



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:47 AM UTC

PDB ID : 2XSZ / pdb_00002xsz
Title : The dodecameric human RuvBL1:RuvBL2 complex with truncated domains II
Authors : Gorynia, S.; Bandejas, T.M.; Matias, P.M.; Pinho, F.G.; McVey, C.E.; Vonrhein, C.; Svergun, D.I.; Round, A.; Donner, P.; Carrondo, M.A.
Deposited on : 2010-10-01
Resolution : 3.00 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

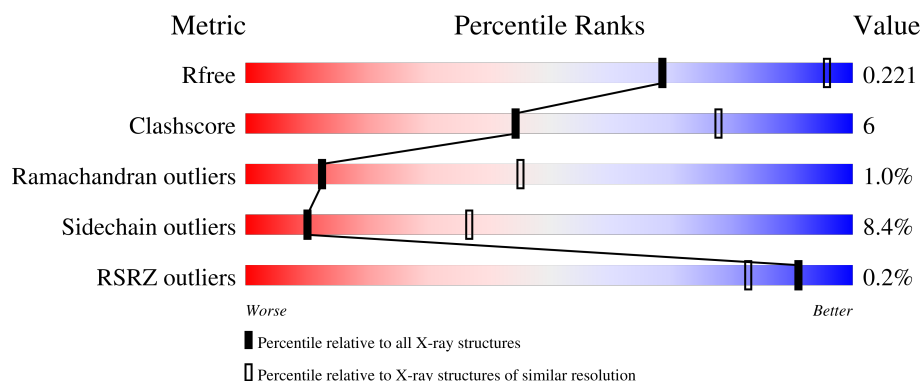
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

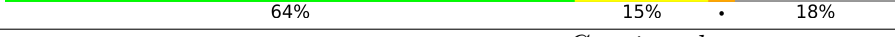
The reported resolution of this entry is 3.00 Å.

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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	367	
1	B	367	
1	C	367	
2	D	378	

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Mol	Chain	Length	Quality of chain
2	E	378	61%20%17%
2	F	378	64%15%18%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14664 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 RUVB-LIKE 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	Se	0	0	0
			2430	1530	421	468	4	7			
1	B	319	Total	C	N	O	S	Se	0	0	0
			2435	1531	422	471	4	7			
1	C	320	Total	C	N	O	S	Se	0	0	0
			2425	1525	420	469	4	7			

There are 45 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	expression tag	UNP Q9Y265
A	2	VAL	-	expression tag	UNP Q9Y265
A	3	HIS	-	expression tag	UNP Q9Y265
A	4	HIS	-	expression tag	UNP Q9Y265
A	5	HIS	-	expression tag	UNP Q9Y265
A	6	HIS	-	expression tag	UNP Q9Y265
A	7	HIS	-	expression tag	UNP Q9Y265
A	8	HIS	-	expression tag	UNP Q9Y265
A	9	LEU	-	expression tag	UNP Q9Y265
A	10	LEU	-	expression tag	UNP Q9Y265
A	11	VAL	-	expression tag	UNP Q9Y265
A	12	PRO	-	expression tag	UNP Q9Y265
A	13	ARG	-	expression tag	UNP Q9Y265
A	14	GLY	-	expression tag	UNP Q9Y265
A	15	SER	-	expression tag	UNP Q9Y265
B	1	MSE	-	expression tag	UNP Q9Y265
B	2	VAL	-	expression tag	UNP Q9Y265
B	3	HIS	-	expression tag	UNP Q9Y265
B	4	HIS	-	expression tag	UNP Q9Y265
B	5	HIS	-	expression tag	UNP Q9Y265
B	6	HIS	-	expression tag	UNP Q9Y265
B	7	HIS	-	expression tag	UNP Q9Y265
B	8	HIS	-	expression tag	UNP Q9Y265

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Chain	Residue	Modelled	Actual	Comment	Reference
B	9	LEU	-	expression tag	UNP Q9Y265
B	10	LEU	-	expression tag	UNP Q9Y265
B	11	VAL	-	expression tag	UNP Q9Y265
B	12	PRO	-	expression tag	UNP Q9Y265
B	13	ARG	-	expression tag	UNP Q9Y265
B	14	GLY	-	expression tag	UNP Q9Y265
B	15	SER	-	expression tag	UNP Q9Y265
C	1	MSE	-	expression tag	UNP Q9Y265
C	2	VAL	-	expression tag	UNP Q9Y265
C	3	HIS	-	expression tag	UNP Q9Y265
C	4	HIS	-	expression tag	UNP Q9Y265
C	5	HIS	-	expression tag	UNP Q9Y265
C	6	HIS	-	expression tag	UNP Q9Y265
C	7	HIS	-	expression tag	UNP Q9Y265
C	8	HIS	-	expression tag	UNP Q9Y265
C	9	LEU	-	expression tag	UNP Q9Y265
C	10	LEU	-	expression tag	UNP Q9Y265
C	11	VAL	-	expression tag	UNP Q9Y265
C	12	PRO	-	expression tag	UNP Q9Y265
C	13	ARG	-	expression tag	UNP Q9Y265
C	14	GLY	-	expression tag	UNP Q9Y265
C	15	SER	-	expression tag	UNP Q9Y265

- Molecule 2 is a protein called RUVB-LIKE 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	D	310	Total	C	N	O	S	Se	0	0	0
			2377	1487	421	457	2	10			
2	E	315	Total	C	N	O	S	Se	0	0	0
			2426	1520	428	466	2	10			
2	F	311	Total	C	N	O	S	Se	0	0	0
			2385	1493	422	458	2	10			

There are 48 discrepancies between the modelled and reference sequences:

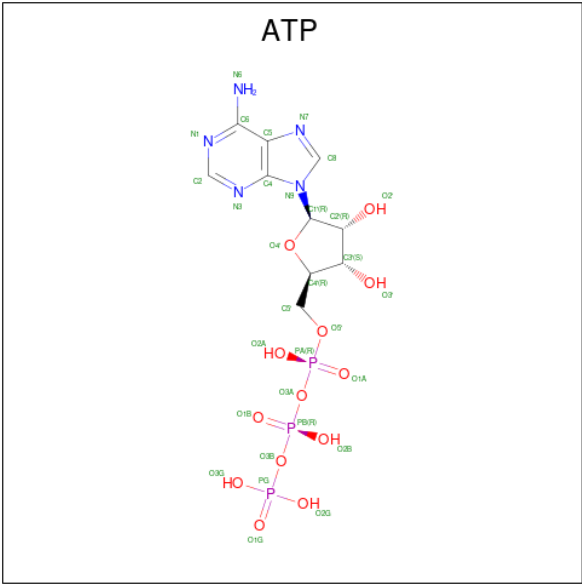
Chain	Residue	Modelled	Actual	Comment	Reference
D	1	MSE	-	expression tag	UNP Q9Y230
D	2	ASP	-	expression tag	UNP Q9Y230
D	3	TYR	-	expression tag	UNP Q9Y230
D	4	LYS	-	expression tag	UNP Q9Y230
D	5	ASP	-	expression tag	UNP Q9Y230
D	6	ASP	-	expression tag	UNP Q9Y230

