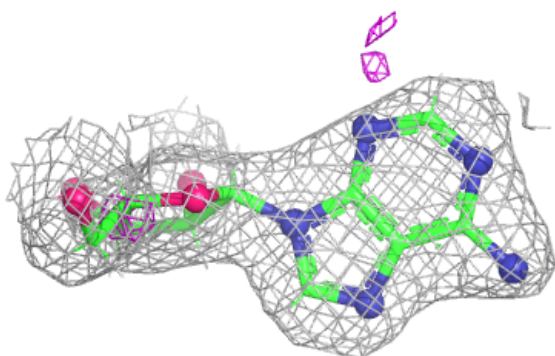
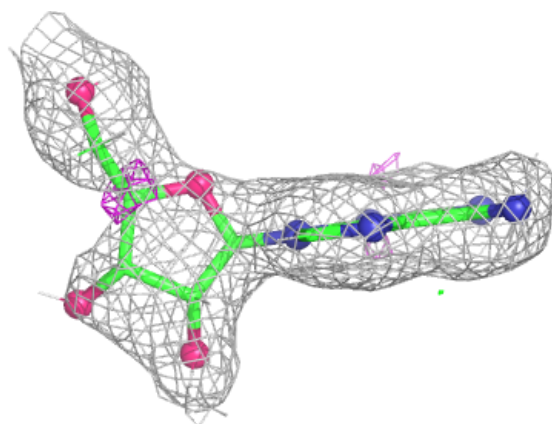
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	ADN	A	502	19/19	0.88	0.11	21,30,39,46	0
3	ADN	B	502	19/19	0.88	0.12	14,32,44,44	0
3	ADN	C	502	19/19	0.91	0.09	20,31,46,46	0
2	NAD	D	501	44/44	0.94	0.09	20,29,49,57	0
2	NAD	B	501	44/44	0.94	0.09	17,26,31,34	0
2	NAD	A	501	44/44	0.95	0.08	14,24,31,32	0
2	NAD	C	501	44/44	0.95	0.08	17,23,35,37	0

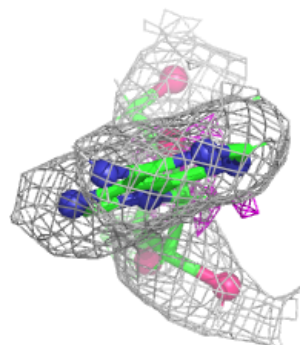
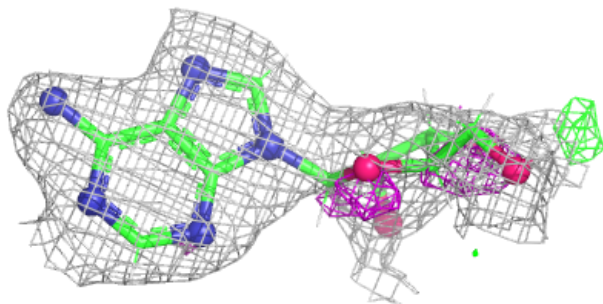
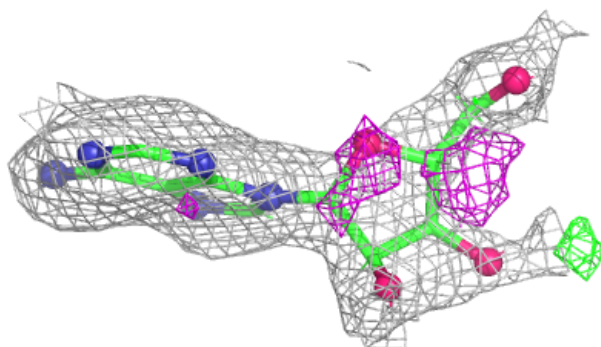
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADN A 502:

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

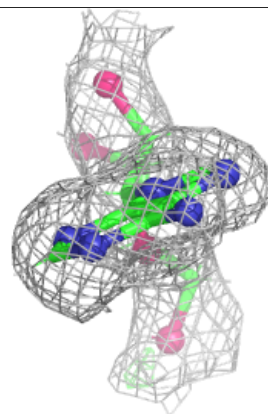
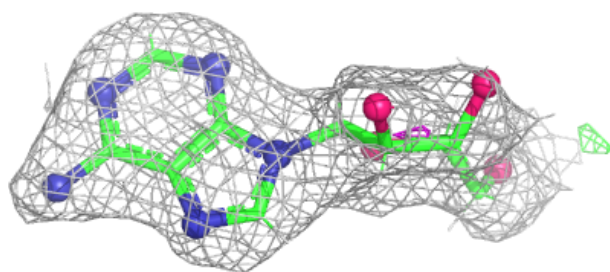
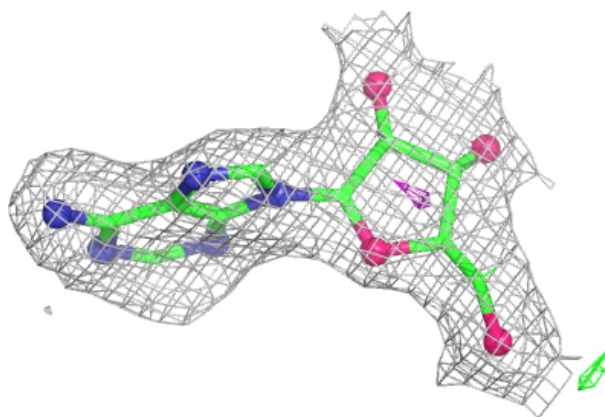
**Electron density around ADN B 502:**

