



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:29 PM UTC

PDB ID : 2NPI / pdb_00002npi
Title : Clp1-ATP-Pcf11 complex
Authors : Noble, C.G.; Beuth, B.; Taylor, I.A.
Deposited on : 2006-10-27
Resolution : 2.95 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

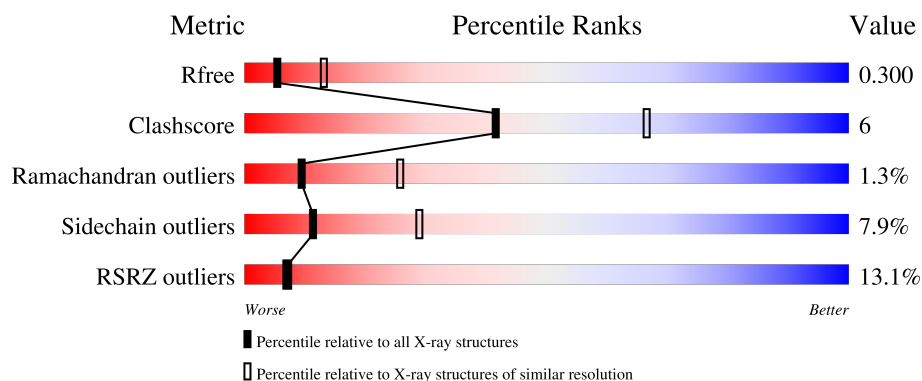
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	460	20% 74% 16% • 7%
1	B	460	4% 72% 20% • 7%
2	C	110	5% 17% • 80%
2	D	110	2% 17% 5% 77%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7437 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 Protein CLP1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	428	Total	C	N	O	S	Se	0	0	0
			3412	2193	571	636	6	6			
1	B	428	Total	C	N	O	S	Se	0	0	0
			3412	2193	571	636	6	6			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	MSE	-	modified residue	UNP Q08685
A	-13	GLY	-	cloning artifact	UNP Q08685
A	-12	SER	-	cloning artifact	UNP Q08685
A	-11	SER	-	cloning artifact	UNP Q08685
A	-10	HIS	-	expression tag	UNP Q08685
A	-9	HIS	-	expression tag	UNP Q08685
A	-8	HIS	-	expression tag	UNP Q08685
A	-7	HIS	-	expression tag	UNP Q08685
A	-6	HIS	-	expression tag	UNP Q08685
A	-5	HIS	-	expression tag	UNP Q08685
A	-4	SER	-	cloning artifact	UNP Q08685
A	-3	GLN	-	cloning artifact	UNP Q08685
A	-2	ASP	-	cloning artifact	UNP Q08685
A	-1	PRO	-	cloning artifact	UNP Q08685
A	0	ASN	-	cloning artifact	UNP Q08685
A	1	SER	-	cloning artifact	UNP Q08685
A	103	MSE	MET	modified residue	UNP Q08685
A	112	MSE	MET	modified residue	UNP Q08685
A	118	MSE	MET	modified residue	UNP Q08685
A	206	MSE	MET	modified residue	UNP Q08685
A	277	MSE	MET	modified residue	UNP Q08685
A	436	MSE	MET	modified residue	UNP Q08685
B	-14	MSE	-	modified residue	UNP Q08685
B	-13	GLY	-	cloning artifact	UNP Q08685
B	-12	SER	-	cloning artifact	UNP Q08685

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-11	SER	-	cloning artifact	UNP Q08685
B	-10	HIS	-	expression tag	UNP Q08685
B	-9	HIS	-	expression tag	UNP Q08685
B	-8	HIS	-	expression tag	UNP Q08685
B	-7	HIS	-	expression tag	UNP Q08685
B	-6	HIS	-	expression tag	UNP Q08685
B	-5	HIS	-	expression tag	UNP Q08685
B	-4	SER	-	cloning artifact	UNP Q08685
B	-3	GLN	-	cloning artifact	UNP Q08685
B	-2	ASP	-	cloning artifact	UNP Q08685
B	-1	PRO	-	cloning artifact	UNP Q08685
B	0	ASN	-	cloning artifact	UNP Q08685
B	1	SER	-	cloning artifact	UNP Q08685
B	103	MSE	MET	modified residue	UNP Q08685
B	112	MSE	MET	modified residue	UNP Q08685
B	118	MSE	MET	modified residue	UNP Q08685
B	206	MSE	MET	modified residue	UNP Q08685
B	277	MSE	MET	modified residue	UNP Q08685
B	436	MSE	MET	modified residue	UNP Q08685

- Molecule 2 is a protein called Protein PCF11.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	22	Total	C	N	O	0	0	0
			189	120	31	38			
2	D	25	Total	C	N	O	0	0	0
			212	133	36	43			

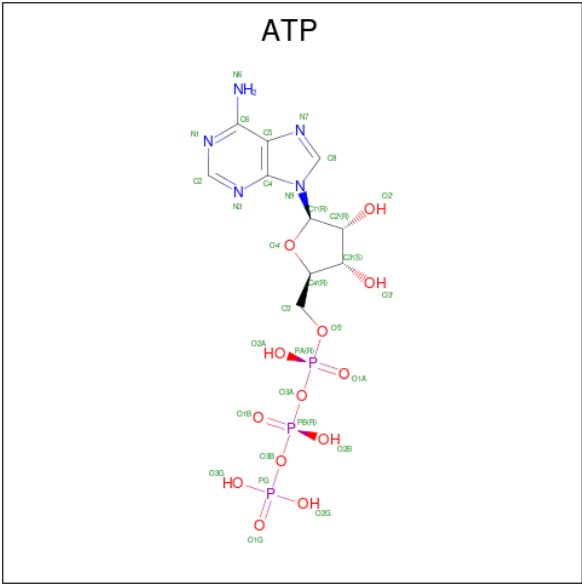
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	560	MSE	MET	modified residue	UNP P39081
D	560	MSE	MET	modified residue	UNP P39081

