



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:11 AM UTC

PDB ID : 3LGX / pdb_00003lgx
Title : Structure of probable D-alanine-poly(phosphoribitol) ligase subunit-1 from Streptococcus pyogenes with ATP
Authors : Ramagopal, U.A.; Toro, R.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2010-01-21
Resolution : 2.60 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

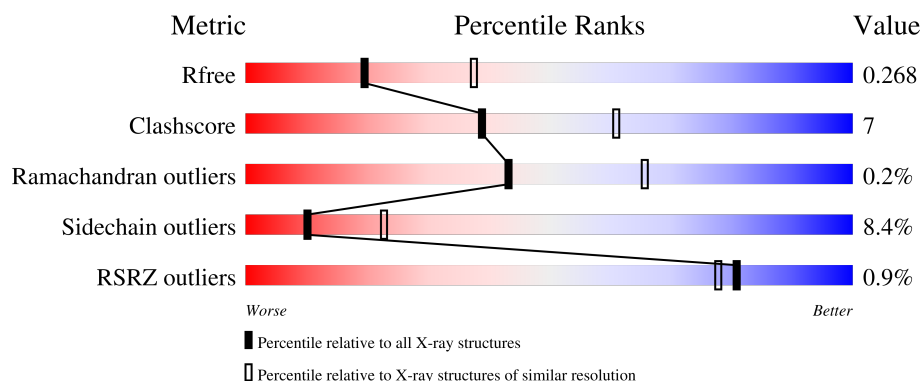
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	521	% 79% 17% ..
1	B	521	% 75% 21% ..
1	C	521	% 78% 17% ..
1	D	521	% 77% 18% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 16060 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 D-alanine--poly(phosphoribitol) ligase subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	508	Total	C	N	O	S	0	0	0
			3969	2554	632	766	17			
1	B	507	Total	C	N	O	S	0	3	0
			3986	2564	640	765	17			
1	C	508	Total	C	N	O	S	0	1	0
			3969	2553	636	763	17			
1	D	508	Total	C	N	O	S	0	0	0
			3964	2551	632	764	17			

There are 44 discrepancies between the modelled and reference sequences:

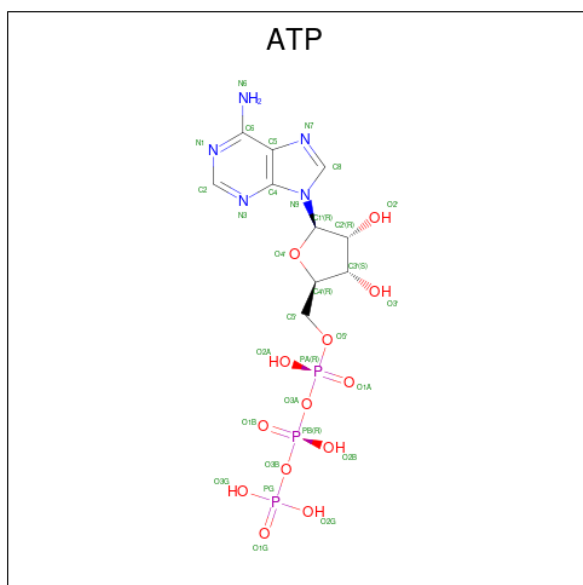
Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	MET	-	expression tag	UNP Q99ZA6
A	-3	SER	-	expression tag	UNP Q99ZA6
A	-2	LEU	-	expression tag	UNP Q99ZA6
A	509	GLU	-	expression tag	UNP Q99ZA6
A	510	GLY	-	expression tag	UNP Q99ZA6
A	511	HIS	-	expression tag	UNP Q99ZA6
A	512	HIS	-	expression tag	UNP Q99ZA6
A	513	HIS	-	expression tag	UNP Q99ZA6
A	514	HIS	-	expression tag	UNP Q99ZA6
A	515	HIS	-	expression tag	UNP Q99ZA6
A	516	HIS	-	expression tag	UNP Q99ZA6
B	-4	MET	-	expression tag	UNP Q99ZA6
B	-3	SER	-	expression tag	UNP Q99ZA6
B	-2	LEU	-	expression tag	UNP Q99ZA6
B	509	GLU	-	expression tag	UNP Q99ZA6
B	510	GLY	-	expression tag	UNP Q99ZA6
B	511	HIS	-	expression tag	UNP Q99ZA6
B	512	HIS	-	expression tag	UNP Q99ZA6
B	513	HIS	-	expression tag	UNP Q99ZA6
B	514	HIS	-	expression tag	UNP Q99ZA6
B	515	HIS	-	expression tag	UNP Q99ZA6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	516	HIS	-	expression tag	UNP Q99ZA6
C	-4	MET	-	expression tag	UNP Q99ZA6
C	-3	SER	-	expression tag	UNP Q99ZA6
C	-2	LEU	-	expression tag	UNP Q99ZA6
C	509	GLU	-	expression tag	UNP Q99ZA6
C	510	GLY	-	expression tag	UNP Q99ZA6
C	511	HIS	-	expression tag	UNP Q99ZA6
C	512	HIS	-	expression tag	UNP Q99ZA6
C	513	HIS	-	expression tag	UNP Q99ZA6
C	514	HIS	-	expression tag	UNP Q99ZA6
C	515	HIS	-	expression tag	UNP Q99ZA6
C	516	HIS	-	expression tag	UNP Q99ZA6
D	-4	MET	-	expression tag	UNP Q99ZA6
D	-3	SER	-	expression tag	UNP Q99ZA6
D	-2	LEU	-	expression tag	UNP Q99ZA6
D	509	GLU	-	expression tag	UNP Q99ZA6
D	510	GLY	-	expression tag	UNP Q99ZA6
D	511	HIS	-	expression tag	UNP Q99ZA6
D	512	HIS	-	expression tag	UNP Q99ZA6
D	513	HIS	-	expression tag	UNP Q99ZA6
D	514	HIS	-	expression tag	UNP Q99ZA6
D	515	HIS	-	expression tag	UNP Q99ZA6
D	516	HIS	-	expression tag	UNP Q99ZA6

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



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

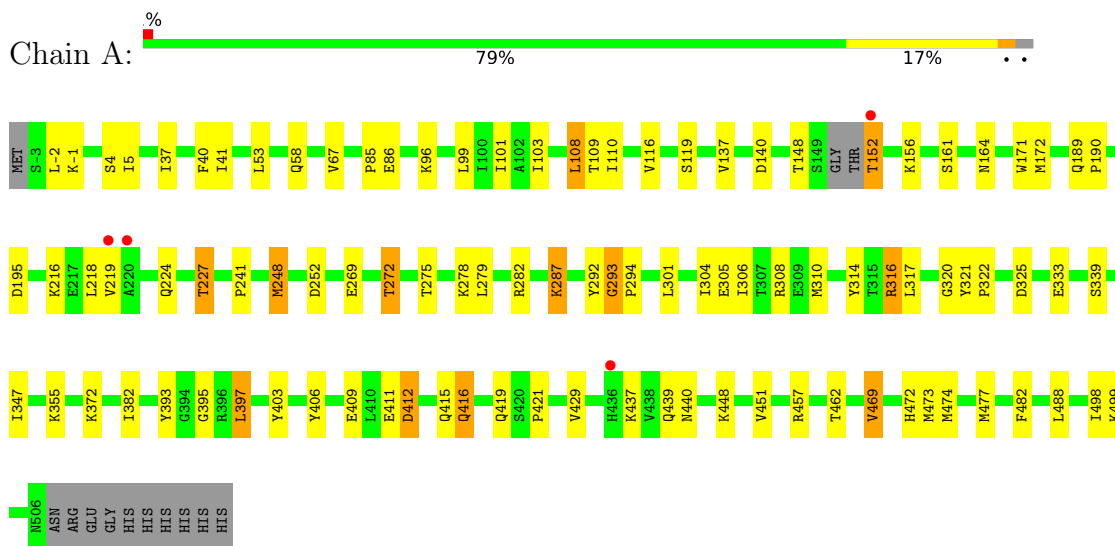
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		
3	B	17	Total	O	0	0
			17	17		
3	C	9	Total	O	0	0
			9	9		
3	D	11	Total	O	0	0
			11	11		

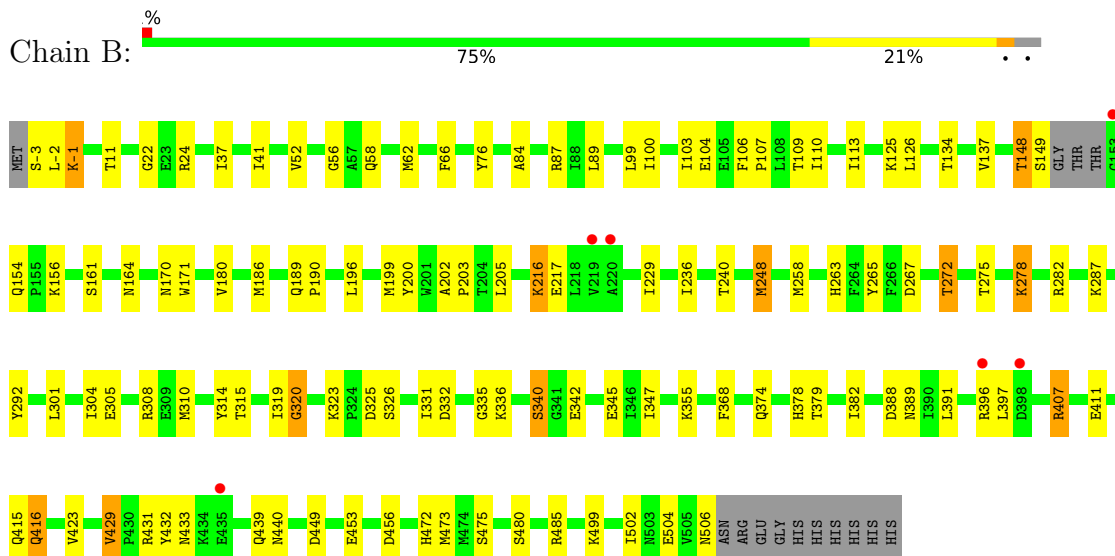
3 Residue-property plots

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.

