



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 04:01 AM UTC

PDB ID : 3GBU / pdb_00003gbu
Title : Crystal structure of an uncharacterized sugar kinase PH1459 from *Pyrococcus horikoshii* in complex with ATP
Authors : Eswaramoorthy, S.; Kumar, G.; Zhang, Z.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2009-02-20
Resolution : 2.20 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

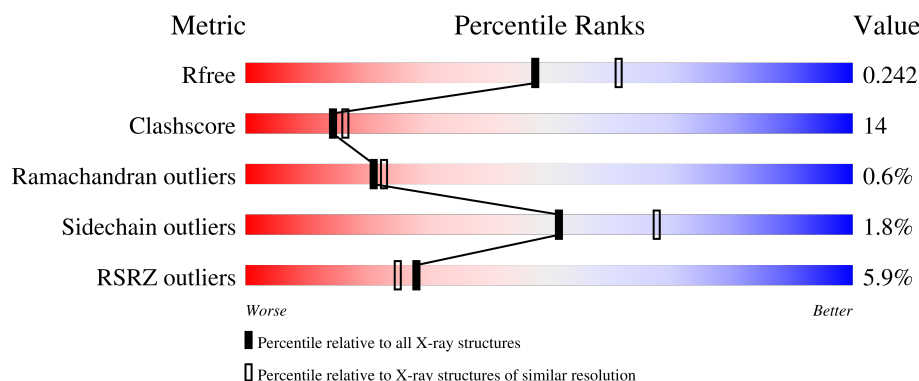
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	313	4% 73% 21% • •
1	B	313	2% 67% 27% • •
1	C	313	12% 66% 29% • •
1	D	313	6% 66% 28% • •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9963 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 Uncharacterized sugar kinase PH1459.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	302	Total	C	N	O	Se	0	0	0
			2336	1499	388	444	5			
1	B	302	Total	C	N	O	Se	0	0	0
			2336	1499	388	444	5			
1	C	302	Total	C	N	O	Se	0	0	0
			2336	1499	388	444	5			
1	D	302	Total	C	N	O	Se	0	0	0
			2336	1499	388	444	5			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	expression tag	UNP O59128
A	2	SER	-	expression tag	UNP O59128
A	3	LEU	-	expression tag	UNP O59128
A	221	PHE	CYS	engineered mutation	UNP O59128
A	306	GLU	-	expression tag	UNP O59128
A	307	GLY	-	expression tag	UNP O59128
A	308	HIS	-	expression tag	UNP O59128
A	309	HIS	-	expression tag	UNP O59128
A	310	HIS	-	expression tag	UNP O59128
A	311	HIS	-	expression tag	UNP O59128
A	312	HIS	-	expression tag	UNP O59128
A	313	HIS	-	expression tag	UNP O59128
B	1	MSE	-	expression tag	UNP O59128
B	2	SER	-	expression tag	UNP O59128
B	3	LEU	-	expression tag	UNP O59128
B	221	PHE	CYS	engineered mutation	UNP O59128
B	306	GLU	-	expression tag	UNP O59128
B	307	GLY	-	expression tag	UNP O59128
B	308	HIS	-	expression tag	UNP O59128
B	309	HIS	-	expression tag	UNP O59128
B	310	HIS	-	expression tag	UNP O59128

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Chain	Residue	Modelled	Actual	Comment	Reference
B	311	HIS	-	expression tag	UNP O59128
B	312	HIS	-	expression tag	UNP O59128
B	313	HIS	-	expression tag	UNP O59128
C	1	MSE	-	expression tag	UNP O59128
C	2	SER	-	expression tag	UNP O59128
C	3	LEU	-	expression tag	UNP O59128
C	221	PHE	CYS	engineered mutation	UNP O59128
C	306	GLU	-	expression tag	UNP O59128
C	307	GLY	-	expression tag	UNP O59128
C	308	HIS	-	expression tag	UNP O59128
C	309	HIS	-	expression tag	UNP O59128
C	310	HIS	-	expression tag	UNP O59128
C	311	HIS	-	expression tag	UNP O59128
C	312	HIS	-	expression tag	UNP O59128
C	313	HIS	-	expression tag	UNP O59128
D	1	MSE	-	expression tag	UNP O59128
D	2	SER	-	expression tag	UNP O59128
D	3	LEU	-	expression tag	UNP O59128
D	221	PHE	CYS	engineered mutation	UNP O59128
D	306	GLU	-	expression tag	UNP O59128
D	307	GLY	-	expression tag	UNP O59128
D	308	HIS	-	expression tag	UNP O59128
D	309	HIS	-	expression tag	UNP O59128
D	310	HIS	-	expression tag	UNP O59128
D	311	HIS	-	expression tag	UNP O59128
D	312	HIS	-	expression tag	UNP O59128
D	313	HIS	-	expression tag	UNP O59128

- 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 31	C 10	N 5	O 13	P 3	0	0
2	B	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	C	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	D	1	Total 31	C 10	N 5	O 13	P 3	0	0

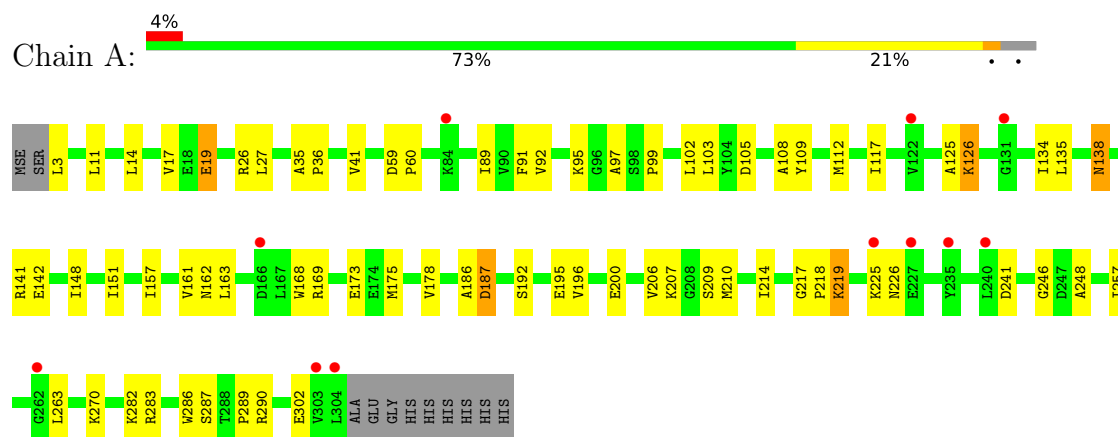
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	152	Total O 152 152	0	0
3	B	143	Total O 143 143	0	0
3	C	101	Total O 101 101	0	0
3	D	99	Total O 99 99	0	0

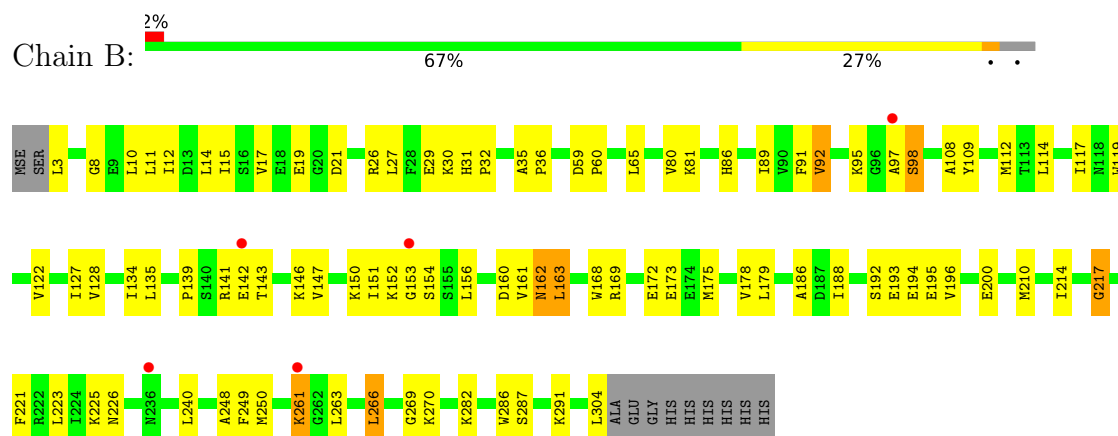
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

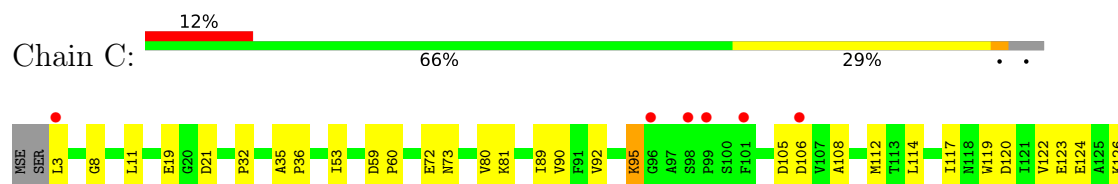
• Molecule 1: Uncharacterized sugar kinase PH1459

