



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 18, 2024 – 11:29 AM EDT

PDB ID : 4A5A
Title : Crystal structure of the C258S/C268S variant of Toxoplasma gondii nucleoside triphosphate diphosphohydrolase 3 (NTPDase3) in complex with magnesium and AMPPNP
Authors : Krug, U.; Zebisch, M.; Straeter, N.
Deposited on : 2011-10-24
Resolution : 2.85 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	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

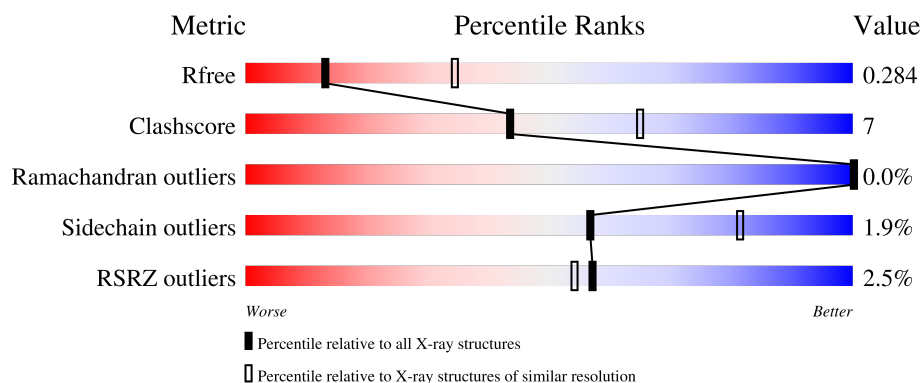
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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}	130704	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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	611	<div> <div>4%</div> <div>80%</div> <div>16%</div> <div>..</div> </div>
1	B	611	<div> <div>3%</div> <div>78%</div> <div>16%</div> <div>..</div> </div>
1	C	611	<div> <div>%</div> <div>78%</div> <div>17%</div> <div>..</div> </div>
1	D	611	<div> <div>%</div> <div>79%</div> <div>17%</div> <div>..</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 18420 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 protein called NUCLEOSIDE-TRIPHOSPHATASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	591	Total	C	N	O	S	0	0	0
			4583	2882	810	866	25			
1	B	585	Total	C	N	O	S	0	0	0
			4531	2851	799	856	25			
1	C	592	Total	C	N	O	S	0	0	0
			4587	2885	807	870	25			
1	D	592	Total	C	N	O	S	0	0	0
			4591	2886	811	869	25			

There are 40 discrepancies between the modelled and reference sequences:

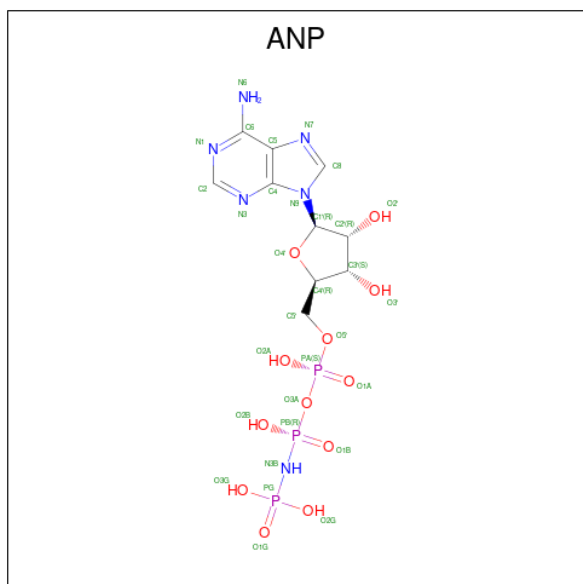
Chain	Residue	Modelled	Actual	Comment	Reference
A	25	MET	-	expression tag	UNP Q27893
A	629	GLU	-	expression tag	UNP Q27893
A	630	HIS	-	expression tag	UNP Q27893
A	631	HIS	-	expression tag	UNP Q27893
A	632	HIS	-	expression tag	UNP Q27893
A	633	HIS	-	expression tag	UNP Q27893
A	634	HIS	-	expression tag	UNP Q27893
A	635	HIS	-	expression tag	UNP Q27893
A	258	SER	CYS	engineered mutation	UNP Q27893
A	268	SER	CYS	engineered mutation	UNP Q27893
B	25	MET	-	expression tag	UNP Q27893
B	629	GLU	-	expression tag	UNP Q27893
B	630	HIS	-	expression tag	UNP Q27893
B	631	HIS	-	expression tag	UNP Q27893
B	632	HIS	-	expression tag	UNP Q27893
B	633	HIS	-	expression tag	UNP Q27893
B	634	HIS	-	expression tag	UNP Q27893
B	635	HIS	-	expression tag	UNP Q27893
B	258	SER	CYS	engineered mutation	UNP Q27893
B	268	SER	CYS	engineered mutation	UNP Q27893
C	25	MET	-	expression tag	UNP Q27893

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Chain	Residue	Modelled	Actual	Comment	Reference
C	629	GLU	-	expression tag	UNP Q27893
C	630	HIS	-	expression tag	UNP Q27893
C	631	HIS	-	expression tag	UNP Q27893
C	632	HIS	-	expression tag	UNP Q27893
C	633	HIS	-	expression tag	UNP Q27893
C	634	HIS	-	expression tag	UNP Q27893
C	635	HIS	-	expression tag	UNP Q27893
C	258	SER	CYS	engineered mutation	UNP Q27893
C	268	SER	CYS	engineered mutation	UNP Q27893
D	25	MET	-	expression tag	UNP Q27893
D	629	GLU	-	expression tag	UNP Q27893
D	630	HIS	-	expression tag	UNP Q27893
D	631	HIS	-	expression tag	UNP Q27893
D	632	HIS	-	expression tag	UNP Q27893
D	633	HIS	-	expression tag	UNP Q27893
D	634	HIS	-	expression tag	UNP Q27893
D	635	HIS	-	expression tag	UNP Q27893
D	258	SER	CYS	engineered mutation	UNP Q27893
D	268	SER	CYS	engineered mutation	UNP Q27893

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: $C_{10}H_{17}N_6O_{12}P_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	
			31	10	6	12	3	

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
2	C	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
2	D	1	Total	C	N	O	P	0	0
			31	10	6	12	3		

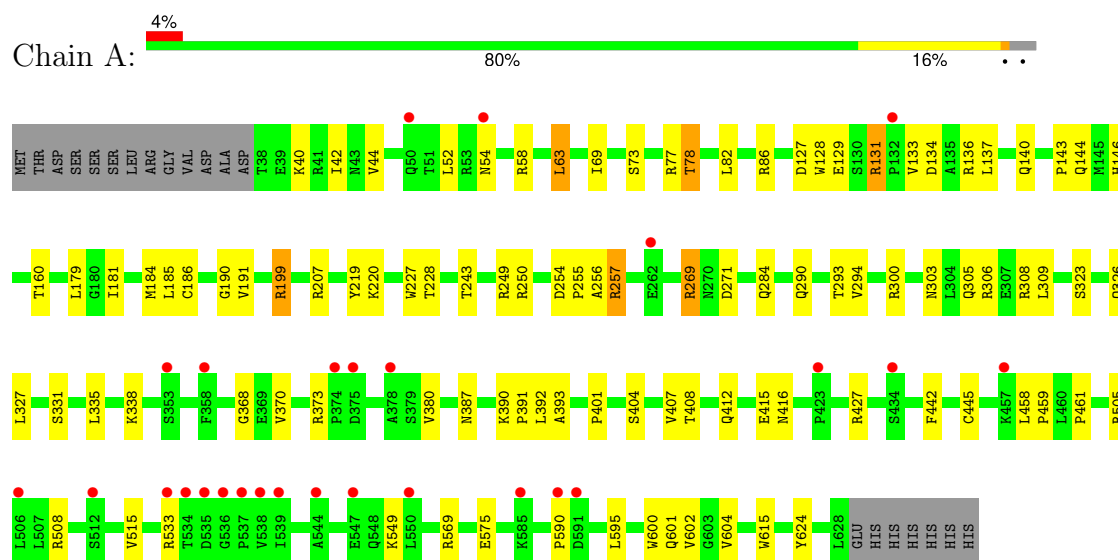
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		

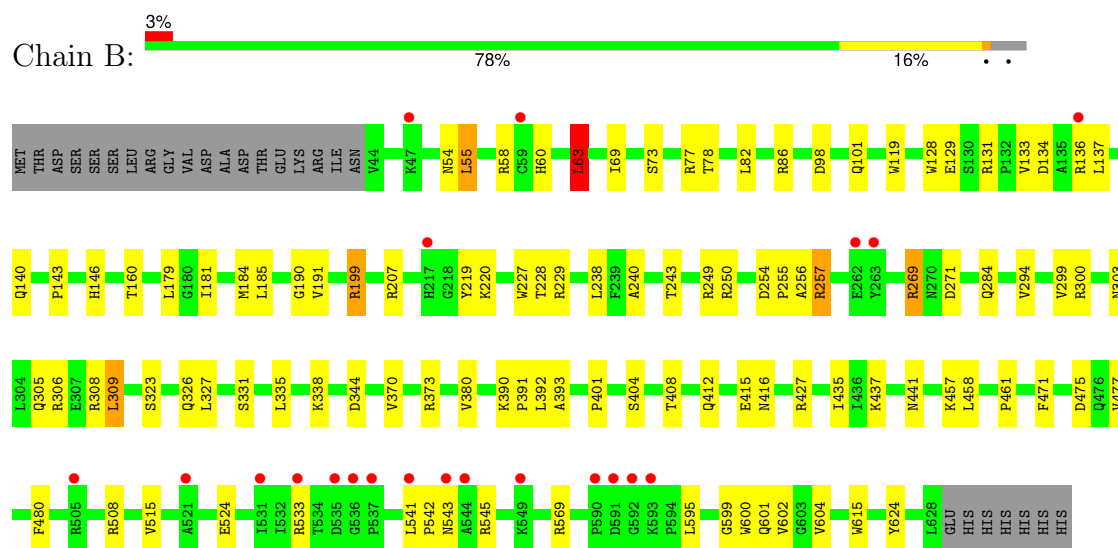
3 Residue-property plots [i](#)

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.