- Molecule 3 is MAGNESIUM ION (CCD ID: 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		

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



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

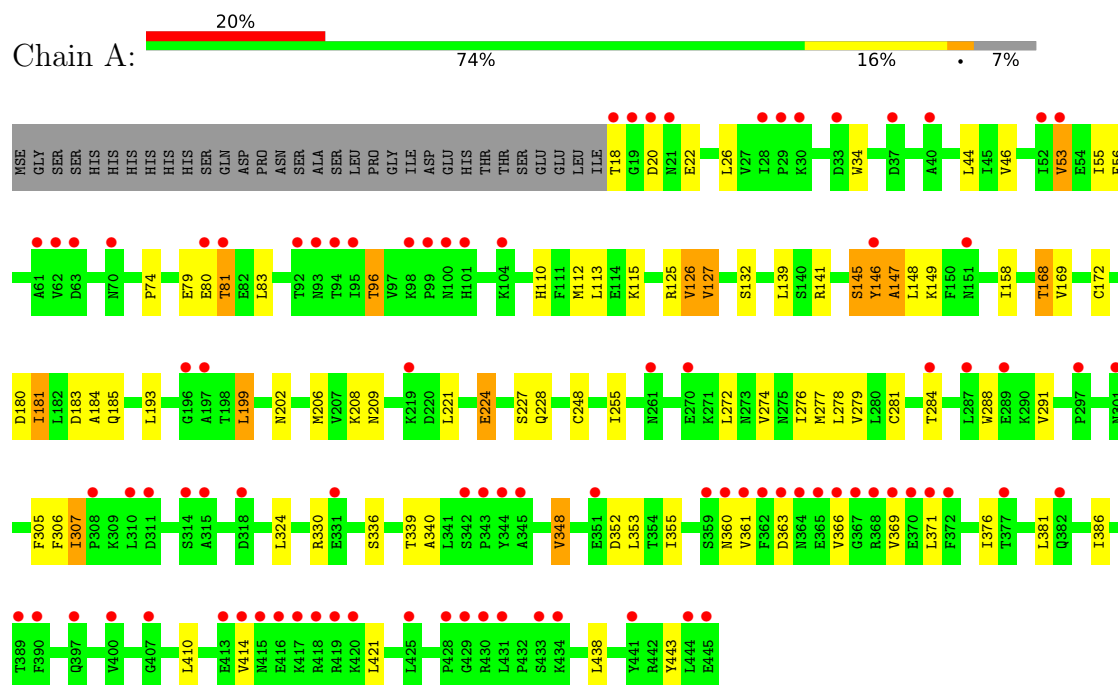
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	57	Total	O	0	0
			57	57		
5	B	77	Total	O	0	0
			77	77		
5	C	7	Total	O	0	0
			7	7		
5	D	7	Total	O	0	0
			7	7		

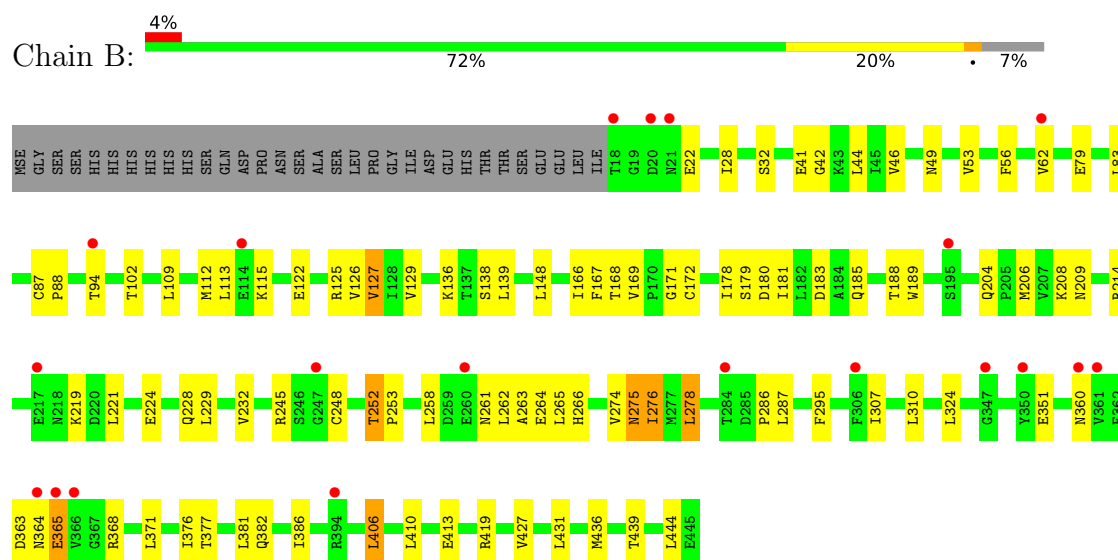
3 Residue-property plots

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.