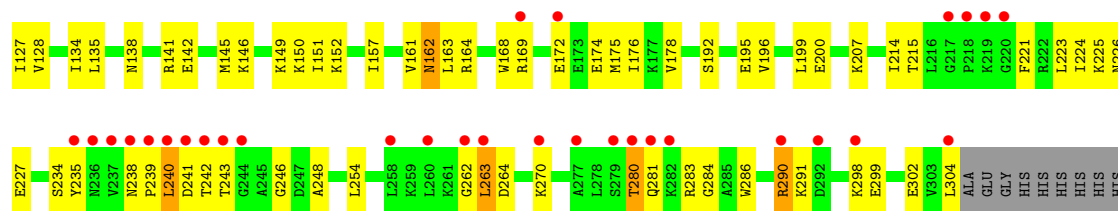


• Molecule 1: Uncharacterized sugar kinase PH1459

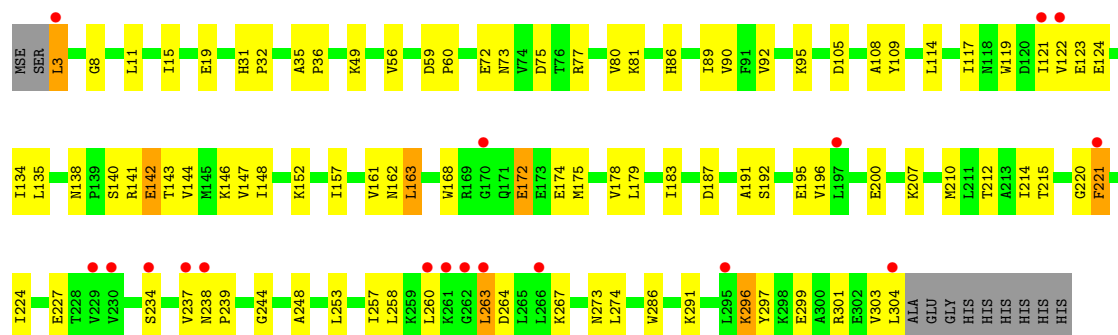


• Molecule 1: Uncharacterized sugar kinase PH1459





● Molecule 1: Uncharacterized sugar kinase PH1459



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.15Å 119.42Å 178.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.66 – 2.20 49.66 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.66-2.20) 100.0 (49.66-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.64 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.216 , 0.242 0.214 , 0.242	Depositor DCC
R_{free} test set	3946 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	33.2	Xtriage
Anisotropy	0.290	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9963	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.42	0/2363	0.93	10/3181 (0.3%)
1	B	0.45	1/2363 (0.0%)	0.92	9/3181 (0.3%)
1	C	0.39	0/2363	0.92	11/3181 (0.3%)
1	D	0.37	0/2363	0.91	8/3181 (0.3%)
All	All	0.41	1/9452 (0.0%)	0.92	38/12724 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	153	GLY	C-O	5.66	1.30	1.23

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	108	ALA	N-CA-C	7.97	120.67	111.11
1	D	108	ALA	N-CA-C	7.75	119.73	111.28
1	C	105	ASP	N-CA-C	7.34	122.41	113.38
1	A	105	ASP	N-CA-C	7.11	122.05	113.23
1	B	108	ALA	N-CA-C	6.87	119.35	111.11
1	C	108	ALA	N-CA-C	6.67	119.40	111.33
1	C	92	VAL	N-CA-C	6.29	117.54	108.48
1	C	240	LEU	N-CA-C	-6.00	107.78	114.62
1	C	126	LYS	N-CA-C	-5.95	105.11	112.90
1	A	126	LYS	N-CA-C	-5.90	106.25	113.50
1	D	105	ASP	N-CA-C	5.88	120.52	113.23
1	A	125	ALA	N-CA-C	5.83	118.94	110.42
1	B	117	ILE	N-CA-C	5.82	117.16	108.54
1	C	163	LEU	N-CA-C	5.82	118.15	109.59
1	C	280	THR	N-CA-C	-5.82	106.52	113.97
1	A	186	ALA	N-CA-C	5.70	119.54	109.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	138	ASN	N-CA-C	-5.61	103.05	110.40
1	A	248	ALA	N-CA-C	-5.61	105.25	111.36
1	B	154	SER	N-CA-C	-5.42	104.45	111.71
1	C	117	ILE	N-CA-C	5.40	116.01	108.89
1	D	238	ASN	CA-C-N	5.35	126.52	119.84
1	D	238	ASN	C-N-CA	5.35	126.52	119.84
1	D	248	ALA	N-CA-C	-5.34	105.37	111.14
1	D	187	ASP	N-CA-C	-5.25	106.02	112.90
1	B	92	VAL	N-CA-C	5.24	116.27	108.46
1	B	217	GLY	CA-C-N	5.21	125.01	119.28
1	B	217	GLY	C-N-CA	5.21	125.01	119.28
1	B	248	ALA	N-CA-C	-5.21	105.68	111.36
1	A	92	VAL	N-CA-C	5.12	116.09	108.46
1	B	186	ALA	N-CA-C	5.10	117.93	109.46
1	A	117	ILE	N-CA-C	5.10	115.69	107.73
1	C	248	ALA	N-CA-C	-5.07	105.83	111.36
1	D	163	LEU	N-CA-C	5.07	117.56	109.96
1	C	238	ASN	CA-C-N	5.06	126.16	119.84
1	C	238	ASN	C-N-CA	5.06	126.16	119.84
1	A	187	ASP	N-CA-C	-5.05	105.86	112.23
1	D	117	ILE	N-CA-C	5.05	115.55	108.89
1	B	240	LEU	N-CA-C	-5.03	105.45	112.45

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	2336	0	2436	56	0
1	B	2336	0	2436	73	0
1	C	2336	0	2436	76	0
1	D	2336	0	2436	75	0
2	A	31	0	12	2	0
2	B	31	0	12	1	0
2	C	31	0	12	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	31	0	12	3	0
3	A	152	0	0	7	0
3	B	143	0	0	6	0
3	C	101	0	0	6	0
3	D	99	0	0	3	0
All	All	9963	0	9792	271	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 14.