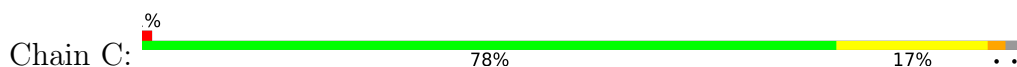
- Molecule 1: D-alanine--poly(phosphoribitol) ligase subunit 1

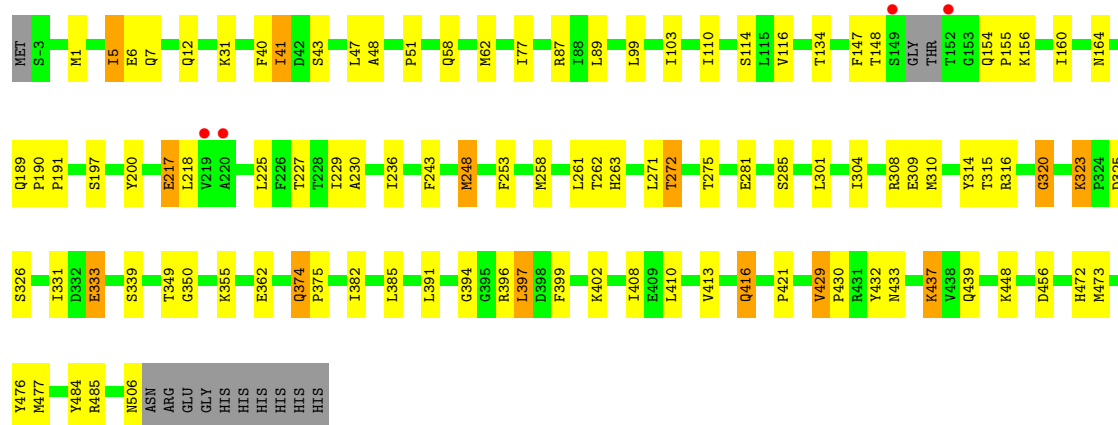


- Molecule 1: D-alanine--poly(phosphoribitol) ligase subunit 1

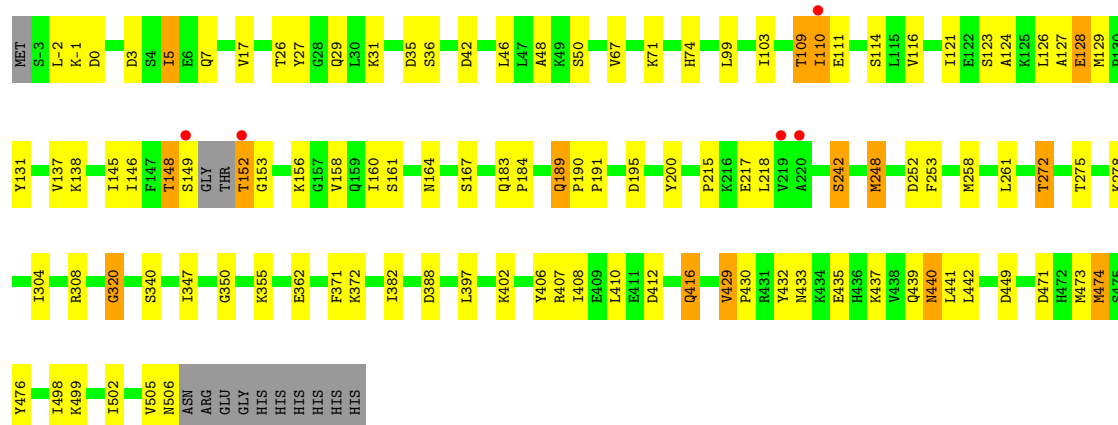
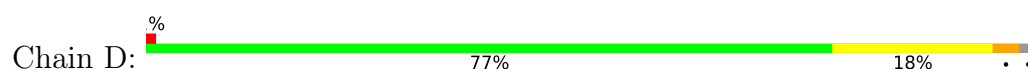


- Molecule 1: D-alanine--poly(phosphoribitol) ligase subunit 1





• Molecule 1: D-alanine-poly(phosphoribitol) ligase subunit 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	174.41Å 174.41Å 176.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.60 50.00 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (50.00-2.60) 99.7 (50.00-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.83 (at 2.61Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.218 , 0.276 0.211 , 0.268	Depositor DCC
R_{free} test set	4149 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	55.2	Xtriage
Anisotropy	0.084	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 21.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.035 for -h,l,k 0.032 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	16060	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.23% 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: 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	0.84	0/4060	1.03	7/5519 (0.1%)
1	B	0.90	0/4077	1.05	14/5540 (0.3%)
1	C	0.85	0/4060	1.04	8/5518 (0.1%)
1	D	0.86	0/4055	1.03	8/5512 (0.1%)
All	All	0.86	0/16252	1.04	37/22089 (0.2%)

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	11	THR	N-CA-C	7.05	119.05	111.36
1	D	320	GLY	N-CA-C	6.63	120.18	111.70
1	A	293	GLY	CA-C-N	6.62	126.79	120.31
1	A	293	GLY	C-N-CA	6.62	126.79	120.31
1	D	189	GLN	CA-C-N	-6.56	114.87	119.66
1	D	189	GLN	C-N-CA	-6.56	114.87	119.66
1	D	350	GLY	CA-C-N	6.54	126.23	119.56
1	D	350	GLY	C-N-CA	6.54	126.23	119.56
1	B	320	GLY	N-CA-C	6.33	119.81	111.52
1	C	331	ILE	N-CA-C	6.07	118.45	108.99
1	D	498	ILE	N-CA-C	6.03	116.77	110.62
1	C	315	THR	N-CA-C	-5.99	106.30	113.97
1	A	451	VAL	N-CA-C	-5.98	106.90	111.62
1	B	84	ALA	CA-C-N	5.95	127.27	119.84
1	B	84	ALA	C-N-CA	5.95	127.27	119.84
1	A	488	LEU	CA-C-N	-5.65	114.15	119.85
1	A	488	LEU	C-N-CA	-5.65	114.15	119.85
1	B	200	TYR	CA-C-N	-5.65	111.73	120.31
1	B	200	TYR	C-N-CA	-5.65	111.73	120.31
1	B	134	THR	N-CA-C	-5.64	107.65	114.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	248	MET	N-CA-C	-5.54	107.10	112.97
1	B	440	ASN	N-CA-C	5.50	115.64	107.88
1	C	134	THR	N-CA-C	-5.36	106.67	113.43
1	B	240	THR	N-CA-C	-5.35	102.69	110.08
1	C	271	LEU	CA-C-N	5.32	128.27	120.71
1	C	271	LEU	C-N-CA	5.32	128.27	120.71
1	D	440	ASN	N-CA-C	5.30	117.73	109.52
1	A	457	ARG	N-CA-C	5.28	118.03	109.06
1	B	62	MET	N-CA-C	-5.23	105.48	111.07
1	A	440	ASN	N-CA-C	5.17	116.38	108.42
1	C	62	MET	N-CA-C	-5.08	105.65	111.14
1	B	-1	LYS	N-CA-C	-5.04	107.79	114.04
1	B	374	GLN	CA-C-N	-5.02	114.59	119.76
1	B	374	GLN	C-N-CA	-5.02	114.59	119.76
1	B	423	VAL	N-CA-C	5.01	115.31	107.99
1	C	320	GLY	N-CA-C	5.00	120.21	112.45
1	D	242	SER	N-CA-C	5.00	116.73	111.28

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	3969	0	3928	43	0
1	B	3986	0	3944	58	0
1	C	3969	0	3921	56	0
1	D	3964	0	3920	58	0
2	A	31	0	12	1	0
2	B	31	0	12	0	0
2	C	31	0	12	0	0
2	D	31	0	12	0	0
3	A	11	0	0	0	0
3	B	17	0	0	0	0
3	C	9	0	0	0	0
3	D	11	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	16060	0	15761	215	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 7.