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Chain	Residue	Modelled	Actual	Comment	Reference
D	7	ASP	-	expression tag	UNP Q9Y230
D	8	ASP	-	expression tag	UNP Q9Y230
D	9	LYS	-	expression tag	UNP Q9Y230
D	10	GLU	-	expression tag	UNP Q9Y230
D	11	ASN	-	expression tag	UNP Q9Y230
D	12	LEU	-	expression tag	UNP Q9Y230
D	13	TYR	-	expression tag	UNP Q9Y230
D	14	PHE	-	expression tag	UNP Q9Y230
D	15	GLN	-	expression tag	UNP Q9Y230
D	16	GLY	-	expression tag	UNP Q9Y230
E	1	MSE	-	expression tag	UNP Q9Y230
E	2	ASP	-	expression tag	UNP Q9Y230
E	3	TYR	-	expression tag	UNP Q9Y230
E	4	LYS	-	expression tag	UNP Q9Y230
E	5	ASP	-	expression tag	UNP Q9Y230
E	6	ASP	-	expression tag	UNP Q9Y230
E	7	ASP	-	expression tag	UNP Q9Y230
E	8	ASP	-	expression tag	UNP Q9Y230
E	9	LYS	-	expression tag	UNP Q9Y230
E	10	GLU	-	expression tag	UNP Q9Y230
E	11	ASN	-	expression tag	UNP Q9Y230
E	12	LEU	-	expression tag	UNP Q9Y230
E	13	TYR	-	expression tag	UNP Q9Y230
E	14	PHE	-	expression tag	UNP Q9Y230
E	15	GLN	-	expression tag	UNP Q9Y230
E	16	GLY	-	expression tag	UNP Q9Y230
F	1	MSE	-	expression tag	UNP Q9Y230
F	2	ASP	-	expression tag	UNP Q9Y230
F	3	TYR	-	expression tag	UNP Q9Y230
F	4	LYS	-	expression tag	UNP Q9Y230
F	5	ASP	-	expression tag	UNP Q9Y230
F	6	ASP	-	expression tag	UNP Q9Y230
F	7	ASP	-	expression tag	UNP Q9Y230
F	8	ASP	-	expression tag	UNP Q9Y230
F	9	LYS	-	expression tag	UNP Q9Y230
F	10	GLU	-	expression tag	UNP Q9Y230
F	11	ASN	-	expression tag	UNP Q9Y230
F	12	LEU	-	expression tag	UNP Q9Y230
F	13	TYR	-	expression tag	UNP Q9Y230
F	14	PHE	-	expression tag	UNP Q9Y230
F	15	GLN	-	expression tag	UNP Q9Y230
F	16	GLY	-	expression tag	UNP Q9Y230

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



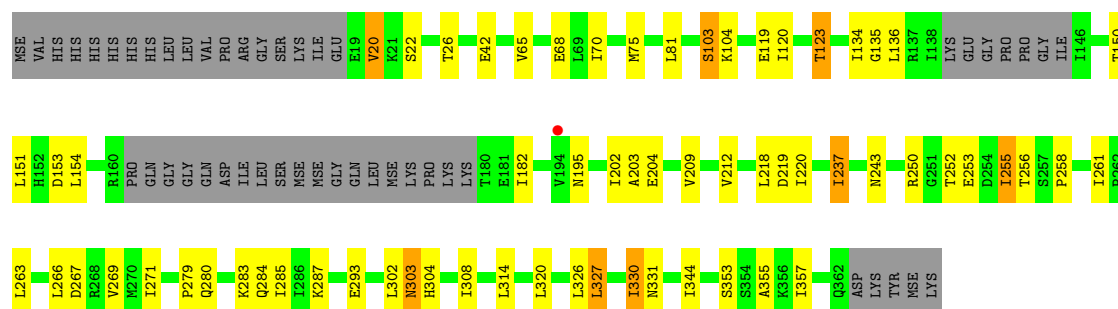
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	C	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	D	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	E	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	F	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

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.

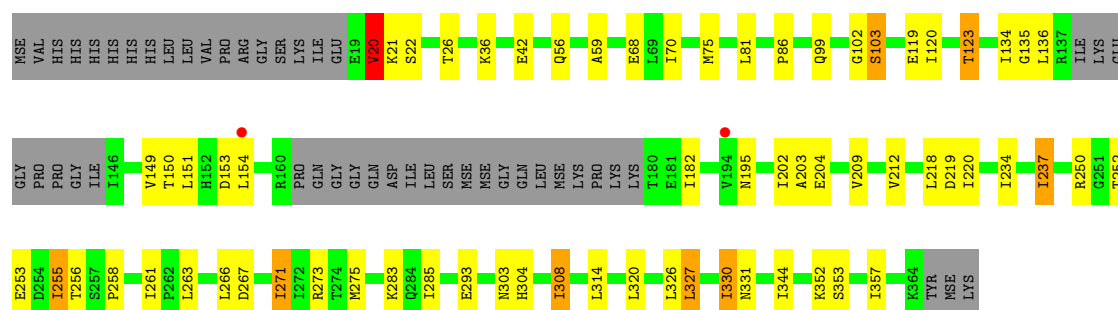
• Molecule 1: RUVB-LIKE 1

Chain A: 



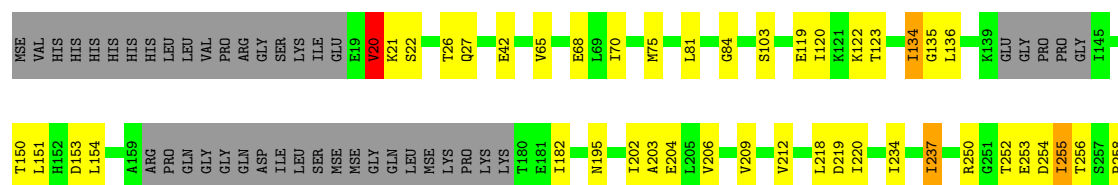
• Molecule 1: RUVB-LIKE 1

Chain B: 



• Molecule 1: RUVB-LIKE 1

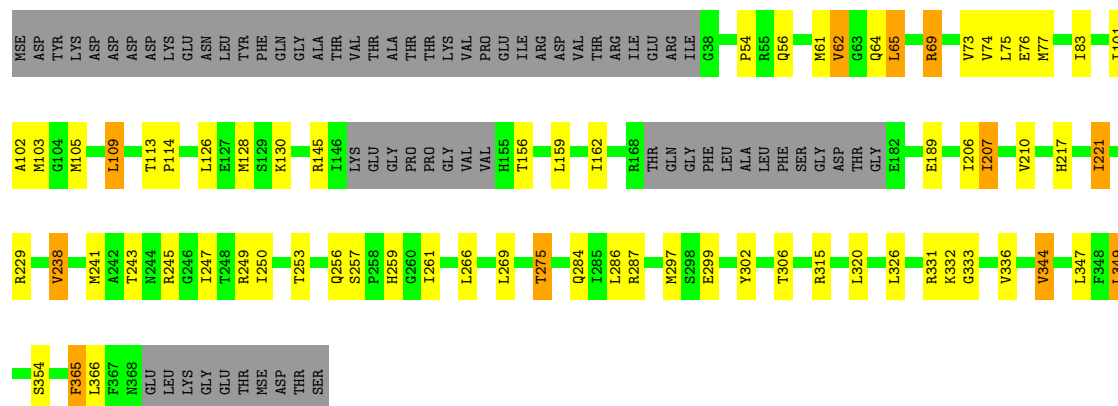
Chain C: 





• Molecule 2: RUVB-LIKE 2

Chain D: 64% 15% 18%



R353	F365	L366	F367	N368	GLU	LEU	LYS	GLY	GLU	THR	MSE	ASP	THR	SER
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4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	111.82Å 187.93Å 244.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.14 – 3.00 46.14 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (46.14-3.00) 99.5 (46.14-3.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 3.01Å)	Xtriage
Refinement program	BUSTER 2.11.0	Depositor
R, R_{free}	0.178 , 0.205 0.196 , 0.221	Depositor DCC
R_{free} test set	2632 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	85.2	Xtriage
Anisotropy	0.619	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 85.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.022 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.034 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14664	wwPDB-VP
Average B, all atoms (Å ²)	111.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% 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	0.80	0/2450	1.39	5/3294 (0.2%)
1	B	0.81	1/2455 (0.0%)	1.38	5/3301 (0.2%)
1	C	0.82	1/2445 (0.0%)	1.39	3/3289 (0.1%)
2	D	0.80	0/2395	1.42	9/3214 (0.3%)
2	E	0.79	0/2445	1.40	5/3280 (0.2%)
2	F	0.83	0/2403	1.43	7/3225 (0.2%)
All	All	0.81	2/14593 (0.0%)	1.40	34/19603 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	134	ILE	C-O	-7.56	1.15	1.24
1	B	275	MSE	SE-CE	5.19	2.11	1.95