All (271) 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:163:LEU:HD11	1:D:175:MSE:HE2	1.41	1.02
1:D:237:VAL:O	1:D:239:PRO:HD3	1.66	0.95
1:C:290:ARG:HH11	1:C:290:ARG:HB3	1.33	0.93
1:B:128:VAL:HG21	1:B:151:ILE:HD12	1.52	0.90
1:D:168:TRP:CD1	1:D:175:MSE:HE3	2.08	0.88
1:C:138:ASN:ND2	1:C:142:GLU:HG2	1.92	0.84
1:B:168:TRP:CD1	1:B:175:MSE:HE3	2.12	0.84
1:C:270:LYS:HE2	1:C:270:LYS:HA	1.58	0.84
1:B:21:ASP:HB3	3:B:327:HOH:O	1.75	0.83
1:C:138:ASN:HD21	1:C:142:GLU:HG2	1.42	0.82
1:C:152:LYS:HA	1:C:157:ILE:HD11	1.62	0.80
1:C:168:TRP:CD1	1:C:175:MSE:HE3	2.18	0.79
1:C:291:LYS:HG3	1:C:304:LEU:HD13	1.64	0.79
1:C:112:MSE:HE3	1:C:134:ILE:HD11	1.66	0.76
1:C:284:GLY:HA2	3:C:336:HOH:O	1.87	0.75
1:D:163:LEU:CD1	1:D:175:MSE:HE2	2.17	0.75
1:C:128:VAL:HG21	1:C:151:ILE:HD12	1.68	0.74
1:B:135:LEU:O	1:B:141:ARG:HD3	1.87	0.74
1:B:139:PRO:O	1:B:142:GLU:HG2	1.87	0.74
1:A:138:ASN:HD21	1:A:142:GLU:HG3	1.51	0.73
1:D:95:LYS:HE3	3:D:402:HOH:O	1.89	0.72
1:B:152:LYS:O	1:B:152:LYS:HG3	1.88	0.71
1:C:240:LEU:HD23	1:C:281:GLN:O	1.90	0.71
1:A:141:ARG:HD2	1:A:178:VAL:HG11	1.71	0.71
1:D:135:LEU:O	1:D:141:ARG:HD3	1.90	0.71
1:C:21:ASP:HB3	3:C:335:HOH:O	1.90	0.70
1:C:119:TRP:O	1:C:123:GLU:HG3	1.92	0.70
1:B:291:LYS:HE3	1:B:304:LEU:HB3	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:LEU:HD11	1:B:175:MSE:HE2	1.75	0.69
1:D:3:LEU:HG	1:D:258:LEU:HD21	1.74	0.69
1:D:274:LEU:HD22	1:D:299:GLU:HB2	1.74	0.68
1:C:135:LEU:O	1:C:141:ARG:HD3	1.92	0.68
1:A:35:ALA:HB3	1:A:36:PRO:HD3	1.77	0.67
1:B:109:TYR:O	1:B:112:MSE:HE3	1.94	0.67
1:A:218:PRO:HG2	1:A:219:LYS:HE2	1.77	0.67
1:D:239:PRO:HG2	2:D:900:ATP:HN62	1.60	0.67
1:A:217:GLY:HA2	2:A:900:ATP:O2A	1.96	0.66
1:B:29:GLU:CD	1:B:31:HIS:HE2	2.04	0.66
1:D:196:VAL:HG21	1:D:214:ILE:HD12	1.78	0.66
1:D:148:ILE:HG23	1:D:157:ILE:HD13	1.77	0.65
1:B:173:GLU:HB2	3:B:369:HOH:O	1.95	0.65
1:C:192:SER:OG	1:C:195:GLU:HG3	1.95	0.65
1:D:172:GLU:CD	1:D:172:GLU:H	2.04	0.65
1:A:148:ILE:HD13	1:A:151:ILE:HD11	1.77	0.65
1:B:266:LEU:HD12	1:B:270:LYS:HE3	1.79	0.65
1:C:246:GLY:HA2	2:C:900:ATP:H5'2	1.78	0.64
1:D:210:MSE:HE1	1:D:257:ILE:HG21	1.78	0.64
1:B:119:TRP:HA	1:B:122:VAL:HG22	1.79	0.64
1:D:239:PRO:HG2	2:D:900:ATP:N6	2.14	0.63
1:D:192:SER:OG	1:D:195:GLU:HG3	1.98	0.63
1:A:168:TRP:CH2	1:A:178:VAL:HG21	2.34	0.63
1:C:290:ARG:HH11	1:C:290:ARG:CB	2.09	0.63
1:B:89:ILE:HD11	1:C:89:ILE:HD11	1.80	0.62
1:D:119:TRP:O	1:D:123:GLU:HG3	2.00	0.62
1:C:241:ASP:O	1:C:280:THR:HA	2.00	0.61
1:A:19:GLU:OE1	1:A:95:LYS:HD3	2.00	0.61
1:B:19:GLU:CD	1:D:19:GLU:HG3	2.26	0.61
1:B:221:PHE:HE1	1:B:223:LEU:HG	1.65	0.61
1:C:119:TRP:HA	1:C:122:VAL:HG22	1.83	0.61
1:A:196:VAL:HG21	1:A:214:ILE:HD12	1.83	0.61
1:B:17:VAL:CG2	1:B:27:LEU:HB2	2.31	0.61
1:A:163:LEU:CD1	1:A:175:MSE:HE2	2.30	0.60
1:B:97:ALA:O	1:B:98:SER:HB2	2.01	0.60
1:C:262:GLY:O	1:C:263:LEU:C	2.45	0.60
1:B:163:LEU:CD1	1:B:175:MSE:HE2	2.31	0.60
1:C:169:ARG:HA	3:C:353:HOH:O	2.02	0.59
1:C:242:THR:HG23	1:C:280:THR:HG23	1.84	0.59
1:D:138:ASN:ND2	1:D:142:GLU:H	1.99	0.59
1:B:3:LEU:N	3:B:360:HOH:O	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:246:GLY:HA3	2:C:900:ATP:O2B	2.02	0.59
1:C:112:MSE:HE3	1:C:134:ILE:CD1	2.31	0.59
1:C:95:LYS:H	1:C:95:LYS:HD3	1.67	0.58
1:D:138:ASN:ND2	1:D:141:ARG:HB3	2.18	0.58
1:C:242:THR:CG2	1:C:280:THR:HG23	2.34	0.58
1:A:163:LEU:HD11	1:A:175:MSE:HE2	1.85	0.58
1:C:11:LEU:HD13	1:C:90:VAL:HG12	1.86	0.58
1:D:221:PHE:CZ	1:D:273:ASN:HB2	2.38	0.58
1:A:218:PRO:HG2	1:A:219:LYS:CE	2.34	0.58
1:C:225:LYS:HD3	1:C:263:LEU:HD12	1.86	0.58
1:C:239:PRO:HB3	1:C:280:THR:HG21	1.83	0.58
1:B:19:GLU:HG2	1:D:19:GLU:OE2	2.04	0.58
1:D:297:TYR:O	1:D:301:ARG:HG3	2.03	0.58
1:A:270:LYS:HD2	3:A:411:HOH:O	2.04	0.57
1:B:128:VAL:HG21	1:B:151:ILE:CD1	2.31	0.57
1:D:49:LYS:HD2	1:D:49:LYS:N	2.17	0.57
1:D:221:PHE:HZ	1:D:273:ASN:HB2	1.69	0.57
1:A:282:LYS:HE3	1:A:289:PRO:HA	1.86	0.57
1:C:196:VAL:O	1:C:200:GLU:HG3	2.03	0.57
1:D:237:VAL:C	1:D:239:PRO:HD3	2.28	0.57
1:C:174:GLU:O	1:C:178:VAL:HG23	2.03	0.57
1:D:291:LYS:HG3	1:D:304:LEU:HD13	1.86	0.57
1:D:35:ALA:HB3	1:D:36:PRO:HD3	1.86	0.57
1:B:196:VAL:O	1:B:200:GLU:HG3	2.05	0.56
1:A:290:ARG:HD3	3:A:412:HOH:O	2.04	0.56
1:B:196:VAL:HG21	1:B:214:ILE:HD12	1.88	0.56
1:B:19:GLU:OE1	1:D:19:GLU:HG3	2.05	0.56
1:C:221:PHE:HE1	1:C:223:LEU:HG	1.71	0.55
1:A:151:ILE:HG13	1:A:157:ILE:HD11	1.87	0.55
1:C:19:GLU:HG2	1:C:95:LYS:CB	2.36	0.55
1:C:291:LYS:HE3	1:C:304:LEU:HB3	1.89	0.55
1:A:135:LEU:O	1:A:141:ARG:HD3	2.06	0.55
1:D:291:LYS:HE3	1:D:304:LEU:HB3	1.89	0.55
1:A:14:LEU:HD13	1:A:91:PHE:CE1	2.41	0.55
1:B:161:VAL:O	1:B:162:ASN:C	2.48	0.55
1:B:160:ASP:HB2	1:B:250:MSE:HE2	1.89	0.55
1:A:19:GLU:CD	1:A:95:LYS:HD3	2.31	0.54
1:C:128:VAL:HG21	1:C:151:ILE:CD1	2.36	0.54
1:D:196:VAL:O	1:D:200:GLU:HG3	2.07	0.54
1:D:210:MSE:HE1	1:D:257:ILE:CG2	2.37	0.54
1:A:26:ARG:NH1	1:C:283:ARG:HH11	2.05	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:GLY:HA2	2:B:900:ATP:O2A	2.08	0.54
1:A:196:VAL:O	1:A:200:GLU:HG3	2.07	0.54
1:A:246:GLY:HA3	2:A:900:ATP:O2B	2.08	0.54
1:B:10:LEU:HG	1:B:65:LEU:HD13	1.89	0.54
1:C:192:SER:HA	1:C:215:THR:OG1	2.06	0.54
1:A:59:ASP:HB2	1:A:60:PRO:CD	2.38	0.53
1:A:17:VAL:CG2	1:A:27:LEU:HB2	2.39	0.53
1:A:11:LEU:C	1:A:11:LEU:HD12	2.33	0.53
1:C:298:LYS:HG3	1:C:299:GLU:N	2.24	0.53
1:B:17:VAL:HG22	1:B:27:LEU:HB2	1.91	0.53
1:A:282:LYS:HD2	1:A:287:SER:O	2.09	0.53
1:D:72:GLU:O	1:D:73:ASN:HB2	2.09	0.53
1:B:89:ILE:HD11	1:C:89:ILE:CD1	2.40	0.52
1:A:169:ARG:HA	3:A:426:HOH:O	2.09	0.52
1:D:163:LEU:HD11	1:D:179:LEU:HD11	1.90	0.52
1:B:80:VAL:HG12	1:B:81:LYS:N	2.25	0.52
1:C:207:LYS:NZ	1:C:224:ILE:HG21	2.24	0.52
1:C:35:ALA:HB3	1:C:36:PRO:HD3	1.91	0.52
1:D:59:ASP:HB2	1:D:60:PRO:CD	2.40	0.52
1:D:161:VAL:O	1:D:162:ASN:C	2.52	0.52
1:A:3:LEU:N	3:A:359:HOH:O	2.43	0.52
1:C:152:LYS:O	1:C:152:LYS:HG2	2.08	0.52
1:A:17:VAL:HG22	1:A:27:LEU:HB2	1.92	0.51
1:D:142:GLU:HB3	3:D:628:HOH:O	2.09	0.51
1:C:299:GLU:N	1:C:299:GLU:OE1	2.36	0.51
1:B:35:ALA:HB3	1:B:36:PRO:HD3	1.93	0.51
1:C:8:GLY:HA3	1:C:36:PRO:HG2	1.92	0.51
1:D:221:PHE:N	1:D:221:PHE:CD2	2.79	0.51