All (215) 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:D:474:MET:HG2	1:D:476:TYR:CE1	1.83	1.13
1:C:382:ILE:HD11	1:C:397:LEU:HD13	1.19	1.12
1:B:407:ARG:HG3	1:B:407:ARG:HH11	1.10	1.09
1:C:485[B]:ARG:HH11	1:C:485[B]:ARG:HG2	1.16	1.07
1:D:382:ILE:HD11	1:D:397:LEU:HD12	1.39	1.03
1:D:474:MET:HG2	1:D:476:TYR:CZ	2.05	0.92
1:C:272:THR:HG22	1:C:275:THR:H	1.35	0.90
1:B:272:THR:HG22	1:B:275:THR:H	1.36	0.89
1:A:161:SER:H	1:A:164:ASN:HD22	1.22	0.86
1:B:407:ARG:HG3	1:B:407:ARG:NH1	1.89	0.86
1:A:416:GLN:HG2	1:A:473:MET:HG2	1.58	0.85
1:B:161:SER:H	1:B:164:ASN:HD22	1.21	0.84
1:C:485[B]:ARG:HG2	1:C:485[B]:ARG:NH1	1.94	0.82
1:D:272:THR:HG22	1:D:275:THR:H	1.45	0.81
1:C:154:GLN:HE21	1:C:155:PRO:HD2	1.47	0.79
1:A:272:THR:HG22	1:A:275:THR:H	1.48	0.79
1:D:148:THR:HG22	1:D:156:LYS:H	1.47	0.78
1:D:253:PHE:HA	1:D:258:MET:CE	2.14	0.77
1:B:161:SER:H	1:B:164:ASN:ND2	1.82	0.77
1:D:161:SER:H	1:D:164:ASN:ND2	1.83	0.77
1:D:382:ILE:CD1	1:D:397:LEU:HD12	2.14	0.76
1:C:382:ILE:CD1	1:C:397:LEU:HD13	2.09	0.76
1:D:161:SER:H	1:D:164:ASN:HD22	1.31	0.76
1:D:31:LYS:O	1:D:35:ASP:OD1	2.06	0.74
1:A:248:MET:HE3	1:A:278:LYS:HD2	1.72	0.72
1:A:152:THR:N	1:A:499:LYS:HE3	2.04	0.71
1:D:253:PHE:HA	1:D:258:MET:HE1	1.74	0.69
1:B:416:GLN:HG2	1:B:473:MET:HG2	1.74	0.69
1:A:241:PRO:HG2	1:A:269:GLU:HG3	1.75	0.69
1:B:382:ILE:HD11	1:B:397:LEU:HD13	1.73	0.69
1:C:413:VAL:HG22	1:C:477:MET:HG2	1.75	0.69
1:D:416:GLN:HG3	1:D:473:MET:HG2	1.73	0.68
1:C:272:THR:CG2	1:C:275:THR:H	2.06	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:36:SER:OG	1:D:124:ALA:O	2.13	0.67
1:A:161:SER:H	1:A:164:ASN:ND2	1.90	0.66
1:C:160:ILE:HA	1:C:164:ASN:HD22	1.61	0.66
1:D:148:THR:HG23	1:D:149:SER:N	2.08	0.66
1:C:350:GLY:HA2	1:C:374:GLN:HG2	1.78	0.65
1:C:421:PRO:O	1:C:448:LYS:HE2	1.96	0.65
1:B:433:ASN:HD22	1:B:439:GLN:HE21	1.45	0.65
1:D:432:TYR:H	1:D:506:ASN:HD21	1.44	0.65
1:A:189:GLN:HB3	1:A:190:PRO:HD3	1.78	0.64
1:C:382:ILE:HD11	1:C:397:LEU:CD1	2.12	0.64
1:D:304:ILE:HB	1:D:320:GLY:HA2	1.80	0.63
1:C:58:GLN:HE21	1:C:191:PRO:HA	1.64	0.62
1:B:485[A]:ARG:NH2	1:B:504:GLU:OE1	2.32	0.62
1:C:333:GLU:CD	1:C:333:GLU:H	2.08	0.62
1:C:304:ILE:HB	1:C:320:GLY:HA2	1.82	0.61
1:C:99:LEU:HD11	1:C:116:VAL:HG23	1.83	0.61
1:B:379:THR:O	1:B:396[B]:ARG:NH1	2.34	0.61
1:B:407:ARG:HH11	1:B:407:ARG:CG	1.98	0.61
1:A:316:ARG:HD3	1:A:415:GLN:HG3	1.83	0.60
1:D:429:VAL:HG22	1:D:502:ILE:HG12	1.84	0.59
1:A:304:ILE:HB	1:A:320:GLY:HA2	1.85	0.59
1:D:17:VAL:HG23	1:D:27:TYR:CZ	2.38	0.59
1:D:347:ILE:C	1:D:347:ILE:HD12	2.28	0.59
1:A:310:MET:HG3	1:A:314:TYR:CE2	2.38	0.58
1:D:99:LEU:HD11	1:D:116:VAL:HG23	1.84	0.58
1:B:432:TYR:H	1:B:506:ASN:HD21	1.52	0.58
1:D:50:SER:O	1:D:74:HIS:HD2	1.86	0.58
1:B:304:ILE:HB	1:B:320:GLY:HA2	1.86	0.57
1:C:323:LYS:HG3	1:C:326:SER:HB2	1.85	0.57
1:C:408:ILE:HD11	1:C:476:TYR:CE2	2.39	0.57
1:D:0:ASP:HB3	1:D:3:ASP:HB2	1.86	0.57
1:B:52:VAL:HG13	1:B:99:LEU:HB3	1.86	0.57
1:D:190:PRO:HD2	1:D:200:TYR:CE2	2.40	0.57
1:D:252:ASP:O	1:D:258:MET:HE2	2.06	0.56
1:D:416:GLN:CG	1:D:473:MET:HG2	2.35	0.56
1:A:248:MET:O	1:A:282:ARG:NH2	2.38	0.56
1:D:26:THR:OG1	1:D:29:GLN:HG3	2.05	0.55
1:D:474:MET:CG	1:D:476:TYR:CE1	2.76	0.55
1:B:287:LYS:HD2	1:B:305:GLU:OE2	2.07	0.55
1:B:189:GLN:HB3	1:B:190:PRO:HD3	1.89	0.54
1:D:110:ILE:N	1:D:110:ILE:HD13	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:248:MET:HE3	1:B:278:LYS:HB3	1.89	0.54
1:B:433:ASN:HB3	1:B:439:GLN:HG2	1.88	0.54
1:C:6:GLU:OE1	1:C:31:LYS:NZ	2.27	0.54
1:C:316:ARG:NH2	1:C:394:GLY:O	2.41	0.54
1:B:22:GLY:O	1:B:24:ARG:HG3	2.08	0.54
1:D:161:SER:N	1:D:164:ASN:HD22	2.04	0.53
1:C:154:GLN:HE21	1:C:155:PRO:CD	2.21	0.53
1:D:5:ILE:HD11	1:D:67:VAL:HG11	1.89	0.53
1:A:101:ILE:HG22	1:A:103:ILE:HG12	1.89	0.53
1:B:301:LEU:C	1:B:301:LEU:HD12	2.34	0.53
1:B:310:MET:HG2	1:B:314:TYR:CE2	2.44	0.53
1:C:432:TYR:HA	1:C:437:LYS:O	2.08	0.52
1:C:154:GLN:NE2	1:C:155:PRO:HD2	2.21	0.52
1:C:301:LEU:HD12	1:C:301:LEU:C	2.33	0.52
1:C:408:ILE:HD11	1:C:476:TYR:HE2	1.71	0.52
1:C:484:TYR:O	1:C:485[B]:ARG:HG2	2.09	0.52
1:D:402:LYS:HA	1:D:406:TYR:O	2.10	0.52
1:A:85:PRO:HB3	1:A:108:LEU:HD21	1.92	0.52
1:C:225:LEU:HD21	1:C:229:ILE:HD11	1.91	0.52
1:C:325:ASP:C	1:C:325:ASP:OD1	2.53	0.51
1:D:190:PRO:HD2	1:D:200:TYR:HE2	1.74	0.51
1:A:409:GLU:O	1:A:412:ASP:HB2	2.11	0.51
1:B:171:TRP:CE3	1:B:323:LYS:HD3	2.46	0.51
1:B:416:GLN:CG	1:B:473:MET:HG2	2.40	0.51
1:C:416:GLN:HG3	1:C:472:HIS:HD2	1.75	0.50
1:B:196:LEU:HD13	1:B:267:ASP:HB3	1.93	0.50
1:B:248:MET:HG2	1:B:278:LYS:HG2	1.92	0.50
1:B:229:ILE:HG22	1:B:258:MET:HE2	1.94	0.50
1:B:432:TYR:H	1:B:506:ASN:ND2	2.09	0.50
1:B:248:MET:O	1:B:282:ARG:NH2	2.39	0.49
1:B:161:SER:N	1:B:164:ASN:HD22	1.99	0.49
1:C:225:LEU:C	1:C:225:LEU:HD23	2.37	0.49
1:C:189:GLN:HB3	1:C:190:PRO:HD3	1.93	0.49
1:D:50:SER:O	1:D:74:HIS:CD2	2.64	0.48
1:D:145:ILE:HA	1:D:158:VAL:O	2.13	0.48
1:C:432:TYR:H	1:C:506:ASN:HD21	1.61	0.48