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	255	ILE	N-CA-C	8.48	119.48	108.12
1	C	255	ILE	N-CA-C	8.33	119.29	108.12
1	B	255	ILE	N-CA-C	8.09	118.96	108.12
2	F	65	LEU	N-CA-C	8.00	119.69	110.97
2	D	365	PHE	CA-CB-CG	6.87	120.67	113.80
2	F	365	PHE	CA-CB-CG	6.68	120.48	113.80
2	E	365	PHE	CA-CB-CG	6.51	120.31	113.80
2	E	128	MSE	N-CA-C	6.17	117.78	108.46
1	A	103	SER	CA-C-N	6.03	133.06	121.54
1	A	103	SER	C-N-CA	6.03	133.06	121.54
2	D	275	THR	CB-CA-C	5.98	117.48	108.86
2	D	128	MSE	N-CA-C	5.84	117.29	108.46
2	E	275	THR	CB-CA-C	5.82	117.84	109.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	275	THR	CB-CA-C	5.71	117.35	108.61
1	C	153	ASP	CA-CB-CG	5.68	118.28	112.60
1	B	308	ILE	CA-C-N	5.65	126.25	119.98
1	B	308	ILE	C-N-CA	5.65	126.25	119.98
1	B	153	ASP	CA-CB-CG	5.57	118.17	112.60
2	E	65	LEU	N-CA-C	5.56	117.02	111.07
1	A	153	ASP	CA-CB-CG	5.53	118.13	112.60
2	F	128	MSE	N-CA-C	5.50	117.47	108.55
2	D	65	LEU	N-CA-C	5.49	117.70	111.11
2	D	64	GLN	CA-C-N	5.28	127.67	120.54
2	D	64	GLN	C-N-CA	5.28	127.67	120.54
1	B	271	ILE	N-CA-C	5.25	116.26	108.23
2	F	221	ILE	N-CA-CB	5.25	117.68	110.54
1	A	303	ASN	CA-CB-CG	5.21	117.81	112.60
2	D	221	ILE	N-CA-CB	5.20	117.61	110.54
2	D	103	MSE	CA-C-N	5.13	125.67	119.98
2	D	103	MSE	C-N-CA	5.13	125.67	119.98
2	E	365	PHE	O-C-N	5.07	129.46	122.46
2	F	184	LYS	CA-C-N	5.03	127.02	120.28
2	F	184	LYS	C-N-CA	5.03	127.02	120.28
1	C	254	ASP	CA-CB-CG	5.03	117.63	112.60