1:B:221:PHE:CE1	1:B:223:LEU:HG	2.45	0.51
1:C:298:LYS:O	1:C:302:GLU:HG3	2.10	0.50
1:B:112:MSE:HE3	1:B:134:ILE:HD11	1.93	0.50
1:B:221:PHE:HZ	1:B:269:GLY:HA2	1.77	0.50
1:C:196:VAL:HG21	1:C:214:ILE:HD12	1.94	0.49
1:D:192:SER:HA	1:D:215:THR:OG1	2.12	0.49
1:B:225:LYS:O	1:B:226:ASN:HB2	2.13	0.49
1:A:173:GLU:HG3	3:A:329:HOH:O	2.11	0.49
1:A:225:LYS:HD2	1:A:263:LEU:HD12	1.94	0.49
1:D:207:LYS:NZ	1:D:224:ILE:HG21	2.28	0.49
1:A:126:LYS:HE3	3:A:388:HOH:O	2.12	0.49
1:B:86:HIS:HD2	3:C:528:HOH:O	1.96	0.49
1:B:29:GLU:HG2	1:B:31:HIS:NE2	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:TRP:CD1	1:A:175:MSE:HE3	2.48	0.48
1:D:220:GLY:HA2	1:D:234:SER:OG	2.12	0.48
1:D:75:ASP:OD1	1:D:77:ARG:HG3	2.13	0.48
1:B:282:LYS:HD2	1:B:287:SER:O	2.14	0.48
1:A:26:ARG:NH1	1:C:283:ARG:NH1	2.61	0.48
1:A:109:TYR:O	1:A:134:ILE:HD11	2.13	0.48
1:A:17:VAL:O	1:A:95:LYS:NZ	2.47	0.48
1:A:102:LEU:HD11	1:D:86:HIS:CD2	2.49	0.48
1:A:206:VAL:O	1:A:207:LYS:HG3	2.14	0.48
1:B:196:VAL:CG2	1:B:214:ILE:HD12	2.44	0.48
1:A:112:MSE:HE2	1:A:134:ILE:HD11	1.96	0.47
1:A:241:ASP:OD2	1:A:283:ARG:HG2	2.15	0.47
1:D:11:LEU:HD13	1:D:90:VAL:HG12	1.96	0.47
1:A:97:ALA:C	1:A:99:PRO:HD3	2.38	0.47
1:B:19:GLU:OE2	1:B:95:LYS:HD3	2.14	0.47
1:B:59:ASP:HB2	1:B:60:PRO:HD2	1.96	0.47
1:C:161:VAL:O	1:C:162:ASN:C	2.57	0.47
1:B:14:LEU:HD13	1:B:91:PHE:CE1	2.49	0.47
1:B:59:ASP:HB2	1:B:60:PRO:CD	2.45	0.47
1:C:192:SER:O	1:C:196:VAL:HG23	2.14	0.47
1:D:239:PRO:CG	2:D:900:ATP:N6	2.78	0.47
1:C:172:GLU:O	1:C:176:ILE:HG12	2.15	0.47
1:A:196:VAL:CG2	1:A:214:ILE:HD12	2.45	0.46
1:B:188:ILE:HG12	1:B:210:MSE:HE2	1.97	0.46
1:D:260:LEU:HD11	1:D:303:VAL:HG22	1.97	0.46
1:B:17:VAL:HG21	1:B:27:LEU:HB2	1.96	0.46
1:C:150:LYS:HB2	3:C:341:HOH:O	2.16	0.46
1:D:152:LYS:HD3	1:D:157:ILE:HD12	1.98	0.46
1:B:114:LEU:HD11	1:B:146:LYS:HD2	1.97	0.46
1:A:59:ASP:HB2	1:A:60:PRO:HD2	1.98	0.46
1:A:103:LEU:C	1:A:103:LEU:HD23	2.41	0.46
1:A:210:MSE:HG2	1:A:263:LEU:HD11	1.97	0.46
1:B:163:LEU:CD1	1:B:179:LEU:HD11	2.45	0.46
1:C:119:TRP:CZ3	1:C:150:LYS:HD3	2.50	0.46
1:C:270:LYS:HE2	1:C:270:LYS:CA	2.39	0.45
1:A:192:SER:OG	1:A:195:GLU:HG3	2.17	0.45
1:B:19:GLU:CD	1:B:19:GLU:H	2.25	0.45
1:A:138:ASN:ND2	1:A:142:GLU:HG3	2.25	0.45
1:C:119:TRP:O	1:C:122:VAL:HG22	2.17	0.45
1:D:148:ILE:CG2	1:D:157:ILE:HD13	2.43	0.45
1:C:145:MSE:O	1:C:149:LYS:HG3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:114:LEU:HD11	1:D:146:LYS:HD3	1.98	0.45
1:D:264:ASP:CG	1:D:267:LYS:HG3	2.42	0.45
1:C:11:LEU:C	1:C:11:LEU:HD12	2.41	0.45
1:B:194:GLU:OE1	1:B:194:GLU:N	2.43	0.45
1:C:72:GLU:O	1:C:73:ASN:HB2	2.17	0.45
1:D:11:LEU:HD23	1:D:109:TYR:CE2	2.52	0.44
1:A:210:MSE:HE1	1:A:257:ILE:HG21	1.99	0.44
1:D:210:MSE:HE3	1:D:263:LEU:HD21	2.00	0.44
1:D:119:TRP:HA	1:D:122:VAL:HG22	2.00	0.44
1:B:192:SER:OG	1:B:195:GLU:HG3	2.17	0.44
1:C:59:ASP:HB2	1:C:60:PRO:CD	2.48	0.44
1:D:174:GLU:O	1:D:178:VAL:HG23	2.18	0.44
1:B:80:VAL:CG1	1:B:81:LYS:N	2.80	0.44
1:C:120:ASP:O	1:C:124:GLU:HG3	2.17	0.44
1:A:89:ILE:HD11	1:D:89:ILE:HD11	2.00	0.44
1:B:36:PRO:HG3	3:B:333:HOH:O	2.18	0.43
1:A:302:GLU:HB3	3:A:343:HOH:O	2.19	0.43
1:D:121:ILE:O	1:D:124:GLU:HG2	2.18	0.43
1:C:196:VAL:CG2	1:C:214:ILE:HD12	2.48	0.43
1:B:11:LEU:HD12	1:B:11:LEU:C	2.43	0.43
1:C:270:LYS:HA	1:C:270:LYS:CE	2.38	0.43
1:D:253:LEU:O	1:D:257:ILE:HG13	2.18	0.43
1:A:187:ASP:HA	1:A:209:SER:OG	2.19	0.43
1:C:221:PHE:CE1	1:C:223:LEU:HG	2.51	0.43
1:B:12:ILE:HD13	1:B:65:LEU:HD21	2.01	0.43
1:B:261:LYS:HD2	1:B:261:LYS:O	2.19	0.43
1:D:296:LYS:NZ	1:D:296:LYS:HB3	2.34	0.43
1:B:15:ILE:HG12	1:B:92:VAL:CG2	2.49	0.43
1:D:8:GLY:HA3	1:D:36:PRO:HG2	2.00	0.43
1:A:41:VAL:HG21	1:A:286:TRP:CE2	2.54	0.43
1:D:56:VAL:O	1:D:81:LYS:HA	2.19	0.43
1:C:59:ASP:HB2	1:C:60:PRO:HD2	2.01	0.42
1:B:172:GLU:CD	1:B:172:GLU:H	2.26	0.42
1:B:221:PHE:CD2	1:B:249:PHE:HE1	2.37	0.42
1:C:32:PRO:HB2	1:C:286:TRP:CZ2	2.54	0.42
1:D:179:LEU:O	1:D:183:ILE:HG13	2.20	0.42
1:D:138:ASN:HD21	1:D:142:GLU:H	1.67	0.42
1:C:114:LEU:HD11	1:C:146:LYS:CD	2.50	0.42
1:D:212:THR:HB	1:D:224:ILE:HB	2.01	0.42
1:D:15:ILE:HD12	1:D:31:HIS:CD2	2.54	0.42
1:D:80:VAL:HG12	1:D:81:LYS:N	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:244:GLY:HA2	3:D:363:HOH:O	2.19	0.42
1:B:29:GLU:CD	1:B:31:HIS:NE2	2.75	0.42
1:B:32:PRO:HB2	1:B:286:TRP:CZ2	2.55	0.42
1:B:169:ARG:O	1:B:169:ARG:HG3	2.19	0.42
1:C:80:VAL:HG12	1:C:81:LYS:N	2.35	0.42
1:C:199:LEU:HD23	1:C:199:LEU:HA	1.90	0.42
1:B:29:GLU:HG3	3:B:377:HOH:O	2.19	0.42
1:B:127:ILE:HG12	1:B:156:LEU:HB2	2.01	0.42
1:D:59:ASP:HB2	1:D:60:PRO:HD2	2.00	0.42
1:A:225:LYS:HD2	1:A:263:LEU:CD1	2.50	0.42
1:D:264:ASP:OD2	1:D:267:LYS:HG3	2.19	0.42
1:A:225:LYS:O	1:A:226:ASN:HB2	2.20	0.42
1:B:143:THR:O	1:B:147:VAL:HG23	2.20	0.42
1:D:15:ILE:HD12	1:D:31:HIS:HD2	1.84	0.42
1:D:140:SER:O	1:D:144:VAL:HG23	2.19	0.41
1:D:191:ALA:O	1:D:214:ILE:HA	2.20	0.41
1:C:138:ASN:ND2	1:C:142:GLU:H	2.18	0.41
1:C:227:GLU:O	1:C:227:GLU:HG2	2.20	0.41
1:A:161:VAL:O	1:A:162:ASN:C	2.63	0.41
1:C:162:ASN:OD1	1:C:164:ARG:NH1	2.54	0.41
1:C:127:ILE:HG21	1:C:254:LEU:HD22	2.01	0.41
1:A:263:LEU:HD23	1:A:263:LEU:HA	1.94	0.41
1:B:26:ARG:NE	3:B:370:HOH:O	2.45	0.41
1:C:234:SER:O	1:C:235:TYR:C	2.62	0.41
1:B:193:GLU:HG2	1:B:194:GLU:OE1	2.20	0.41
1:B:168:TRP:CH2	1:B:178:VAL:HG21	2.55	0.41
1:B:221:PHE:HE1	1:B:223:LEU:CG	2.32	0.41
1:C:168:TRP:CE3	1:C:175:MSE:HB2	2.56	0.41
1:D:32:PRO:HB2	1:D:286:TRP:CZ2	2.55	0.41
1:D:163:LEU:CD1	1:D:179:LEU:HD11	2.51	0.41
1:D:163:LEU:HD12	1:D:163:LEU:HA	1.89	0.41
1:D:109:TYR:O	1:D:134:ILE:HD11	2.21	0.41
1:C:53:ILE:HD12	1:C:122:VAL:HG12	2.03	0.40
1:B:8:GLY:HA3	1:B:36:PRO:HG2	2.03	0.40
1:C:225:LYS:O	1:C:226:ASN:HB3	2.21	0.40
1:D:143:THR:O	1:D:147:VAL:HG23	2.21	0.40
1:B:86:HIS:CD2	3:C:528:HOH:O	2.72	0.40
1:B:168:TRP:CE3	1:B:175:MSE:HB2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	300/313 (96%)	290 (97%)	10 (3%)	0	100	100
1	B	300/313 (96%)	289 (96%)	8 (3%)	3 (1%)	12	11
1	C	300/313 (96%)	280 (93%)	17 (6%)	3 (1%)	12	11
1	D	300/313 (96%)	291 (97%)	8 (3%)	1 (0%)	36	42
All	All	1200/1252 (96%)	1150 (96%)	43 (4%)	7 (1%)	21	23