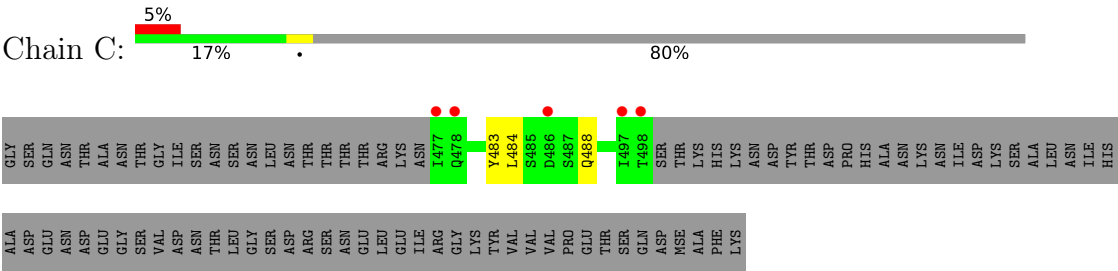
• Molecule 1: Protein CLP1



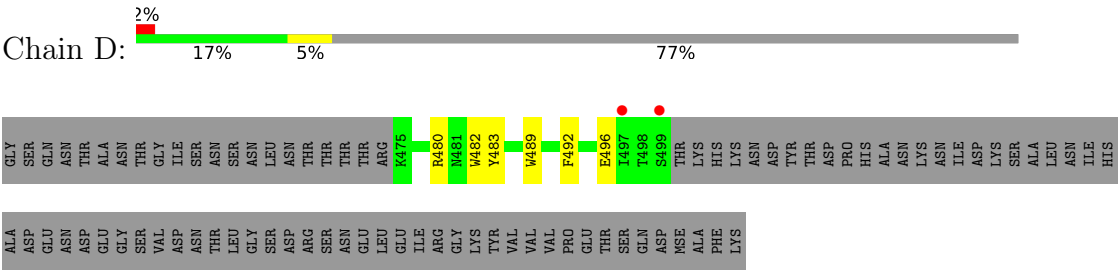
• Molecule 1: Protein CLP1



● Molecule 2: Protein PCF11



● Molecule 2: Protein PCF11



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.97Å 94.99Å 181.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.95 15.00 – 2.95	Depositor EDS
% Data completeness (in resolution range)	90.0 (15.00-2.95) 89.2 (15.00-2.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 2.96Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.248 , 0.303 0.257 , 0.300	Depositor DCC
R_{free} test set	1493 reflections (4.55%)	wwPDB-VP
Wilson B-factor (Å ²)	53.1	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 63.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	7437	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% 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, 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	0.43	0/3483	0.82	2/4729 (0.0%)
1	B	0.43	0/3483	0.80	1/4729 (0.0%)
2	C	0.37	0/194	0.70	0/263
2	D	0.44	0/217	0.74	0/293
All	All	0.42	0/7377	0.80	3/10014 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	147	ALA	N-CA-C	-6.15	105.26	112.89
1	B	172	CYS	N-CA-C	5.77	118.96	109.85
1	A	284	THR	N-CA-C	-5.34	107.43	114.31

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	3412	0	3451	44	0
1	B	3412	0	3451	48	0
2	C	189	0	168	2	0
2	D	212	0	192	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	31	0	12	0	0
4	B	31	0	12	0	0
5	A	57	0	0	0	0
5	B	77	0	0	7	0
5	C	7	0	0	0	0
5	D	7	0	0	0	0
All	All	7437	0	7286	93	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 6.

