



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:32 AM UTC

PDB ID : 1ESQ / pdb_00001esq
Title : CRYSTAL STRUCTURE OF THIAZOLE KINASE MUTANT (C198S)
WITH ATP AND THIAZOLE PHOSPHATE.
Authors : Campobasso, N.; Mathews, I.I.; Begley, T.P.; Ealick, S.E.
Deposited on : 2000-04-10
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 Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

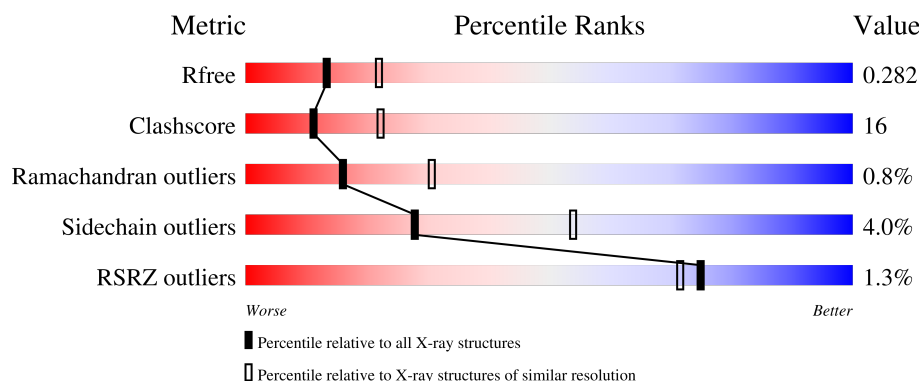
1 Overall quality at a glance

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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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	284	65% 24% • 9%
1	B	284	% 66% 23% • 9%
1	C	284	2% 64% 30% • •

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6061 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 HYDROXYETHYLTHIAZOLE KINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	0	0
			1881	1176	331	368	6			
1	B	258	Total	C	N	O	S	0	0	0
			1876	1174	331	365	6			
1	C	274	Total	C	N	O	S	0	0	0
			1982	1230	362	383	7			

There are 39 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	expression tag	UNP P39593
A	-10	ARG	-	expression tag	UNP P39593
A	-9	GLY	-	expression tag	UNP P39593
A	-8	SER	-	expression tag	UNP P39593
A	-7	HIS	-	expression tag	UNP P39593
A	-6	HIS	-	expression tag	UNP P39593
A	-5	HIS	-	expression tag	UNP P39593
A	-4	HIS	-	expression tag	UNP P39593
A	-3	HIS	-	expression tag	UNP P39593
A	-2	HIS	-	expression tag	UNP P39593
A	-1	GLY	-	expression tag	UNP P39593
A	0	SER	-	expression tag	UNP P39593
A	198	SER	CYS	engineered mutation	UNP P39593
B	-11	MET	-	expression tag	UNP P39593
B	-10	ARG	-	expression tag	UNP P39593
B	-9	GLY	-	expression tag	UNP P39593
B	-8	SER	-	expression tag	UNP P39593
B	-7	HIS	-	expression tag	UNP P39593
B	-6	HIS	-	expression tag	UNP P39593
B	-5	HIS	-	expression tag	UNP P39593
B	-4	HIS	-	expression tag	UNP P39593
B	-3	HIS	-	expression tag	UNP P39593
B	-2	HIS	-	expression tag	UNP P39593

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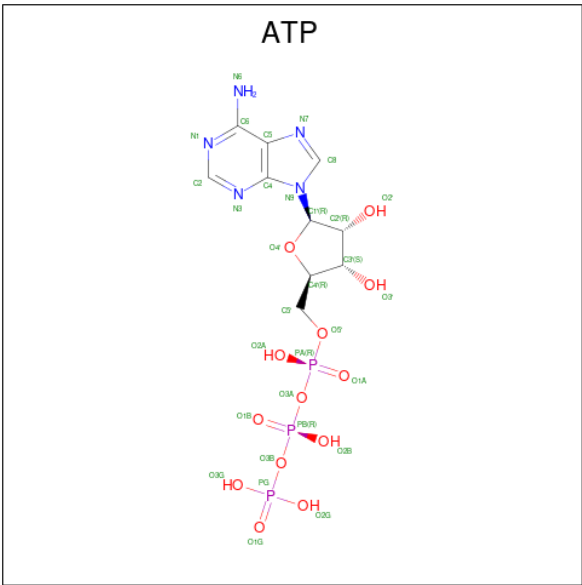
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	GLY	-	expression tag	UNP P39593
B	0	SER	-	expression tag	UNP P39593
B	198	SER	CYS	engineered mutation	UNP P39593
C	-11	MET	-	expression tag	UNP P39593
C	-10	ARG	-	expression tag	UNP P39593
C	-9	GLY	-	expression tag	UNP P39593
C	-8	SER	-	expression tag	UNP P39593
C	-7	HIS	-	expression tag	UNP P39593
C	-6	HIS	-	expression tag	UNP P39593
C	-5	HIS	-	expression tag	UNP P39593
C	-4	HIS	-	expression tag	UNP P39593
C	-3	HIS	-	expression tag	UNP P39593
C	-2	HIS	-	expression tag	UNP P39593
C	-1	GLY	-	expression tag	UNP P39593
C	0	SER	-	expression tag	UNP P39593
C	198	SER	CYS	engineered mutation	UNP P39593

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

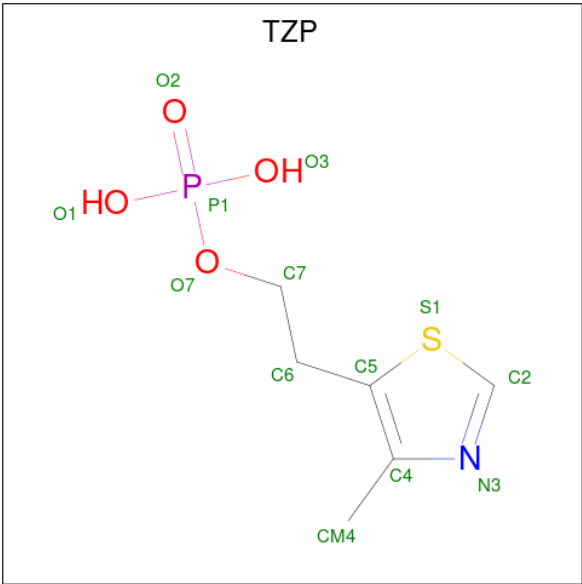
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total 2 Mg 2	0	0
2	B	2	Total 2 Mg 2	0	0
2	C	2	Total 2 Mg 2	0	0

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



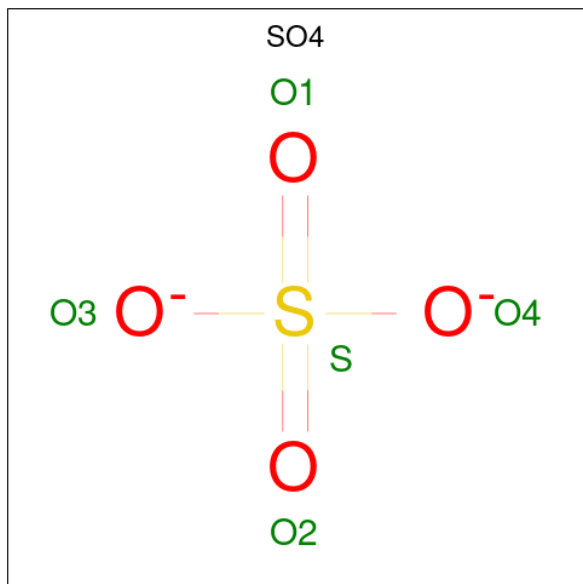
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	C	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 4 is 4-METHYL-5-HYDROXYETHYLTHIAZOLE PHOSPHATE (CCD ID: TZP) (formula: C₆H₁₀NO₄PS).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	S	0	0
			13	6	1	4	1	1		
4	B	1	Total	C	N	O	P	S	0	0
			13	6	1	4	1	1		
4	C	1	Total	C	N	O	P	S	0	0
			13	6	1	4	1	1		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	S	0	0
			5	4	1		