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	98	SER
1	C	243	THR
1	C	263	LEU
1	D	263	LEU
1	B	162	ASN
1	B	263	LEU
1	C	162	ASN

5.3.2 Protein sidechains [i](#)

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	257/260 (99%)	255 (99%)	2 (1%)	73	85
1	B	257/260 (99%)	252 (98%)	5 (2%)	50	66
1	C	257/260 (99%)	252 (98%)	5 (2%)	50	66

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	D	257/260 (99%)	250 (97%)	7 (3%)	39 53
All	All	1028/1040 (99%)	1009 (98%)	19 (2%)	51 68

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	GLU
1	A	219	LYS
1	B	30	LYS
1	B	150	LYS
1	B	163	LEU
1	B	261	LYS
1	B	266	LEU
1	C	3	LEU
1	C	95	LYS
1	C	106	ASP
1	C	264	ASP
1	C	290	ARG
1	D	3	LEU
1	D	92	VAL
1	D	142	GLU
1	D	172	GLU
1	D	221	PHE
1	D	227	GLU
1	D	296	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	86	HIS
1	A	111	ASN
1	A	115	ASN
1	A	138	ASN
1	A	238	ASN
1	A	281	GLN
1	B	115	ASN
1	B	129	ASN
1	B	138	ASN
1	B	201	ASN
1	B	226	ASN
1	B	238	ASN

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Mol	Chain	Res	Type
1	C	129	ASN
1	C	138	ASN
1	C	201	ASN
1	C	202	GLN
1	C	273	ASN
1	D	111	ASN
1	D	138	ASN
1	D	202	GLN
1	D	226	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](#)

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	C	900	-	32,33,33	1.96	8 (25%)	48,52,52	1.87	11 (22%)
2	ATP	D	900	-	32,33,33	1.71	5 (15%)	48,52,52	1.84	12 (25%)
2	ATP	B	900	-	32,33,33	1.71	6 (18%)	48,52,52	1.76	10 (20%)
2	ATP	A	900	-	32,33,33	1.89	8 (25%)	48,52,52	1.77	11 (22%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ATP	C	900	-	-	7/22/38/38	0/3/3/3
2	ATP	D	900	-	-	4/22/38/38	0/3/3/3
2	ATP	B	900	-	-	7/22/38/38	0/3/3/3
2	ATP	A	900	-	-	7/22/38/38	0/3/3/3

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	900	ATP	C5-C4	5.60	1.49	1.39
2	A	900	ATP	C5-C4	5.60	1.49	1.39
2	C	900	ATP	C5-C4	5.41	1.48	1.39
2	C	900	ATP	PA-O3A	5.11	1.65	1.59
2	D	900	ATP	C5-C4	5.02	1.48	1.39
2	D	900	ATP	PA-O3A	4.61	1.64	1.59
2	A	900	ATP	PA-O3A	4.53	1.64	1.59
2	B	900	ATP	PA-O3A	3.86	1.63	1.59
2	C	900	ATP	PB-O3B	3.78	1.63	1.59
2	A	900	ATP	PB-O3B	3.52	1.63	1.59
2	C	900	ATP	C5-C6	3.00	1.49	1.41
2	B	900	ATP	C5-N7	-2.91	1.33	1.39
2	C	900	ATP	C5-N7	-2.82	1.33	1.39
2	A	900	ATP	C5-C6	2.82	1.48	1.41
2	D	900	ATP	C5-N7	-2.75	1.34	1.39
2	D	900	ATP	C5-C6	2.65	1.48	1.41
2	A	900	ATP	PB-O3A	2.56	1.62	1.59
2	A	900	ATP	C8-N7	2.56	1.36	1.31
2	D	900	ATP	PB-O3B	2.52	1.62	1.59
2	B	900	ATP	PB-O3B	2.50	1.62	1.59
2	C	900	ATP	PB-O3A	2.43	1.62	1.59
2	B	900	ATP	C5-C6	2.37	1.47	1.41
2	A	900	ATP	C5-N7	-2.36	1.34	1.39
2	C	900	ATP	C8-N7	2.34	1.36	1.31
2	C	900	ATP	C4-N3	2.25	1.38	1.34
2	B	900	ATP	C8-N7	2.23	1.36	1.31
2	A	900	ATP	C4-N3	2.11	1.38	1.34

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	900	ATP	C5-C4-N3	-5.48	119.18	126.72
2	B	900	ATP	C5-C4-N3	-5.10	119.70	126.72
2	D	900	ATP	C5-C4-N3	-4.94	119.92	126.72
2	A	900	ATP	C5-C4-N3	-4.78	120.14	126.72
2	D	900	ATP	N3-C2-N1	-4.52	121.73	128.58
2	C	900	ATP	N3-C4-N9	4.43	134.70	127.17
2	A	900	ATP	N3-C2-N1	-4.42	121.89	128.58
2	C	900	ATP	N3-C2-N1	-4.37	121.97	128.58
2	B	900	ATP	N3-C2-N1	-4.37	121.97	128.58
2	D	900	ATP	N3-C4-N9	4.25	134.40	127.17
2	B	900	ATP	N3-C4-N9	4.14	134.21	127.17
2	C	900	ATP	C2-N3-C4	3.96	121.51	111.83
2	D	900	ATP	C2-N3-C4	3.89	121.32	111.83
2	A	900	ATP	N3-C4-N9	3.84	133.70	127.17
2	B	900	ATP	C2-N3-C4	3.78	121.07	111.83
2	A	900	ATP	C2-N3-C4	3.78	121.06	111.83
2	D	900	ATP	C2'-C1'-N9	-3.58	104.40	113.30
2	C	900	ATP	C4-C5-N7	-3.48	106.61	110.58
2	C	900	ATP	O4'-C1'-N9	3.48	114.76	108.09
2	A	900	ATP	C4-C5-N7	-3.01	107.14	110.58
2	C	900	ATP	C5-N7-C8	2.95	108.08	103.45
2	B	900	ATP	C2'-C1'-N9	-2.85	106.22	113.30
2	B	900	ATP	C4-C5-N7	-2.85	107.32	110.58
2	A	900	ATP	C2'-C1'-N9	-2.83	106.28	113.30
2	A	900	ATP	C6-C5-N7	2.65	137.21	132.09
2	D	900	ATP	C4-C5-N7	-2.65	107.56	110.58
2	C	900	ATP	C6-C5-N7	2.61	137.12	132.09
2	A	900	ATP	O4'-C1'-N9	2.60	113.08	108.09
2	A	900	ATP	C2-N1-C6	2.47	122.78	118.73
2	D	900	ATP	C6-C5-N7	2.45	136.81	132.09
2	B	900	ATP	C6-C5-N7	2.43	136.77	132.09
2	D	900	ATP	O4'-C1'-N9	2.42	112.73	108.09
2	B	900	ATP	C2-N1-C6	2.39	122.66	118.73
2	D	900	ATP	C2-N1-C6	2.34	122.58	118.73
2	A	900	ATP	C5-N7-C8	2.27	107.02	103.45
2	D	900	ATP	C4-N9-C8	2.25	108.11	105.74
2	B	900	ATP	C5-N7-C8	2.25	106.99	103.45
2	A	900	ATP	C4-N9-C8	2.22	108.07	105.74
2	C	900	ATP	C2'-C1'-N9	-2.17	107.92	113.30
2	B	900	ATP	O4'-C1'-N9	2.15	112.22	108.09
2	C	900	ATP	C2-N1-C6	2.13	122.23	118.73
2	D	900	ATP	C5-N7-C8	2.11	106.76	103.45
2	D	900	ATP	O4'-C4'-C5'	-2.02	102.87	109.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	900	ATP	C4-N9-C8	2.01	107.85	105.74

There are no chirality outliers.