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

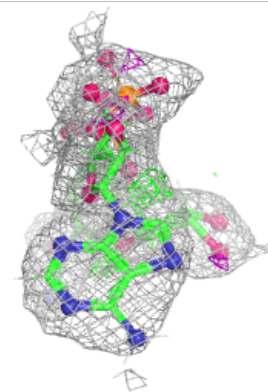
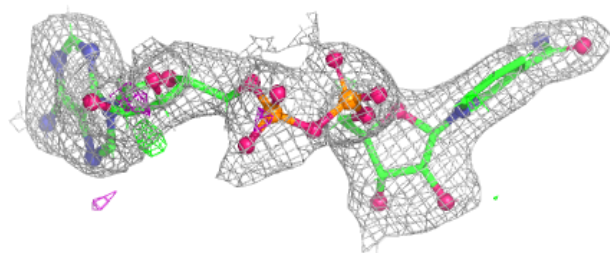
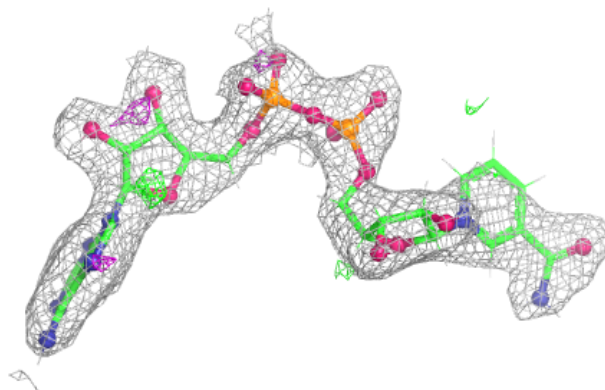


Electron density around ADN C 502:

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

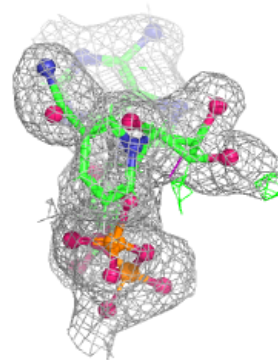
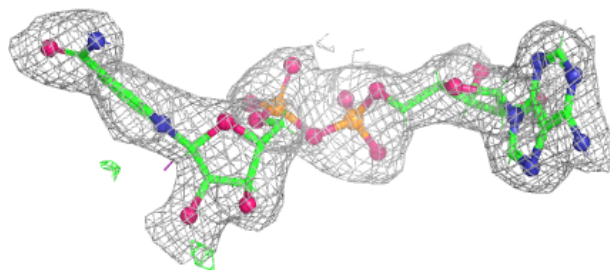
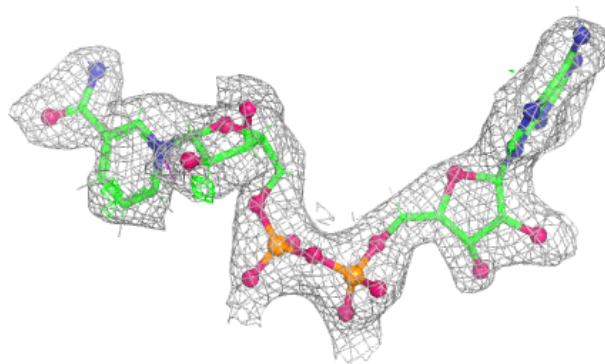
**Electron density around NAD D 501:**