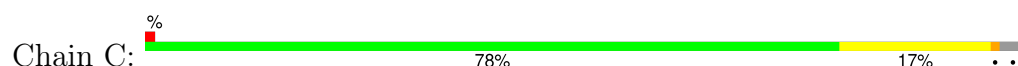
• Molecule 1: NUCLEOSIDE-TRIPHOSPHATASE 1

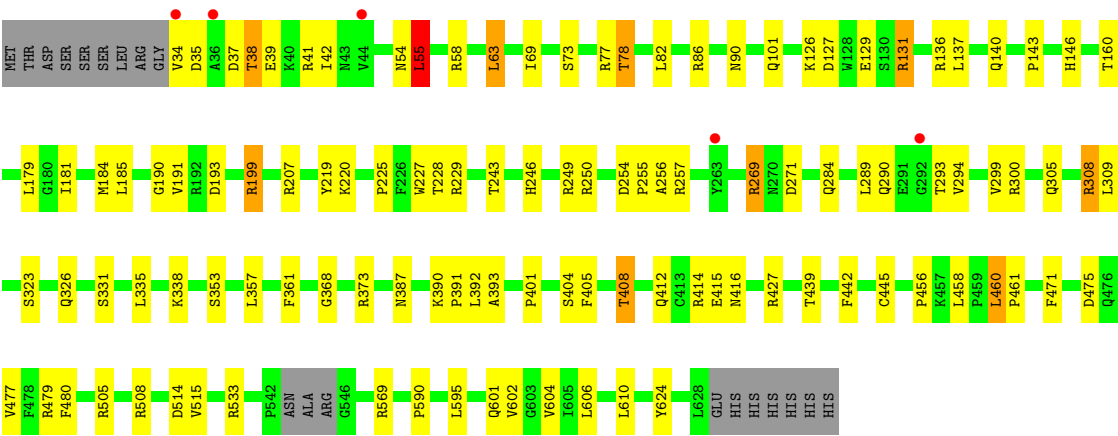


• Molecule 1: NUCLEOSIDE-TRIPHOSPHATASE 1

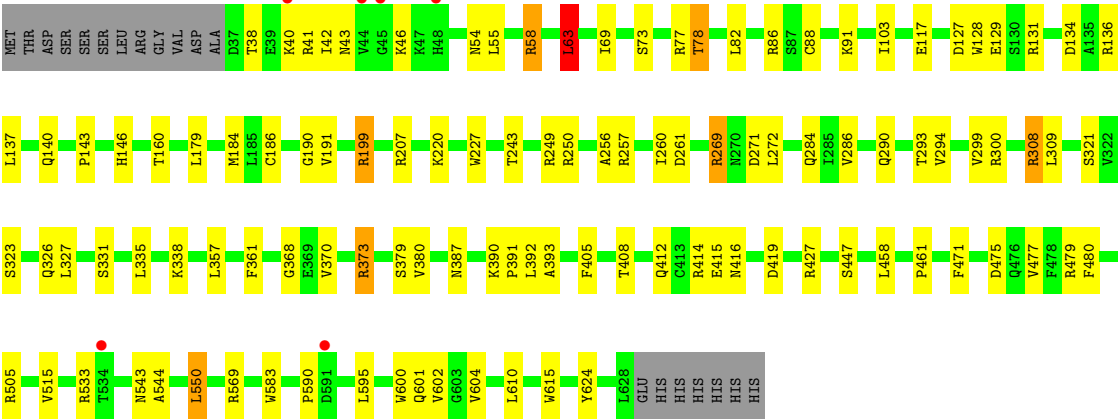
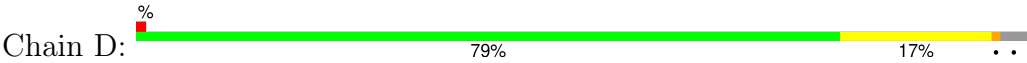


• Molecule 1: NUCLEOSIDE-TRIPHOSPHATASE 1





● Molecule 1: NUCLEOSIDE-TRIPHOSPHATASE 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	71.95Å 150.67Å 486.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.85 38.96 – 2.85	Depositor EDS
% Data completeness (in resolution range)	97.8 (40.00-2.85) 97.8 (38.96-2.85)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 2.86Å)	Xtriage
Refinement program	REFMAC 5.6.0116	Depositor
R, R_{free}	0.234 , 0.284 0.233 , 0.284	Depositor DCC
R_{free} test set	3124 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	36.3	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 26.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	18420	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.46% 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

5.1 Standard geometry

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

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	0.54	2/4672 (0.0%)	0.84	24/6323 (0.4%)
1	B	0.55	4/4620 (0.1%)	0.94	27/6254 (0.4%)
1	C	0.66	2/4675 (0.0%)	1.03	32/6327 (0.5%)
1	D	0.64	5/4680 (0.1%)	0.97	30/6334 (0.5%)
All	All	0.60	13/18647 (0.1%)	0.95	113/25238 (0.4%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	58	ARG	CZ-NH1	-10.98	1.18	1.33
1	C	229	ARG	CZ-NH2	-9.00	1.21	1.33
1	C	229	ARG	CZ-NH1	-8.23	1.22	1.33
1	B	128	TRP	CD2-CE2	5.68	1.48	1.41
1	D	583	TRP	CD2-CE2	5.42	1.47	1.41
1	D	615	TRP	CD2-CE2	5.39	1.47	1.41
1	B	615	TRP	CD2-CE2	5.38	1.47	1.41
1	D	128	TRP	CD2-CE2	5.19	1.47	1.41
1	A	128	TRP	CD2-CE2	5.14	1.47	1.41
1	D	58	ARG	CZ-NH2	-5.12	1.26	1.33
1	B	119	TRP	CD2-CE2	5.09	1.47	1.41
1	B	600	TRP	CD2-CE2	5.05	1.47	1.41
1	A	615	TRP	CD2-CE2	5.00	1.47	1.41