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	2430	0	2514	30	0
1	B	2435	0	2509	31	0
1	C	2425	0	2482	29	0
2	D	2377	0	2392	32	0
2	E	2426	0	2449	40	0
2	F	2385	0	2403	36	0
3	A	31	0	12	0	0
3	B	31	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	31	0	12	0	0
3	D	31	0	12	0	0
3	E	31	0	12	0	0
3	F	31	0	12	0	0
All	All	14664	0	14821	177	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 (177) 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:68:GLU:HB2	2:D:326:LEU:HD21	1.53	0.88
1:B:68:GLU:HB2	2:E:326:LEU:HD21	1.55	0.85
2:E:243:THR:HG21	2:E:259:HIS:HB3	1.60	0.83
2:D:243:THR:HG21	2:D:259:HIS:HB3	1.62	0.82
2:F:243:THR:HG21	2:F:259:HIS:HB3	1.61	0.81
1:A:283:LYS:HD3	1:A:303:ASN:HA	1.66	0.76
1:C:68:GLU:HB2	2:F:326:LEU:HD21	1.71	0.73
2:F:350:ASP:OD1	2:F:353:ARG:HB2	1.91	0.70
2:E:320:LEU:HD22	2:E:344:VAL:HG13	1.76	0.67
1:B:273:ARG:HB2	2:E:349:LEU:HD11	1.77	0.65
2:E:287:ARG:NH2	2:E:299:GLU:OE1	2.30	0.65
2:E:302:TYR:O	2:E:306:THR:HG23	1.99	0.62
1:A:252:THR:HB	1:A:255:ILE:HD12	1.82	0.62
2:E:210:VAL:HG13	2:E:240:ILE:HD12	1.81	0.62
1:C:252:THR:HB	1:C:255:ILE:HD12	1.82	0.61
2:F:302:TYR:O	2:F:306:THR:HG23	2.01	0.61
1:B:252:THR:HB	1:B:255:ILE:HD12	1.81	0.61
1:C:70:ILE:HG13	1:C:75:MSE:HG3	1.83	0.61
2:D:302:TYR:O	2:D:306:THR:HG23	2.01	0.60
1:B:36:LYS:HA	1:B:99:GLN:HE21	1.66	0.60
1:C:135:GLY:HA2	1:C:150:THR:HA	1.84	0.60
2:F:286:LEU:HD12	2:F:306:THR:HG22	1.83	0.60
2:D:286:LEU:HD12	2:D:306:THR:HG22	1.84	0.60
2:D:145:ARG:HG2	2:D:156:THR:HG22	1.85	0.59
1:A:135:GLY:HA2	1:A:150:THR:HA	1.85	0.59
1:B:135:GLY:HA2	1:B:150:THR:HA	1.84	0.59
2:E:365:PHE:O	2:E:368:ASN:HB2	2.02	0.59
2:E:145:ARG:HG2	2:E:156:THR:HG22	1.85	0.59
2:F:145:ARG:HG2	2:F:156:THR:HG22	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:102:ALA:HA	2:D:105:MSE:HE3	1.86	0.57
2:E:73:VAL:HG12	2:E:77:MSE:HE2	1.87	0.57
2:F:210:VAL:HG13	2:F:240:ILE:HD12	1.85	0.57
1:B:70:ILE:HG13	1:B:75:MSE:HG3	1.86	0.56
2:D:74:VAL:HA	2:D:77:MSE:HE3	1.86	0.56
2:D:73:VAL:HG12	2:D:77:MSE:HE2	1.86	0.56
2:E:102:ALA:HA	2:E:105:MSE:HE3	1.87	0.56
2:F:102:ALA:HA	2:F:105:MSE:HE3	1.87	0.56
2:E:259:HIS:HB2	2:E:261:ILE:HD12	1.88	0.56
2:D:56:GLN:HA	2:D:65:LEU:HD21	1.87	0.56
1:A:81:LEU:HD23	1:A:271:ILE:HG12	1.88	0.56
1:A:279:PRO:O	1:A:283:LYS:HG3	2.06	0.55
2:F:56:GLN:HA	2:F:65:LEU:HD21	1.89	0.55
2:E:56:GLN:HA	2:E:65:LEU:HD21	1.89	0.54
2:D:101:ILE:HG22	2:D:105:MSE:HE2	1.88	0.54
2:D:217:HIS:CG	2:D:245:ARG:HD2	2.43	0.54
2:E:286:LEU:HD12	2:E:306:THR:HG22	1.90	0.54
2:E:101:ILE:HG22	2:E:105:MSE:HE2	1.89	0.54
2:F:217:HIS:CG	2:F:245:ARG:HD2	2.42	0.54
2:F:349:LEU:CD2	2:F:353:ARG:NH1	2.72	0.53
2:E:217:HIS:CG	2:E:245:ARG:HD2	2.43	0.53
2:F:259:HIS:HB2	2:F:261:ILE:HD12	1.89	0.53
2:D:259:HIS:HB2	2:D:261:ILE:HD12	1.91	0.53
2:F:101:ILE:HG22	2:F:105:MSE:HE2	1.91	0.53
1:C:81:LEU:HD23	1:C:271:ILE:HG12	1.91	0.52
1:A:355:ALA:HB2	2:F:258:PRO:HB3	1.91	0.52
2:D:266:LEU:HD23	2:D:269:LEU:HD12	1.92	0.52
2:E:89:LEU:HB2	2:E:269:LEU:HD13	1.91	0.52
1:B:327:LEU:HD12	1:B:344:ILE:HG12	1.90	0.52
2:E:213:ILE:HD12	2:E:239:LEU:HD11	1.91	0.52
1:A:327:LEU:HD12	1:A:344:ILE:HG12	1.92	0.52
2:F:349:LEU:HD22	2:F:353:ARG:NH1	2.25	0.52
1:A:70:ILE:HG13	1:A:75:MSE:HG3	1.91	0.51
1:A:258:PRO:HB2	1:A:266:LEU:HD21	1.93	0.51
1:B:134:ILE:HB	1:B:151:LEU:HD12	1.92	0.51
2:E:74:VAL:HG22	2:E:270:LEU:HD13	1.93	0.51
2:F:89:LEU:HD11	2:F:243:THR:HG22	1.92	0.51
1:B:209:VAL:HG22	1:B:237:ILE:HG23	1.92	0.51
1:B:81:LEU:HD23	1:B:271:ILE:HG12	1.93	0.50
1:A:209:VAL:HG22	1:A:237:ILE:HG23	1.92	0.50
2:E:250:ILE:HD11	2:E:257:SER:HB2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:297:MSE:HG3	2:E:336:VAL:HB	1.94	0.50
2:D:297:MSE:HG3	2:D:336:VAL:HB	1.93	0.50
1:C:258:PRO:HB2	1:C:266:LEU:HD21	1.94	0.50
1:C:327:LEU:HD12	1:C:344:ILE:HG12	1.92	0.50
2:D:54:PRO:HD3	2:D:69:ARG:HB2	1.94	0.49
1:B:283:LYS:HD3	1:B:303:ASN:HA	1.94	0.49
1:A:134:ILE:HB	1:A:151:LEU:HD12	1.94	0.49
1:C:209:VAL:HG22	1:C:237:ILE:HG23	1.94	0.49
2:E:159:LEU:HA	2:E:162:ILE:HD12	1.95	0.49
2:F:249:ARG:HA	2:F:256:GLN:HA	1.95	0.49
1:A:219:ASP:HA	1:A:250:ARG:HB3	1.95	0.49
1:B:20:VAL:HG12	1:B:21:LYS:H	1.77	0.49
2:F:159:LEU:HA	2:F:162:ILE:HD12	1.95	0.49
1:C:134:ILE:HB	1:C:151:LEU:HD12	1.95	0.48
1:C:219:ASP:HA	1:C:250:ARG:HB3	1.94	0.48
2:E:212:PHE:HA	2:E:240:ILE:O	2.13	0.48
2:F:297:MSE:HG3	2:F:336:VAL:HB	1.93	0.48
1:B:330:ILE:HD11	2:D:76:GLU:HG3	1.95	0.48
2:D:159:LEU:HA	2:D:162:ILE:HD12	1.95	0.48
2:E:109:LEU:HB3	2:E:113:THR:HG21	1.95	0.48
1:B:258:PRO:HB2	1:B:266:LEU:HD21	1.94	0.48
2:D:241:MSE:HE1	2:D:269:LEU:HD21	1.96	0.48
2:E:61:MSE:HE2	2:E:68:ARG:HD2	1.94	0.48
1:B:219:ASP:HA	1:B:250:ARG:HB3	1.95	0.48
1:B:326:LEU:HD23	2:D:76:GLU:HB3	1.96	0.48
1:A:285:ILE:HG21	1:A:314:LEU:HD21	1.96	0.47
2:F:350:ASP:OD1	2:F:350:ASP:N	2.30	0.47
1:A:252:THR:O	1:A:253:GLU:HG2	2.14	0.47
2:F:250:ILE:HD11	2:F:257:SER:HB2	1.94	0.47
2:F:210:VAL:CG1	2:F:240:ILE:HD12	2.45	0.47
2:F:54:PRO:HD3	2:F:69:ARG:HB2	1.96	0.47
1:C:267:ASP:HA	2:F:315:ARG:HD2	1.95	0.47
1:B:267:ASP:HA	2:E:315:ARG:HD2	1.96	0.47
1:C:65:VAL:HG13	2:F:326:LEU:HD23	1.96	0.47
2:E:249:ARG:HA	2:E:256:GLN:HA	1.96	0.47
1:B:353:SER:O	1:B:357:ILE:HG12	2.14	0.47
2:D:250:ILE:HD11	2:D:257:SER:HB2	1.95	0.47
2:D:349:LEU:HD12	2:D:354:SER:HA	1.97	0.47
1:A:320:LEU:HD22	1:A:344:ILE:HG22	1.97	0.46
1:A:353:SER:O	1:A:357:ILE:HG12	2.16	0.46
1:C:84:GLY:HA3	1:C:274:THR:OG1	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:109:LEU:HB3	2:D:113:THR:HG21	1.96	0.46
1:A:123:THR:HG23	2:D:126:LEU:HB3	1.98	0.46
2:D:249:ARG:HA	2:D:256:GLN:HA	1.96	0.46
1:B:252:THR:O	1:B:253:GLU:HG2	2.15	0.46
2:E:216:VAL:HG21	2:E:241:MSE:SE	2.66	0.46
1:C:353:SER:O	1:C:357:ILE:HG12	2.16	0.45
1:C:320:LEU:HD22	1:C:344:ILE:HG22	1.98	0.45
1:B:320:LEU:HD22	1:B:344:ILE:HG22	1.98	0.45
1:C:212:VAL:HG11	1:C:218:LEU:HD11	1.99	0.45
1:C:326:LEU:HD23	2:E:76:GLU:HB3	1.99	0.45
1:A:330:ILE:HG22	1:A:331:ASN:HD22	1.82	0.45
1:C:252:THR:O	1:C:253:GLU:HG2	2.17	0.45
1:C:330:ILE:HG22	1:C:331:ASN:HD22	1.82	0.45
1:C:283:LYS:HD3	1:C:303:ASN:HA	1.98	0.45
2:E:366:LEU:O	2:E:369:GLU:HB2	2.16	0.45
2:E:54:PRO:HD3	2:E:69:ARG:HB2	2.00	0.44
1:C:20:VAL:HG12	1:C:21:LYS:H	1.83	0.44
1:B:330:ILE:HG22	1:B:331:ASN:HD22	1.83	0.44
1:C:122:LYS:HD2	2:F:126:LEU:HD21	2.00	0.43
1:B:123:THR:HG23	2:E:126:LEU:HB3	2.01	0.43
1:A:287:LYS:HB2	1:A:302:LEU:HD21	1.99	0.43
1:A:330:ILE:HD11	2:F:76:GLU:HG3	1.99	0.43
2:F:109:LEU:HB3	2:F:113:THR:HG21	1.99	0.43
1:A:151:LEU:HA	1:A:154:LEU:HD12	1.99	0.43
1:B:136:LEU:HD22	1:B:203:ALA:HB1	2.01	0.43
1:B:261:ILE:HB	1:B:266:LEU:HD13	2.01	0.43
1:C:151:LEU:HA	1:C:154:LEU:HD12	2.01	0.43
2:F:320:LEU:HD22	2:F:344:VAL:HG13	2.00	0.43
1:A:261:ILE:HB	1:A:266:LEU:HD13	2.01	0.43
2:F:114:PRO:HD3	2:F:207:ILE:HD12	2.01	0.43
2:D:114:PRO:HD3	2:D:207:ILE:HD12	2.01	0.43
1:B:151:LEU:HA	1:B:154:LEU:HD12	2.00	0.42
1:B:212:VAL:HG11	1:B:218:LEU:HD11	2.01	0.42
1:B:304:HIS:O	1:B:308:ILE:HG12	2.19	0.42
2:E:114:PRO:HD3	2:E:207:ILE:HD12	2.02	0.42
1:A:65:VAL:HG13	2:D:326:LEU:HD23	2.00	0.42
1:A:136:LEU:HD22	1:A:203:ALA:HB1	2.01	0.42
2:F:349:LEU:HD22	2:F:353:ARG:CZ	2.49	0.42
1:B:285:ILE:HG21	1:B:314:LEU:HD21	2.00	0.42
2:E:210:VAL:HG22	2:E:238:VAL:HG13	2.00	0.42
1:A:267:ASP:HA	2:D:315:ARG:HD2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:62:VAL:HG21	2:E:284:GLN:HB3	2.02	0.42
1:B:56:GLN:HG3	1:B:59:ALA:HB3	2.02	0.42
1:C:304:HIS:O	1:C:308:ILE:HG12	2.20	0.42
1:B:86:PRO:HD2	2:E:361:TYR:OH	2.20	0.42
2:E:75:LEU:HD21	2:E:108:ALA:HB3	2.02	0.42
1:A:212:VAL:HG11	1:A:218:LEU:HD11	2.01	0.41
2:D:284:GLN:HE22	2:D:287:ARG:HD2	1.85	0.41
1:C:285:ILE:HG21	1:C:314:LEU:HD21	2.02	0.41
1:C:136:LEU:HD22	1:C:203:ALA:HB1	2.02	0.41
1:C:355:ALA:HB2	2:E:258:PRO:HB3	2.01	0.41
2:D:320:LEU:HD22	2:D:344:VAL:HG13	2.02	0.41
2:D:62:VAL:HG21	2:D:284:GLN:HB3	2.03	0.41
1:A:326:LEU:HD23	2:F:76:GLU:HB3	2.02	0.41
1:C:347:LEU:HB3	2:E:70:ALA:HB2	2.03	0.41
2:D:287:ARG:NH2	2:D:299:GLU:OE1	2.54	0.41
2:F:62:VAL:HG21	2:F:284:GLN:HB3	2.02	0.41
2:F:210:VAL:HG22	2:F:238:VAL:HG13	2.02	0.41
1:A:280:GLN:HE21	1:A:284:GLN:HG3	1.86	0.41
1:C:261:ILE:HB	1:C:266:LEU:HD13	2.02	0.41
1:B:20:VAL:HG13	1:B:149:VAL:HG13	2.03	0.40
1:A:304:HIS:O	1:A:308:ILE:HG12	2.21	0.40
2:E:284:GLN:HE22	2:E:287:ARG:HD2	1.87	0.40
2:F:297:MSE:HE2	2:F:302:TYR:HA	2.03	0.40
2:F:40:HIS:HB2	2:F:103:MSE:HE3	2.04	0.40
2:D:210:VAL:HG22	2:D:238:VAL:HG13	2.01	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	312/367 (85%)	296 (95%)	12 (4%)	4 (1%)	9 38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	313/367 (85%)	296 (95%)	12 (4%)	5 (2%)	7	34
1	C	314/367 (86%)	299 (95%)	12 (4%)	3 (1%)	12	45
2	D	304/378 (80%)	295 (97%)	7 (2%)	2 (1%)	18	53
2	E	309/378 (82%)	300 (97%)	7 (2%)	2 (1%)	21	56
2	F	305/378 (81%)	297 (97%)	6 (2%)	2 (1%)	18	53
All	All	1857/2235 (83%)	1783 (96%)	56 (3%)	18 (1%)	12	45