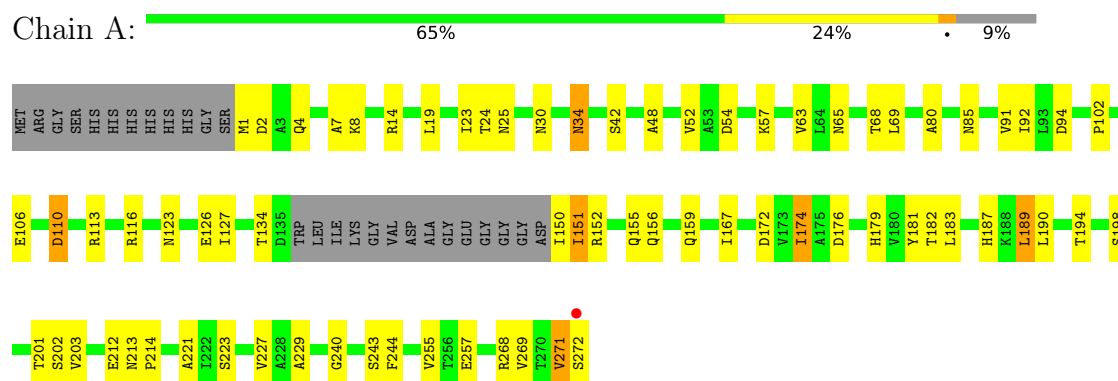
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	56	Total	O	0	0
			56	56		
6	B	73	Total	O	0	0
			73	73		
6	C	50	Total	O	0	0
			50	50		

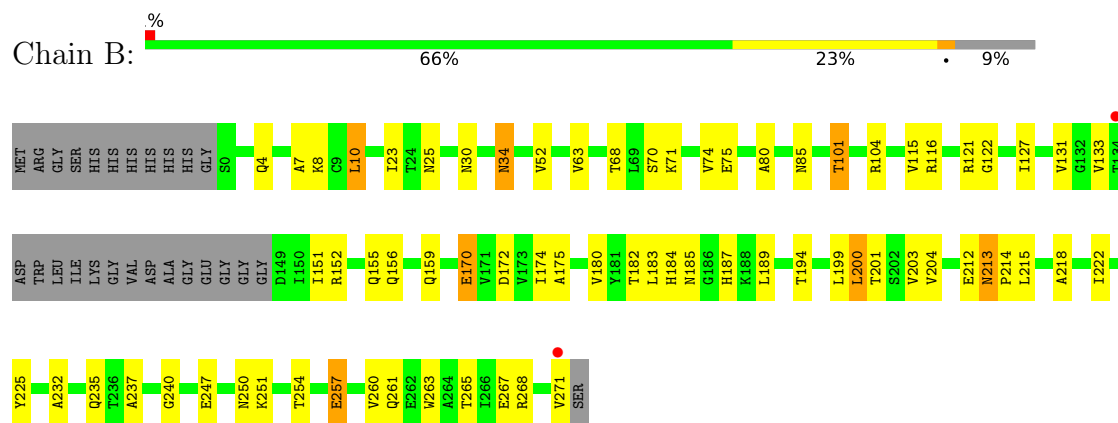
3 Residue-property plots

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.

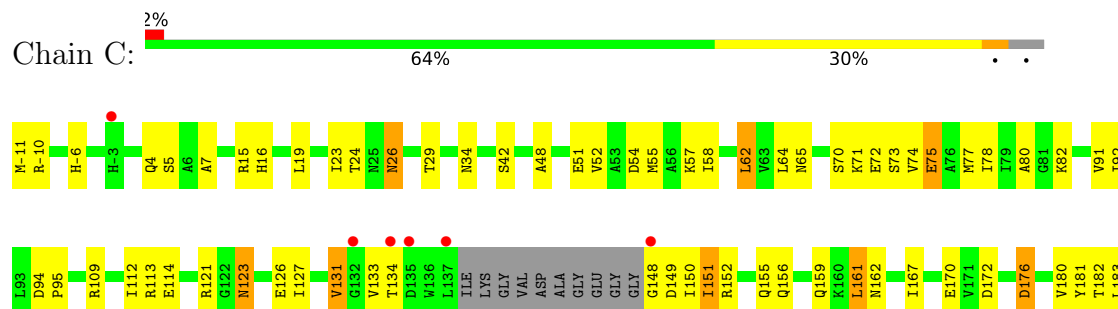
• Molecule 1: HYDROXYETHYLTHIAZOLE KINASE

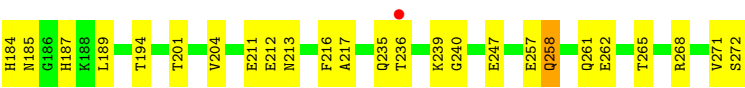


• Molecule 1: HYDROXYETHYLTHIAZOLE KINASE



• Molecule 1: HYDROXYETHYLTHIAZOLE KINASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.08Å 100.84Å 72.51Å 90.00° 95.10° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	94.9 (20.00-2.50) 99.6 (20.00-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.16 (at 2.50Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.209 , 0.282 0.210 , 0.282	Depositor DCC
R_{free} test set	1300 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.474	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 34.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6061	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.28% 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: SO4, MG, TZP, ATP

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	0.46	0/1903	0.90	4/2589 (0.2%)
1	B	0.47	0/1898	0.97	7/2582 (0.3%)
1	C	0.46	0/2010	0.94	5/2727 (0.2%)
All	All	0.46	0/5811	0.94	16/7898 (0.2%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	23	ILE	N-CA-C	-7.17	96.66	106.85
1	A	240	GLY	CA-C-N	7.11	127.43	119.32
1	A	240	GLY	C-N-CA	7.11	127.43	119.32
1	B	257	GLU	N-CA-C	-7.06	103.58	111.28
1	A	23	ILE	N-CA-C	-6.36	96.60	106.72
1	C	240	GLY	CA-C-N	5.96	126.11	119.32
1	C	240	GLY	C-N-CA	5.96	126.11	119.32
1	B	133	VAL	N-CA-C	5.52	116.25	108.36
1	B	115	VAL	N-CA-C	5.42	115.70	108.11
1	C	23	ILE	N-CA-C	-5.42	98.65	106.88
1	C	131	VAL	N-CA-C	5.35	117.25	111.58
1	C	176	ASP	N-CA-C	-5.32	101.63	109.18
1	B	237	ALA	CB-CA-C	-5.14	110.23	117.23
1	B	240	GLY	CA-C-N	5.05	124.65	119.05
1	B	240	GLY	C-N-CA	5.05	124.65	119.05
1	A	69	LEU	N-CA-C	5.02	117.01	110.43

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	1881	0	1921	73	0
1	B	1876	0	1913	54	0
1	C	1982	0	1980	73	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
3	A	31	0	12	0	0
3	B	31	0	12	0	0
3	C	31	0	12	1	0
4	A	13	0	8	1	0
4	B	13	0	8	0	0
4	C	13	0	8	0	0
5	B	5	0	0	0	0
6	A	56	0	0	4	0
6	B	73	0	0	2	0
6	C	50	0	0	1	0
All	All	6061	0	5874	192	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 16.