All (25) torsion outliers are listed below:

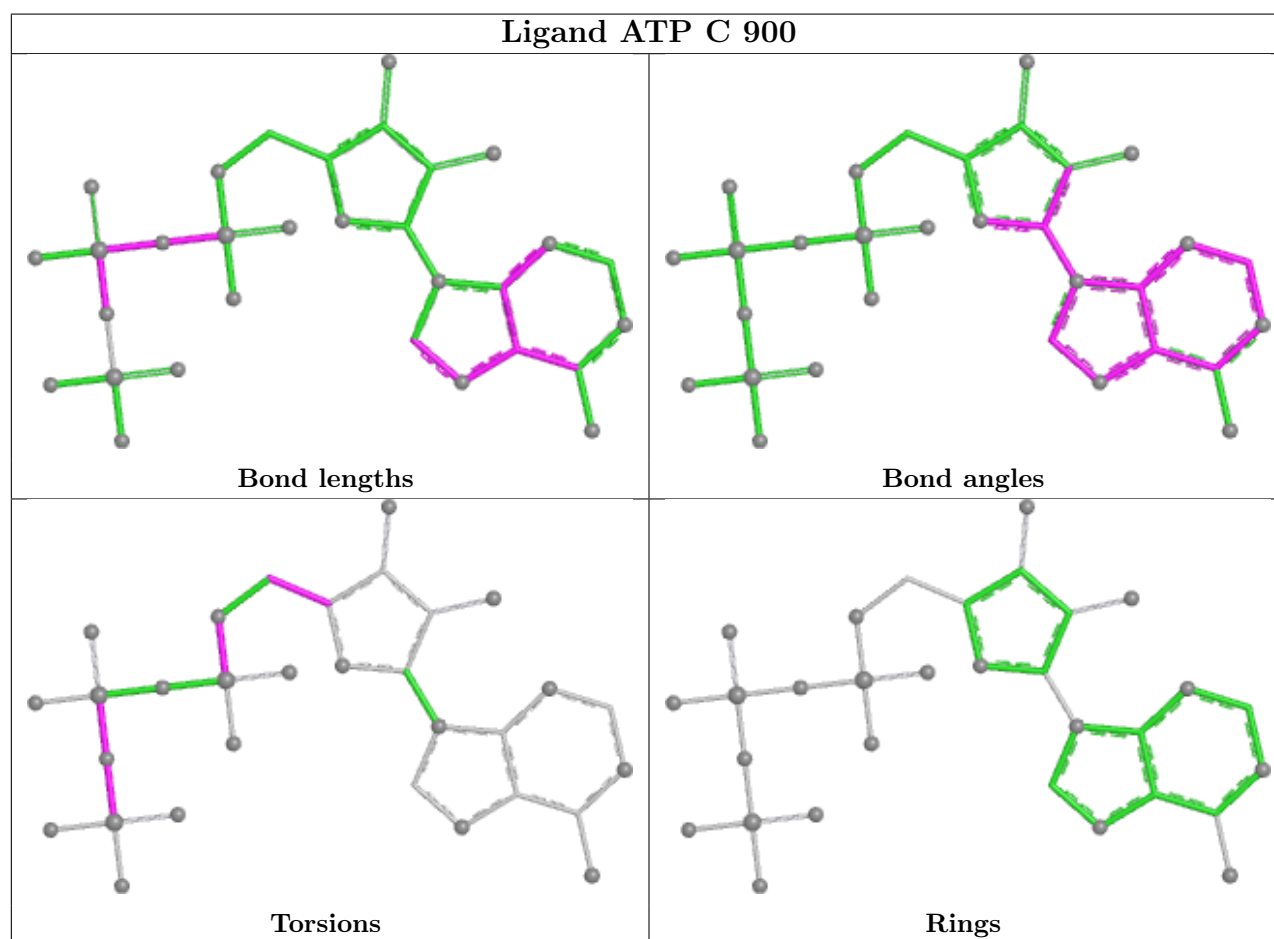
Mol	Chain	Res	Type	Atoms
2	A	900	ATP	C5'-O5'-PA-O1A
2	A	900	ATP	C5'-O5'-PA-O3A
2	B	900	ATP	C5'-O5'-PA-O1A
2	C	900	ATP	C5'-O5'-PA-O1A
2	C	900	ATP	C5'-O5'-PA-O2A
2	C	900	ATP	C5'-O5'-PA-O3A
2	A	900	ATP	C3'-C4'-C5'-O5'
2	B	900	ATP	C3'-C4'-C5'-O5'
2	C	900	ATP	C3'-C4'-C5'-O5'
2	D	900	ATP	C3'-C4'-C5'-O5'
2	A	900	ATP	O4'-C4'-C5'-O5'
2	B	900	ATP	O4'-C4'-C5'-O5'
2	C	900	ATP	O4'-C4'-C5'-O5'
2	D	900	ATP	O4'-C4'-C5'-O5'
2	A	900	ATP	PB-O3B-PG-O1G
2	A	900	ATP	PB-O3B-PG-O3G
2	A	900	ATP	C5'-O5'-PA-O2A
2	B	900	ATP	C5'-O5'-PA-O3A
2	D	900	ATP	PA-O3A-PB-O2B
2	C	900	ATP	PB-O3B-PG-O1G
2	B	900	ATP	PG-O3B-PB-O2B
2	B	900	ATP	PA-O3A-PB-O1B
2	C	900	ATP	PG-O3B-PB-O2B
2	D	900	ATP	PG-O3B-PB-O1B
2	B	900	ATP	PG-O3B-PB-O1B

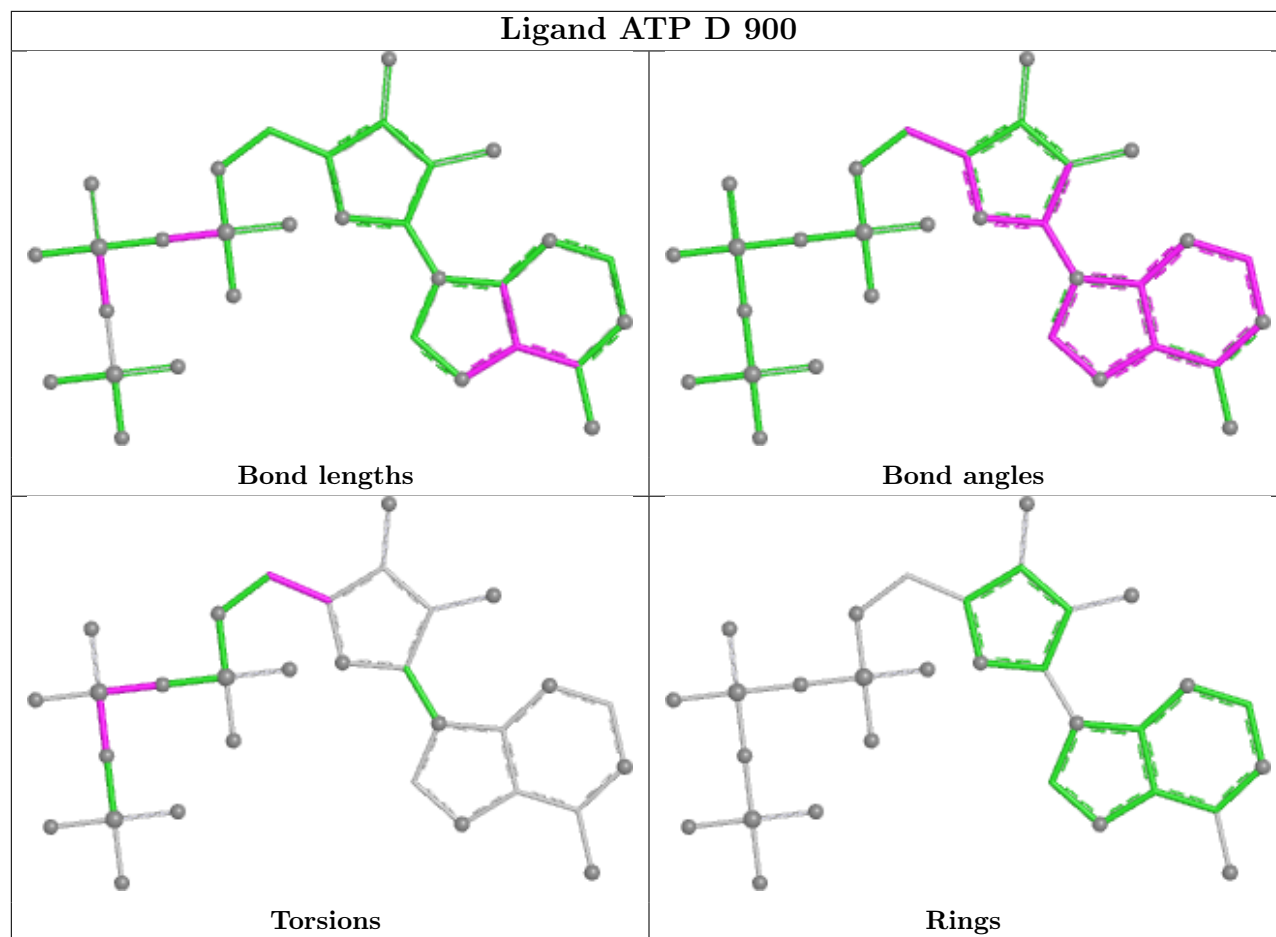
There are no ring outliers.