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	20	VAL
1	A	104	LYS
1	B	20	VAL
1	C	20	VAL
2	D	332	LYS
2	E	332	LYS
2	F	332	LYS
1	B	103	SER
2	F	333	GLY
1	A	22	SER
1	B	22	SER
1	B	102	GLY
1	C	22	SER
2	D	333	GLY
2	E	333	GLY
1	A	182	ILE
1	C	182	ILE
1	B	182	ILE

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	267/299 (89%)	248 (93%)	19 (7%)	13	43

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	267/299 (89%)	248 (93%)	19 (7%)	13	43
1	C	263/299 (88%)	243 (92%)	20 (8%)	12	41
2	D	250/305 (82%)	228 (91%)	22 (9%)	9	35
2	E	257/305 (84%)	230 (90%)	27 (10%)	6	27
2	F	251/305 (82%)	228 (91%)	23 (9%)	8	33
All	All	1555/1812 (86%)	1425 (92%)	130 (8%)	10	37

All (130) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	VAL
1	A	26	THR
1	A	42	GLU
1	A	103	SER
1	A	119	GLU
1	A	120	ILE
1	A	123	THR
1	A	195	ASN
1	A	202	ILE
1	A	204	GLU
1	A	220	ILE
1	A	237	ILE
1	A	243	ASN
1	A	256	THR
1	A	263	LEU
1	A	269	VAL
1	A	293	GLU
1	A	327	LEU
1	A	330	ILE
1	B	20	VAL
1	B	26	THR
1	B	42	GLU
1	B	103	SER
1	B	119	GLU
1	B	120	ILE
1	B	123	THR
1	B	195	ASN
1	B	202	ILE
1	B	204	GLU
1	B	220	ILE

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Mol	Chain	Res	Type
1	B	234	ILE
1	B	237	ILE
1	B	256	THR
1	B	263	LEU
1	B	293	GLU
1	B	327	LEU
1	B	330	ILE
1	B	352	LYS
1	C	20	VAL
1	C	26	THR
1	C	27	GLN
1	C	42	GLU
1	C	103	SER
1	C	119	GLU
1	C	120	ILE
1	C	123	THR
1	C	195	ASN
1	C	202	ILE
1	C	204	GLU
1	C	206	VAL
1	C	220	ILE
1	C	234	ILE
1	C	237	ILE
1	C	256	THR
1	C	263	LEU
1	C	293	GLU
1	C	327	LEU
1	C	330	ILE
2	D	61	MSE
2	D	62	VAL
2	D	69	ARG
2	D	75	LEU
2	D	83	ILE
2	D	109	LEU
2	D	130	LYS
2	D	189	GLU
2	D	206	ILE
2	D	207	ILE
2	D	221	ILE
2	D	229	ARG
2	D	238	VAL
2	D	247	ILE

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Mol	Chain	Res	Type
2	D	253	THR
2	D	275	THR
2	D	331	ARG
2	D	344	VAL
2	D	347	LEU
2	D	349	LEU
2	D	365	PHE
2	D	366	LEU
2	E	48	LEU
2	E	58	SER
2	E	61	MSE
2	E	62	VAL
2	E	83	ILE
2	E	86	ARG
2	E	109	LEU
2	E	121	SER
2	E	130	LYS
2	E	189	GLU
2	E	206	ILE
2	E	207	ILE
2	E	221	ILE
2	E	229	ARG
2	E	238	VAL
2	E	251	ARG
2	E	253	THR
2	E	272	VAL
2	E	275	THR
2	E	331	ARG
2	E	344	VAL
2	E	347	LEU
2	E	349	LEU
2	E	352	SER
2	E	355	THR
2	E	365	PHE
2	E	366	LEU
2	F	48	LEU
2	F	62	VAL
2	F	69	ARG
2	F	75	LEU
2	F	83	ILE
2	F	109	LEU
2	F	121	SER

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Mol	Chain	Res	Type
2	F	130	LYS
2	F	189	GLU
2	F	206	ILE
2	F	207	ILE
2	F	221	ILE
2	F	229	ARG
2	F	235	MSE
2	F	238	VAL
2	F	253	THR
2	F	273	SER
2	F	275	THR
2	F	331	ARG
2	F	344	VAL
2	F	347	LEU
2	F	365	PHE
2	F	366	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	48	GLN
1	A	280	GLN
1	A	284	GLN
1	A	325	ASN
1	A	331	ASN
1	B	27	GLN
1	B	48	GLN
1	B	99	GLN
1	B	147	GLN
1	B	216	HIS
1	B	284	GLN
1	B	303	ASN
1	B	325	ASN
1	B	331	ASN
1	C	48	GLN
1	C	58	ASN
1	C	147	GLN
1	C	243	ASN
1	C	284	GLN
1	C	331	ASN
2	D	56	GLN

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Mol	Chain	Res	Type
2	D	284	GLN
2	D	337	GLN
2	E	56	GLN
2	E	284	GLN
2	E	337	GLN
2	F	56	GLN
2	F	284	GLN
2	F	337	GLN

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](#)

6 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	$\# Z > 2$
3	ATP	A	501	-	32,33,33	1.29	5 (15%)	48,52,52	0.93	2 (4%)
3	ATP	E	501	-	32,33,33	1.46	5 (15%)	48,52,52	0.87	2 (4%)
3	ATP	B	501	-	32,33,33	1.39	7 (21%)	48,52,52	0.92	2 (4%)
3	ATP	C	501	-	32,33,33	1.25	5 (15%)	48,52,52	0.95	2 (4%)
3	ATP	F	501	-	32,33,33	1.39	6 (18%)	48,52,52	0.92	2 (4%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ATP	D	501	-	32,33,33	1.28	5 (15%)	48,52,52	0.84	1 (2%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ATP	A	501	-	-	3/22/38/38	0/3/3/3
3	ATP	E	501	-	-	4/22/38/38	0/3/3/3
3	ATP	B	501	-	-	3/22/38/38	0/3/3/3
3	ATP	C	501	-	-	3/22/38/38	0/3/3/3
3	ATP	F	501	-	-	4/22/38/38	0/3/3/3
3	ATP	D	501	-	-	4/22/38/38	0/3/3/3