All (192) 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:185:ASN:HD21	1:C:265:THR:H	1.14	0.90
1:A:190:LEU:HD21	1:A:229:ALA:HB1	1.57	0.86
1:B:127:ILE:O	1:B:131:VAL:HG12	1.78	0.84
1:A:271:VAL:HG12	1:A:272:SER:H	1.46	0.80
1:C:71:LYS:O	1:C:75:GLU:HG2	1.82	0.80
1:B:170:GLU:H	1:B:170:GLU:CD	1.87	0.79
1:C:123:ASN:HD22	1:C:123:ASN:C	1.92	0.78
1:A:187:HIS:HD2	1:A:189:LEU:HB2	1.48	0.77
1:C:-11:MET:HB2	1:C:16:HIS:ND1	2.02	0.75
1:B:182:THR:HG22	1:B:267:GLU:HB3	1.67	0.74
1:C:123:ASN:HD21	1:C:126:GLU:HG3	1.52	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:187:HIS:HD2	1:C:189:LEU:H	1.37	0.72
1:A:94:ASP:HB2	1:A:201:THR:HG21	1.70	0.71
1:A:34:ASN:HD21	1:C:194:THR:H	1.39	0.70
1:A:151:ILE:HD11	1:A:152:ARG:NH1	2.05	0.70
1:A:243:SER:H	1:B:250:ASN:HD21	1.39	0.70
1:B:172:ASP:HB2	1:B:183:LEU:HB2	1.72	0.70
1:B:213:ASN:C	1:B:213:ASN:HD22	2.00	0.70
1:B:71:LYS:O	1:B:75:GLU:HG3	1.93	0.69
1:A:91:VAL:C	1:A:92:ILE:HD12	2.18	0.69
1:C:156:GLN:HA	1:C:159:GLN:HE21	1.57	0.69
1:A:34:ASN:ND2	1:C:194:THR:H	1.90	0.68
1:C:187:HIS:CD2	1:C:189:LEU:H	2.10	0.68
1:C:123:ASN:ND2	1:C:126:GLU:H	1.90	0.68
1:A:187:HIS:CD2	1:A:189:LEU:HB2	2.29	0.67
1:B:101:THR:HG21	1:C:51:GLU:OE1	1.94	0.67
1:B:185:ASN:HD21	1:B:265:THR:H	1.41	0.67
1:C:172:ASP:O	1:C:182:THR:HA	1.94	0.66
1:A:151:ILE:HG12	1:A:152:ARG:N	2.10	0.66
1:A:187:HIS:HD2	1:A:189:LEU:H	1.42	0.66
1:A:213:ASN:C	1:A:213:ASN:HD22	2.04	0.66
1:B:52:VAL:HB	1:B:80:ALA:HB2	1.78	0.65
1:A:187:HIS:CD2	1:A:189:LEU:H	2.14	0.65
1:B:199:LEU:O	1:B:203:VAL:HG23	1.95	0.65
1:B:7:ALA:HB2	1:B:260:VAL:HB	1.79	0.65
1:A:151:ILE:HG12	1:A:152:ARG:H	1.60	0.65
1:C:271:VAL:HG22	1:C:272:SER:N	2.12	0.65
1:B:101:THR:HG22	1:B:104:ARG:H	1.61	0.65
1:B:182:THR:HG23	1:B:184:HIS:NE2	2.12	0.65
1:A:85:ASN:HD21	1:A:116:ARG:H	1.43	0.64
1:B:4:GLN:HA	1:B:261:GLN:HG2	1.79	0.64
1:C:123:ASN:ND2	1:C:126:GLU:HG3	2.12	0.64
1:C:55:MET:CE	1:C:58:ILE:HD11	2.27	0.64
1:A:212:GLU:HG2	6:A:453:HOH:O	1.97	0.64
1:B:7:ALA:HB1	1:B:257:GLU:HG2	1.80	0.63
1:A:172:ASP:O	1:A:182:THR:HA	1.98	0.63
1:B:4:GLN:O	1:B:8:LYS:HG2	1.98	0.63
1:B:187:HIS:HD2	1:B:189:LEU:H	1.47	0.63
1:B:152:ARG:CG	1:B:156:GLN:HE21	2.12	0.63
1:B:213:ASN:HD22	1:B:214:PRO:N	1.97	0.63
1:C:72:GLU:CD	1:C:72:GLU:H	2.06	0.63
1:A:85:ASN:ND2	1:A:116:ARG:H	1.97	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:ASN:OD1	1:A:126:GLU:HG3	1.99	0.62
1:A:172:ASP:HB2	1:A:183:LEU:HB2	1.81	0.61
1:A:174:ILE:N	1:A:174:ILE:HD12	2.15	0.61
1:C:123:ASN:HD21	1:C:126:GLU:H	1.48	0.61
1:B:174:ILE:HD11	1:B:222:ILE:HD12	1.82	0.61
1:C:185:ASN:ND2	1:C:265:THR:H	1.92	0.60
1:B:30:ASN:ND2	6:B:539:HOH:O	2.32	0.59
1:B:85:ASN:HD21	1:B:116:ARG:H	1.49	0.59
1:B:121:ARG:HD2	1:B:122:GLY:N	2.18	0.59
1:C:127:ILE:HG13	1:C:167:ILE:HG12	1.84	0.59
1:A:190:LEU:HD21	1:A:229:ALA:CB	2.31	0.59
1:C:161:LEU:N	1:C:161:LEU:HD12	2.18	0.59
1:A:151:ILE:N	1:A:151:ILE:HD13	2.18	0.59
1:C:185:ASN:HD21	1:C:265:THR:N	1.94	0.59
1:B:218:ALA:O	1:B:222:ILE:HG13	2.04	0.58
1:C:24:THR:HA	1:C:65:ASN:HB3	1.83	0.58
1:C:91:VAL:C	1:C:92:ILE:HD12	2.29	0.58
1:C:183:LEU:HB3	6:C:416:HOH:O	2.04	0.58
1:A:7:ALA:HB1	1:A:257:GLU:HG2	1.85	0.57
1:C:26:ASN:HD22	1:C:26:ASN:C	2.09	0.57
1:B:187:HIS:CD2	1:B:189:LEU:H	2.22	0.57
1:B:213:ASN:HD22	1:B:214:PRO:CD	2.17	0.57
1:B:194:THR:H	1:C:34:ASN:ND2	2.02	0.56
1:A:271:VAL:HG12	1:A:272:SER:N	2.18	0.56
1:C:48:ALA:O	1:C:52:VAL:HG13	2.06	0.55
1:B:194:THR:H	1:C:34:ASN:HD21	1.55	0.55
1:B:194:THR:HA	6:B:574:HOH:O	2.07	0.55
1:C:271:VAL:HG22	1:C:272:SER:H	1.72	0.55
1:C:4:GLN:HA	1:C:261:GLN:HG3	1.90	0.54
1:A:151:ILE:HD13	1:A:151:ILE:H	1.73	0.54
1:A:194:THR:H	1:B:34:ASN:ND2	2.06	0.54
1:A:213:ASN:HD22	1:A:214:PRO:N	2.04	0.54
1:B:200:LEU:HG	1:B:225:TYR:CD2	2.43	0.54
1:A:30:ASN:HB3	6:A:401:HOH:O	2.08	0.54
1:B:152:ARG:HG3	1:B:156:GLN:HE21	1.73	0.54
1:C:94:ASP:HB2	1:C:201:THR:HG21	1.90	0.54
1:C:156:GLN:HA	1:C:159:GLN:NE2	2.23	0.54
1:C:258:GLN:HE21	1:C:258:GLN:H	1.57	0.53
1:C:92:ILE:HD12	1:C:92:ILE:N	2.24	0.53
1:B:170:GLU:CD	1:B:170:GLU:N	2.64	0.53
1:C:26:ASN:C	1:C:26:ASN:ND2	2.66	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:ASN:HD21	1:A:116:ARG:N	2.05	0.53
1:C:62:LEU:HD11	1:C:64:LEU:HD21	1.90	0.53
1:A:127:ILE:HG13	1:A:167:ILE:HG12	1.91	0.53
1:C:149:ASP:O	1:C:151:ILE:N	2.36	0.52
1:A:181:TYR:CZ	1:A:268:ARG:HD3	2.44	0.51