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

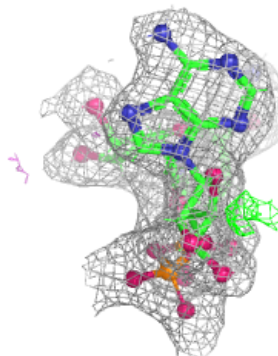
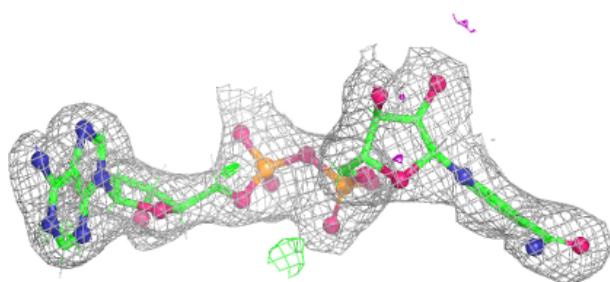
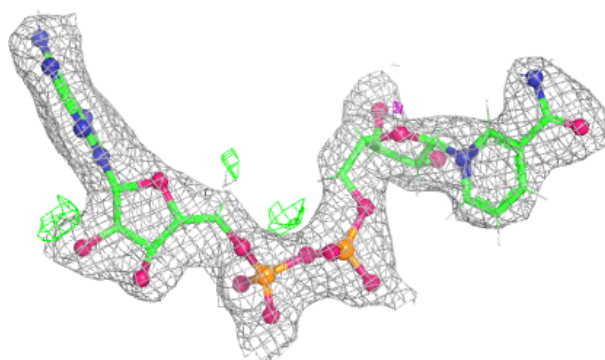


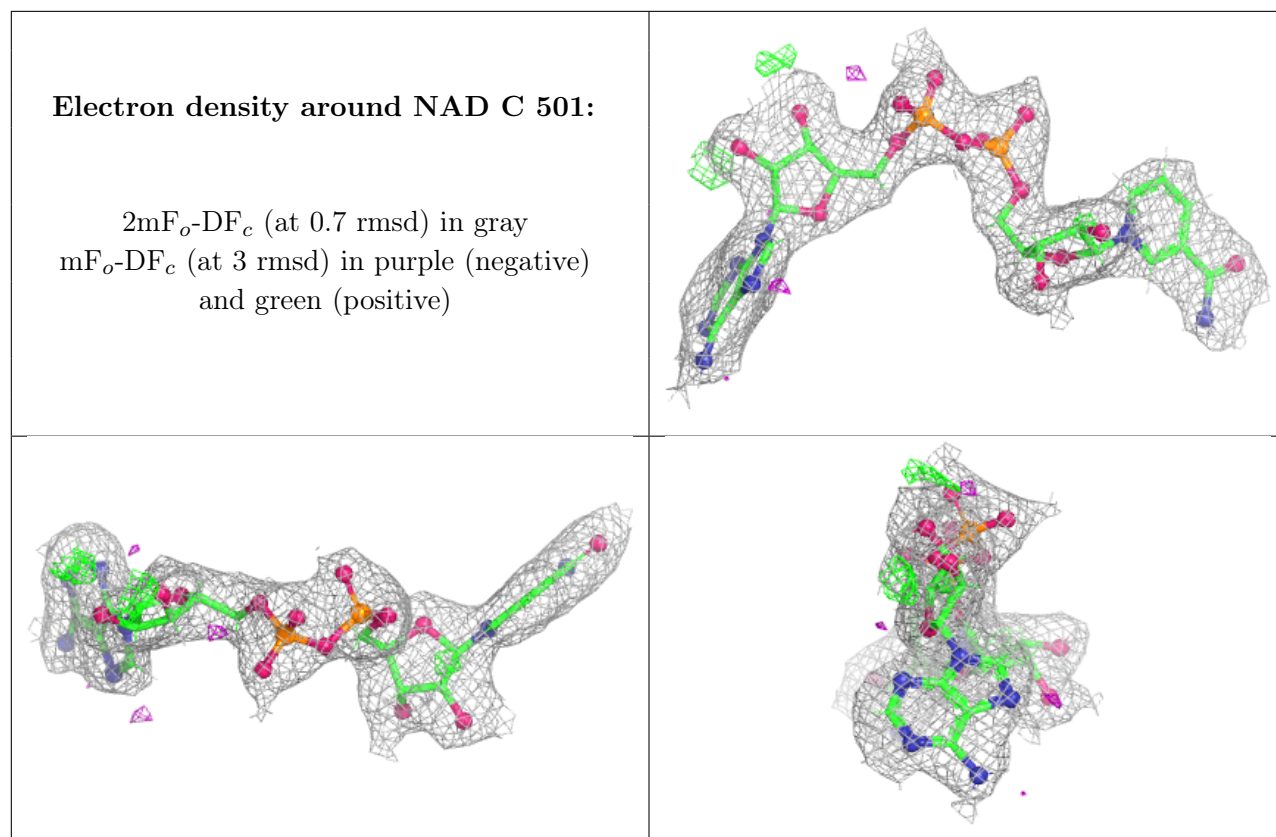
Electron density around NAD B 501:

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

**Electron density around NAD A 501:**

$2mF_o-DF_c$ (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.