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	501	ATP	PB-O3A	-4.64	1.54	1.59
3	D	501	ATP	PA-O3A	-3.60	1.55	1.59
3	F	501	ATP	PA-O3A	-3.37	1.55	1.59
3	B	501	ATP	PB-O3A	-3.34	1.55	1.59
3	C	501	ATP	PB-O3A	-3.27	1.56	1.59
3	A	501	ATP	PB-O3A	-3.11	1.56	1.59
3	F	501	ATP	PB-O3A	-2.85	1.56	1.59
3	A	501	ATP	PG-O1G	2.80	1.59	1.50
3	C	501	ATP	PG-O1G	2.75	1.59	1.50
3	A	501	ATP	O4'-C1'	2.75	1.48	1.42
3	B	501	ATP	PG-O1G	2.70	1.58	1.50
3	D	501	ATP	PB-O3A	-2.67	1.56	1.59
3	F	501	ATP	C2-N1	2.57	1.38	1.33
3	F	501	ATP	PG-O1G	2.53	1.58	1.50
3	D	501	ATP	C2-N1	2.50	1.38	1.33
3	E	501	ATP	PG-O1G	2.47	1.58	1.50
3	B	501	ATP	C2-N3	2.45	1.38	1.33
3	A	501	ATP	C2-N3	2.40	1.38	1.33
3	E	501	ATP	C2-N1	2.37	1.38	1.33
3	B	501	ATP	C2-N1	2.37	1.38	1.33
3	F	501	ATP	C4-N3	2.35	1.38	1.34
3	E	501	ATP	C5-C4	2.32	1.43	1.39
3	C	501	ATP	O4'-C1'	2.29	1.47	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	501	ATP	PB-O3B	2.22	1.61	1.59
3	D	501	ATP	PG-O1G	2.22	1.57	1.50
3	E	501	ATP	C2-N3	2.18	1.37	1.33
3	F	501	ATP	PB-O3B	-2.16	1.57	1.59
3	A	501	ATP	PA-O1A	2.16	1.58	1.50
3	B	501	ATP	O4'-C1'	2.15	1.47	1.42
3	C	501	ATP	C5-N7	-2.13	1.35	1.39
3	B	501	ATP	PA-O1A	2.09	1.58	1.50
3	D	501	ATP	O4'-C1'	2.06	1.46	1.42
3	C	501	ATP	PG-O2G	2.03	1.62	1.54

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	501	ATP	O2G-PG-O3B	3.33	115.81	104.64
3	A	501	ATP	O2G-PG-O3B	3.10	115.02	104.64
3	B	501	ATP	O2G-PG-O3B	3.01	114.73	104.64
3	C	501	ATP	O3A-PB-O1B	-2.56	103.00	110.70
3	B	501	ATP	O3A-PB-O1B	-2.55	103.05	110.70
3	D	501	ATP	O2G-PG-O3B	2.54	113.16	104.64
3	F	501	ATP	O2G-PG-O3B	2.46	112.87	104.64
3	F	501	ATP	C5-N7-C8	2.42	107.26	103.45
3	A	501	ATP	O3A-PB-O1B	-2.33	103.70	110.70
3	E	501	ATP	O2G-PG-O3B	2.17	111.91	104.64
3	E	501	ATP	C5-N7-C8	2.09	106.74	103.45

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	ATP	C5'-O5'-PA-O1A
3	B	501	ATP	C5'-O5'-PA-O1A
3	B	501	ATP	C5'-O5'-PA-O3A
3	C	501	ATP	C5'-O5'-PA-O1A
3	D	501	ATP	PB-O3B-PG-O3G
3	E	501	ATP	PB-O3B-PG-O3G
3	F	501	ATP	PB-O3B-PG-O3G
3	D	501	ATP	PA-O3A-PB-O1B
3	E	501	ATP	PA-O3A-PB-O1B
3	F	501	ATP	PA-O3A-PB-O1B
3	A	501	ATP	C5'-O5'-PA-O2A
3	A	501	ATP	C5'-O5'-PA-O3A

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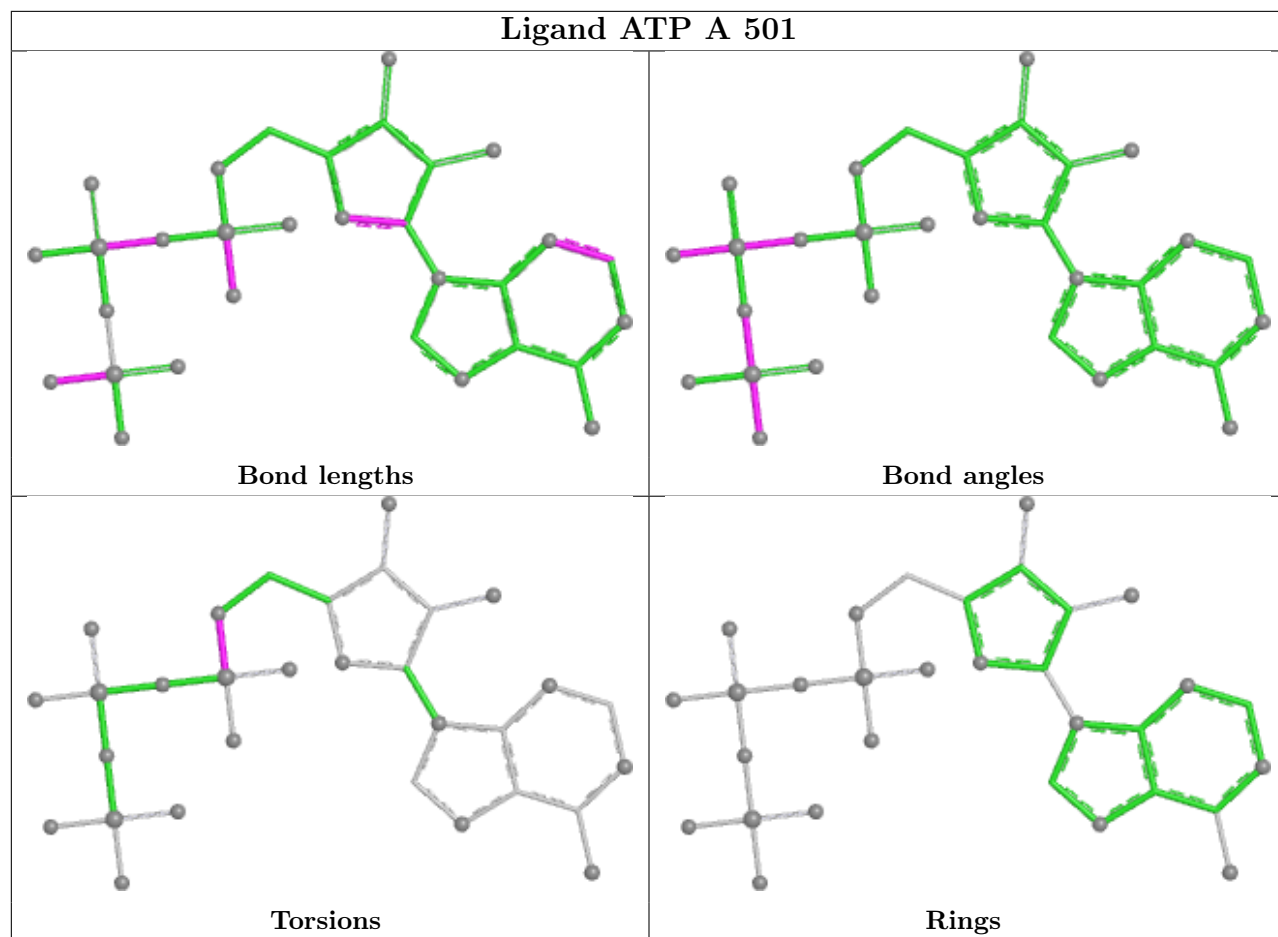
Mol	Chain	Res	Type	Atoms
3	B	501	ATP	C5'-O5'-PA-O2A
3	C	501	ATP	C5'-O5'-PA-O2A
3	C	501	ATP	C5'-O5'-PA-O3A
3	D	501	ATP	PA-O3A-PB-O2B
3	E	501	ATP	PA-O3A-PB-O2B
3	D	501	ATP	PB-O3B-PG-O1G
3	E	501	ATP	PB-O3B-PG-O1G
3	F	501	ATP	PB-O3B-PG-O1G
3	F	501	ATP	PA-O3A-PB-O2B

There are no ring outliers.