1:B:70:SER:O	1:B:74:VAL:HG23	2.11	0.51
1:C:7:ALA:HB1	1:C:257:GLU:HB2	1.92	0.51
1:A:194:THR:H	1:B:34:ASN:HD21	1.57	0.51
1:C:62:LEU:CD1	1:C:64:LEU:HD21	2.41	0.51
1:B:213:ASN:C	1:B:213:ASN:ND2	2.65	0.51
1:C:55:MET:HE3	1:C:55:MET:HA	1.93	0.51
1:A:54:ASP:HA	1:A:57:LYS:HE3	1.92	0.51
1:C:-11:MET:HB2	1:C:16:HIS:HD1	1.74	0.51
1:A:183:LEU:HB3	6:A:405:HOH:O	2.11	0.50
1:B:232:ALA:HB1	1:B:247:GLU:HB3	1.93	0.50
1:A:203:VAL:HG12	1:A:221:ALA:HB1	1.92	0.50
1:A:182:THR:HG23	1:A:269:VAL:CG2	2.42	0.50
1:B:263:TRP:O	1:B:265:THR:HG23	2.12	0.49
1:A:110:ASP:O	1:A:113:ARG:HG2	2.12	0.49
1:B:174:ILE:CD1	1:B:222:ILE:HD12	2.41	0.49
1:A:2:ASP:OD2	1:A:4:GLN:HG2	2.13	0.49
1:C:155:GLN:OE1	1:C:271:VAL:HG21	2.11	0.49
1:A:52:VAL:HB	1:A:80:ALA:HB2	1.94	0.49
1:A:187:HIS:O	1:A:190:LEU:HD23	2.12	0.49
1:C:149:ASP:C	1:C:151:ILE:H	2.20	0.48
1:A:223:SER:O	1:A:227:VAL:HG13	2.13	0.48
1:C:236:THR:HG22	1:C:239:LYS:HD2	1.94	0.48
1:B:235:GLN:HG2	1:B:247:GLU:OE2	2.14	0.48
1:C:123:ASN:C	1:C:123:ASN:ND2	2.65	0.47
1:A:102:PRO:O	1:A:106:GLU:HG3	2.15	0.47
1:A:198:SER:N	4:A:320:TZP:O3	2.44	0.47
1:C:109:ARG:O	1:C:113:ARG:HG3	2.15	0.47
1:A:2:ASP:CG	1:A:4:GLN:HG2	2.39	0.47
1:A:151:ILE:HD11	1:A:152:ARG:HH12	1.77	0.47
1:A:189:LEU:HD23	1:A:244:PHE:CG	2.49	0.47
1:A:268:ARG:HH11	1:A:268:ARG:HG2	1.80	0.47
1:B:10:LEU:HD12	1:B:260:VAL:HG21	1.96	0.47
1:C:95:PRO:HD2	1:C:121:ARG:O	2.14	0.47
1:A:48:ALA:O	1:A:52:VAL:HG13	2.15	0.47
1:A:155:GLN:O	1:A:159:GLN:HG3	2.15	0.46
1:B:155:GLN:HE21	1:B:159:GLN:CD	2.23	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:213:ASN:HD22	1:B:214:PRO:HD2	1.80	0.46
1:C:78:ILE:HG22	1:C:82:LYS:HE2	1.98	0.46
1:C:236:THR:CG2	1:C:247:GLU:HG3	2.46	0.46
1:B:271:VAL:O	1:B:271:VAL:HG12	2.15	0.46
1:A:187:HIS:H	1:A:190:LEU:HD23	1.81	0.45
1:B:182:THR:HG22	1:B:267:GLU:CB	2.40	0.45
1:C:112:ILE:HD13	1:C:161:LEU:HD23	1.97	0.45
1:A:151:ILE:H	1:A:151:ILE:CD1	2.28	0.45
1:A:150:ILE:CA	6:A:467:HOH:O	2.64	0.45
1:B:175:ALA:CB	1:B:180:VAL:HG22	2.46	0.45
1:B:201:THR:O	1:B:204:VAL:HB	2.15	0.45
1:C:271:VAL:CG2	1:C:272:SER:N	2.78	0.45
1:C:161:LEU:O	1:C:162:ASN:C	2.60	0.45
1:C:-6:HIS:HD2	1:C:5:SER:OG	2.00	0.44
1:C:15:ARG:HG2	1:C:16:HIS:CD2	2.52	0.44
1:C:70:SER:O	1:C:74:VAL:HG23	2.17	0.44
1:C:258:GLN:H	1:C:258:GLN:NE2	2.15	0.44
1:C:54:ASP:HA	1:C:57:LYS:HE2	1.99	0.44
1:A:25:ASN:HB2	1:A:68:THR:HG23	2.00	0.44
1:A:19:LEU:HA	1:A:42:SER:O	2.18	0.43
1:A:151:ILE:CG1	1:A:152:ARG:H	2.30	0.43
1:C:26:ASN:HA	1:C:29:THR:HG23	2.00	0.43
1:A:92:ILE:HD12	1:A:92:ILE:N	2.33	0.43
1:A:268:ARG:HG2	1:A:268:ARG:NH1	2.34	0.43
1:C:201:THR:O	1:C:204:VAL:HB	2.18	0.43
1:B:151:ILE:HG23	1:B:180:VAL:HG11	2.01	0.43
1:A:113:ARG:HB3	1:A:113:ARG:NH1	2.34	0.42
1:A:113:ARG:HH11	1:A:113:ARG:CB	2.32	0.42
1:C:211:GLU:HB3	1:C:217:ALA:HB2	2.01	0.42
1:A:63:VAL:HG21	1:A:202:SER:HA	2.01	0.42
1:A:176:ASP:OD1	1:A:179:HIS:N	2.48	0.42
1:A:8:LYS:HA	1:A:8:LYS:HD2	1.84	0.42
1:C:52:VAL:HB	1:C:80:ALA:HB2	2.00	0.42
1:A:151:ILE:CG1	1:A:152:ARG:N	2.80	0.42
1:A:179:HIS:ND1	1:A:268:ARG:NH2	2.68	0.42
1:A:34:ASN:HD22	1:A:34:ASN:HA	1.67	0.41
1:A:151:ILE:N	1:A:151:ILE:CD1	2.83	0.41
1:A:152:ARG:O	1:A:156:GLN:HG3	2.20	0.41
1:B:175:ALA:HB2	1:B:180:VAL:HG13	2.03	0.41
1:A:271:VAL:CG1	1:A:272:SER:H	2.17	0.41
1:C:213:ASN:HB3	1:C:216:PHE:CD1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:19:LEU:HA	1:C:42:SER:HB2	2.02	0.41
1:B:212:GLU:OE2	1:B:212:GLU:N	2.41	0.41
1:B:213:ASN:ND2	1:B:215:LEU:H	2.19	0.41
1:C:161:LEU:HD12	1:C:161:LEU:H	1.84	0.41
1:C:181:TYR:CZ	1:C:268:ARG:HD3	2.56	0.41
1:A:14:ARG:HH22	1:A:255:VAL:HG12	1.86	0.40
1:C:-10:ARG:HG2	1:C:212:GLU:OE1	2.21	0.40
1:C:148:GLY:O	1:C:152:ARG:HB2	2.20	0.40
1:C:170:GLU:HA	3:C:310:ATP:C6	2.56	0.40
1:C:184:HIS:HB2	1:C:265:THR:OG1	2.19	0.40
1:C:176:ASP:OD1	1:C:176:ASP:C	2.64	0.40
1:C:131:VAL:HG12	1:C:161:LEU:HD21	2.03	0.40
1:A:24:THR:HA	1:A:65:ASN:HB3	2.03	0.40
1:B:25:ASN:HB2	1:B:68:THR:HG23	2.03	0.40
1:A:189:LEU:HD23	1:A:244:PHE:CD1	2.57	0.40
1:B:251:LYS:HA	1:B:254:THR:HB	2.03	0.40
1:C:73:SER:HB3	1:C:77:MET:HE3	2.03	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	254/284 (89%)	243 (96%)	9 (4%)	2 (1%)	16	31
1	B	254/284 (89%)	242 (95%)	12 (5%)	0	100	100
1	C	270/284 (95%)	254 (94%)	12 (4%)	4 (2%)	8	16
All	All	778/852 (91%)	739 (95%)	33 (4%)	6 (1%)	16	31