All (113) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	269	ARG	NE-CZ-NH2	-18.67	110.97	120.30
1	C	249	ARG	NE-CZ-NH2	-17.98	111.31	120.30
1	B	207	ARG	NE-CZ-NH2	-17.64	111.48	120.30
1	C	414	ARG	NE-CZ-NH1	17.22	128.91	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	249	ARG	NE-CZ-NH1	16.88	128.74	120.30
1	C	414	ARG	NE-CZ-NH2	-16.60	112.00	120.30
1	D	58	ARG	NE-CZ-NH2	16.14	128.37	120.30
1	D	373	ARG	NE-CZ-NH2	-16.09	112.25	120.30
1	B	207	ARG	NE-CZ-NH1	15.91	128.26	120.30
1	D	373	ARG	NE-CZ-NH1	15.84	128.22	120.30
1	C	229	ARG	NE-CZ-NH1	15.50	128.05	120.30
1	B	533	ARG	NE-CZ-NH2	-15.05	112.78	120.30
1	D	269	ARG	NE-CZ-NH1	14.85	127.72	120.30
1	B	533	ARG	NE-CZ-NH1	14.67	127.64	120.30
1	C	505	ARG	NE-CZ-NH2	-13.96	113.32	120.30
1	B	131	ARG	NE-CZ-NH2	-13.95	113.33	120.30
1	A	136	ARG	NE-CZ-NH1	13.37	126.98	120.30
1	C	505	ARG	NE-CZ-NH1	13.36	126.98	120.30
1	D	86	ARG	NE-CZ-NH2	-12.99	113.81	120.30
1	A	136	ARG	NE-CZ-NH2	-12.87	113.87	120.30
1	B	131	ARG	NE-CZ-NH1	12.62	126.61	120.30
1	B	86	ARG	NE-CZ-NH2	-12.41	114.10	120.30
1	B	86	ARG	NE-CZ-NH1	12.01	126.30	120.30
1	D	86	ARG	NE-CZ-NH1	11.79	126.20	120.30
1	A	308	ARG	NE-CZ-NH2	-11.41	114.60	120.30
1	A	308	ARG	NE-CZ-NH1	11.23	125.92	120.30
1	C	229	ARG	NH1-CZ-NH2	-10.78	107.54	119.40
1	D	308	ARG	NE-CZ-NH2	-10.70	114.95	120.30
1	D	308	ARG	NE-CZ-NH1	10.57	125.59	120.30
1	C	308	ARG	NE-CZ-NH2	-10.21	115.19	120.30
1	A	86	ARG	NE-CZ-NH1	10.03	125.31	120.30
1	C	308	ARG	NE-CZ-NH1	9.84	125.22	120.30
1	B	308	ARG	NE-CZ-NH2	-9.79	115.40	120.30
1	A	86	ARG	NE-CZ-NH2	-9.51	115.55	120.30
1	B	257	ARG	NE-CZ-NH1	9.35	124.98	120.30
1	A	199	ARG	NE-CZ-NH2	-9.33	115.64	120.30
1	B	308	ARG	NE-CZ-NH1	9.30	124.95	120.30
1	C	199	ARG	NE-CZ-NH1	9.22	124.91	120.30
1	B	257	ARG	NE-CZ-NH2	-9.09	115.76	120.30
1	C	249	ARG	CD-NE-CZ	8.84	135.98	123.60
1	B	199	ARG	NE-CZ-NH2	-8.77	115.92	120.30
1	C	199	ARG	NE-CZ-NH2	-8.76	115.92	120.30
1	D	199	ARG	NE-CZ-NH1	8.75	124.67	120.30
1	A	199	ARG	NE-CZ-NH1	8.69	124.64	120.30
1	A	257	ARG	NE-CZ-NH1	8.44	124.52	120.30
1	D	199	ARG	NE-CZ-NH2	-8.40	116.10	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	257	ARG	NE-CZ-NH2	-8.40	116.10	120.30
1	C	414	ARG	CD-NE-CZ	8.25	135.15	123.60
1	C	86	ARG	NE-CZ-NH2	-8.20	116.20	120.30
1	C	229	ARG	NE-CZ-NH2	8.20	124.40	120.30
1	D	58	ARG	NH1-CZ-NH2	-8.03	110.57	119.40
1	B	199	ARG	NE-CZ-NH1	7.95	124.27	120.30
1	D	207	ARG	NE-CZ-NH2	-7.89	116.36	120.30
1	D	249	ARG	NE-CZ-NH2	-7.78	116.41	120.30
1	C	269	ARG	NE-CZ-NH2	-7.60	116.50	120.30
1	A	207	ARG	NE-CZ-NH1	7.57	124.08	120.30
1	D	373	ARG	CD-NE-CZ	7.56	134.18	123.60
1	C	207	ARG	NE-CZ-NH1	7.47	124.03	120.30
1	B	533	ARG	CD-NE-CZ	7.38	133.94	123.60
1	D	207	ARG	NE-CZ-NH1	7.37	123.99	120.30
1	B	207	ARG	CD-NE-CZ	7.36	133.90	123.60
1	B	249	ARG	NE-CZ-NH2	-7.11	116.75	120.30
1	B	229	ARG	NE-CZ-NH1	7.07	123.83	120.30
1	D	257	ARG	NE-CZ-NH1	7.06	123.83	120.30
1	C	257	ARG	NE-CZ-NH2	-7.05	116.77	120.30
1	D	257	ARG	NE-CZ-NH2	-7.01	116.79	120.30
1	C	86	ARG	NE-CZ-NH1	7.00	123.80	120.30
1	C	257	ARG	NE-CZ-NH1	6.96	123.78	120.30
1	C	505	ARG	CD-NE-CZ	6.95	133.32	123.60
1	C	131	ARG	NE-CZ-NH2	-6.62	116.99	120.30
1	B	131	ARG	CD-NE-CZ	6.45	132.63	123.60
1	B	269	ARG	NE-CZ-NH2	-6.42	117.09	120.30
1	A	207	ARG	NE-CZ-NH2	-6.40	117.10	120.30
1	A	136	ARG	CD-NE-CZ	6.39	132.54	123.60
1	A	131	ARG	NE-CZ-NH2	-6.33	117.13	120.30
1	D	86	ARG	CD-NE-CZ	6.25	132.35	123.60
1	A	373	ARG	NE-CZ-NH2	-6.16	117.22	120.30
1	C	373	ARG	NE-CZ-NH1	6.10	123.35	120.30
1	B	86	ARG	CD-NE-CZ	6.05	132.07	123.60
1	D	505	ARG	NE-CZ-NH1	6.03	123.32	120.30
1	A	533	ARG	NE-CZ-NH2	-5.99	117.31	120.30
1	A	308	ARG	CD-NE-CZ	5.95	131.92	123.60
1	A	249	ARG	NE-CZ-NH2	-5.94	117.33	120.30
1	D	550	LEU	CA-CB-CG	5.93	128.94	115.30
1	C	207	ARG	NE-CZ-NH2	-5.91	117.35	120.30
1	C	131	ARG	NE-CZ-NH1	5.86	123.23	120.30
1	C	55	LEU	CB-CG-CD1	5.82	120.90	111.00
1	A	249	ARG	NE-CZ-NH1	5.75	123.17	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	269	ARG	NE-CZ-NH2	-5.69	117.46	120.30
1	A	373	ARG	NE-CZ-NH1	5.60	123.10	120.30
1	D	308	ARG	CD-NE-CZ	5.59	131.43	123.60
1	C	136	ARG	NE-CZ-NH1	5.59	123.09	120.30
1	D	533	ARG	NE-CZ-NH1	5.57	123.09	120.30
1	B	249	ARG	NE-CZ-NH1	5.56	123.08	120.30
1	C	193	ASP	CB-CG-OD1	5.50	123.25	118.30
1	D	505	ARG	NE-CZ-NH2	-5.40	117.60	120.30
1	C	533	ARG	NE-CZ-NH1	5.39	123.00	120.30
1	D	63	LEU	CA-CB-CG	5.37	127.64	115.30
1	C	373	ARG	NE-CZ-NH2	-5.35	117.63	120.30
1	B	63	LEU	CA-CB-CG	5.33	127.56	115.30
1	B	595	LEU	CA-CB-CG	5.30	127.49	115.30
1	D	414	ARG	NE-CZ-NH2	-5.30	117.65	120.30
1	D	269	ARG	CD-NE-CZ	5.29	131.01	123.60
1	A	86	ARG	CB-CG-CD	-5.27	97.90	111.60
1	D	533	ARG	NE-CZ-NH2	-5.23	117.68	120.30
1	D	136	ARG	NE-CZ-NH1	5.19	122.90	120.30
1	B	373	ARG	NE-CZ-NH2	-5.18	117.71	120.30
1	B	136	ARG	NE-CZ-NH2	-5.13	117.73	120.30
1	B	373	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	D	63	LEU	CB-CG-CD2	5.10	119.67	111.00
1	C	136	ARG	NE-CZ-NH2	-5.09	117.76	120.30
1	A	131	ARG	NE-CZ-NH1	5.05	122.82	120.30
1	A	505	ARG	NE-CZ-NH2	-5.01	117.79	120.30

There are no chirality outliers.

There are no planarity outliers.

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	A	4583	0	4564	58	1
1	B	4531	0	4508	61	2
1	C	4587	0	4561	69	4
1	D	4591	0	4568	60	5

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	31	0	13	0	0
2	B	31	0	13	0	0
2	C	31	0	13	0	0
2	D	31	0	13	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
All	All	18420	0	18253	244	6

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 7.