1:C:433:ASN:HB3	1:C:439:GLN:HG2	1.95	0.48
1:C:473:MET:HE2	1:C:477:MET:HB3	1.94	0.48
1:D:253:PHE:HD1	1:D:258:MET:HE1	1.79	0.48
1:B:407:ARG:NH1	1:B:407:ARG:CG	2.66	0.48
1:A:53:LEU:C	1:A:53:LEU:HD23	2.38	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:382:ILE:HG13	1:A:397:LEU:HD13	1.96	0.48
1:B:99:LEU:HG	1:B:100:ILE:N	2.28	0.48
1:C:310:MET:HG2	1:C:314:TYR:CE2	2.49	0.48
1:D:429:VAL:CG1	1:D:505:VAL:HG21	2.44	0.47
1:D:433:ASN:HD22	1:D:439:GLN:HE21	1.62	0.47
1:D:126:LEU:C	1:D:128:GLU:H	2.22	0.47
1:A:469:VAL:HG12	1:A:472:HIS:HD2	1.80	0.47
1:D:42:ASP:OD2	1:D:131:TYR:HE1	1.98	0.47
1:B:148:THR:CG2	1:B:149:SER:N	2.77	0.47
1:B:199:MET:O	1:B:203:PRO:HG3	2.15	0.47
1:C:253:PHE:HA	1:C:258:MET:HE3	1.97	0.46
1:B:382:ILE:HD11	1:B:397:LEU:CD1	2.40	0.46
1:D:215:PRO:HD2	1:D:218:LEU:HD12	1.97	0.46
1:B:388:ASP:O	1:B:389:ASN:HB3	2.16	0.46
1:A:403:TYR:O	1:A:406:TYR:HB2	2.15	0.46
1:C:189:GLN:HG3	1:C:243:PHE:CE1	2.51	0.46
1:A:393:TYR:CZ	1:A:395:GLY:HA2	2.50	0.46
1:A:473:MET:CE	1:A:477:MET:HB3	2.45	0.46
1:C:217:GLU:O	1:C:218:LEU:HD12	2.16	0.46
1:D:109:THR:C	1:D:110:ILE:HD13	2.41	0.46
1:C:349:THR:HA	1:C:375:PRO:O	2.16	0.45
1:B:37:ILE:HD11	1:B:125:LYS:HG2	1.98	0.45
1:B:202:ALA:N	1:B:203:PRO:HD2	2.32	0.45
1:A:287:LYS:HD3	1:A:305:GLU:OE2	2.16	0.45
1:A:421:PRO:O	1:A:448:LYS:HE2	2.16	0.45
1:B:382:ILE:CD1	1:B:397:LEU:HD13	2.43	0.45
1:B:416:GLN:HG3	1:B:472:HIS:HD2	1.82	0.45
1:C:51:PRO:HB2	1:C:77:ILE:HD11	1.98	0.45
1:A:325:ASP:OD1	1:A:325:ASP:C	2.59	0.45
1:B:325:ASP:OD1	1:B:325:ASP:C	2.60	0.45
1:B:331:ILE:HD12	1:B:335:GLY:HA2	1.97	0.45
1:B:319:ILE:HG13	1:B:391:LEU:HB3	1.98	0.44
1:B:180:VAL:HG22	1:B:265:TYR:OH	2.18	0.44
1:C:236:ILE:HG12	1:C:263:HIS:HB2	1.99	0.44
1:B:-3:SER:HB2	1:B:170:ASN:OD1	2.17	0.44
1:B:429:VAL:HG22	1:B:502:ILE:HG12	1.99	0.44
1:D:355:LYS:HB3	1:D:355:LYS:HE3	1.78	0.44
1:C:484:TYR:O	1:C:485[B]:ARG:NH1	2.50	0.44
1:B:345:GLU:OE2	1:B:378:HIS:HB3	2.17	0.43
1:C:304:ILE:HB	1:C:320:GLY:CA	2.46	0.43
1:D:473:MET:HE2	1:D:473:MET:HB3	1.91	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:236:ILE:HG12	1:B:263:HIS:HB2	2.00	0.43
1:D:253:PHE:CD1	1:D:258:MET:HE1	2.53	0.43
1:A:292:TYR:CG	1:A:293:GLY:N	2.86	0.43
1:C:147:PHE:HA	1:C:156:LYS:O	2.17	0.43
1:A:416:GLN:CG	1:A:473:MET:HG2	2.38	0.43
1:B:332:ASP:OD2	1:B:336:LYS:HB2	2.18	0.43
1:C:396:ARG:O	1:C:399:PHE:N	2.47	0.43
1:D:190:PRO:HA	1:D:191:PRO:HD3	1.84	0.43
1:D:433:ASN:HB3	1:D:439:GLN:HG2	2.00	0.43
1:B:106:PHE:HA	1:B:107:PRO:HD3	1.78	0.43
1:C:103:ILE:N	1:C:103:ILE:HD12	2.34	0.43
1:C:103:ILE:HD12	1:C:103:ILE:H	1.84	0.43
1:A:411:GLU:O	1:A:415:GLN:HG2	2.18	0.42
1:A:416:GLN:O	1:A:419:GLN:HB2	2.19	0.42
1:B:205:LEU:HD23	1:B:205:LEU:HA	1.76	0.42
1:A:224:GLN:OE1	1:A:224:GLN:HA	2.18	0.42
1:C:429:VAL:HA	1:C:430:PRO:HD3	1.94	0.42
1:C:1:MET:O	1:C:5:ILE:HG23	2.20	0.42
1:C:484:TYR:C	1:C:485[B]:ARG:HG2	2.45	0.42
1:D:46:LEU:HD11	1:D:99:LEU:HD22	2.01	0.42
1:B:433:ASN:HB3	1:B:439:GLN:HE21	1.85	0.42
1:A:227:THR:HG23	1:A:252:ASP:OD1	2.20	0.42
1:B:66:PHE:HB3	1:B:76:TYR:CE1	2.54	0.42
1:B:323:LYS:HE3	1:B:326:SER:OG	2.20	0.42
1:C:197:SER:HA	1:C:200:TYR:CE2	2.54	0.42
1:C:385:LEU:HD13	1:C:391:LEU:HD13	2.01	0.42
1:A:140:ASP:O	1:A:355:LYS:NZ	2.52	0.42
1:A:306:ILE:HD12	1:A:317:LEU:HD13	2.02	0.42
1:D:408:ILE:HD11	1:D:476:TYR:HE2	1.85	0.42
1:B:186:MET:HE3	1:B:236:ILE:HG21	2.02	0.42
1:D:160:ILE:HA	1:D:164:ASN:HD22	1.85	0.42
1:A:171:TRP:O	1:A:172:MET:C	2.62	0.41
1:C:40:PHE:HD2	1:C:41:ILE:HD12	1.84	0.41
1:A:-2:LEU:HD12	1:A:-2:LEU:HA	1.90	0.41
1:D:164:ASN:O	1:D:167:SER:HB3	2.19	0.41
1:A:333:GLU:H	1:A:333:GLU:HG3	1.73	0.41
1:A:310:MET:HG2	1:A:314:TYR:CD2	2.55	0.41
1:B:368:PHE:N	1:B:368:PHE:CD1	2.88	0.41
1:A:321:TYR:HA	1:A:322:PRO:HD3	1.93	0.41
1:C:227:THR:O	1:C:230:ALA:HB3	2.20	0.41
1:C:410:LEU:HD23	1:C:410:LEU:HA	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:152:THR:HB	1:D:153:GLY:H	1.56	0.41
1:A:294:PRO:HA	2:A:600:ATP:H5'1	2.03	0.41
1:A:301:LEU:HD12	1:A:301:LEU:C	2.46	0.41
1:D:248:MET:HE3	1:D:278:LYS:HB2	2.01	0.41
1:A:37:ILE:O	1:A:40:PHE:HB3	2.21	0.41
1:D:410:LEU:HD23	1:D:410:LEU:HA	1.80	0.41
1:D:371:PHE:O	1:D:372:LYS:C	2.65	0.40
1:A:5:ILE:HD11	1:A:67:VAL:HG11	2.02	0.40
1:A:99:LEU:HD11	1:A:116:VAL:HG23	2.03	0.40
1:D:429:VAL:HA	1:D:430:PRO:HD3	1.92	0.40
1:B:56:GLY:HA2	1:B:104:GLU:HG2	2.04	0.40
1:B:411:GLU:O	1:B:415:GLN:HG2	2.20	0.40
1:B:431:ARG:HA	1:B:506:ASN:HD21	1.87	0.40
1:D:67:VAL:HG12	1:D:71:LYS:HE3	2.04	0.40
1:A:462:THR:HG23	1:A:482:PHE:CD2	2.57	0.40
1:C:47:LEU:O	1:C:48:ALA:C	2.63	0.40
1:D:127:ALA:O	1:D:129:MET:N	2.54	0.40
1:D:183:GLN:N	1:D:184:PRO:CD	2.85	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	504/521 (97%)	477 (95%)	27 (5%)	0	100	100
1	B	506/521 (97%)	485 (96%)	19 (4%)	2 (0%)	30	51
1	C	505/521 (97%)	478 (95%)	27 (5%)	0	100	100
1	D	504/521 (97%)	474 (94%)	28 (6%)	2 (0%)	30	51
All	All	2019/2084 (97%)	1914 (95%)	101 (5%)	4 (0%)	43	66