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	133	VAL
1	C	134	THR
1	C	150	ILE
1	C	151	ILE
1	A	134	THR
1	A	271	VAL

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	197/216 (91%)	191 (97%)	6 (3%)	36	64
1	B	195/216 (90%)	187 (96%)	8 (4%)	27	53
1	C	203/216 (94%)	193 (95%)	10 (5%)	22	45
All	All	595/648 (92%)	571 (96%)	24 (4%)	28	54

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	34	ASN
1	A	110	ASP
1	A	151	ILE
1	A	174	ILE
1	A	189	LEU
1	B	10	LEU
1	B	34	ASN
1	B	63	VAL
1	B	101	THR
1	B	170	GLU
1	B	200	LEU
1	B	213	ASN
1	B	268	ARG
1	C	26	ASN
1	C	62	LEU
1	C	75	GLU

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Mol	Chain	Res	Type
1	C	114	GLU
1	C	123	ASN
1	C	161	LEU
1	C	180	VAL
1	C	235	GLN
1	C	258	GLN
1	C	262	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	ASN
1	A	85	ASN
1	A	162	ASN
1	A	187	HIS
1	A	213	ASN
1	A	235	GLN
1	A	261	GLN
1	B	16	HIS
1	B	34	ASN
1	B	85	ASN
1	B	155	GLN
1	B	156	GLN
1	B	159	GLN
1	B	179	HIS
1	B	185	ASN
1	B	187	HIS
1	B	213	ASN
1	B	230	GLN
1	B	250	ASN
1	B	261	GLN
1	C	-6	HIS
1	C	4	GLN
1	C	26	ASN
1	C	34	ASN
1	C	123	ASN
1	C	156	GLN
1	C	159	GLN
1	C	179	HIS
1	C	185	ASN
1	C	187	HIS
1	C	234	GLN

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Mol	Chain	Res	Type
1	C	258	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 ⓘ

Of 13 ligands modelled in this entry, 6 are monoatomic - leaving 7 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$
3	ATP	B	305	2	32,33,33	3.21	14 (43%)	48,52,52	1.72	9 (18%)
5	SO4	B	370	-	4,4,4	0.31	0	6,6,6	0.12	0
4	TZP	C	330	2	12,13,13	3.18	5 (41%)	15,18,18	2.99	7 (46%)
4	TZP	A	320	2	12,13,13	3.05	5 (41%)	15,18,18	3.11	6 (40%)
3	ATP	C	310	2	32,33,33	2.91	12 (37%)	48,52,52	1.68	9 (18%)
3	ATP	A	300	2	32,33,33	3.01	13 (40%)	48,52,52	1.64	8 (16%)
4	TZP	B	325	2	12,13,13	3.07	6 (50%)	15,18,18	3.11	6 (40%)

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
3	ATP	B	305	2	-	9/22/38/38	0/3/3/3
4	TZP	C	330	2	-	1/7/7/7	0/1/1/1
4	TZP	A	320	2	-	1/7/7/7	0/1/1/1
3	ATP	C	310	2	-	9/22/38/38	0/3/3/3
3	ATP	A	300	2	-	9/22/38/38	0/3/3/3
4	TZP	B	325	2	-	3/7/7/7	0/1/1/1

All (55) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	305	ATP	PB-O3B	-11.18	1.47	1.59
3	A	300	ATP	PB-O3B	-10.41	1.48	1.59
3	C	310	ATP	PB-O3B	-9.77	1.48	1.59
4	C	330	TZP	C4-N3	6.65	1.47	1.38
4	B	325	TZP	C4-N3	6.52	1.47	1.38
4	A	320	TZP	C4-N3	6.39	1.46	1.38
4	C	330	TZP	P1-O7	-5.99	1.41	1.60
3	B	305	ATP	PA-O3A	5.56	1.65	1.59
4	B	325	TZP	P1-O7	-5.42	1.43	1.60
3	C	310	ATP	C1'-N9	-5.40	1.31	1.46
4	A	320	TZP	P1-O7	-5.40	1.43	1.60
3	B	305	ATP	C1'-N9	-5.37	1.31	1.46
3	B	305	ATP	PB-O3A	5.21	1.65	1.59
3	A	300	ATP	C1'-N9	-4.93	1.32	1.46
3	B	305	ATP	C4-N3	4.87	1.43	1.34
3	B	305	ATP	C8-N9	4.85	1.45	1.37
3	A	300	ATP	C8-N9	4.82	1.45	1.37
3	A	300	ATP	PB-O3A	4.78	1.64	1.59
3	A	300	ATP	PA-O3A	4.70	1.64	1.59
3	C	310	ATP	PB-O3A	4.69	1.64	1.59
3	C	310	ATP	C4-N3	4.64	1.43	1.34
3	C	310	ATP	C8-N9	4.56	1.45	1.37
3	A	300	ATP	C4-N3	4.36	1.42	1.34
3	C	310	ATP	PA-O3A	4.13	1.64	1.59
3	B	305	ATP	PG-O3G	3.91	1.69	1.54
3	A	300	ATP	PG-O3G	3.68	1.68	1.54
3	C	310	ATP	PG-O3G	3.66	1.68	1.54
4	B	325	TZP	P1-O3	3.63	1.68	1.54
4	A	320	TZP	P1-O3	3.62	1.68	1.54
4	C	330	TZP	P1-O3	3.49	1.67	1.54
3	C	310	ATP	PA-O1A	3.32	1.62	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	300	ATP	PA-O1A	3.28	1.62	1.50
3	B	305	ATP	PA-O1A	3.25	1.62	1.50
4	B	325	TZP	C2-N3	3.24	1.38	1.30
4	C	330	TZP	C2-N3	3.17	1.38	1.30
4	A	320	TZP	C2-N3	3.13	1.38	1.30
4	C	330	TZP	C2-S1	2.87	1.77	1.72
4	B	325	TZP	C2-S1	2.70	1.77	1.72
3	A	300	ATP	C5'-C4'	2.68	1.59	1.51
3	B	305	ATP	C5-C4	2.58	1.43	1.39
3	C	310	ATP	C2-N1	2.57	1.38	1.33
3	A	300	ATP	C2-N1	2.51	1.38	1.33
3	B	305	ATP	C2-N1	2.49	1.38	1.33
4	A	320	TZP	C2-S1	2.41	1.76	1.72
3	A	300	ATP	C5-C4	2.39	1.43	1.39
3	C	310	ATP	C2-N3	2.31	1.38	1.33
3	C	310	ATP	PG-O2G	-2.28	1.46	1.54
3	A	300	ATP	PG-O2G	-2.23	1.46	1.54
3	B	305	ATP	PG-O2G	-2.21	1.46	1.54
3	C	310	ATP	C6-N1	2.16	1.41	1.35
3	B	305	ATP	C6-N1	2.11	1.41	1.35
3	B	305	ATP	C5'-C4'	2.10	1.57	1.51
3	A	300	ATP	C5-C6	2.08	1.46	1.41
3	B	305	ATP	C5-C6	2.06	1.46	1.41
4	B	325	TZP	CM4-C4	2.06	1.53	1.49

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	325	TZP	P1-O7-C7	7.25	137.93	118.21
4	A	320	TZP	P1-O7-C7	6.91	137.00	118.21
4	C	330	TZP	S1-C2-N3	-6.35	110.62	116.08
4	A	320	TZP	S1-C2-N3	-6.25	110.71	116.08
4	C	330	TZP	P1-O7-C7	6.16	134.97	118.21
4	B	325	TZP	S1-C2-N3	-6.01	110.92	116.08
3	C	310	ATP	O4'-C4'-C5'	-5.61	91.37	109.33
3	B	305	ATP	O4'-C4'-C5'	-5.44	91.91	109.33
3	A	300	ATP	O4'-C4'-C5'	-5.18	92.75	109.33
3	A	300	ATP	C5'-C4'-C3'	4.91	132.87	115.21
3	B	305	ATP	C5'-C4'-C3'	4.90	132.87	115.21
3	C	310	ATP	C5'-C4'-C3'	4.87	132.75	115.21
4	A	320	TZP	O3-P1-O7	4.38	118.10	106.67
4	A	320	TZP	C2-N3-C4	3.88	115.07	110.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	325	TZP	O3-P1-O7	3.83	116.65	106.67
4	C	330	TZP	C2-N3-C4	3.76	114.94	110.71
4	C	330	TZP	O3-P1-O7	3.68	116.26	106.67
4	B	325	TZP	C2-N3-C4	3.61	114.77	110.71
4	B	325	TZP	C6-C5-S1	-3.51	112.69	120.90
4	C	330	TZP	C6-C5-S1	-3.26	113.28	120.90
3	B	305	ATP	O2A-PA-O3A	3.24	116.03	107.27
3	B	305	ATP	O3A-PA-O1A	-3.19	101.10	110.70
3	B	305	ATP	C2'-C1'-N9	3.19	121.23	113.30
3	A	300	ATP	C2'-C1'-N9	3.15	121.14	113.30
3	C	310	ATP	O3A-PA-O1A	-3.12	101.31	110.70
3	B	305	ATP	O4'-C1'-N9	-3.11	102.12	108.09
3	C	310	ATP	O2A-PA-O3A	3.03	115.46	107.27
4	A	320	TZP	C6-C5-S1	-2.96	113.98	120.90
4	A	320	TZP	O3-P1-O2	-2.96	99.32	110.83
4	C	330	TZP	O3-P1-O2	-2.95	99.34	110.83
3	C	310	ATP	O4'-C1'-N9	-2.91	102.51	108.09
3	A	300	ATP	O2A-PA-O3A	2.88	115.06	107.27
3	A	300	ATP	O4'-C1'-N9	-2.79	102.73	108.09
3	A	300	ATP	O3A-PA-O1A	-2.78	102.35	110.70
4	B	325	TZP	O3-P1-O2	-2.74	100.17	110.83
3	B	305	ATP	C6-C5-C4	-2.68	113.52	117.18
3	C	310	ATP	C2'-C1'-N9	2.59	119.75	113.30
3	A	300	ATP	C6-C5-C4	-2.48	113.79	117.18
3	C	310	ATP	C6-C5-C4	-2.38	113.94	117.18
3	A	300	ATP	O3G-PG-O2G	-2.37	98.91	107.80
3	B	305	ATP	O3G-PG-O2G	-2.25	99.35	107.80
3	C	310	ATP	O3G-PG-O2G	-2.22	99.48	107.80
3	C	310	ATP	N3-C2-N1	-2.18	125.28	128.58
4	C	330	TZP	O1-P1-O7	2.15	112.26	106.67
3	B	305	ATP	O3B-PB-O1B	2.03	116.82	110.70