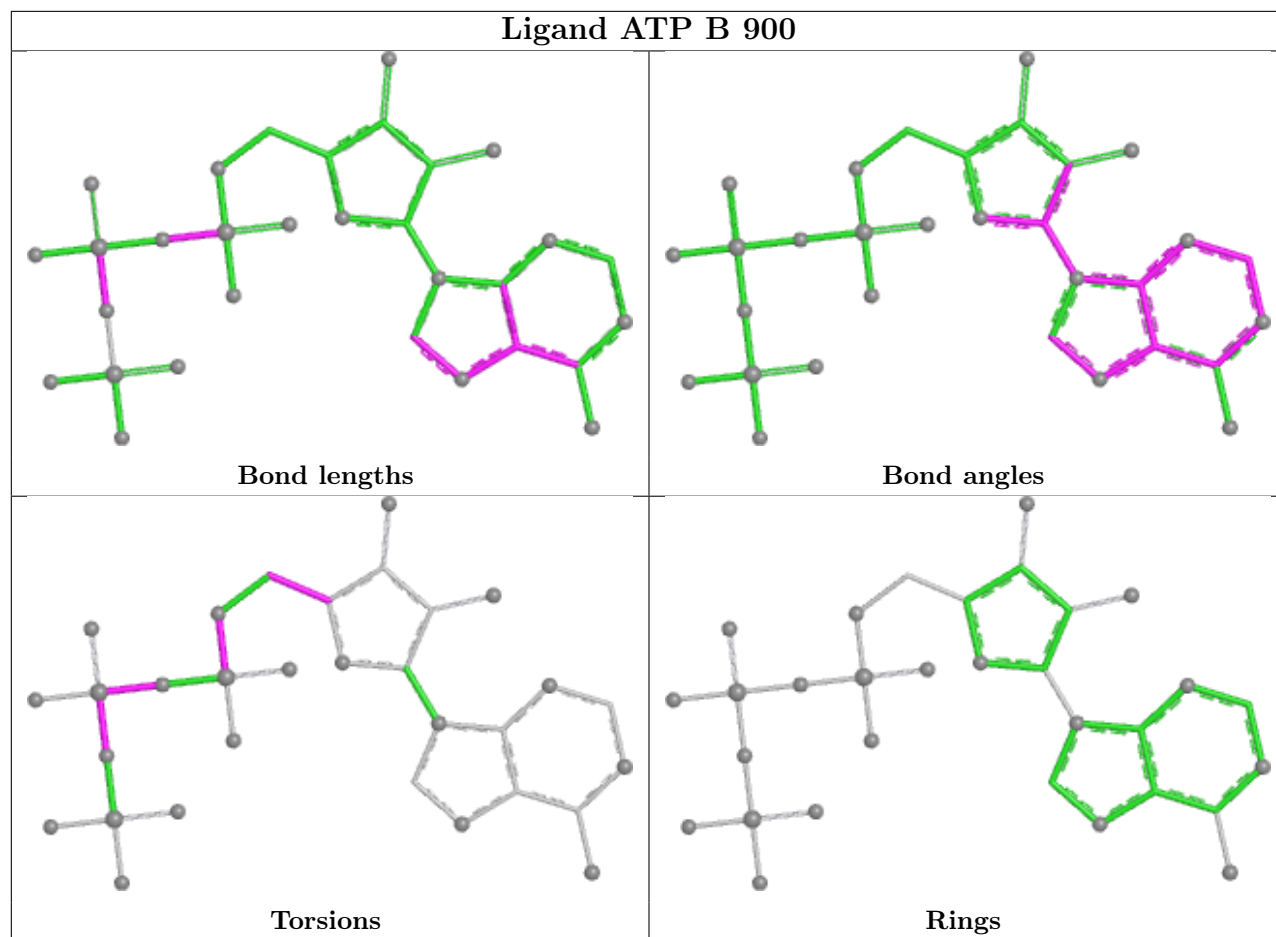
4 monomers are involved in 8 short contacts:

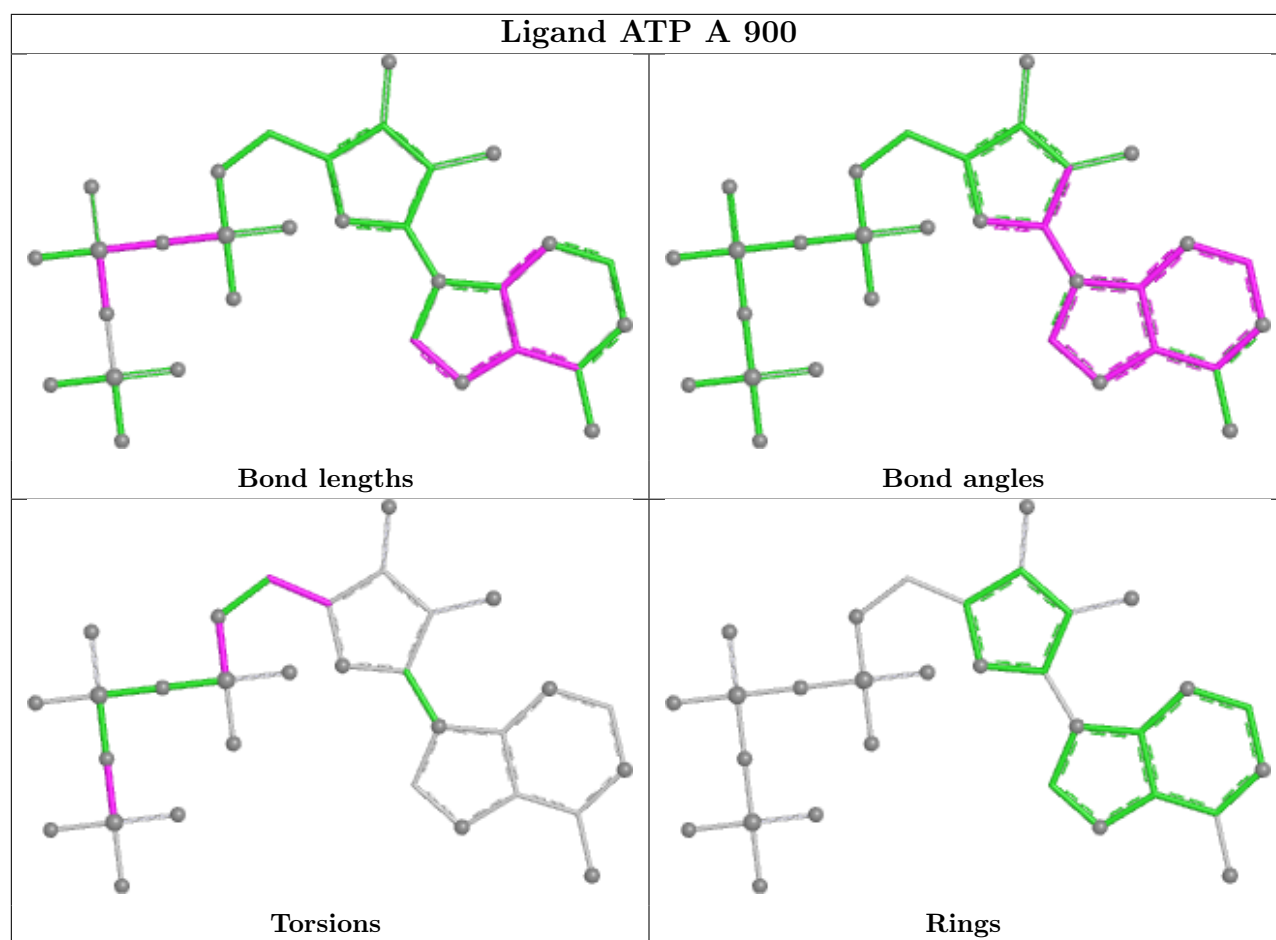
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	900	ATP	2	0
2	D	900	ATP	3	0
2	B	900	ATP	1	0
2	A	900	ATP	2	0