All (93) 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:A:148:LEU:HD11	1:A:180:ASP:HA	1.46	0.95
1:B:109:LEU:HD21	1:B:278:LEU:HD11	1.55	0.89
1:B:148:LEU:HD11	1:B:180:ASP:HA	1.56	0.85
1:B:125:ARG:H	1:B:275:ASN:HD21	1.26	0.84
1:B:376:ILE:HG23	1:B:381:LEU:HD12	1.63	0.78
1:A:172:CYS:HB3	1:A:209:ASN:HA	1.69	0.73
1:A:168:THR:HG22	1:A:169:VAL:H	1.53	0.73
1:B:127:VAL:HG22	1:B:274:VAL:HG11	1.73	0.69
1:B:180:ASP:HB2	5:B:2508:HOH:O	1.93	0.67
1:B:406:LEU:HD22	2:D:482:TRP:HB3	1.82	0.60
1:A:278:LEU:HD22	1:A:307:ILE:HD11	1.84	0.59
1:A:360:ASN:HB2	1:A:363:ASP:HB2	1.84	0.59
1:A:168:THR:HG23	2:C:483:TYR:OH	2.01	0.59
1:A:53:VAL:HG13	1:A:81:THR:HG21	1.84	0.58
1:A:376:ILE:HG23	1:A:381:LEU:HD12	1.86	0.57
1:B:360:ASN:HB2	1:B:363:ASP:HB2	1.87	0.56
1:B:214:ARG:HD3	5:B:2527:HOH:O	2.05	0.56
1:A:127:VAL:HG22	1:A:274:VAL:HG11	1.87	0.56
1:A:279:VAL:HG21	1:A:291:VAL:HG11	1.87	0.55
1:A:34:TRP:O	1:A:96:THR:O	2.25	0.55
1:A:278:LEU:HD23	1:A:305:PHE:HB2	1.87	0.55
1:A:279:VAL:HG12	1:A:281:CYS:HB2	1.87	0.55
1:A:353:LEU:HD23	1:A:355:ILE:HD11	1.87	0.55
1:B:168:THR:HG23	2:D:483:TYR:OH	2.06	0.55
1:A:110:HIS:HE1	1:A:147:ALA:HA	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:427:VAL:HG22	2:D:480:ARG:O	2.07	0.53
2:C:484:LEU:HD22	2:C:488:GLN:HE21	1.74	0.53
1:B:382:GLN:HE21	1:B:413:GLU:HA	1.73	0.52
1:A:183:ASP:OD1	1:A:185:GLN:HG2	2.10	0.52
1:B:125:ARG:N	1:B:275:ASN:HD21	2.02	0.52
1:B:263:ALA:HA	1:B:266:HIS:HD2	1.75	0.51
1:B:129:VAL:HA	5:B:2509:HOH:O	2.11	0.50
1:B:208:LYS:HD2	1:B:228:GLN:HE21	1.76	0.50
1:B:360:ASN:HB3	1:B:363:ASP:H	1.75	0.50
1:B:46:VAL:HG13	1:B:83:LEU:HD13	1.94	0.50
1:A:336:SER:H	1:A:339:THR:HB	1.78	0.49
1:A:56:PHE:CZ	1:A:74:PRO:HD2	2.49	0.48
1:A:125:ARG:HG2	1:A:248:CYS:HB3	1.95	0.48
1:A:46:VAL:HG13	1:A:83:LEU:HD13	1.95	0.48
1:A:224:GLU:HG3	1:A:369:VAL:HG11	1.96	0.48
1:A:20:ASP:OD2	1:A:22:GLU:HG2	2.14	0.48
1:A:410:LEU:CD1	1:A:443:TYR:HD2	2.26	0.47
1:B:208:LYS:HD2	1:B:228:GLN:HG3	1.95	0.47
1:B:148:LEU:HD11	1:B:180:ASP:CA	2.36	0.47
1:B:166:ILE:HG22	1:B:167:PHE:CD1	2.49	0.47
1:B:56:PHE:CE2	1:B:138:SER:HB3	2.50	0.46
1:A:208:LYS:HD2	1:A:228:GLN:HG3	1.98	0.46
1:B:439:THR:HG21	5:B:2517:HOH:O	2.14	0.46
1:B:406:LEU:CD2	2:D:482:TRP:HB3	2.46	0.46
1:B:204:GLN:HG3	2:D:492:PHE:CE2	2.51	0.46
1:A:348:VAL:O	1:A:421:LEU:N	2.48	0.45
1:B:102:THR:HG21	1:B:310:LEU:HA	1.99	0.45
1:A:110:HIS:CE1	1:A:147:ALA:HA	2.52	0.45
1:B:206:MSE:HB2	2:D:489:TRP:CE2	2.52	0.45
1:B:381:LEU:HD22	1:B:386:ILE:HD11	1.98	0.45
1:B:102:THR:HB	5:B:2544:HOH:O	2.18	0.44
1:B:252:THR:HG22	1:B:253:PRO:HD2	1.99	0.44
1:B:262:LEU:HB3	1:B:265:LEU:HD12	1.99	0.44
1:B:109:LEU:HD22	1:B:307:ILE:HD11	2.00	0.44
1:A:146:TYR:CB	1:A:149:LYS:HB3	2.48	0.44
1:B:87:CYS:HA	1:B:88:PRO:HD3	1.89	0.44
1:B:126:VAL:HG13	1:B:276:ILE:HG12	2.01	0.43
1:B:168:THR:HG22	1:B:169:VAL:HG22	2.00	0.43
1:A:141:ARG:O	1:A:145:SER:HB2	2.19	0.43
1:B:28:ILE:HD11	1:B:32:SER:O	2.18	0.42
1:B:360:ASN:H	1:B:364:ASN:HD22	1.65	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:THR:O	1:B:189:TRP:HB2	2.18	0.42
1:B:183:ASP:OD1	1:B:185:GLN:HG2	2.19	0.42
1:A:110:HIS:CE1	1:A:146:TYR:O	2.73	0.42
1:B:148:LEU:HD13	1:B:178:ILE:HG22	2.02	0.42
1:A:79:GLU:HG3	1:A:80:GLU:H	1.85	0.42
1:A:148:LEU:CD1	1:A:180:ASP:HA	2.32	0.42
1:B:125:ARG:HA	1:B:248:CYS:HB3	2.02	0.42
1:A:146:TYR:HB3	1:A:149:LYS:HB3	2.01	0.42
1:A:127:VAL:HG23	1:A:277:MSE:HG3	2.02	0.41
1:B:386:ILE:CG2	1:B:436:MSE:HB3	2.50	0.41
1:A:181:ILE:H	1:A:181:ILE:HG13	1.58	0.41
1:A:55:ILE:HD12	1:A:184:ALA:HB1	2.03	0.41
1:B:148:LEU:CD1	1:B:180:ASP:HA	2.41	0.41
1:A:126:VAL:HA	1:A:276:ILE:O	2.21	0.41
1:A:199:LEU:H	1:A:199:LEU:HD13	1.86	0.41
1:A:348:VAL:HG13	1:A:352:ASP:HB2	2.03	0.41
1:B:351:GLU:CD	1:B:351:GLU:H	2.29	0.41
1:A:288:TRP:CZ3	1:A:306:PHE:HB3	2.56	0.41
1:A:414:VAL:HA	1:A:421:LEU:HD23	2.04	0.40
1:B:258:LEU:HD22	1:B:264:GLU:HG2	2.04	0.40
1:A:193:LEU:HD23	1:A:202:ASN:HB2	2.02	0.40
1:B:171:GLY:HA3	5:B:2517:HOH:O	2.20	0.40
1:B:245:ARG:HD3	5:B:2513:HOH:O	2.20	0.40
1:A:125:ARG:HG2	1:A:248:CYS:CB	2.51	0.40
1:A:126:VAL:HG13	1:A:276:ILE:HG12	2.03	0.40
1:B:265:LEU:HB2	1:B:295:PHE:CZ	2.57	0.40
1:A:158:ILE:HD13	1:A:272:LEU:HD22	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	426/460 (93%)	380 (89%)	42 (10%)	4 (1%)	14	34
1	B	426/460 (93%)	394 (92%)	24 (6%)	8 (2%)	6	17
2	C	20/110 (18%)	20 (100%)	0	0	100	100
2	D	23/110 (21%)	23 (100%)	0	0	100	100
All	All	895/1140 (78%)	817 (91%)	66 (7%)	12 (1%)	9	26