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	300	ATP	C5'-O5'-PA-O1A
3	A	300	ATP	C5'-O5'-PA-O3A
3	B	305	ATP	C5'-O5'-PA-O1A
3	B	305	ATP	C5'-O5'-PA-O3A
3	C	310	ATP	C5'-O5'-PA-O1A
3	C	310	ATP	C5'-O5'-PA-O3A

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Mol	Chain	Res	Type	Atoms
4	A	320	TZP	C5-C6-C7-O7
4	B	325	TZP	C7-O7-P1-O1
4	C	330	TZP	C5-C6-C7-O7
3	A	300	ATP	C3'-C4'-C5'-O5'
3	B	305	ATP	C3'-C4'-C5'-O5'
3	C	310	ATP	C3'-C4'-C5'-O5'
3	A	300	ATP	O4'-C4'-C5'-O5'
3	B	305	ATP	O4'-C4'-C5'-O5'
3	C	310	ATP	O4'-C4'-C5'-O5'
3	A	300	ATP	PA-O3A-PB-O3B
3	B	305	ATP	PA-O3A-PB-O3B
3	C	310	ATP	PA-O3A-PB-O3B
4	B	325	TZP	C5-C6-C7-O7
3	A	300	ATP	PG-O3B-PB-O2B
3	B	305	ATP	PG-O3B-PB-O2B
3	C	310	ATP	PG-O3B-PB-O2B
3	A	300	ATP	C5'-O5'-PA-O2A
3	B	305	ATP	C5'-O5'-PA-O2A
3	C	310	ATP	C5'-O5'-PA-O2A
4	B	325	TZP	C7-O7-P1-O2
3	A	300	ATP	PG-O3B-PB-O1B
3	A	300	ATP	PA-O3A-PB-O1B
3	B	305	ATP	PG-O3B-PB-O1B
3	B	305	ATP	PA-O3A-PB-O1B
3	C	310	ATP	PG-O3B-PB-O1B
3	C	310	ATP	PA-O3A-PB-O1B

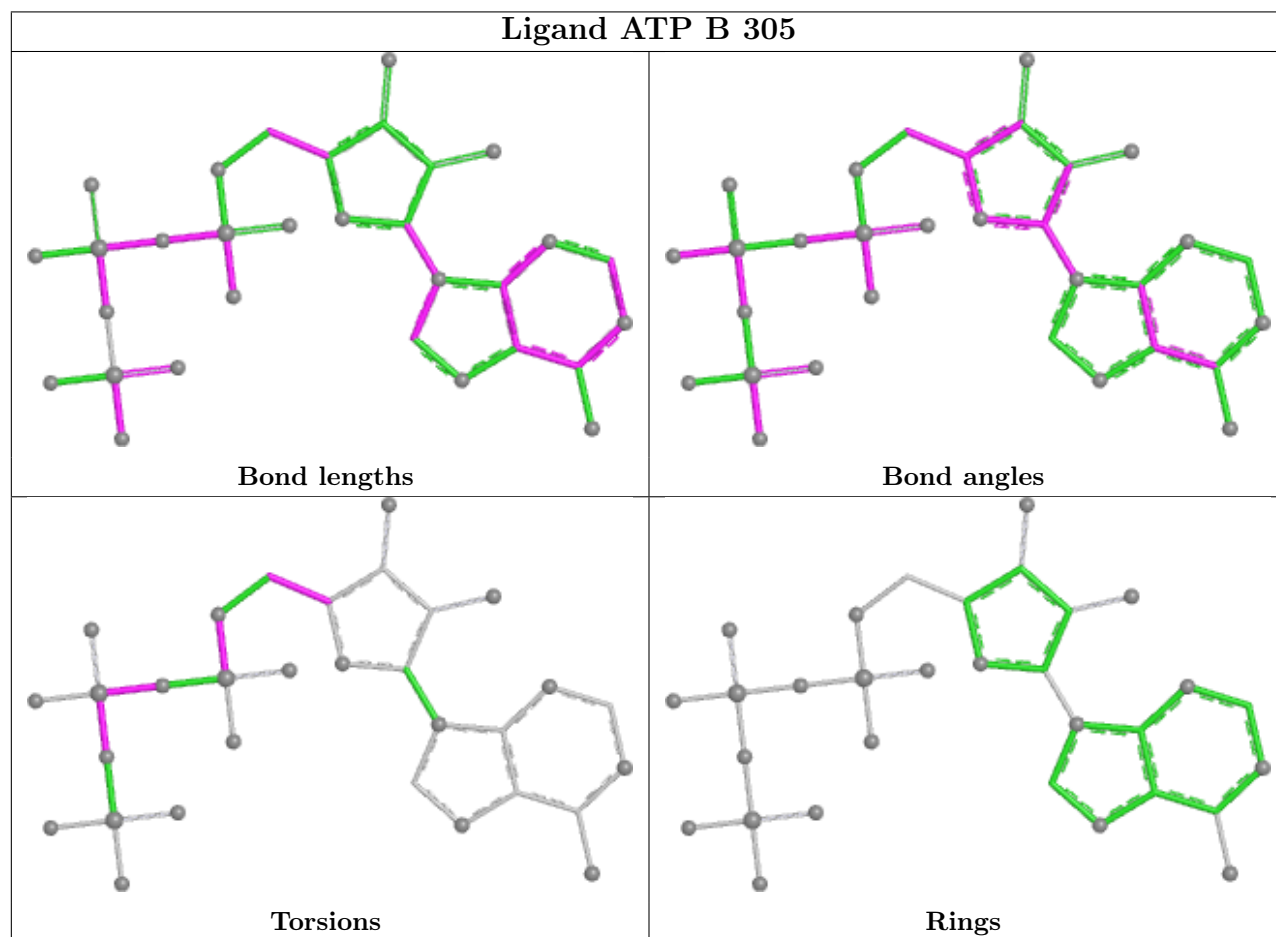
There are no ring outliers.