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	128	GLU
1	B	216	LYS
1	B	340	SER
1	D	48	ALA

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	432/447 (97%)	395 (91%)	37 (9%)	10	22
1	B	432/447 (97%)	396 (92%)	36 (8%)	10	23
1	C	430/447 (96%)	399 (93%)	31 (7%)	13	30
1	D	431/447 (96%)	391 (91%)	40 (9%)	8	18
All	All	1725/1788 (96%)	1581 (92%)	144 (8%)	10	23

All (144) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-1	LYS
1	A	4	SER
1	A	41	ILE
1	A	58	GLN
1	A	86	GLU
1	A	96	LYS
1	A	108	LEU
1	A	109	THR
1	A	110	ILE
1	A	119	SER
1	A	137	VAL
1	A	148	THR
1	A	152	THR
1	A	156	LYS
1	A	195	ASP
1	A	216	LYS
1	A	218	LEU
1	A	219	VAL

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Mol	Chain	Res	Type
1	A	227	THR
1	A	248	MET
1	A	272	THR
1	A	279	LEU
1	A	287	LYS
1	A	308	ARG
1	A	316	ARG
1	A	339	SER
1	A	347	ILE
1	A	372	LYS
1	A	397	LEU
1	A	412	ASP
1	A	416	GLN
1	A	429	VAL
1	A	437	LYS
1	A	439	GLN
1	A	469	VAL
1	A	474	MET
1	A	498	ILE
1	B	-2	LEU
1	B	-1	LYS
1	B	41	ILE
1	B	58	GLN
1	B	87	ARG
1	B	89	LEU
1	B	103	ILE
1	B	109	THR
1	B	110	ILE
1	B	113	ILE
1	B	126	LEU
1	B	137	VAL
1	B	148	THR
1	B	154	GLN
1	B	156	LYS
1	B	216	LYS
1	B	217	GLU
1	B	248	MET
1	B	272	THR
1	B	278	LYS
1	B	292	TYR
1	B	308	ARG
1	B	315	THR

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Mol	Chain	Res	Type
1	B	340	SER
1	B	342	GLU
1	B	347	ILE
1	B	355	LYS
1	B	407	ARG
1	B	416	GLN
1	B	429	VAL
1	B	449	ASP
1	B	453	GLU
1	B	456	ASP
1	B	475	SER
1	B	480	SER
1	B	499	LYS
1	C	5	ILE
1	C	7	GLN
1	C	12	GLN
1	C	41	ILE
1	C	43	SER
1	C	87	ARG
1	C	89	LEU
1	C	110	ILE
1	C	114	SER
1	C	148	THR
1	C	217	GLU
1	C	248	MET
1	C	261	LEU
1	C	262	THR
1	C	272	THR
1	C	281	GLU
1	C	285	SER
1	C	308	ARG
1	C	309	GLU
1	C	323	LYS
1	C	333	GLU
1	C	339	SER
1	C	355	LYS
1	C	362	GLU
1	C	374	GLN
1	C	397	LEU
1	C	402	LYS
1	C	416	GLN
1	C	429	VAL

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Mol	Chain	Res	Type
1	C	437	LYS
1	C	456	ASP
1	D	-2	LEU
1	D	-1	LYS
1	D	5	ILE
1	D	7	GLN
1	D	103	ILE
1	D	109	THR
1	D	110	ILE
1	D	111	GLU
1	D	114	SER
1	D	121	ILE
1	D	123	SER
1	D	137	VAL
1	D	138	LYS
1	D	146	ILE
1	D	148	THR
1	D	152	THR
1	D	189	GLN
1	D	195	ASP
1	D	217	GLU
1	D	242	SER
1	D	248	MET
1	D	261	LEU
1	D	272	THR
1	D	308	ARG
1	D	340	SER
1	D	362	GLU
1	D	388	ASP
1	D	407	ARG
1	D	412	ASP
1	D	416	GLN
1	D	429	VAL
1	D	435	GLU
1	D	437	LYS
1	D	440	ASN
1	D	441	LEU
1	D	442	LEU
1	D	449	ASP
1	D	471	ASP
1	D	474	MET
1	D	499	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	58	GLN
1	A	154	GLN
1	A	164	ASN
1	A	183	GLN
1	A	185	GLN
1	A	189	GLN
1	A	433	ASN
1	A	472	HIS
1	A	503	ASN
1	B	12	GLN
1	B	154	GLN
1	B	164	ASN
1	B	183	GLN
1	B	185	GLN
1	B	189	GLN
1	B	313	ASN
1	B	416	GLN
1	B	439	GLN
1	B	472	HIS
1	B	503	ASN
1	B	506	ASN
1	C	58	GLN
1	C	74	HIS
1	C	154	GLN
1	C	164	ASN
1	C	185	GLN
1	C	231	GLN
1	C	255	GLN
1	C	313	ASN
1	C	343	GLN
1	C	359	ASN
1	C	400	GLN
1	C	416	GLN
1	C	440	ASN
1	C	503	ASN
1	C	506	ASN
1	D	74	HIS
1	D	154	GLN
1	D	164	ASN
1	D	189	GLN

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Mol	Chain	Res	Type
1	D	224	GLN
1	D	313	ASN
1	D	359	ASN
1	D	439	GLN
1	D	503	ASN
1	D	506	ASN

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 ⓘ

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	ATP	D	600	-	32,33,33	1.59	6 (18%)	48,52,52	1.64	6 (12%)
2	ATP	A	600	-	32,33,33	1.38	7 (21%)	48,52,52	1.99	11 (22%)
2	ATP	B	600	-	32,33,33	1.28	3 (9%)	48,52,52	1.91	15 (31%)
2	ATP	C	600	-	32,33,33	1.32	5 (15%)	48,52,52	1.95	12 (25%)

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	ATP	D	600	-	-	7/22/38/38	0/3/3/3
2	ATP	A	600	-	-	9/22/38/38	0/3/3/3
2	ATP	B	600	-	-	7/22/38/38	0/3/3/3
2	ATP	C	600	-	-	4/22/38/38	0/3/3/3

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	600	ATP	C5-C4	5.03	1.48	1.39
2	B	600	ATP	C5-C4	4.42	1.47	1.39
2	C	600	ATP	C5-C4	4.10	1.46	1.39
2	D	600	ATP	PA-O3A	3.36	1.63	1.59
2	A	600	ATP	PB-O3A	3.28	1.63	1.59
2	A	600	ATP	C5-C4	3.22	1.44	1.39
2	D	600	ATP	PB-O3A	3.12	1.62	1.59
2	B	600	ATP	C8-N7	2.77	1.37	1.31
2	C	600	ATP	C5-N7	-2.63	1.34	1.39
2	B	600	ATP	C5-C6	2.58	1.48	1.41
2	A	600	ATP	PA-O3A	2.57	1.62	1.59
2	A	600	ATP	C8-N9	-2.56	1.33	1.37
2	D	600	ATP	C5-N7	-2.45	1.34	1.39
2	D	600	ATP	C5-C6	2.43	1.47	1.41
2	C	600	ATP	PB-O3A	2.41	1.62	1.59
2	D	600	ATP	C8-N7	2.30	1.36	1.31
2	A	600	ATP	C8-N7	2.28	1.36	1.31
2	C	600	ATP	C8-N7	2.24	1.36	1.31
2	A	600	ATP	C5-N7	-2.19	1.35	1.39
2	A	600	ATP	C5-C6	2.05	1.46	1.41
2	C	600	ATP	C5-C6	2.04	1.46	1.41

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	600	ATP	C5-C4-N3	-6.42	117.88	126.72
2	D	600	ATP	C5-C4-N3	-6.13	118.28	126.72
2	C	600	ATP	C5-C4-N3	-5.86	118.65	126.72
2	B	600	ATP	C5-C4-N3	-5.19	119.57	126.72
2	D	600	ATP	N3-C4-N9	5.01	135.68	127.17
2	A	600	ATP	N3-C4-N9	4.80	135.33	127.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	600	ATP	N3-C2-N1	-4.66	121.52	128.58
2	B	600	ATP	N3-C4-N9	4.61	135.02	127.17
2	C	600	ATP	N3-C4-N9	4.59	134.98	127.17
2	A	600	ATP	C2-N3-C4	4.30	122.33	111.83
2	A	600	ATP	N3-C2-N1	-4.29	122.09	128.58
2	A	600	ATP	C4-C5-N7	-4.22	105.75	110.58
2	C	600	ATP	C2-N3-C4	4.08	121.80	111.83
2	B	600	ATP	C2-N3-C4	3.89	121.34	111.83
2	B	600	ATP	N3-C2-N1	-3.87	122.72	128.58
2	C	600	ATP	C4-C5-N7	-3.67	106.39	110.58
2	D	600	ATP	C2-N3-C4	3.62	120.68	111.83
2	B	600	ATP	C4-N9-C8	3.58	109.49	105.74
2	B	600	ATP	O3A-PA-O1A	-3.42	100.41	110.70
2	A	600	ATP	C5-N7-C8	3.24	108.54	103.45
2	C	600	ATP	O2B-PB-O3A	-3.19	98.66	107.27
2	A	600	ATP	C4-N9-C8	3.13	109.03	105.74
2	C	600	ATP	C5-N7-C8	2.93	108.05	103.45
2	A	600	ATP	N9-C8-N7	-2.86	109.88	113.94
2	D	600	ATP	N3-C2-N1	-2.68	124.53	128.58
2	C	600	ATP	C2-N1-C6	2.60	123.01	118.73
2	C	600	ATP	C4-N9-C8	2.59	108.46	105.74
2	A	600	ATP	C6-C5-N7	2.52	136.94	132.09
2	B	600	ATP	N9-C8-N7	-2.41	110.52	113.94
2	B	600	ATP	C4-C5-N7	-2.39	107.84	110.58
2	B	600	ATP	O2A-PA-O1A	2.39	123.56	112.44
2	A	600	ATP	O2B-PB-O1B	2.38	123.51	112.44
2	D	600	ATP	C4-C5-N7	-2.35	107.89	110.58
2	D	600	ATP	O3A-PA-O1A	-2.33	103.69	110.70
2	C	600	ATP	C6-C5-N7	2.32	136.56	132.09
2	C	600	ATP	N9-C8-N7	-2.27	110.72	113.94
2	A	600	ATP	O3G-PG-O2G	2.26	116.28	107.80
2	B	600	ATP	O3G-PG-O2G	2.09	115.63	107.80
2	B	600	ATP	C5-C6-N6	-2.06	118.18	123.29
2	B	600	ATP	C5-N7-C8	2.06	106.69	103.45
2	B	600	ATP	C6-C5-N7	2.03	136.01	132.09
2	C	600	ATP	O3A-PA-O1A	-2.03	104.60	110.70
2	B	600	ATP	O2B-PB-O1B	2.03	121.89	112.44
2	B	600	ATP	N6-C6-N1	2.02	122.89	118.38