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	419	ARG
1	B	42	GLY
1	B	62	VAL
1	B	261	ASN
1	A	96	THR
1	A	340	ALA
1	A	81	THR
1	B	41	GLU
1	B	209	ASN
1	B	365	GLU
1	B	286	PRO
1	A	366	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	385/407 (95%)	356 (92%)	29 (8%)	12	31
1	B	385/407 (95%)	351 (91%)	34 (9%)	9	24
2	C	20/97 (21%)	20 (100%)	0	100	100
2	D	23/97 (24%)	22 (96%)	1 (4%)	26	52
All	All	813/1008 (81%)	749 (92%)	64 (8%)	11	29

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	THR
1	A	26	LEU
1	A	44	LEU
1	A	53	VAL
1	A	112	MSE
1	A	113	LEU
1	A	115	LYS
1	A	126	VAL
1	A	127	VAL
1	A	132	SER
1	A	139	LEU
1	A	145	SER
1	A	146	TYR
1	A	168	THR
1	A	181	ILE
1	A	199	LEU
1	A	206	MSE
1	A	221	LEU
1	A	224	GLU
1	A	227	SER
1	A	255	ILE
1	A	307	ILE
1	A	324	LEU
1	A	330	ARG
1	A	348	VAL
1	A	361	VAL
1	A	371	LEU
1	A	386	ILE
1	A	438	LEU
1	B	22	GLU
1	B	44	LEU
1	B	49	ASN
1	B	53	VAL
1	B	79	GLU
1	B	94	THR
1	B	112	MSE
1	B	113	LEU
1	B	115	LYS
1	B	122	GLU
1	B	127	VAL
1	B	136	LYS
1	B	139	LEU
1	B	179	SER

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Mol	Chain	Res	Type
1	B	181	ILE
1	B	219	LYS
1	B	221	LEU
1	B	224	GLU
1	B	229	LEU
1	B	232	VAL
1	B	252	THR
1	B	275	ASN
1	B	276	ILE
1	B	278	LEU
1	B	287	LEU
1	B	324	LEU
1	B	365	GLU
1	B	368	ARG
1	B	371	LEU
1	B	377	THR
1	B	406	LEU
1	B	410	LEU
1	B	431	LEU
1	B	444	LEU
2	D	496	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	GLN
1	A	110	HIS
1	A	120	ASN
1	A	133	GLN
1	A	154	GLN
1	A	163	GLN
1	A	209	ASN
1	A	238	HIS
1	A	242	GLN
1	A	273	ASN
1	A	302	ASN
1	A	364	ASN
1	A	380	ASN
1	B	49	ASN
1	B	120	ASN
1	B	154	GLN
1	B	159	ASN

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Mol	Chain	Res	Type
1	B	228	GLN
1	B	266	HIS
1	B	267	HIS
1	B	273	ASN
1	B	275	ASN
1	B	364	ASN
1	B	380	ASN
1	B	382	GLN
2	C	488	GLN
2	D	476	ASN

5.3.3 RNA [i](#)

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	ATP	B	1600	3	32,33,33	1.52	6 (18%)	48,52,52	1.67	8 (16%)
4	ATP	A	600	3	32,33,33	1.39	4 (12%)	48,52,52	1.72	9 (18%)

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
4	ATP	B	1600	3	-	3/22/38/38	0/3/3/3
4	ATP	A	600	3	-	0/22/38/38	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1600	ATP	C5-C4	4.95	1.47	1.39
4	A	600	ATP	C5-C4	4.71	1.47	1.39
4	B	1600	ATP	C5-C6	2.85	1.48	1.41
4	A	600	ATP	C5-C6	2.80	1.48	1.41
4	B	1600	ATP	PA-O3A	2.51	1.62	1.59
4	B	1600	ATP	C8-N7	2.42	1.36	1.31
4	A	600	ATP	C5-N7	-2.41	1.34	1.39
4	A	600	ATP	C8-N7	2.34	1.36	1.31
4	B	1600	ATP	C5-N7	-2.29	1.34	1.39
4	B	1600	ATP	PB-O3A	2.21	1.61	1.59

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	600	ATP	C5-C4-N3	-5.54	119.09	126.72
4	B	1600	ATP	C5-C4-N3	-5.53	119.10	126.72
4	A	600	ATP	N3-C4-N9	4.43	134.69	127.17
4	B	1600	ATP	N3-C4-N9	4.37	134.60	127.17
4	A	600	ATP	C2-N3-C4	3.75	120.99	111.83
4	A	600	ATP	N3-C2-N1	-3.64	123.08	128.58
4	B	1600	ATP	C2-N3-C4	3.57	120.56	111.83
4	A	600	ATP	C4-C5-N7	-3.40	106.69	110.58
4	B	1600	ATP	C4-C5-N7	-3.33	106.78	110.58
4	B	1600	ATP	N3-C2-N1	-3.24	123.67	128.58
4	A	600	ATP	C4-N9-C8	2.75	108.62	105.74
4	B	1600	ATP	C4-N9-C8	2.48	108.35	105.74
4	A	600	ATP	C5-N7-C8	2.48	107.35	103.45
4	B	1600	ATP	C5-N7-C8	2.39	107.21	103.45
4	A	600	ATP	C6-C5-N7	2.20	136.34	132.09
4	B	1600	ATP	C6-C5-N7	2.09	136.12	132.09
4	A	600	ATP	C2-N1-C6	2.02	122.05	118.73

There are no chirality outliers.