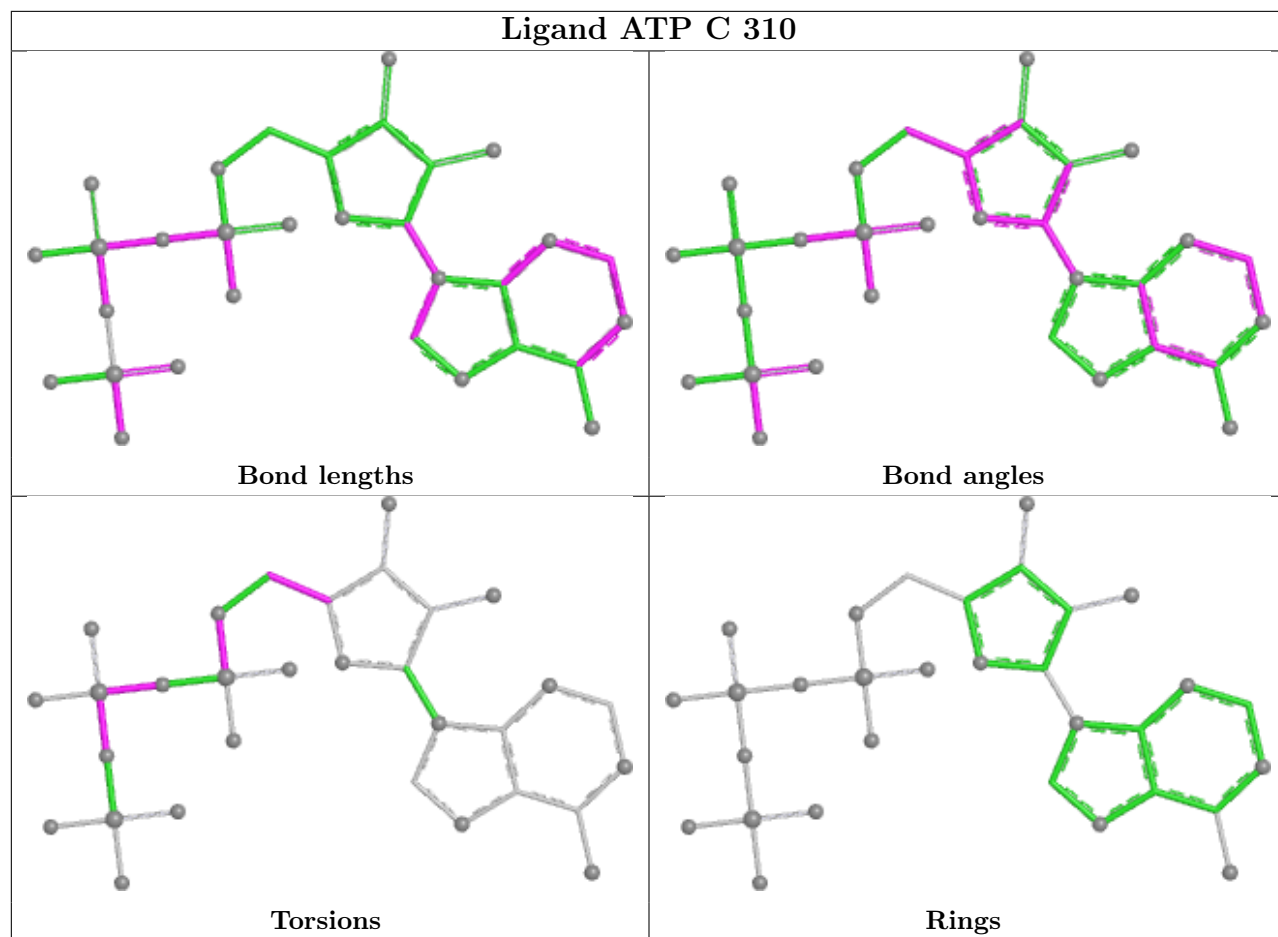
2 monomers are involved in 2 short contacts:

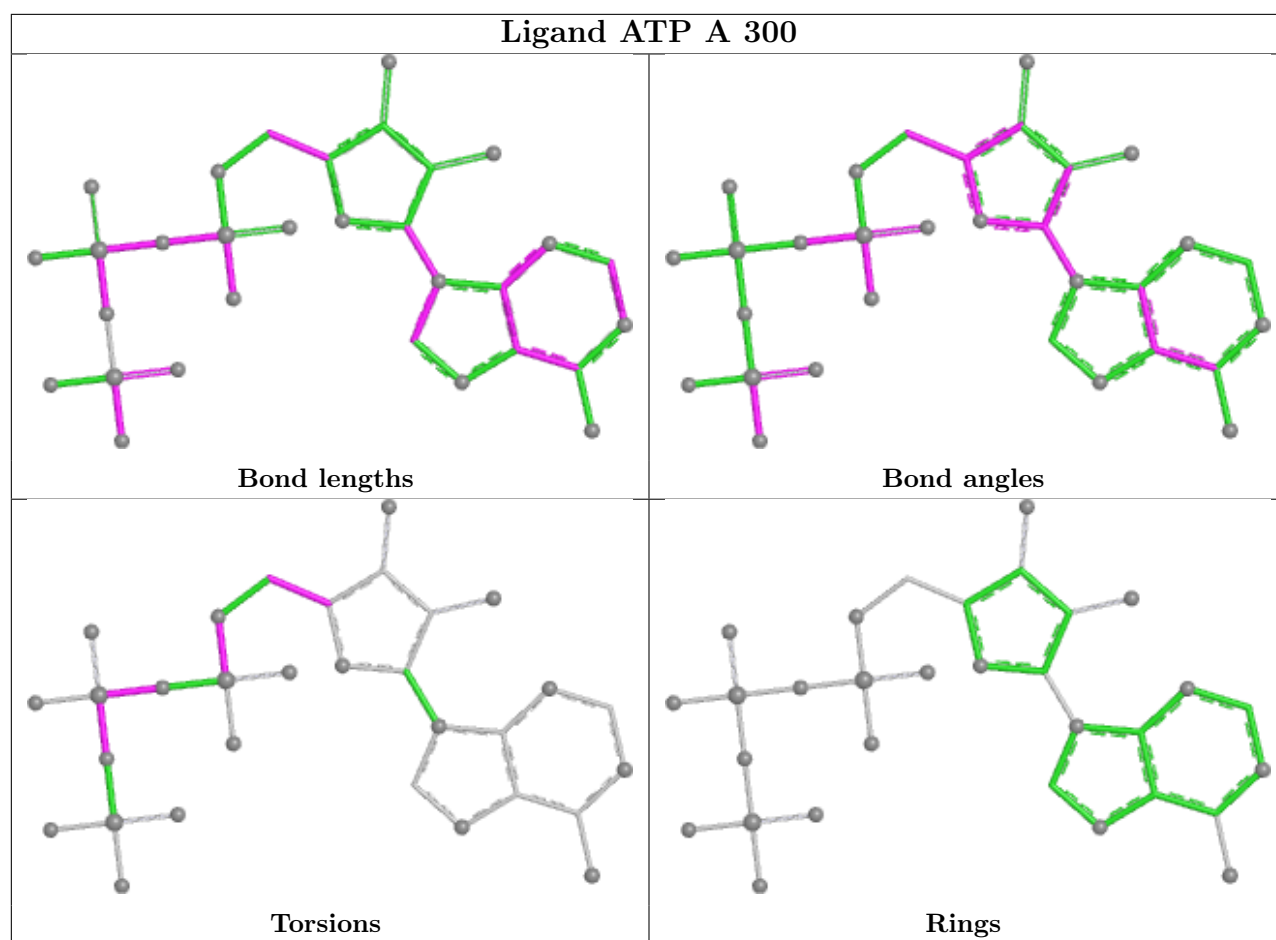
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	320	TZP	1	0
3	C	310	ATP	1	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	258/284 (90%)	-0.15	1 (0%) 88 86	16, 30, 49, 88	0
1	B	258/284 (90%)	-0.14	2 (0%) 82 80	15, 29, 50, 66	0
1	C	274/284 (96%)	-0.06	7 (2%) 57 52	14, 30, 54, 63	0
All	All	790/852 (92%)	-0.12	10 (1%) 75 71	14, 29, 51, 88	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	134	THR	2.8
1	B	271	VAL	2.7
1	C	137	LEU	2.5
1	C	236	THR	2.4
1	C	134	THR	2.3
1	C	132	GLY	2.3
1	A	272	SER	2.3
1	C	-3	HIS	2.1
1	C	148	GLY	2.0
1	C	135	ASP	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 ⓘ