All (244) 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:C:34:VAL:HG13	1:C:37:ASP:HB2	1.33	1.07
1:C:191:VAL:CG1	1:C:199:ARG:HG3	2.12	0.80
1:D:54:ASN:O	1:D:58:ARG:HG2	1.81	0.80
1:A:191:VAL:CG1	1:A:199:ARG:HG3	2.13	0.78
1:B:129:GLU:O	1:B:427:ARG:NH2	2.16	0.78
1:A:191:VAL:HG12	1:A:199:ARG:HG3	1.65	0.77
1:A:129:GLU:O	1:A:427:ARG:NH2	2.16	0.75
1:C:191:VAL:HG12	1:C:199:ARG:HG3	1.69	0.73
1:D:42:ILE:O	1:D:46:LYS:HB2	1.89	0.73
1:C:415:GLU:HG3	1:C:458:LEU:HD22	1.70	0.72
1:D:63:LEU:HD11	1:D:82:LEU:HD13	1.73	0.71
1:D:129:GLU:O	1:D:427:ARG:NH2	2.23	0.70
1:B:415:GLU:HG3	1:B:458:LEU:HD22	1.75	0.69
1:A:408:THR:O	1:A:412:GLN:HG3	1.92	0.68
1:C:34:VAL:HG12	1:C:34:VAL:O	1.93	0.67
1:A:415:GLU:HG3	1:A:458:LEU:HD22	1.75	0.67
1:B:393:ALA:O	1:B:427:ARG:NH1	2.28	0.67
1:C:34:VAL:CG1	1:C:37:ASP:HB2	2.18	0.66
1:C:243:THR:HG23	1:C:602:VAL:HB	1.77	0.66
1:D:256:ALA:O	1:D:269:ARG:HD2	1.95	0.66
1:A:243:THR:HG23	1:A:602:VAL:HB	1.77	0.66
1:D:415:GLU:HG3	1:D:458:LEU:HD22	1.77	0.66
1:A:326:GLN:HA	1:A:331:SER:HB3	1.78	0.66
1:C:69:ILE:HG12	1:C:78:THR:HG22	1.79	0.65
1:A:335:LEU:HD11	1:A:416:ASN:HB3	1.77	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:SER:HA	1:B:190:GLY:HA3	1.79	0.64
1:B:326:GLN:HA	1:B:331:SER:HB3	1.79	0.64
1:B:54:ASN:O	1:B:58:ARG:HG2	1.97	0.64
1:C:129:GLU:O	1:C:427:ARG:NH2	2.26	0.64
1:D:326:GLN:HA	1:D:331:SER:HB3	1.79	0.64
1:C:335:LEU:HD11	1:C:416:ASN:HB3	1.79	0.63
1:C:34:VAL:O	1:C:38:THR:HB	1.99	0.63
1:C:78:THR:HG21	1:C:160:THR:OG1	1.98	0.63
1:D:335:LEU:HD11	1:D:416:ASN:HB3	1.81	0.63
1:C:326:GLN:HA	1:C:331:SER:HB3	1.81	0.62
1:A:73:SER:HA	1:A:190:GLY:HA3	1.81	0.62
1:B:63:LEU:HD11	1:B:82:LEU:HD13	1.81	0.62
1:B:335:LEU:HD11	1:B:416:ASN:HB3	1.81	0.62
1:C:38:THR:CG2	1:C:39:GLU:N	2.63	0.62
1:C:54:ASN:O	1:C:58:ARG:HG2	2.00	0.62
1:A:54:ASN:O	1:A:58:ARG:HG2	2.00	0.62
1:A:69:ILE:HG12	1:A:78:THR:HG22	1.82	0.61
1:A:40:LYS:O	1:A:44:VAL:HG13	2.01	0.61
1:B:284:GLN:HG3	1:B:323:SER:HB2	1.83	0.61
1:B:69:ILE:HG12	1:B:78:THR:HG22	1.83	0.60
1:D:38:THR:HA	1:D:41:ARG:HD2	1.82	0.60
1:D:390:LYS:HB3	1:D:391:PRO:HD3	1.83	0.60
1:C:38:THR:HG23	1:C:39:GLU:N	2.17	0.60
1:B:408:THR:O	1:B:412:GLN:HG3	2.02	0.59
1:C:515:VAL:HG11	1:C:569:ARG:HG3	1.85	0.59
1:B:191:VAL:HG12	1:B:199:ARG:HD3	1.85	0.59
1:A:393:ALA:O	1:A:427:ARG:NH1	2.35	0.58
1:C:408:THR:O	1:C:412:GLN:HG3	2.04	0.58
1:C:127:ASP:O	1:C:131:ARG:HD3	2.05	0.57
1:A:220:LYS:HD3	1:A:624:TYR:CE2	2.38	0.57
1:A:390:LYS:HB3	1:A:391:PRO:HD3	1.86	0.57
1:D:191:VAL:HG12	1:D:199:ARG:HD3	1.87	0.57
1:B:601:GLN:O	1:B:604:VAL:HG12	2.05	0.57
1:A:78:THR:HG21	1:A:160:THR:OG1	2.05	0.56
1:D:73:SER:HA	1:D:190:GLY:HA3	1.85	0.56
1:D:191:VAL:CG1	1:D:199:ARG:HG3	2.34	0.56
1:D:515:VAL:HG11	1:D:569:ARG:HG3	1.86	0.56
1:D:220:LYS:HD3	1:D:624:TYR:CE2	2.40	0.56
1:B:250:ARG:HD3	1:B:271:ASP:O	2.06	0.56
1:D:243:THR:HG23	1:D:602:VAL:HB	1.88	0.56
1:C:73:SER:HA	1:C:190:GLY:HA3	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:38:THR:O	1:C:41:ARG:N	2.36	0.55
1:C:393:ALA:O	1:C:427:ARG:NH1	2.36	0.55
1:A:127:ASP:O	1:A:131:ARG:HD3	2.07	0.55
1:B:542:PRO:O	1:B:545:ARG:HG3	2.07	0.55
1:C:69:ILE:HG12	1:C:78:THR:CG2	2.37	0.55
1:B:238:LEU:HD13	1:B:309:LEU:HD21	1.90	0.54
1:D:117:GLU:OE1	1:D:390:LYS:NZ	2.40	0.54
1:D:408:THR:O	1:D:412:GLN:HG3	2.06	0.54
1:B:137:LEU:O	1:B:140:GLN:HG2	2.07	0.54
1:A:338:LYS:HE2	1:A:427:ARG:O	2.08	0.54
1:D:300:ARG:HB3	1:D:300:ARG:NH1	2.23	0.54
1:D:284:GLN:HG3	1:D:323:SER:HB2	1.90	0.54
1:D:373:ARG:HD2	1:D:379:SER:HB2	1.89	0.54
1:B:243:THR:HG23	1:B:602:VAL:HB	1.90	0.54
1:B:78:THR:HG21	1:B:160:THR:OG1	2.08	0.53
1:B:63:LEU:HD12	1:B:82:LEU:HD22	1.89	0.53
1:C:220:LYS:HD3	1:C:624:TYR:CE2	2.44	0.53
1:A:601:GLN:O	1:A:604:VAL:HG12	2.09	0.52
1:B:69:ILE:HG12	1:B:78:THR:CG2	2.40	0.52
1:A:69:ILE:HG12	1:A:78:THR:CG2	2.39	0.52
1:A:143:PRO:O	1:A:146:HIS:HB3	2.10	0.52
1:D:78:THR:HG21	1:D:160:THR:OG1	2.10	0.52
1:B:191:VAL:HG12	1:B:191:VAL:O	2.09	0.52
1:A:184:MET:HG2	1:A:227:TRP:HB3	1.91	0.52
1:C:35:ASP:O	1:C:38:THR:HG22	2.09	0.51
1:A:515:VAL:HG11	1:A:569:ARG:HG3	1.93	0.51
1:C:405:PHE:CE1	1:D:479:ARG:HD3	2.46	0.51
1:D:393:ALA:O	1:D:427:ARG:NH1	2.42	0.51
1:C:305:GLN:N	1:C:305:GLN:OE1	2.44	0.51
1:B:143:PRO:O	1:B:146:HIS:HB3	2.11	0.50
1:C:39:GLU:HA	1:C:42:ILE:HD12	1.93	0.50
1:D:127:ASP:O	1:D:131:ARG:HD3	2.10	0.50
1:B:508:ARG:HG3	1:B:508:ARG:HH11	1.77	0.50
1:C:143:PRO:O	1:C:146:HIS:HB3	2.11	0.50
1:C:250:ARG:HD3	1:C:271:ASP:O	2.11	0.50
1:D:137:LEU:O	1:D:140:GLN:HG2	2.11	0.50
1:A:185:LEU:HB3	1:A:228:THR:HG23	1.94	0.50
1:D:250:ARG:HD3	1:D:271:ASP:O	2.11	0.50
1:A:392:LEU:HD11	1:A:461:PRO:HG2	1.94	0.50
1:C:101:GLN:O	1:C:101:GLN:HG3	2.11	0.50
1:B:401:PRO:HA	1:B:404:SER:HB3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:471:PHE:CD1	1:B:477:VAL:HA	2.47	0.49
1:C:137:LEU:O	1:C:140:GLN:HG2	2.13	0.49
1:B:390:LYS:HB3	1:B:391:PRO:HD3	1.93	0.49
1:D:63:LEU:CD1	1:D:82:LEU:HD13	2.40	0.49
1:C:254:ASP:HB2	1:C:255:PRO:HD2	1.94	0.49
1:C:300:ARG:NH1	1:C:300:ARG:HB3	2.27	0.48
1:D:40:LYS:O	1:D:43:ASN:HB2	2.14	0.48
1:A:137:LEU:O	1:A:140:GLN:HG2	2.13	0.48
1:A:191:VAL:HG12	1:A:191:VAL:O	2.13	0.48
1:A:250:ARG:HD3	1:A:271:ASP:O	2.13	0.48
1:A:63:LEU:HD21	1:A:82:LEU:HD13	1.96	0.48
1:B:300:ARG:NH1	1:B:300:ARG:HB3	2.29	0.48
1:C:38:THR:CG2	1:C:39:GLU:H	2.27	0.48
1:B:63:LEU:O	1:B:179:LEU:HD12	2.13	0.48
1:C:508:ARG:HG3	1:C:508:ARG:HH11	1.77	0.48
1:A:134:ASP:HB3	1:A:137:LEU:HD12	1.95	0.48
1:B:220:LYS:HD3	1:B:624:TYR:CE2	2.48	0.48
1:D:338:LYS:HE2	1:D:427:ARG:O	2.14	0.48
1:B:98:ASP:OD1	1:B:257:ARG:NH2	2.39	0.48
1:D:392:LEU:HD11	1:D:461:PRO:HG2	1.95	0.48
1:B:338:LYS:HE2	1:B:427:ARG:O	2.15	0.47
1:C:284:GLN:HG3	1:C:323:SER:HB2	1.96	0.47
1:C:442:PHE:O	1:C:445:CYS:HB3	2.14	0.47
1:D:69:ILE:HG12	1:D:78:THR:CG2	2.44	0.47
1:D:590:PRO:HD2	1:D:595:LEU:HD11	1.96	0.47
1:C:390:LYS:HB3	1:C:391:PRO:HD3	1.96	0.47
1:A:254:ASP:OD1	1:A:257:ARG:HD3	2.15	0.47
1:A:256:ALA:O	1:A:269:ARG:HD3	2.15	0.47
1:D:601:GLN:O	1:D:604:VAL:HG12	2.13	0.47
1:B:515:VAL:HG11	1:B:569:ARG:HG3	1.97	0.47
1:C:34:VAL:O	1:C:34:VAL:CG1	2.62	0.47
1:D:184:MET:CG	1:D:227:TRP:HB3	2.44	0.47
1:B:191:VAL:HG12	1:B:199:ARG:CD	2.45	0.46
1:D:69:ILE:HG12	1:D:78:THR:HG22	1.96	0.46
1:B:191:VAL:CG1	1:B:199:ARG:HG3	2.45	0.46
1:C:185:LEU:HB3	1:C:228:THR:HG23	1.97	0.46
1:C:338:LYS:HE2	1:C:427:ARG:O	2.16	0.46
1:C:471:PHE:CD1	1:C:477:VAL:HA	2.51	0.46
1:D:63:LEU:HD12	1:D:82:LEU:HD22	1.98	0.46
1:B:305:GLN:N	1:B:305:GLN:OE1	2.49	0.45
1:C:368:GLY:HA2	1:C:387:ASN:HA	1.96	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:143:PRO:O	1:D:146:HIS:HB3	2.16	0.45
1:C:63:LEU:CD2	1:C:82:LEU:HD13	2.47	0.45
1:D:290:GLN:O	1:D:293:THR:OG1	2.34	0.45
1:D:191:VAL:HG12	1:D:191:VAL:O	2.16	0.45
1:D:191:VAL:HG11	1:D:199:ARG:HG3	1.98	0.45
1:A:590:PRO:HD2	1:A:595:LEU:HD11	1.97	0.45
1:C:254:ASP:HB2	1:C:255:PRO:CD	2.47	0.45
1:D:471:PHE:CD1	1:D:477:VAL:HA	2.51	0.45
1:B:541:LEU:HB3	1:B:545:ARG:HB2	1.99	0.45
1:D:184:MET:HG2	1:D:227:TRP:HB3	1.99	0.45
1:C:184:MET:CG	1:C:227:TRP:HB3	2.47	0.45
1:D:63:LEU:O	1:D:179:LEU:HD12	2.17	0.45
1:C:308:ARG:NH2	1:D:308:ARG:NH1	2.65	0.45
1:A:442:PHE:O	1:A:445:CYS:HB3	2.17	0.44
1:B:134:ASP:HB3	1:B:137:LEU:HD12	1.99	0.44
1:C:601:GLN:O	1:C:604:VAL:HG12	2.16	0.44
1:A:40:LYS:O	1:A:44:VAL:CG1	2.65	0.44
1:A:63:LEU:CD2	1:A:82:LEU:HD13	2.47	0.44
1:C:479:ARG:HD3	1:D:405:PHE:CE1	2.52	0.44
1:D:191:VAL:HG12	1:D:199:ARG:CD	2.46	0.44
1:B:191:VAL:O	1:B:191:VAL:CG1	2.65	0.44
1:B:370:VAL:HG22	1:B:380:VAL:HG22	2.00	0.44
1:C:401:PRO:HA	1:C:404:SER:HB3	1.98	0.44
1:D:368:GLY:HA2	1:D:387:ASN:HA	1.99	0.44
1:A:401:PRO:HA	1:A:404:SER:HB3	2.00	0.44
1:D:286:VAL:HG13	1:D:321:SER:HB2	2.00	0.44
1:D:299:VAL:HG21	1:D:480:PHE:HB3	2.00	0.44
1:C:191:VAL:HG12	1:C:191:VAL:O	2.17	0.44
1:A:63:LEU:O	1:A:179:LEU:HD12	2.18	0.44
1:A:144:GLN:OE1	1:A:144:GLN:N	2.51	0.44
1:C:246:HIS:CG	1:C:606:LEU:HD11	2.53	0.44
1:A:186:CYS:HB2	1:A:600:TRP:CH2	2.53	0.43
1:A:40:LYS:C	1:A:42:ILE:N	2.72	0.43
1:B:303:ASN:HB3	1:B:306:ARG:HB3	2.00	0.43
1:B:55:LEU:HD23	1:B:60:HIS:NE2	2.34	0.43
1:C:357:LEU:HD13	1:C:361:PHE:CE2	2.53	0.43
1:C:256:ALA:O	1:C:269:ARG:HD3	2.18	0.43
1:A:181:ILE:HD12	1:A:219:TYR:HB3	2.01	0.43
1:D:134:ASP:HB3	1:D:137:LEU:HD12	2.01	0.43
1:C:456:PRO:HG3	1:C:460:LEU:HD13	2.00	0.43
1:B:240:ALA:HA	1:B:599:GLY:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:LEU:HD21	1:C:82:LEU:HD13	2.01	0.42
1:D:370:VAL:HG22	1:D:380:VAL:HG22	2.00	0.42
1:D:610:LEU:HD23	1:D:610:LEU:HA	1.90	0.42
1:C:590:PRO:HD2	1:C:595:LEU:HD11	2.01	0.42
1:A:303:ASN:HB3	1:A:306:ARG:HB3	2.01	0.42
1:A:368:GLY:HA2	1:A:387:ASN:HA	2.01	0.42
1:B:184:MET:HG2	1:B:227:TRP:HB3	2.01	0.42
1:B:254:ASP:OD1	1:B:257:ARG:HD3	2.19	0.42
1:B:435:ILE:HD12	1:B:437:LYS:HG3	2.01	0.42
1:B:471:PHE:CE1	1:B:477:VAL:HA	2.55	0.42
1:A:300:ARG:HB3	1:A:300:ARG:NH1	2.35	0.42
1:A:191:VAL:CG1	1:A:191:VAL:O	2.67	0.42
1:A:305:GLN:OE1	1:A:305:GLN:N	2.52	0.42
1:A:407:VAL:HG11	1:B:457:LYS:HE2	2.02	0.42
1:B:542:PRO:O	1:B:545:ARG:CG	2.67	0.42
1:A:284:GLN:HG3	1:A:323:SER:HB2	2.01	0.42
1:C:290:GLN:O	1:C:293:THR:OG1	2.38	0.42
1:B:256:ALA:O	1:B:269:ARG:HD3	2.19	0.42
1:B:54:ASN:HB3	1:B:58:ARG:CZ	2.49	0.42
1:C:55:LEU:HD23	1:C:55:LEU:HA	1.68	0.42
1:A:549:LYS:HE3	1:A:549:LYS:HB2	1.88	0.42
1:A:508:ARG:HG3	1:A:508:ARG:HH11	1.84	0.41
1:B:392:LEU:HD11	1:B:461:PRO:HG2	2.01	0.41
1:C:299:VAL:HG21	1:C:480:PHE:HB3	2.02	0.41
1:C:610:LEU:HD23	1:C:610:LEU:HA	1.90	0.41
1:C:353:SER:HB3	1:C:439:THR:HG21	2.01	0.41
1:A:184:MET:CG	1:A:227:TRP:HB3	2.50	0.41
1:D:543:ASN:O	1:D:544:ALA:C	2.59	0.41
1:B:191:VAL:HG11	1:B:199:ARG:HG3	2.02	0.41
1:C:392:LEU:HD11	1:C:461:PRO:HG2	2.03	0.41
1:B:184:MET:CG	1:B:227:TRP:HB3	2.50	0.41
1:A:42:ILE:N	1:A:42:ILE:HD12	2.36	0.41
1:A:290:GLN:O	1:A:293:THR:OG1	2.38	0.41
1:C:63:LEU:O	1:C:179:LEU:HD12	2.20	0.41
1:C:181:ILE:HD12	1:C:219:TYR:HB3	2.03	0.41
1:D:471:PHE:CE1	1:D:477:VAL:HA	2.56	0.41
1:A:412:GLN:OE1	1:A:459:PRO:HG2	2.20	0.41
1:B:185:LEU:HB3	1:B:228:THR:HG23	2.03	0.41
1:B:524:GLU:OE2	1:B:543:ASN:ND2	2.54	0.41
1:A:370:VAL:HG22	1:A:380:VAL:HG22	2.03	0.41
1:A:254:ASP:HB2	1:A:255:PRO:HD2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:299:VAL:HG21	1:B:480:PHE:HB3	2.01	0.41
1:C:39:GLU:HA	1:C:42:ILE:CG1	2.51	0.41
1:D:82:LEU:HG	1:D:103:ILE:HD13	2.03	0.41
1:D:88:CYS:HB2	1:D:91:LYS:HB2	2.02	0.40
1:D:186:CYS:HB2	1:D:600:TRP:CH2	2.55	0.40
1:B:441:ASN:OD1	1:B:441:ASN:C	2.58	0.40
1:D:419:ASP:C	1:D:419:ASP:OD1	2.59	0.40
1:B:181:ILE:HD12	1:B:219:TYR:HB3	2.02	0.40
1:B:254:ASP:HB2	1:B:255:PRO:HD2	2.02	0.40
1:C:126:LYS:O	1:C:127:ASP:HB3	2.21	0.40
1:C:184:MET:HG2	1:C:227:TRP:HB3	2.04	0.40
1:A:54:ASN:HB3	1:A:58:ARG:CZ	2.52	0.40
1:B:101:GLN:O	1:B:101:GLN:HG3	2.22	0.40
1:D:357:LEU:HD13	1:D:361:PHE:CE2	2.56	0.40