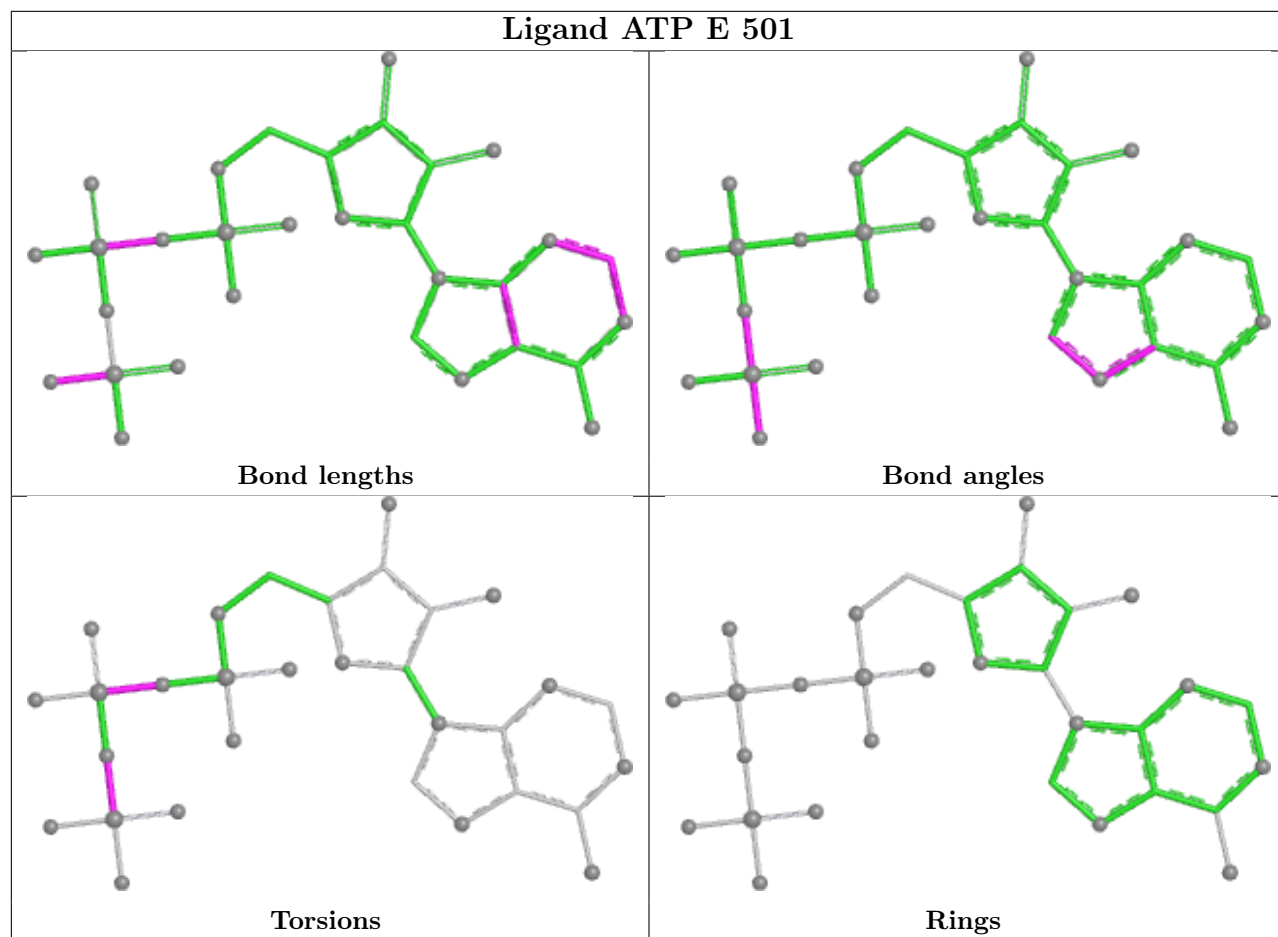
No monomer is involved in short contacts.

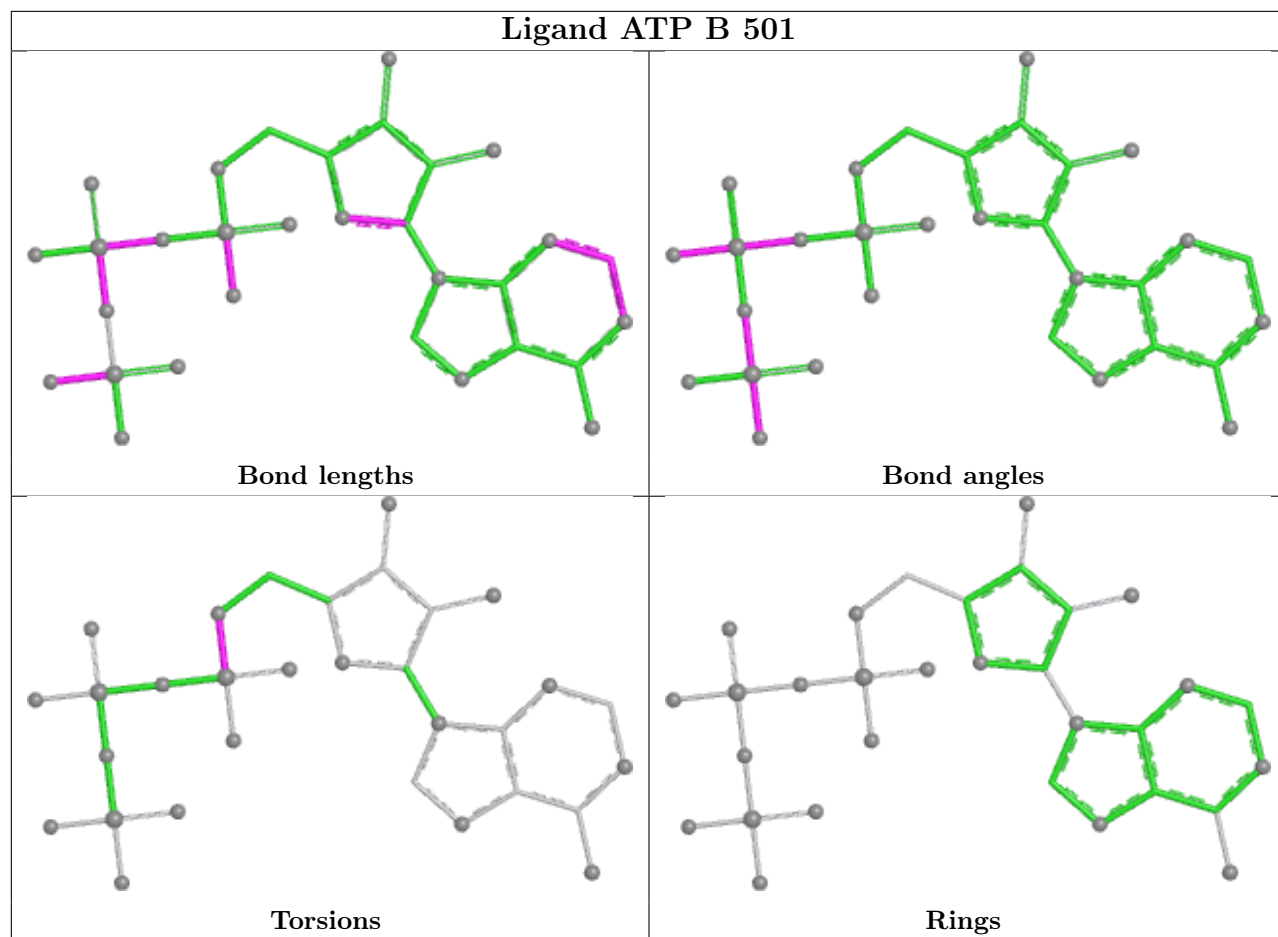
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