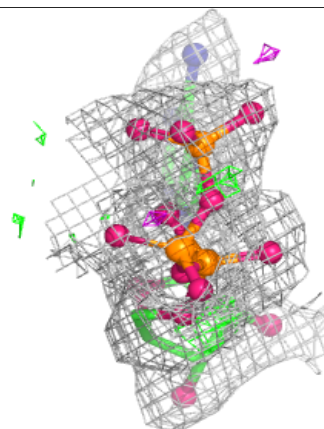
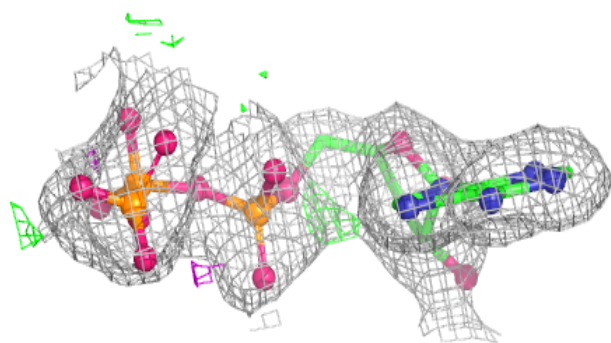
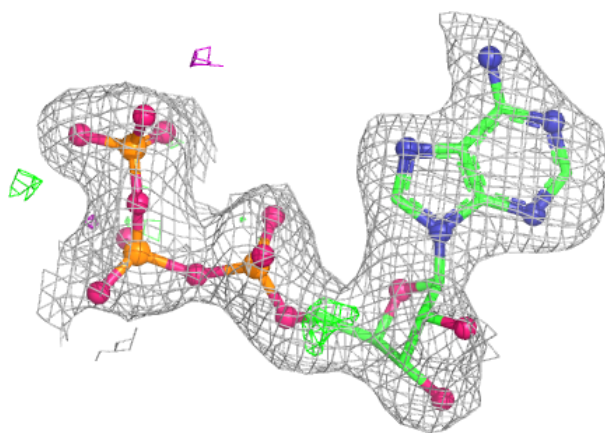
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MG	C	355	1/1	0.74	0.14	40,40,40,40	0
2	MG	A	365	1/1	0.76	0.17	46,46,46,46	0
2	MG	B	360	1/1	0.77	0.16	48,48,48,48	0
2	MG	A	340	1/1	0.91	0.08	32,32,32,32	0
3	ATP	B	305	31/31	0.92	0.09	28,36,42,42	0
2	MG	C	350	1/1	0.94	0.08	41,41,41,41	0
3	ATP	C	310	31/31	0.94	0.08	27,31,33,35	0
2	MG	B	345	1/1	0.95	0.07	22,22,22,22	0
3	ATP	A	300	31/31	0.95	0.08	35,40,44,47	0
4	TZP	B	325	13/13	0.96	0.07	17,22,34,35	0
4	TZP	C	330	13/13	0.97	0.06	20,22,31,32	0
5	SO4	B	370	5/5	0.97	0.08	32,32,34,35	0
4	TZP	A	320	13/13	0.98	0.05	17,20,33,33	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.

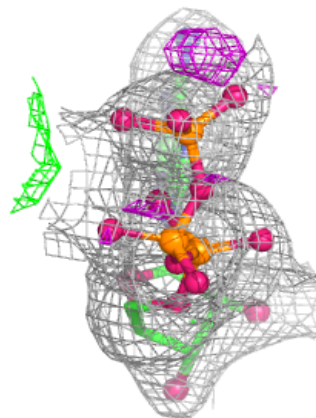
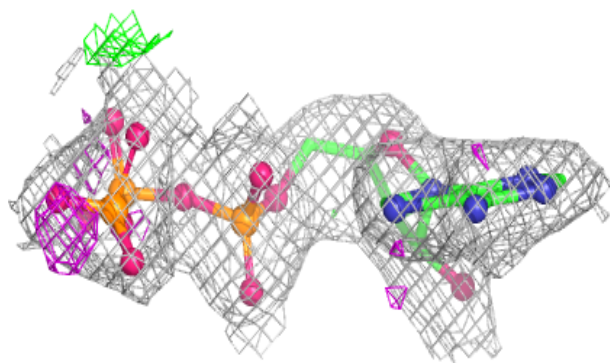
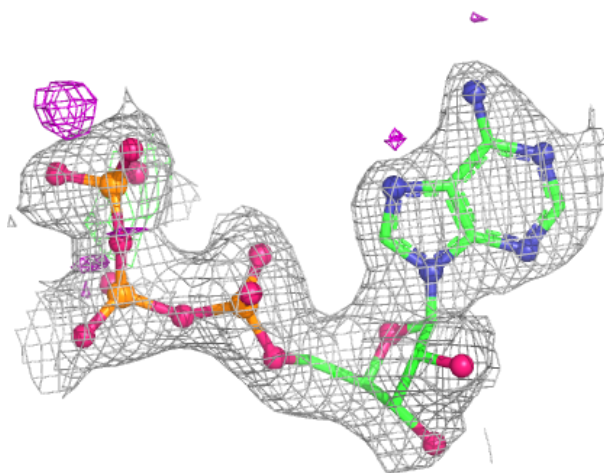
Electron density around ATP B 305:

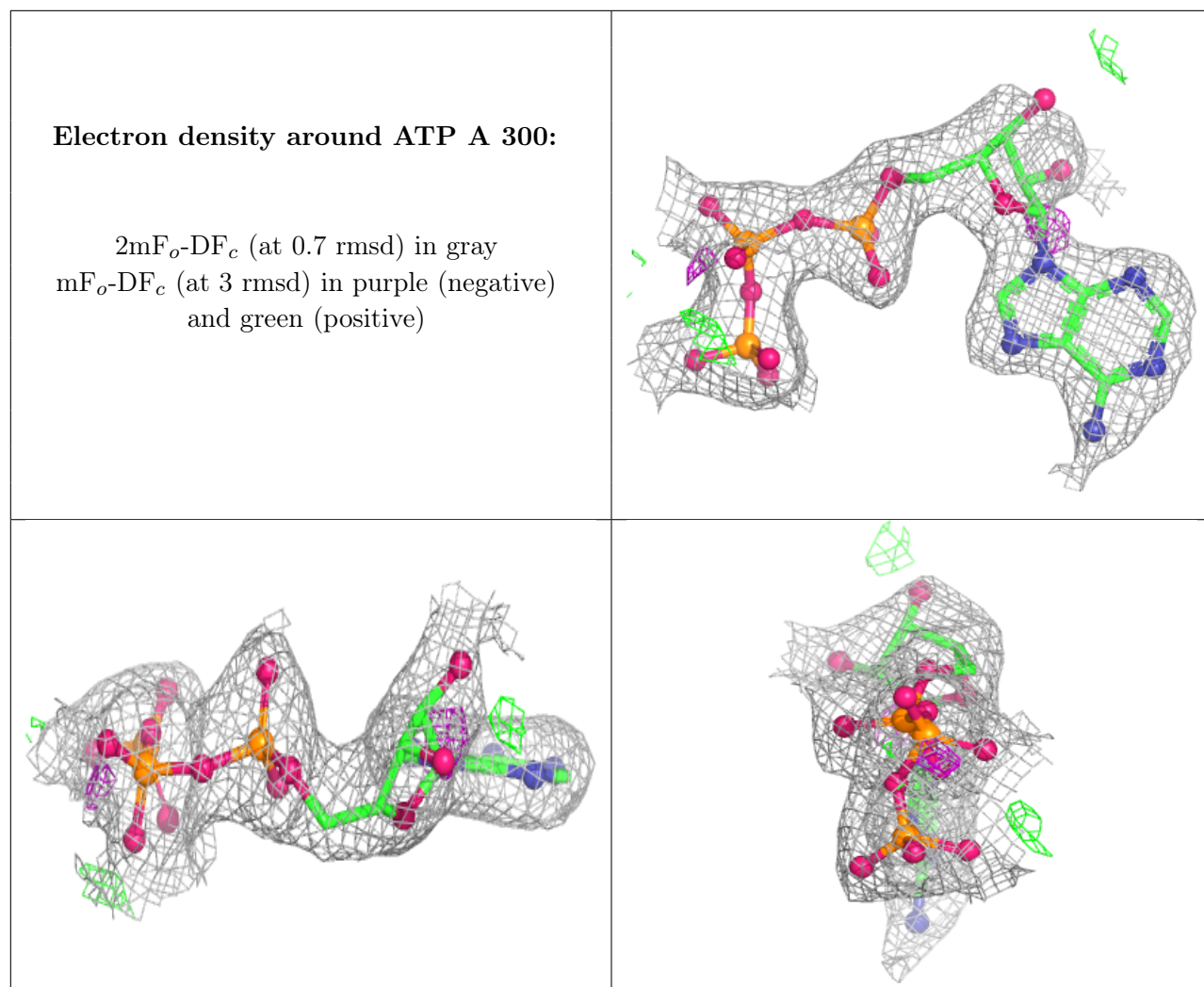
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ATP C 310:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.