All (6) 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)
1:C:514:ASP:OD2	1:D:269:ARG:NH2[1_455]	1.90	0.30
1:C:508:ARG:NH1	1:D:260:ILE:O[1_455]	2.06	0.14
1:C:514:ASP:CG	1:D:269:ARG:NH2[1_455]	2.08	0.12
1:B:344:ASP:OD2	1:D:447:SER:OG[5_545]	2.10	0.10
1:C:508:ARG:NH1	1:D:261:ASP:O[1_455]	2.12	0.08
1:A:575:GLU:O	1:B:508:ARG:NH2[1_455]	2.19	0.01

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	589/611 (96%)	567 (96%)	22 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	583/611 (95%)	560 (96%)	23 (4%)	0	100	100
1	C	588/611 (96%)	564 (96%)	23 (4%)	1 (0%)	47	75
1	D	590/611 (97%)	568 (96%)	22 (4%)	0	100	100
All	All	2350/2444 (96%)	2259 (96%)	90 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	90	ASN

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	503/521 (96%)	495 (98%)	8 (2%)	62	84
1	B	497/521 (95%)	489 (98%)	8 (2%)	62	84
1	C	504/521 (97%)	492 (98%)	12 (2%)	49	77
1	D	504/521 (97%)	494 (98%)	10 (2%)	55	80
All	All	2008/2084 (96%)	1970 (98%)	38 (2%)	57	81

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	LEU
1	A	63	LEU
1	A	77	ARG
1	A	78	THR
1	A	133	VAL
1	A	294	VAL
1	A	309	LEU
1	A	327	LEU
1	B	55	LEU
1	B	63	LEU

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Mol	Chain	Res	Type
1	B	77	ARG
1	B	133	VAL
1	B	294	VAL
1	B	309	LEU
1	B	327	LEU
1	B	475	ASP
1	C	38	THR
1	C	55	LEU
1	C	63	LEU
1	C	77	ARG
1	C	78	THR
1	C	225	PRO
1	C	289	LEU
1	C	294	VAL
1	C	309	LEU
1	C	408	THR
1	C	460	LEU
1	C	475	ASP
1	D	55	LEU
1	D	63	LEU
1	D	77	ARG
1	D	78	THR
1	D	272	LEU
1	D	294	VAL
1	D	309	LEU
1	D	327	LEU
1	D	475	ASP
1	D	550	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	607	HIS
1	B	543	ASN
1	C	208	HIS
1	C	412	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	$\# Z > 2$
2	ANP	D	700	3	29,33,33	1.96	7 (24%)	31,52,52	1.84	7 (22%)
2	ANP	B	700	3	29,33,33	2.08	7 (24%)	31,52,52	1.78	6 (19%)
2	ANP	C	700	3	29,33,33	1.81	7 (24%)	31,52,52	1.87	7 (22%)
2	ANP	A	700	3	29,33,33	1.80	6 (20%)	31,52,52	1.68	6 (19%)

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	ANP	D	700	3	-	8/14/38/38	0/3/3/3
2	ANP	B	700	3	-	8/14/38/38	0/3/3/3
2	ANP	C	700	3	-	8/14/38/38	0/3/3/3
2	ANP	A	700	3	-	8/14/38/38	0/3/3/3

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	700	ANP	PG-N3B	5.18	1.76	1.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	700	ANP	PB-N3B	4.99	1.76	1.63
2	D	700	ANP	PG-N3B	4.88	1.76	1.63
2	A	700	ANP	PB-N3B	4.86	1.76	1.63
2	D	700	ANP	PB-N3B	4.80	1.75	1.63
2	A	700	ANP	PG-N3B	4.71	1.75	1.63
2	C	700	ANP	PG-N3B	4.10	1.74	1.63
2	C	700	ANP	PB-N3B	4.05	1.73	1.63
2	B	700	ANP	PG-O1G	4.01	1.52	1.46
2	B	700	ANP	PA-O3A	3.67	1.63	1.59
2	D	700	ANP	PG-O1G	3.66	1.51	1.46
2	D	700	ANP	PB-O1B	3.52	1.51	1.46
2	C	700	ANP	PG-O1G	3.45	1.51	1.46
2	A	700	ANP	PB-O1B	3.09	1.50	1.46
2	C	700	ANP	PB-O1B	3.09	1.50	1.46
2	C	700	ANP	PA-O3A	2.97	1.62	1.59
2	B	700	ANP	PB-O1B	2.87	1.50	1.46
2	B	700	ANP	PB-O3A	2.79	1.62	1.59
2	A	700	ANP	PG-O1G	2.71	1.50	1.46
2	D	700	ANP	PB-O3A	2.71	1.62	1.59
2	D	700	ANP	PG-O3G	-2.54	1.50	1.56
2	B	700	ANP	C2-N3	2.32	1.35	1.32
2	D	700	ANP	C2-N3	2.25	1.35	1.32
2	C	700	ANP	PB-O3A	2.17	1.61	1.59
2	C	700	ANP	C2-N3	2.10	1.35	1.32
2	A	700	ANP	PB-O2B	-2.06	1.51	1.56
2	A	700	ANP	PG-O3G	-2.05	1.51	1.56