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	297/313 (94%)	0.02	11 (3%) 45 42	17, 29, 52, 66	0
1	B	297/313 (94%)	0.04	5 (1%) 69 66	18, 31, 51, 58	0
1	C	297/313 (94%)	0.56	36 (12%) 8 6	20, 38, 60, 70	0
1	D	297/313 (94%)	0.49	18 (6%) 27 24	22, 41, 60, 70	0
All	All	1188/1252 (94%)	0.28	70 (5%) 28 25	17, 35, 58, 70	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	240	LEU	5.0
1	C	238	ASN	3.9
1	C	237	VAL	3.9
1	C	236	ASN	3.6
1	A	240	LEU	3.6
1	C	282	LYS	3.4
1	C	218	PRO	3.3
1	C	243	THR	3.3
1	C	235	TYR	3.3
1	C	279	SER	3.3
1	C	244	GLY	3.2
1	A	304	LEU	3.2
1	C	262	GLY	3.2
1	D	3	LEU	3.2
1	C	260	LEU	3.2
1	C	304	LEU	3.1
1	D	238	ASN	3.0
1	C	219	LYS	3.0
1	C	101	PHE	2.9
1	D	234	SER	2.9
1	D	122	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	229	VAL	2.9
1	B	261	LYS	2.8
1	D	230	VAL	2.8
1	D	237	VAL	2.8
1	C	281	GLN	2.8
1	C	270	LYS	2.7
1	D	304	LEU	2.7
1	C	106	ASP	2.7
1	A	235	TYR	2.6
1	C	169	ARG	2.6
1	A	225	LYS	2.6
1	C	217	GLY	2.6
1	D	197	LEU	2.6
1	D	295	LEU	2.6
1	D	262	GLY	2.6
1	B	97	ALA	2.5
1	C	99	PRO	2.5
1	C	239	PRO	2.5
1	C	258	LEU	2.5
1	C	220	GLY	2.5
1	C	280	THR	2.5
1	A	122	VAL	2.5
1	C	3	LEU	2.5
1	B	236	ASN	2.5
1	C	241	ASP	2.4
1	C	96	GLY	2.4
1	D	221	PHE	2.4
1	C	172	GLU	2.4
1	A	131	GLY	2.3
1	D	261	LYS	2.3
1	A	303	VAL	2.3
1	A	262	GLY	2.3
1	C	263	LEU	2.3
1	D	266	LEU	2.3
1	C	292	ASP	2.3
1	C	98	SER	2.2
1	A	227	GLU	2.2
1	C	298	LYS	2.2
1	D	121	ILE	2.1
1	A	166	ASP	2.1
1	C	277	ALA	2.1
1	D	260	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	263	LEU	2.1
1	B	153	GLY	2.1
1	D	170	GLY	2.1
1	C	242	THR	2.1
1	A	84	LYS	2.1
1	B	142	GLU	2.1
1	C	290	ARG	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](#)

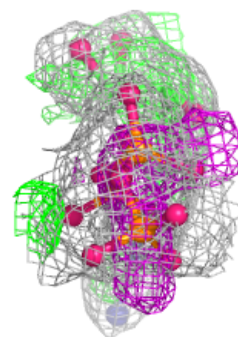
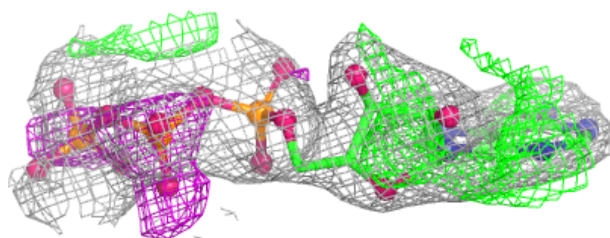
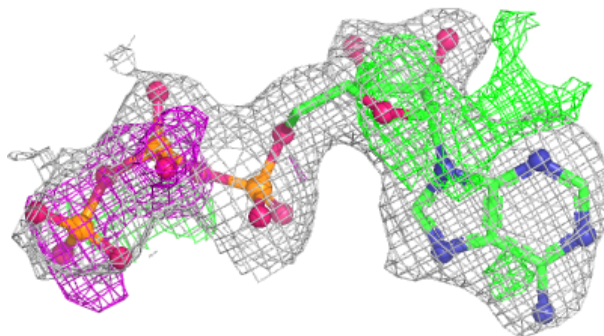
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	ATP	A	900	31/31	0.71	0.22	31,35,74,75	18
2	ATP	C	900	31/31	0.73	0.20	44,47,68,69	18
2	ATP	B	900	31/31	0.77	0.19	21,26,64,65	18
2	ATP	D	900	31/31	0.77	0.18	28,35,70,71	18

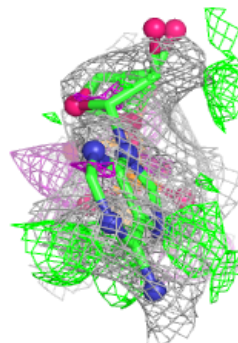
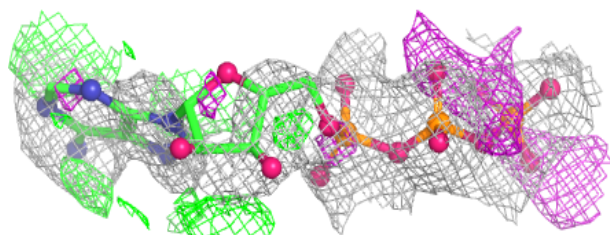
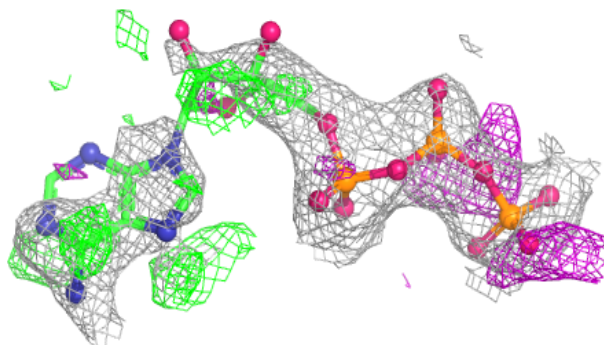
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ATP A 900:

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

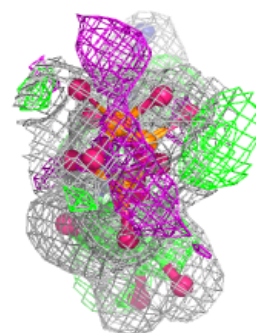
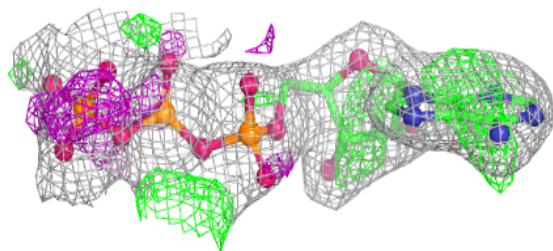
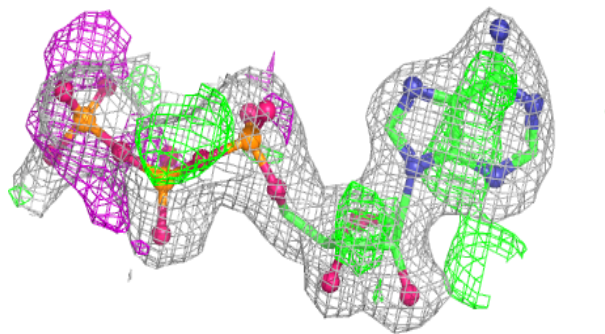
**Electron density around ATP C 900:**

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

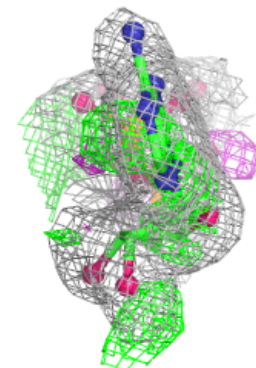
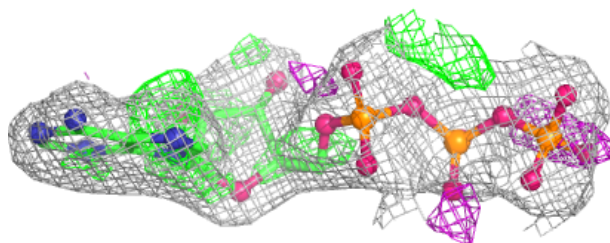
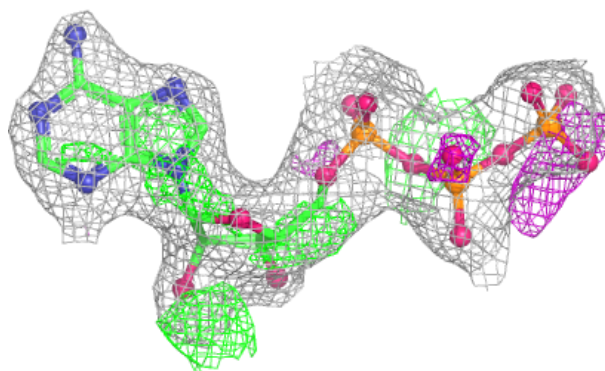


Electron density around ATP B 900:

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

**Electron density around ATP D 900:**

$2mF_o-DF_c$ (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.