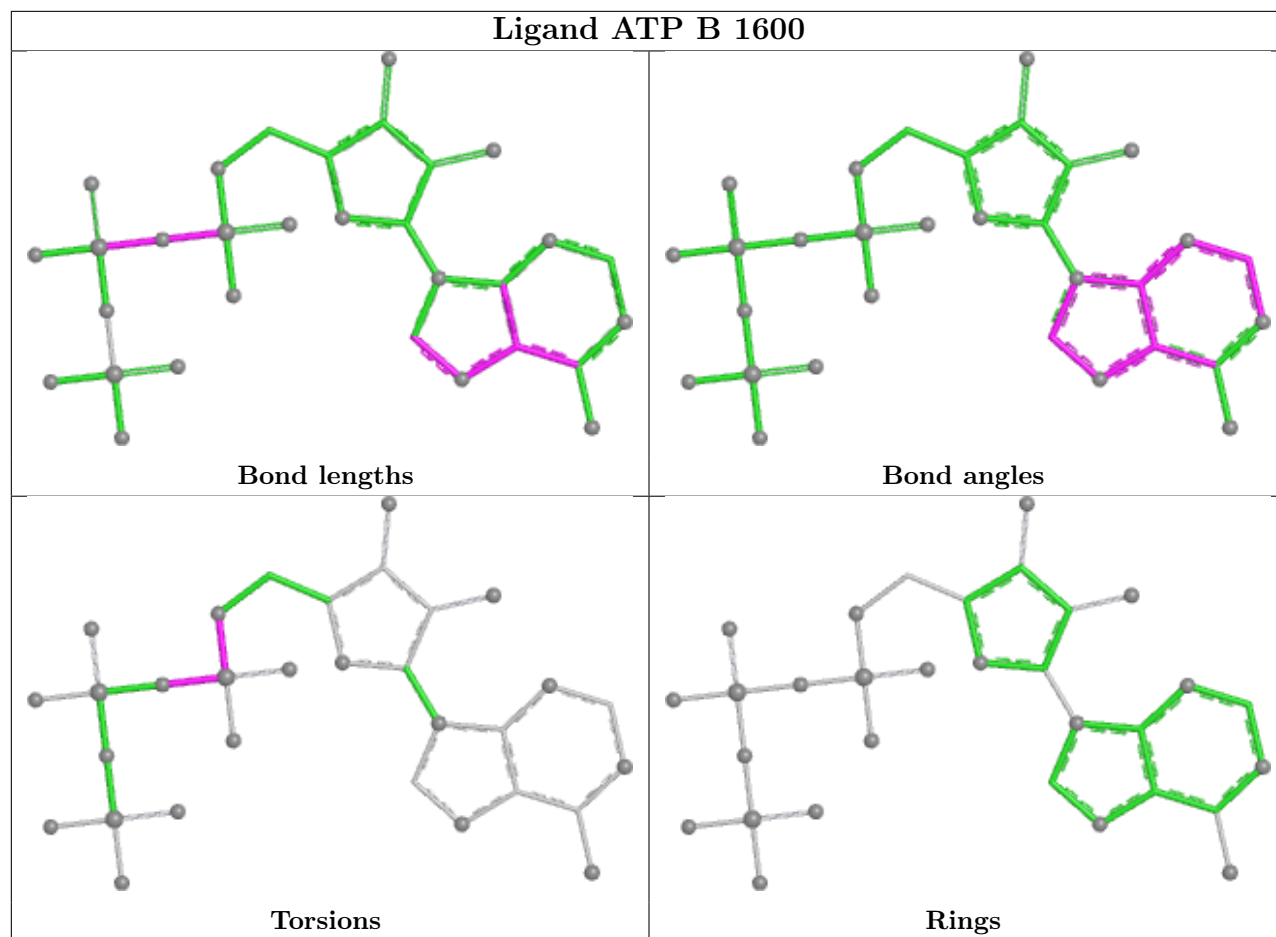
All (3) torsion outliers are listed below:

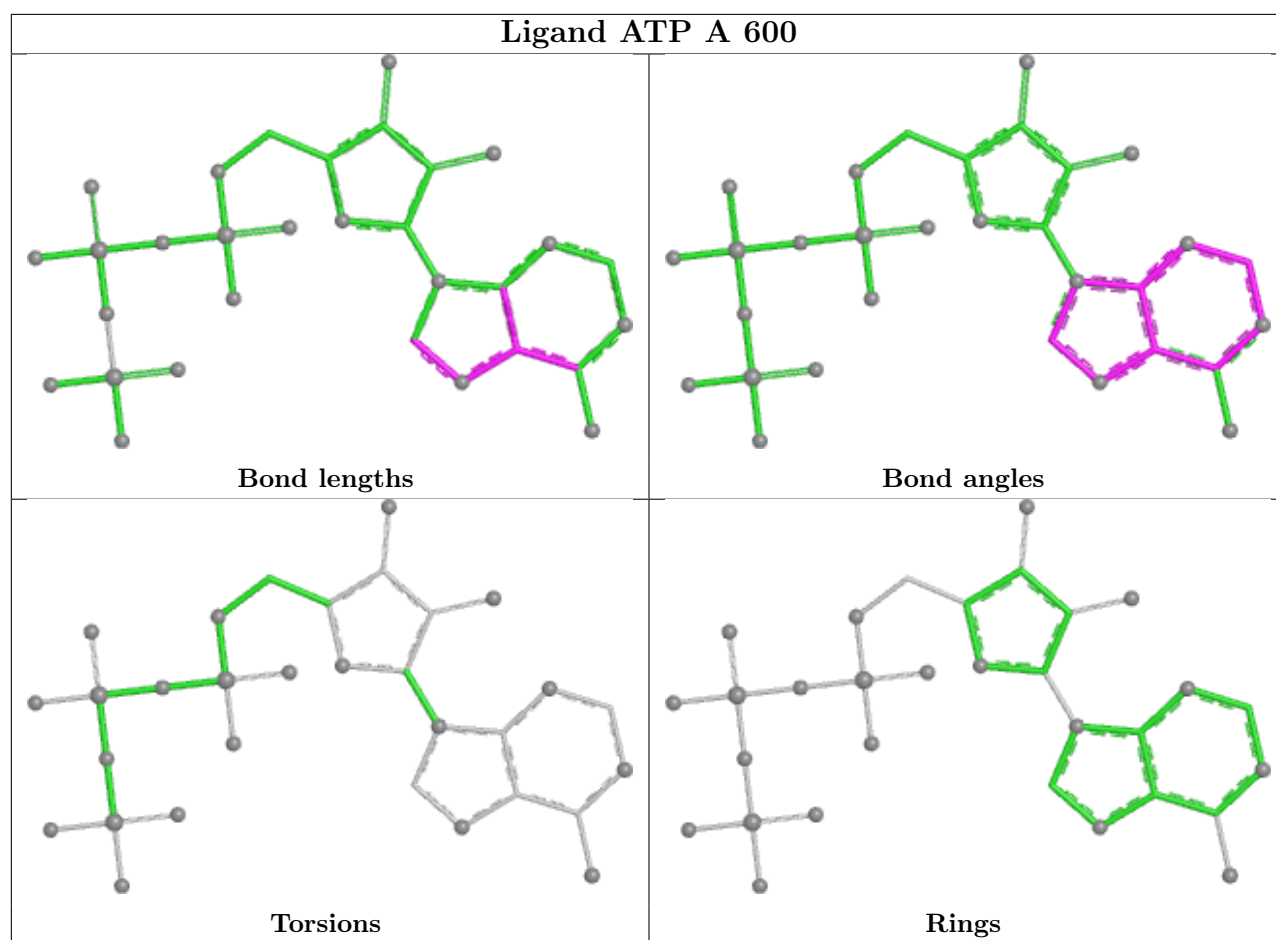
Mol	Chain	Res	Type	Atoms
4	B	1600	ATP	C5'-O5'-PA-O1A
4	B	1600	ATP	PB-O3A-PA-O1A
4	B	1600	ATP	PB-O3A-PA-O2A

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	422/460 (91%)	1.25	90 (21%) 2 2	26, 55, 90, 133	0
1	B	422/460 (91%)	0.60	20 (4%) 36 30	25, 45, 69, 98	0
2	C	22/110 (20%)	0.99	5 (22%) 2 2	31, 42, 62, 77	0
2	D	25/110 (22%)	0.66	2 (8%) 18 16	27, 39, 71, 90	0
All	All	891/1140 (78%)	0.92	117 (13%) 7 7	25, 49, 84, 133	0