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	700	ANP	O1G-PG-N3B	-5.87	103.12	111.77
2	C	700	ANP	O1G-PG-N3B	-4.48	105.17	111.77
2	B	700	ANP	O1G-PG-N3B	-4.46	105.21	111.77
2	C	700	ANP	O1B-PB-N3B	-4.42	105.25	111.77
2	B	700	ANP	O2B-PB-O1B	4.33	119.15	109.87
2	C	700	ANP	O2B-PB-O1B	4.19	118.85	109.87
2	C	700	ANP	N3-C2-N1	-3.99	123.26	128.67
2	A	700	ANP	O2B-PB-O1B	3.96	118.36	109.87
2	A	700	ANP	N3-C2-N1	-3.85	123.44	128.67
2	A	700	ANP	O1G-PG-N3B	-3.85	106.11	111.77
2	D	700	ANP	O2B-PB-O1B	3.79	118.00	109.87
2	B	700	ANP	N3-C2-N1	-3.77	123.55	128.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	700	ANP	C4-C5-N7	-3.66	105.47	109.34
2	D	700	ANP	N3-C2-N1	-3.33	124.16	128.67
2	B	700	ANP	O1B-PB-N3B	-3.26	106.97	111.77
2	D	700	ANP	C4-C5-N7	-3.06	106.10	109.34
2	C	700	ANP	C4-C5-N7	-2.86	106.31	109.34
2	A	700	ANP	C4-C5-N7	-2.77	106.41	109.34
2	D	700	ANP	O1B-PB-N3B	-2.76	107.70	111.77
2	A	700	ANP	O2G-PG-O3G	2.68	114.79	107.59
2	C	700	ANP	O2A-PA-O1A	2.48	123.98	112.44
2	D	700	ANP	O2G-PG-O3G	2.47	114.23	107.59
2	C	700	ANP	O2G-PG-O3G	2.33	113.84	107.59
2	A	700	ANP	O2A-PA-O1A	2.23	122.81	112.44
2	D	700	ANP	O2A-PA-O1A	2.18	122.60	112.44
2	B	700	ANP	O2G-PG-O3G	2.01	113.00	107.59

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	700	ANP	PB-N3B-PG-O1G
2	A	700	ANP	PG-N3B-PB-O1B
2	A	700	ANP	PA-O3A-PB-O2B
2	A	700	ANP	C5'-O5'-PA-O1A
2	A	700	ANP	C3'-C4'-C5'-O5'
2	B	700	ANP	PB-N3B-PG-O1G
2	B	700	ANP	PG-N3B-PB-O1B
2	B	700	ANP	PA-O3A-PB-O2B
2	B	700	ANP	C5'-O5'-PA-O1A
2	B	700	ANP	C3'-C4'-C5'-O5'
2	C	700	ANP	PB-N3B-PG-O1G
2	C	700	ANP	PG-N3B-PB-O1B
2	C	700	ANP	PA-O3A-PB-O2B
2	C	700	ANP	C5'-O5'-PA-O1A
2	C	700	ANP	C3'-C4'-C5'-O5'
2	D	700	ANP	PB-N3B-PG-O1G
2	D	700	ANP	PG-N3B-PB-O1B
2	D	700	ANP	PA-O3A-PB-O2B
2	D	700	ANP	C5'-O5'-PA-O1A
2	D	700	ANP	C5'-O5'-PA-O3A
2	D	700	ANP	C3'-C4'-C5'-O5'
2	A	700	ANP	O4'-C4'-C5'-O5'
2	B	700	ANP	O4'-C4'-C5'-O5'

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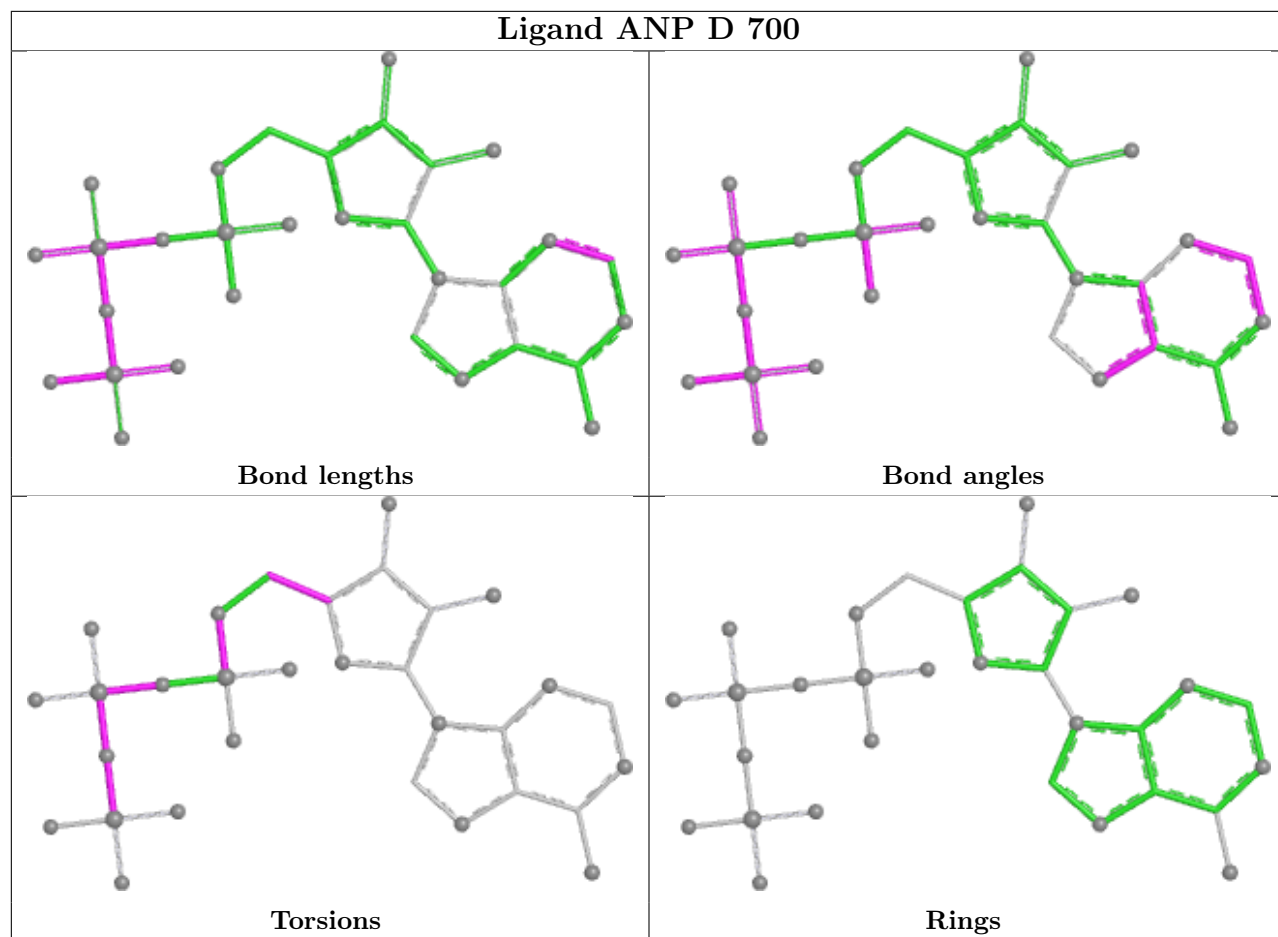
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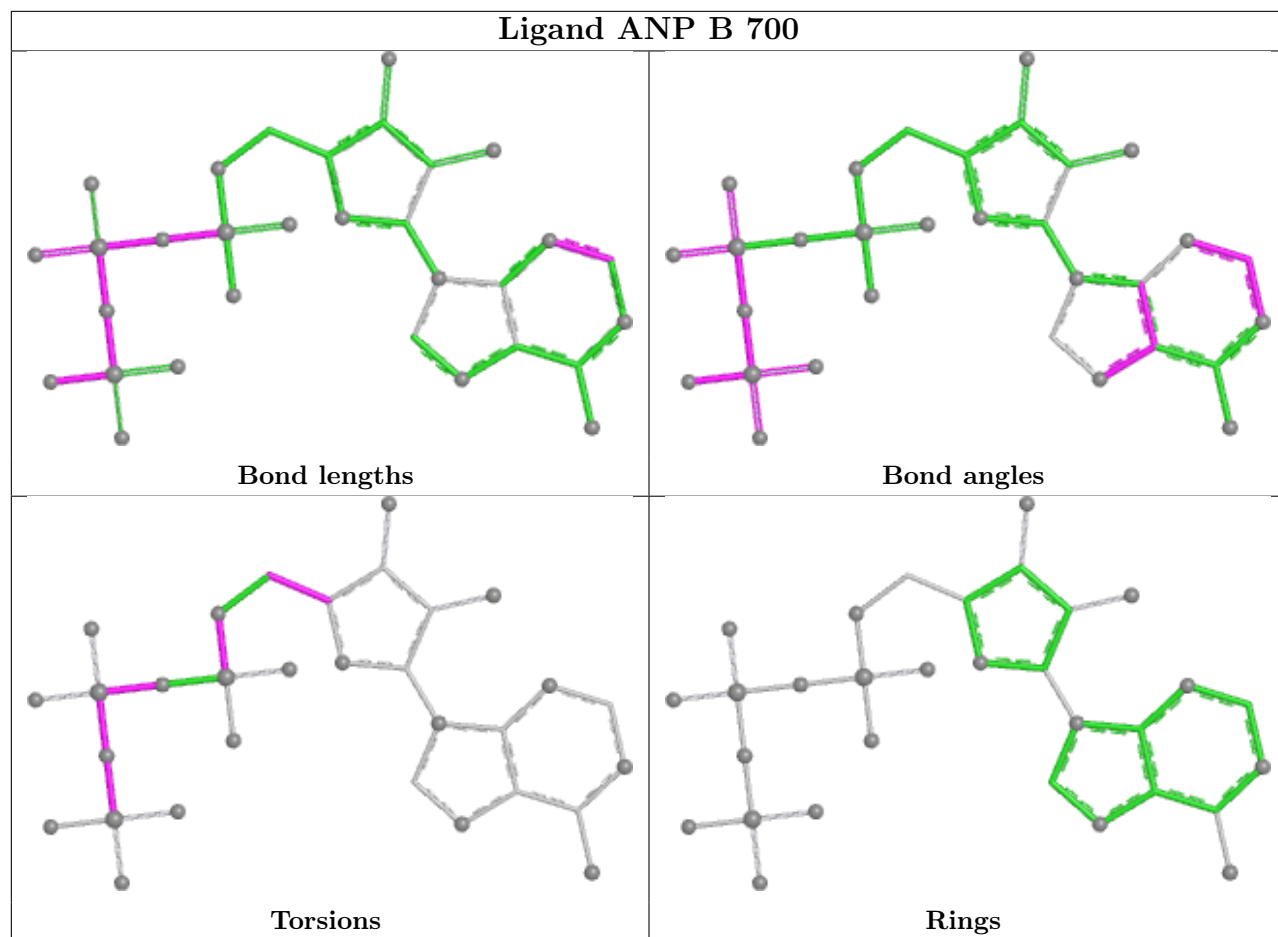
Mol	Chain	Res	Type	Atoms
2	C	700	ANP	O4'-C4'-C5'-O5'
2	D	700	ANP	O4'-C4'-C5'-O5'
2	A	700	ANP	C5'-O5'-PA-O2A
2	A	700	ANP	C5'-O5'-PA-O3A
2	B	700	ANP	C5'-O5'-PA-O2A
2	B	700	ANP	C5'-O5'-PA-O3A
2	C	700	ANP	C5'-O5'-PA-O2A
2	C	700	ANP	C5'-O5'-PA-O3A
2	D	700	ANP	C5'-O5'-PA-O2A