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	600	ATP	C5'-O5'-PA-O2A
2	A	600	ATP	C5'-O5'-PA-O3A
2	B	600	ATP	C5'-O5'-PA-O1A
2	B	600	ATP	C5'-O5'-PA-O2A
2	B	600	ATP	C5'-O5'-PA-O3A
2	C	600	ATP	C5'-O5'-PA-O1A
2	C	600	ATP	C5'-O5'-PA-O2A
2	C	600	ATP	C5'-O5'-PA-O3A
2	D	600	ATP	C5'-O5'-PA-O2A
2	D	600	ATP	C5'-O5'-PA-O3A
2	A	600	ATP	C3'-C4'-C5'-O5'
2	A	600	ATP	PB-O3A-PA-O5'
2	B	600	ATP	PB-O3A-PA-O5'
2	C	600	ATP	PB-O3A-PA-O5'
2	D	600	ATP	PB-O3A-PA-O5'
2	D	600	ATP	C3'-C4'-C5'-O5'
2	A	600	ATP	PA-O3A-PB-O1B
2	B	600	ATP	PB-O3A-PA-O1A
2	D	600	ATP	PA-O3A-PB-O1B
2	B	600	ATP	PB-O3B-PG-O2G
2	D	600	ATP	PB-O3B-PG-O3G
2	A	600	ATP	PG-O3B-PB-O2B
2	A	600	ATP	PA-O3A-PB-O2B
2	D	600	ATP	PA-O3A-PB-O2B
2	A	600	ATP	O4'-C4'-C5'-O5'
2	B	600	ATP	PA-O3A-PB-O2B
2	A	600	ATP	PG-O3B-PB-O3A

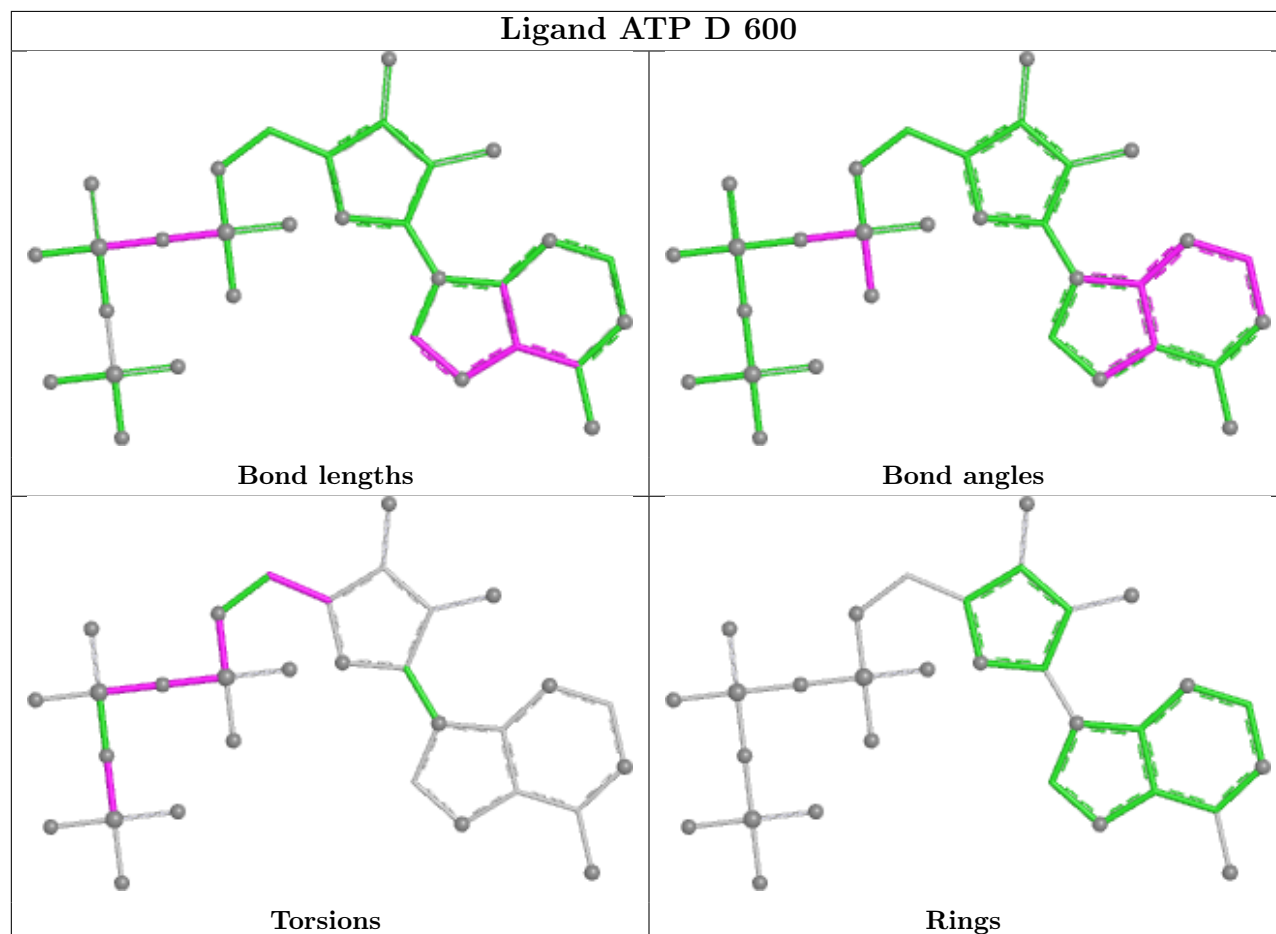
There are no ring outliers.

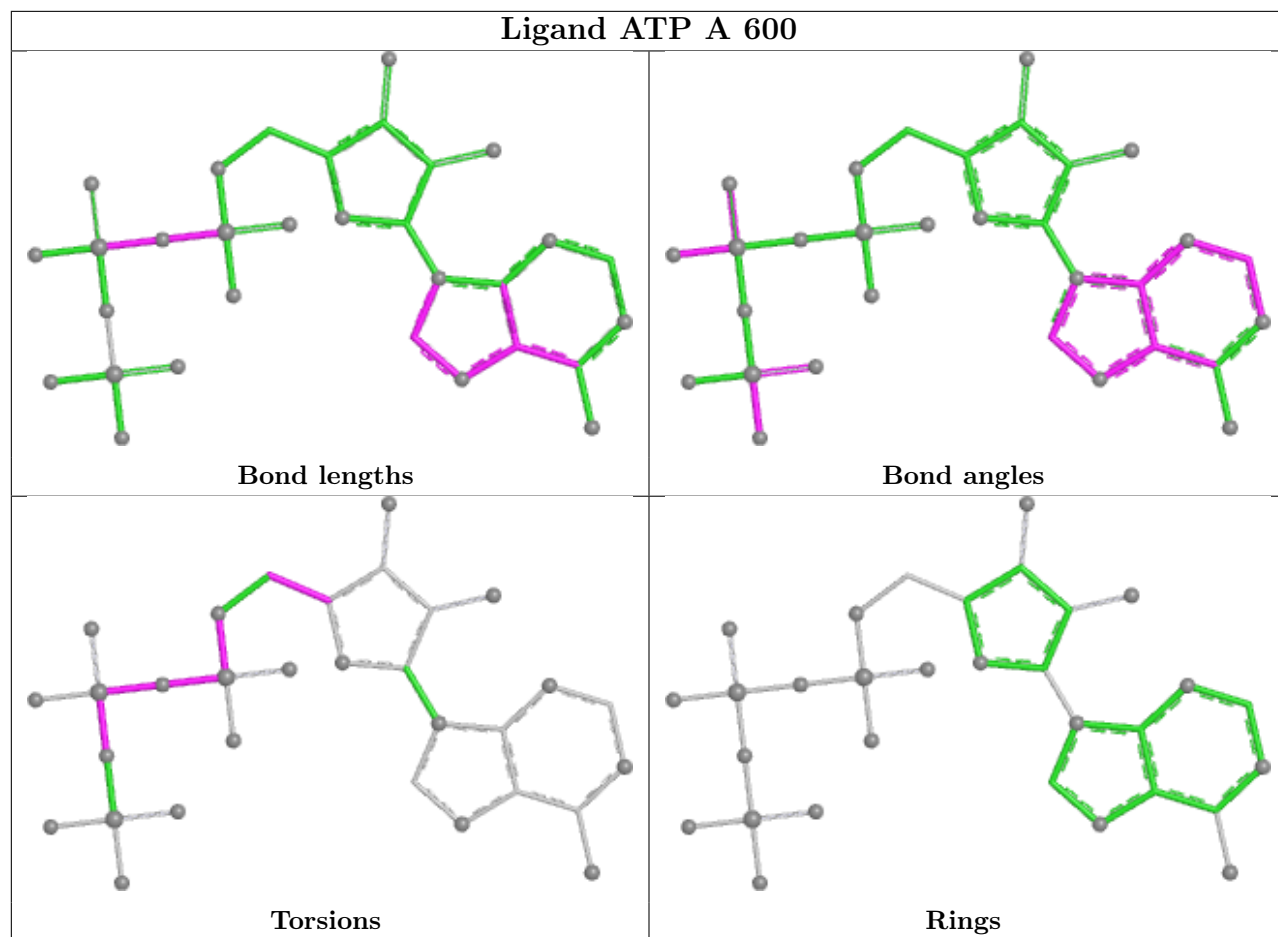
1 monomer is involved in 1 short contact:

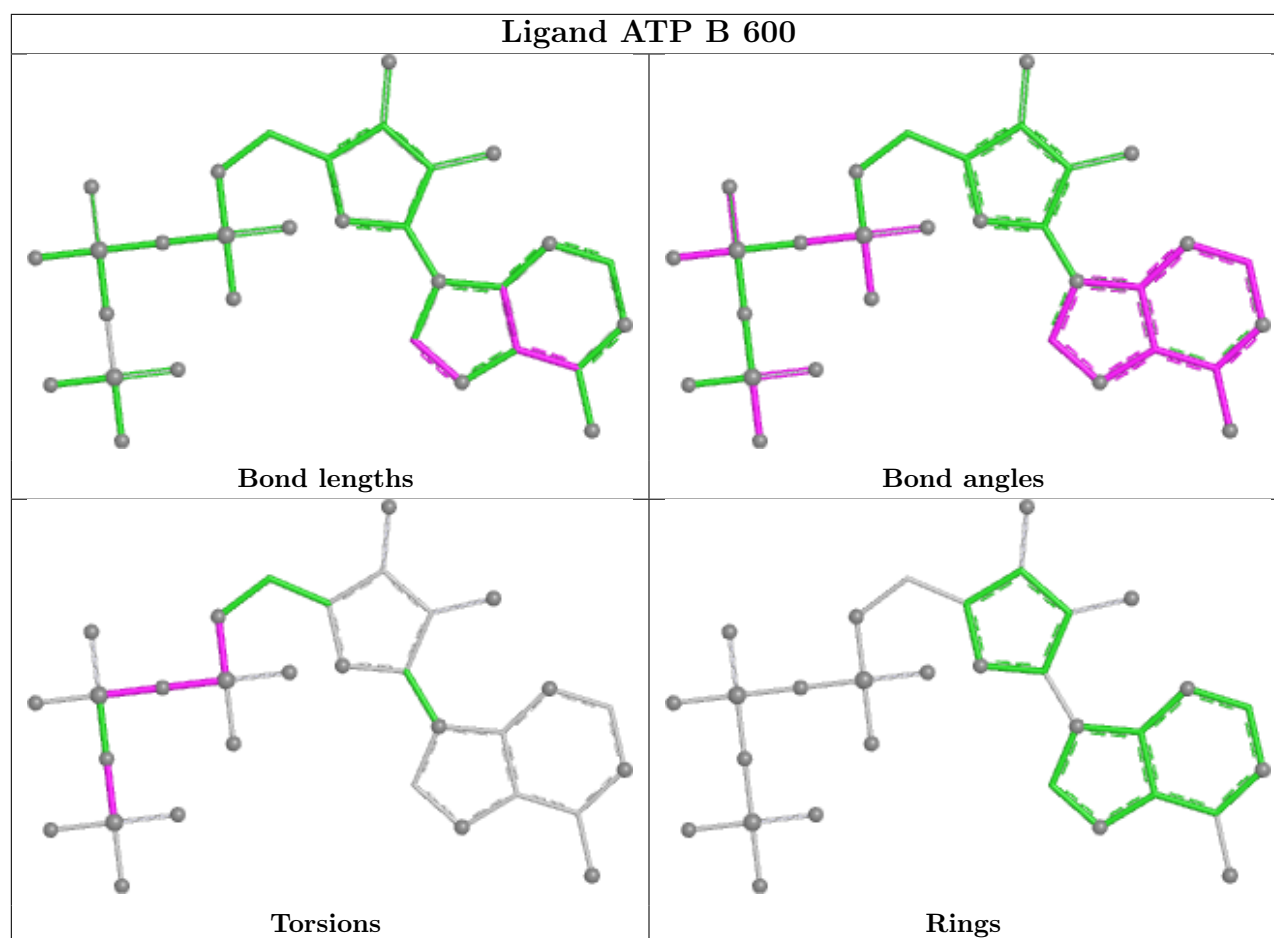
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	ATP	1	0

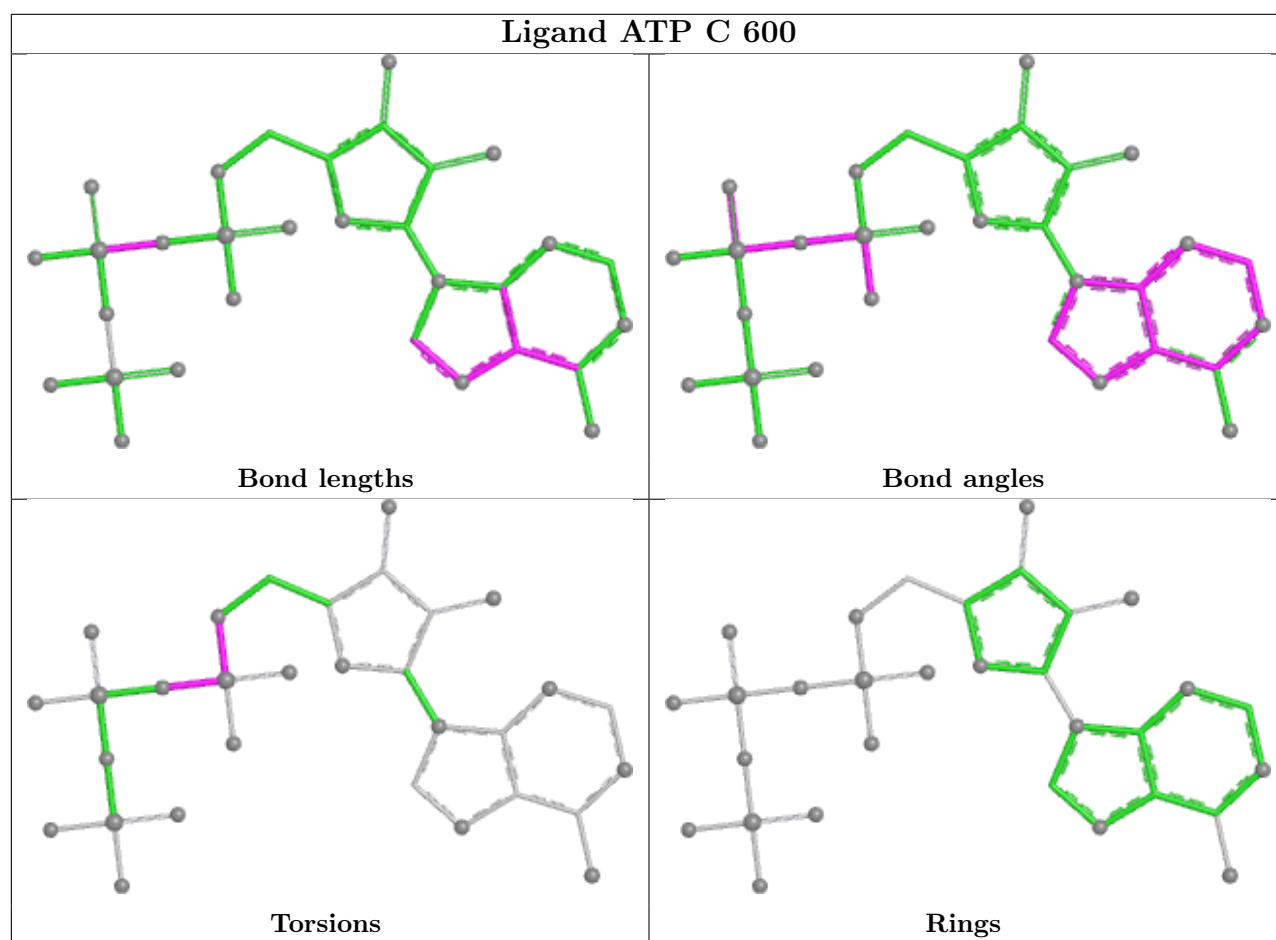
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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