All (117) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	366	VAL	5.1
1	A	364	ASN	5.0
1	A	18	THR	5.0
1	B	18	THR	4.9
1	A	100	ASN	4.7
1	A	367	GLY	4.6
1	A	342	SER	4.5
1	A	359	SER	4.5
1	A	99	PRO	4.4
1	A	361	VAL	4.2
1	A	365	GLU	4.1
2	D	499	SER	4.0
1	A	420	LYS	3.9
1	A	284	THR	3.9
1	A	369	VAL	3.9
1	B	306	PHE	3.7
1	B	365	GLU	3.5
1	A	360	ASN	3.5
1	A	417	LYS	3.4
1	A	400	VAL	3.4
1	A	429	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	80	GLU	3.4
1	A	362	PHE	3.3
1	A	343	PRO	3.2
1	A	444	LEU	3.2
1	A	146	TYR	3.2
1	B	21	ASN	3.1
1	A	413	GLU	3.1
1	A	419	ARG	3.1
1	A	19	GLY	3.0
1	A	29	PRO	3.0
1	A	441	TYR	3.0
1	B	62	VAL	3.0
1	A	344	TYR	3.0
1	B	360	ASN	3.0
1	A	33	ASP	2.9
1	A	61	ALA	2.9
1	A	30	LYS	2.9
1	A	418	ARG	2.8
2	C	477	ILE	2.8
1	A	101	HIS	2.8
1	A	94	THR	2.8
1	A	351	GLU	2.8
1	A	370	GLU	2.8
1	A	416	GLU	2.8
1	A	390	PHE	2.8
1	A	425	LEU	2.7
2	C	498	THR	2.7
1	A	430	ARG	2.7
1	A	397	GLN	2.7
1	A	70	ASN	2.6
1	A	93	ASN	2.6
1	A	98	LYS	2.6
1	A	52	ILE	2.6
1	A	287	LEU	2.6
1	A	289	GLU	2.6
1	A	63	ASP	2.6
2	C	497	ILE	2.5
1	B	247	GLY	2.5
1	A	414	VAL	2.5
1	B	114	GLU	2.5
1	A	331	GLU	2.5
1	A	345	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	382	GLN	2.5
1	B	364	ASN	2.5
1	A	261	ASN	2.5
1	A	431	LEU	2.4
1	A	310	LEU	2.4
1	A	301	ASN	2.4
1	B	361	VAL	2.4
1	B	366	VAL	2.4
1	A	311	ASP	2.4
1	A	40	ALA	2.4
1	A	433	SER	2.4
2	D	497	ILE	2.4
1	A	407	GLY	2.3
1	A	197	ALA	2.3
1	A	104	LYS	2.3
1	A	434	LYS	2.3
1	A	151	ASN	2.3
1	A	270	GLU	2.3
1	A	20	ASP	2.3
1	A	219	LYS	2.3
1	A	389	THR	2.3
1	B	284	THR	2.2
1	B	350	TYR	2.2
1	A	371	LEU	2.2
1	A	92	THR	2.2
1	A	428	PRO	2.2
1	A	314	SER	2.2
1	A	37	ASP	2.2
1	B	20	ASP	2.2
1	B	94	THR	2.2
1	B	347	GLY	2.2
1	A	315	ALA	2.2
1	A	62	VAL	2.1
1	A	377	THR	2.1
1	B	260	GLU	2.1
2	C	486	ASP	2.1
1	A	81	THR	2.1
1	A	196	GLY	2.1
1	A	368	ARG	2.1
1	B	195	SER	2.1
2	C	478	GLN	2.1
1	A	308	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	372	PHE	2.1
1	A	445	GLU	2.1
1	A	297	PRO	2.0
1	A	21	ASN	2.0
1	A	415	ASN	2.0
1	A	95	ILE	2.0
1	A	318	ASP	2.0
1	B	217	GLU	2.0
1	B	394	ARG	2.0
1	A	28	ILE	2.0
1	A	363	ASP	2.0
1	A	53	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](#)

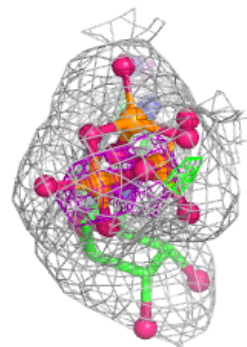
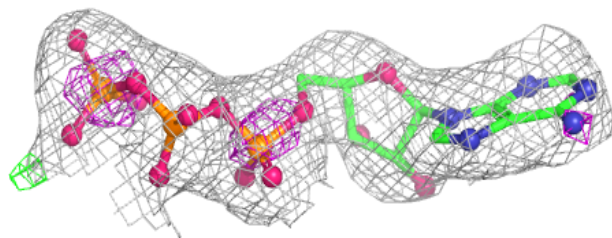
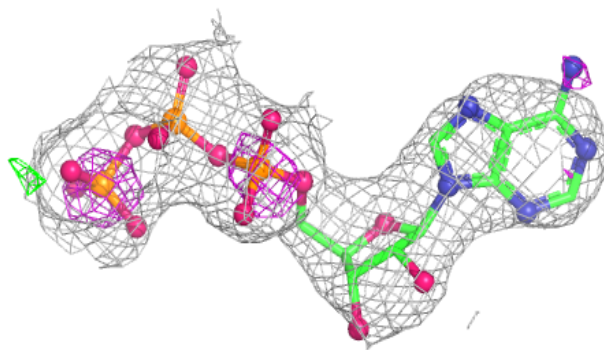
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	ATP	A	600	31/31	0.93	0.10	41,44,45,45	0
4	ATP	B	1600	31/31	0.94	0.09	37,39,40,41	0
3	MG	A	1500	1/1	0.95	0.07	27,27,27,27	0
3	MG	B	2500	1/1	0.96	0.04	15,15,15,15	0

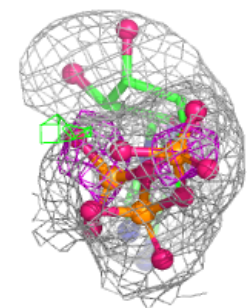
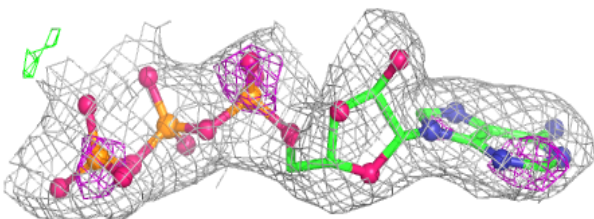
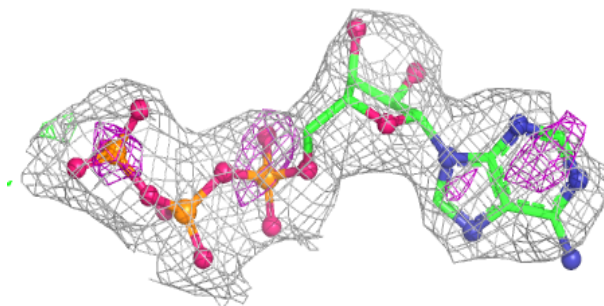
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ATP A 600:

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

**Electron density around ATP B 1600:**

$2mF_o-DF_c$ (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.