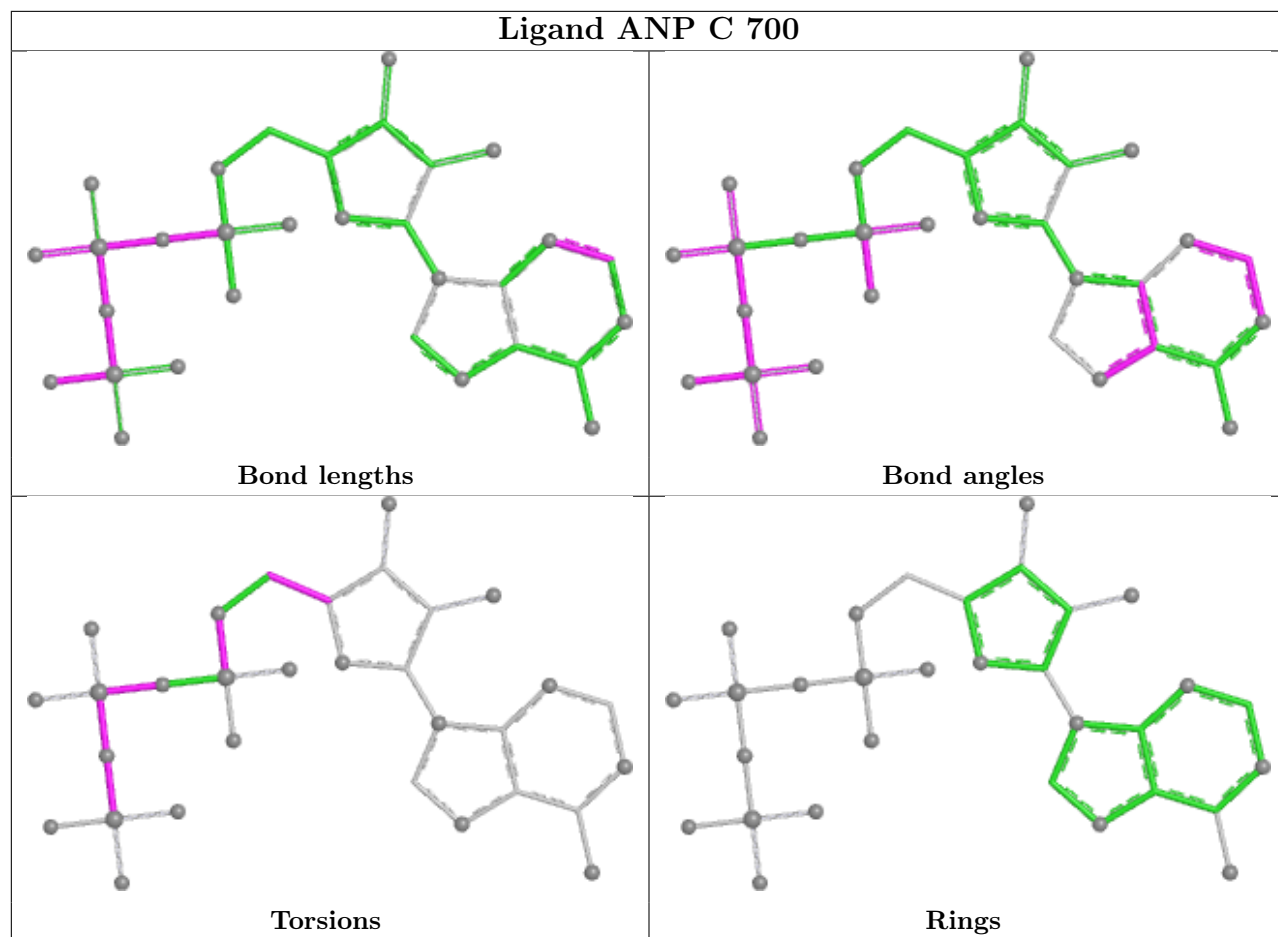
There are no ring outliers.

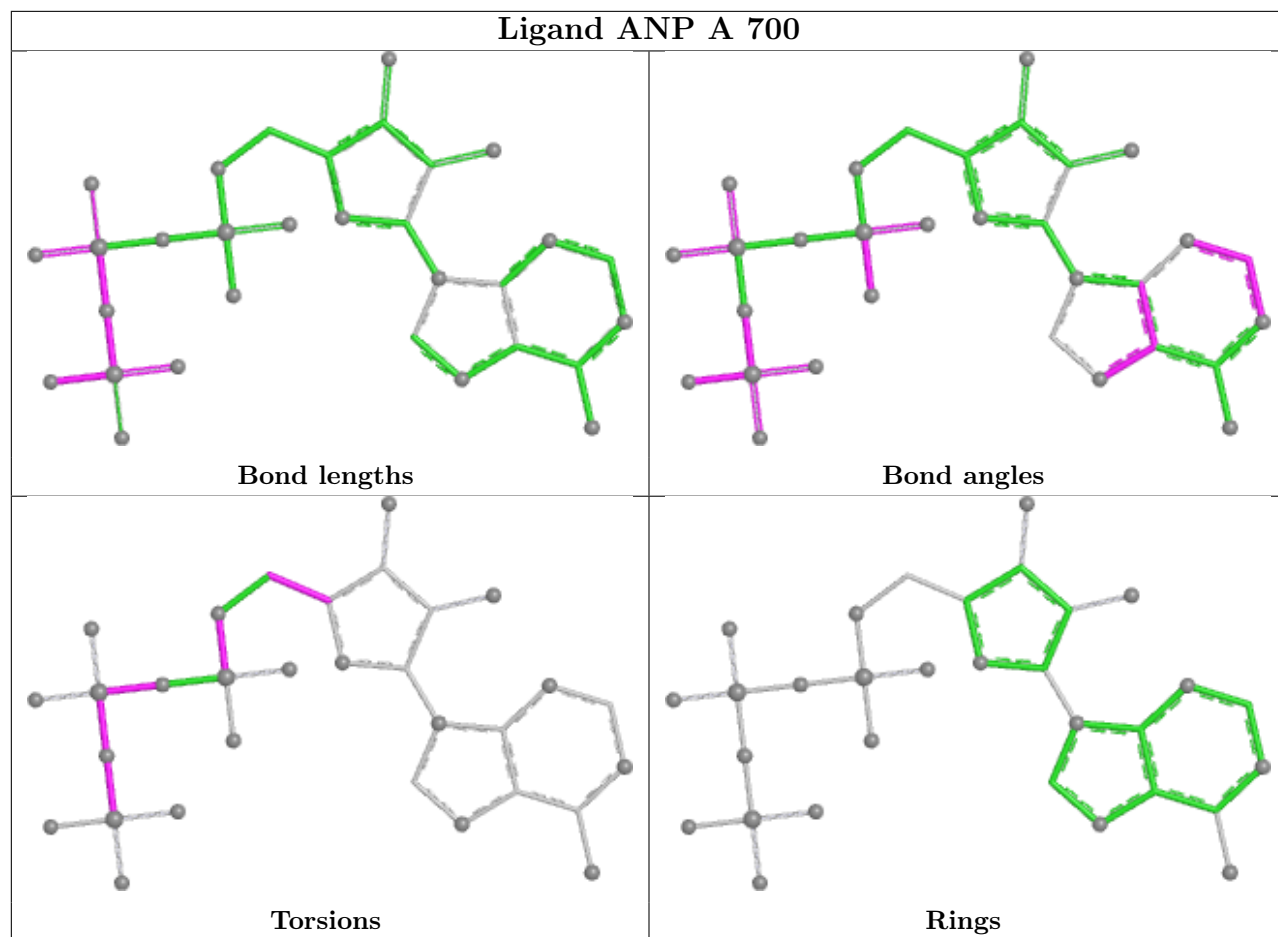
No monomer is involved in short contacts.