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	508/521 (97%)	-0.07	4 (0%) 82 80	30, 52, 80, 97	0
1	B	507/521 (97%)	-0.21	6 (1%) 76 73	18, 45, 66, 82	3 (0%)
1	C	508/521 (97%)	-0.13	4 (0%) 82 80	20, 49, 73, 86	1 (0%)
1	D	508/521 (97%)	-0.11	5 (0%) 79 76	32, 49, 70, 86	0
All	All	2031/2084 (97%)	-0.13	19 (0%) 81 78	18, 49, 74, 97	4 (0%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	219	VAL	3.6
1	B	435	GLU	3.4
1	C	152	THR	3.4
1	D	110	ILE	3.2
1	C	220	ALA	3.1
1	B	219	VAL	3.1
1	A	220	ALA	2.8
1	D	219	VAL	2.7
1	A	152	THR	2.6
1	B	398	ASP	2.6
1	B	396[A]	ARG	2.6
1	B	220	ALA	2.6
1	A	436	HIS	2.4
1	D	149	SER	2.4
1	C	149	SER	2.3
1	D	152	THR	2.3
1	B	153	GLY	2.3
1	C	219	VAL	2.2
1	D	220	ALA	2.1

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](#)

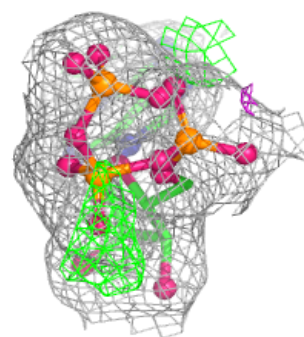
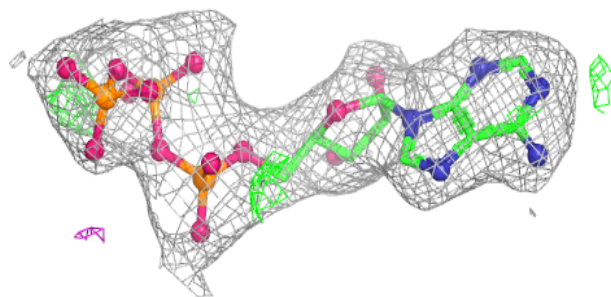
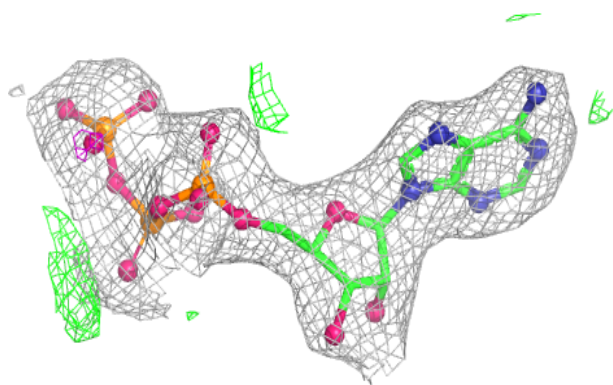
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	ATP	A	600	31/31	0.95	0.07	22,39,67,67	0
2	ATP	C	600	31/31	0.95	0.07	30,41,63,64	0
2	ATP	D	600	31/31	0.95	0.07	26,41,66,67	0
2	ATP	B	600	31/31	0.98	0.05	23,34,58,59	0

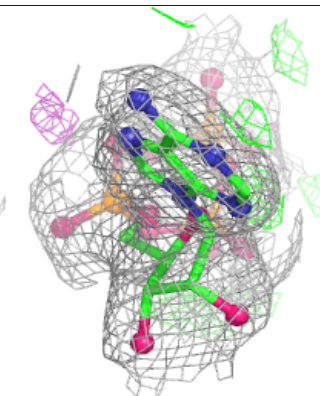
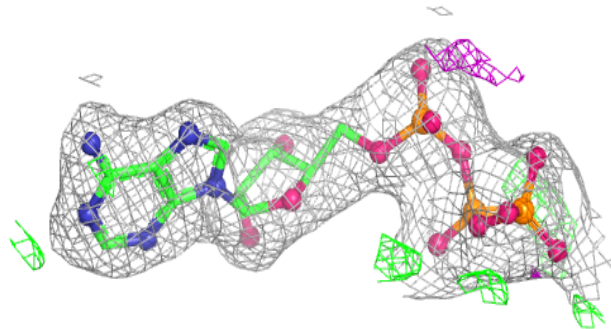
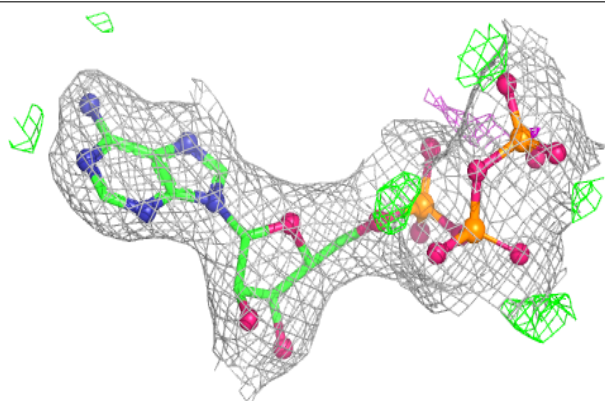
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

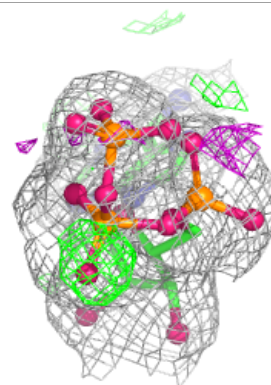
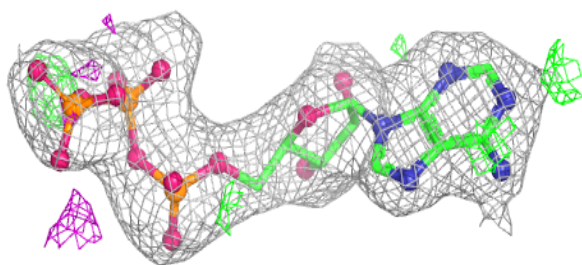
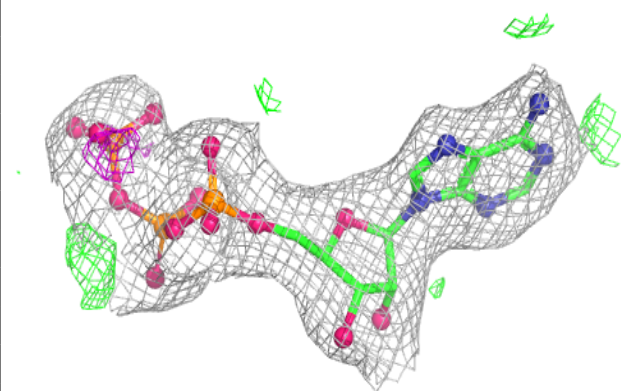
**Electron density around ATP C 600:**

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

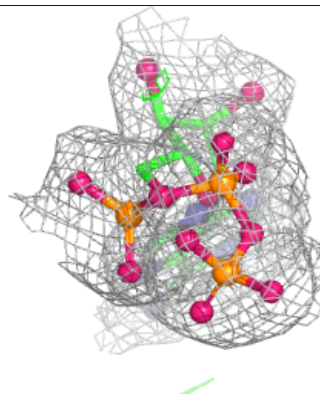
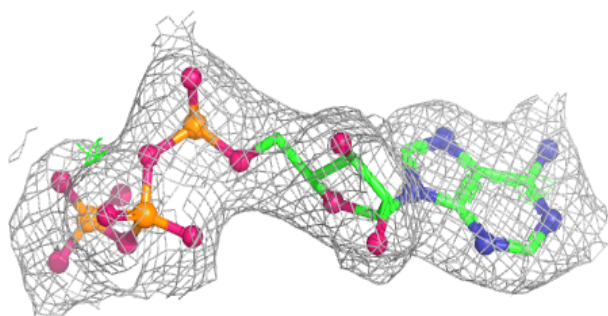
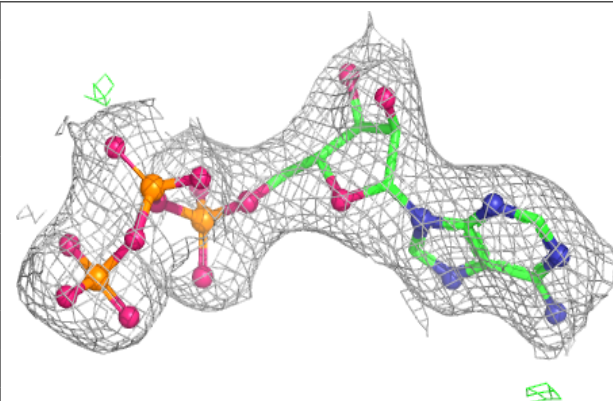


Electron density around ATP D 600:

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

**Electron density around ATP B 600:**

$2mF_o-DF_c$ (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.