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	591/611 (96%)	0.12	27 (4%) 32 27	26, 43, 80, 124	0
1	B	585/611 (95%)	0.13	21 (3%) 42 37	24, 40, 62, 88	0
1	C	592/611 (96%)	-0.25	5 (0%) 86 85	12, 24, 51, 81	0
1	D	592/611 (96%)	-0.20	6 (1%) 82 81	12, 27, 57, 90	0
All	All	2360/2444 (96%)	-0.05	59 (2%) 57 54	12, 34, 64, 124	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	45	GLY	4.1
1	B	262	GLU	4.1
1	C	292	GLY	3.6
1	B	535	ASP	3.6
1	A	544	ALA	3.5
1	A	590	PRO	3.5
1	D	44	VAL	3.3
1	A	457	LYS	3.2
1	A	374	PRO	3.2
1	A	534	THR	3.1
1	D	534	THR	3.1
1	A	506	LEU	3.1
1	A	535	ASP	3.1
1	D	40	LYS	3.1
1	A	262	GLU	2.9
1	B	47	LYS	2.9
1	B	263	TYR	2.8
1	A	50	GLN	2.7
1	B	549	LYS	2.7
1	A	537	PRO	2.7
1	D	591	ASP	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	550	LEU	2.6
1	B	591	ASP	2.6
1	A	538	VAL	2.6
1	A	512	SER	2.6
1	B	593	LYS	2.6
1	A	434	SER	2.5
1	B	217	HIS	2.4
1	A	132	PRO	2.4
1	C	44	VAL	2.4
1	B	592	GLY	2.4
1	C	34	VAL	2.4
1	A	375	ASP	2.4
1	B	544	ALA	2.4
1	A	533	ARG	2.3
1	B	531	ILE	2.3
1	A	353	SER	2.3
1	B	590	PRO	2.3
1	A	423	PRO	2.3
1	D	48	HIS	2.2
1	A	539	ILE	2.2
1	B	136	ARG	2.2
1	C	36	ALA	2.2
1	B	543	ASN	2.2
1	A	536	GLY	2.2
1	C	263	TYR	2.2
1	B	505	ARG	2.1
1	B	533	ARG	2.1
1	B	541	LEU	2.1
1	A	358	PHE	2.1
1	B	521	ALA	2.1
1	B	537	PRO	2.1
1	A	591	ASP	2.1
1	A	54	ASN	2.1
1	A	378	ALA	2.1
1	B	59	CYS	2.0
1	A	585	LYS	2.0
1	A	547	GLU	2.0
1	B	536	GLY	2.0

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

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](#)

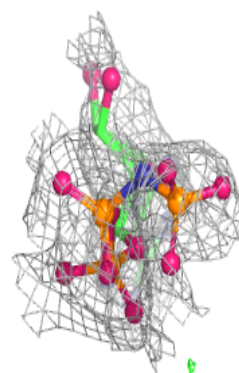
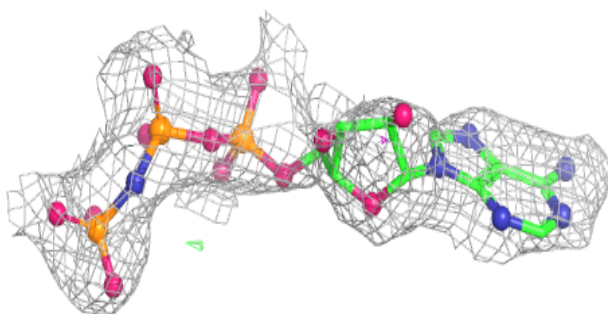
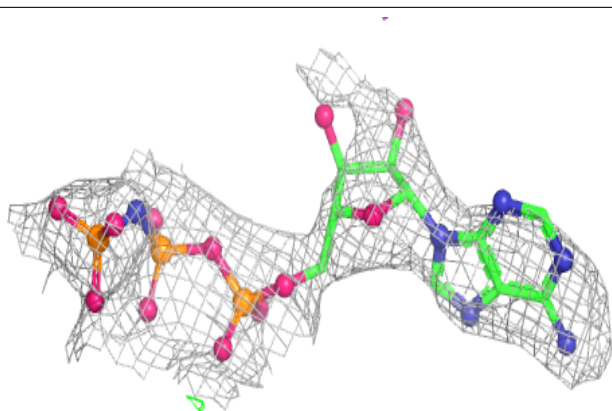
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	ANP	B	700	31/31	0.89	0.20	60,82,95,96	0
2	ANP	D	700	31/31	0.92	0.21	36,52,69,73	0
2	ANP	A	700	31/31	0.93	0.19	47,66,77,79	0
3	MG	A	701	1/1	0.94	0.29	30,30,30,30	0
3	MG	C	701	1/1	0.94	0.25	24,24,24,24	0
2	ANP	C	700	31/31	0.95	0.17	31,41,44,45	0
3	MG	B	701	1/1	0.96	0.19	39,39,39,39	0
3	MG	D	701	1/1	0.98	0.40	25,25,25,25	0

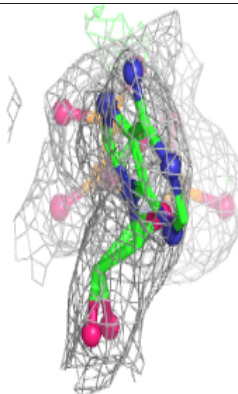
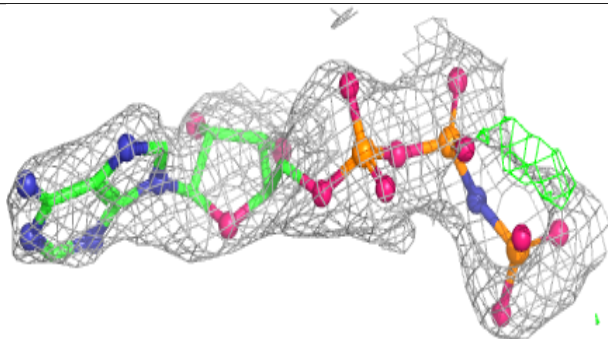
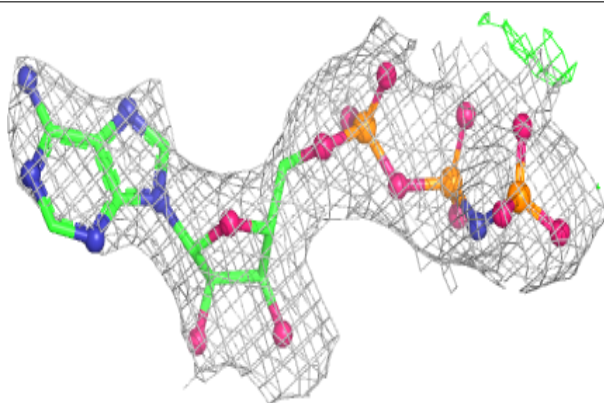
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ANP B 700:

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

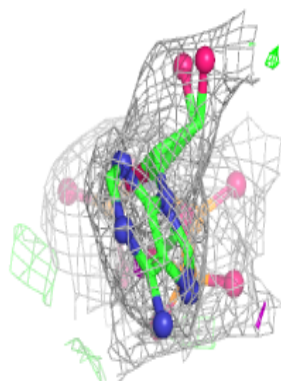
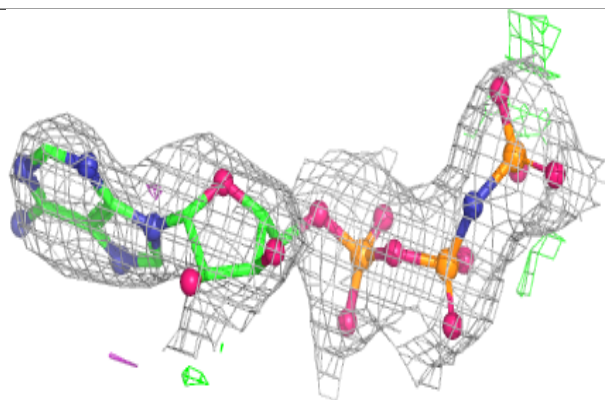
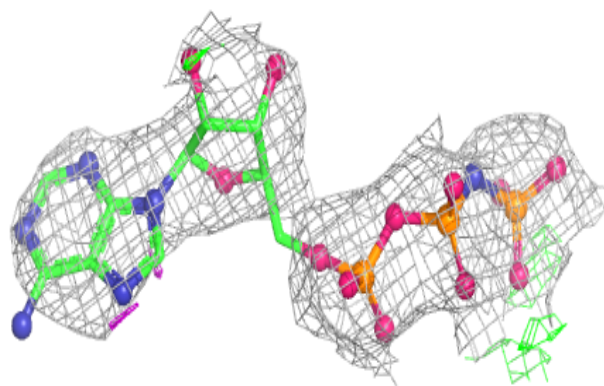
**Electron density around ANP D 700:**

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

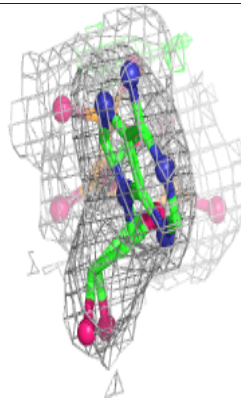
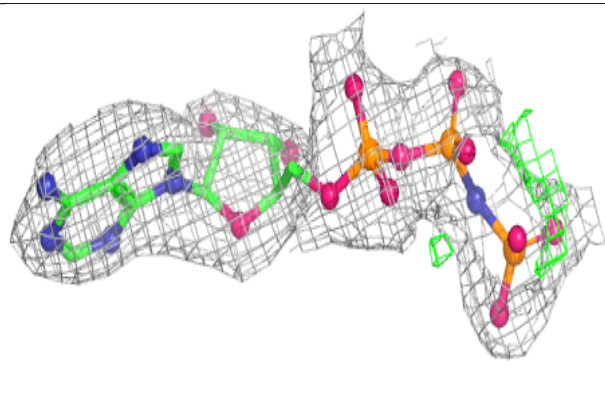
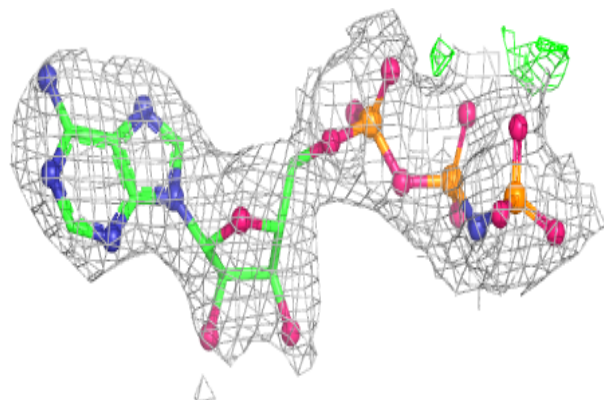


Electron density around ANP A 700:

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

**Electron density around ANP C 700:**

$2mF_o-DF_c$ (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.