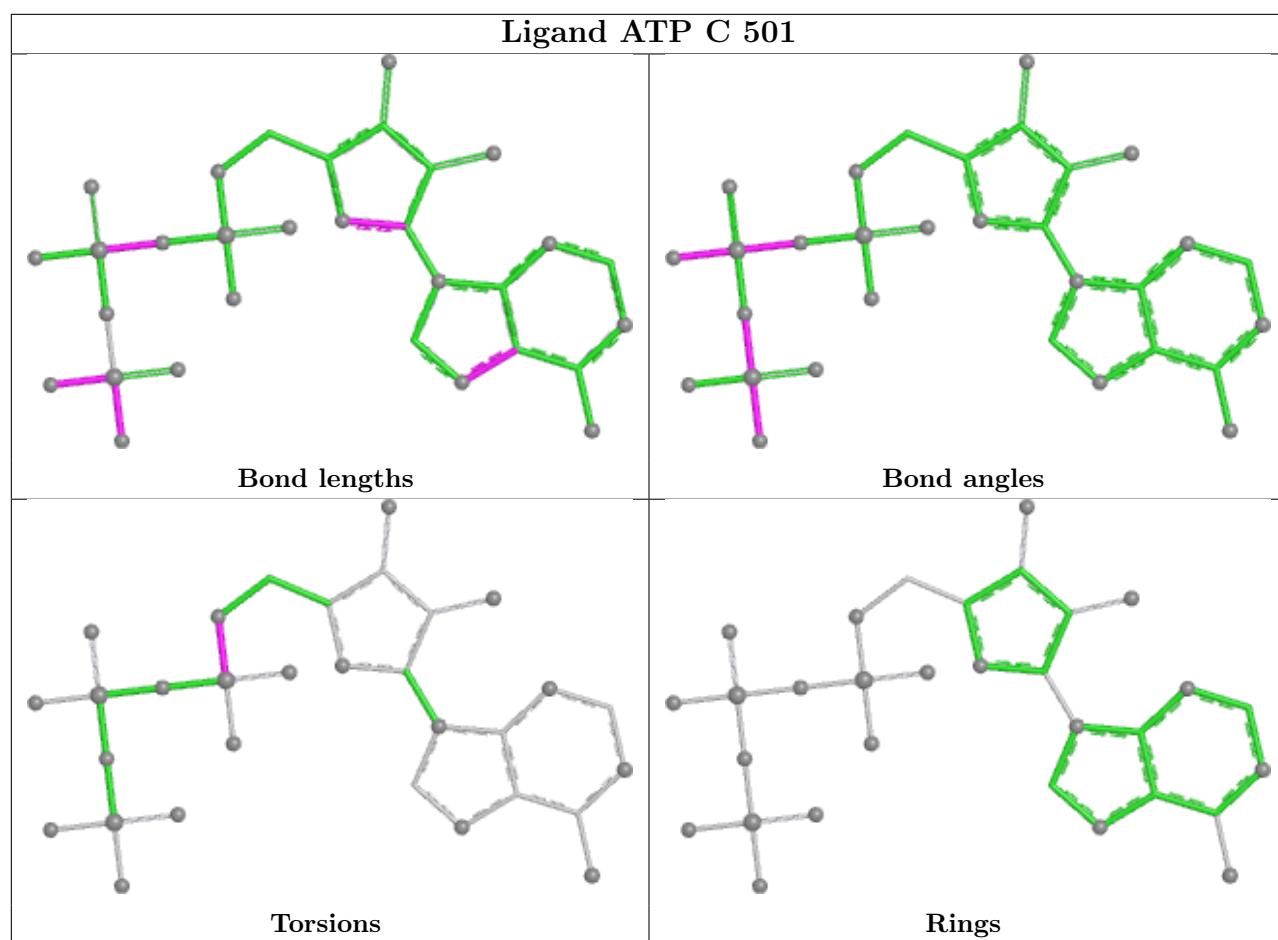
Ligand ATP A 501



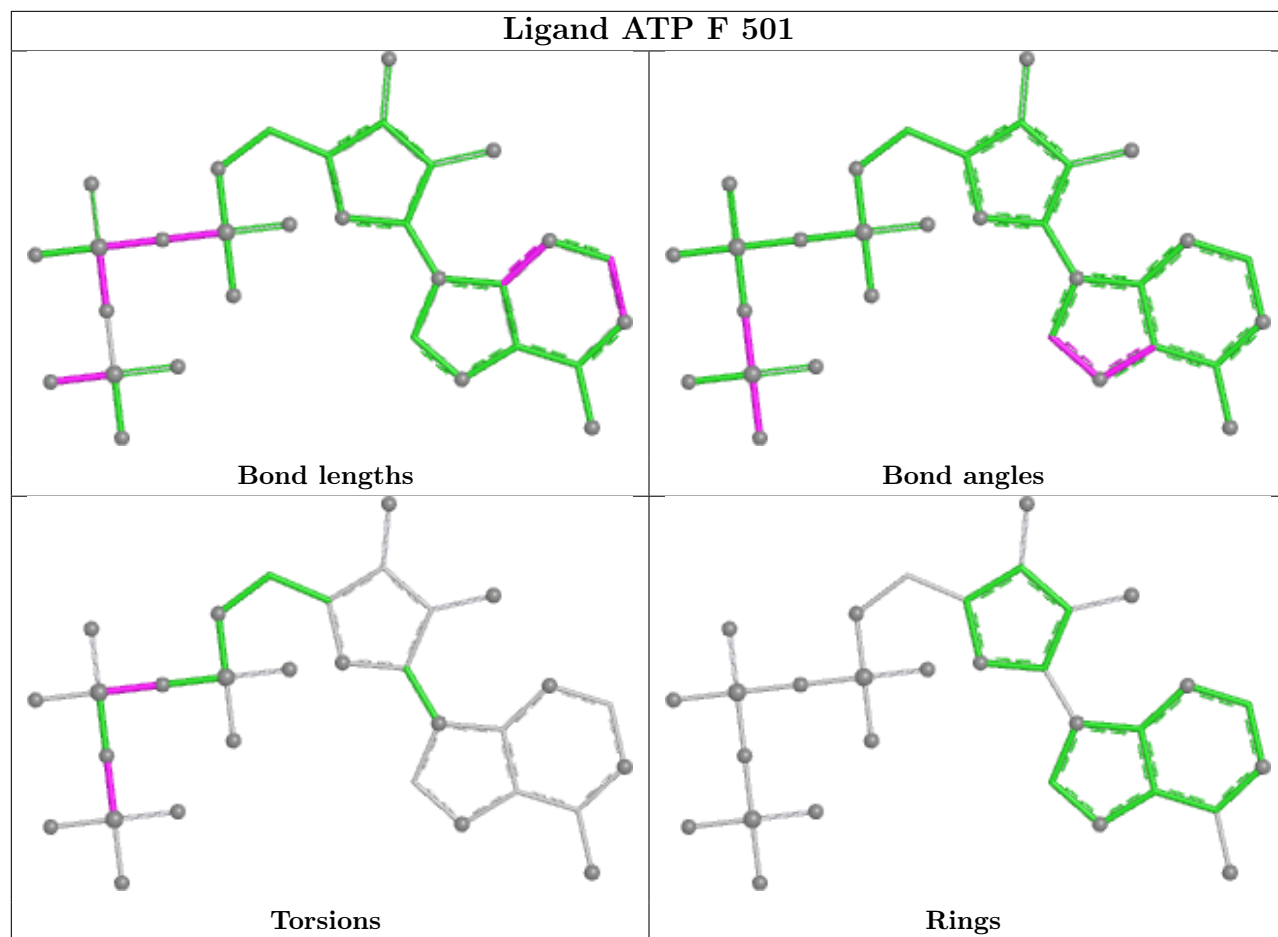
Ligand ATP E 501

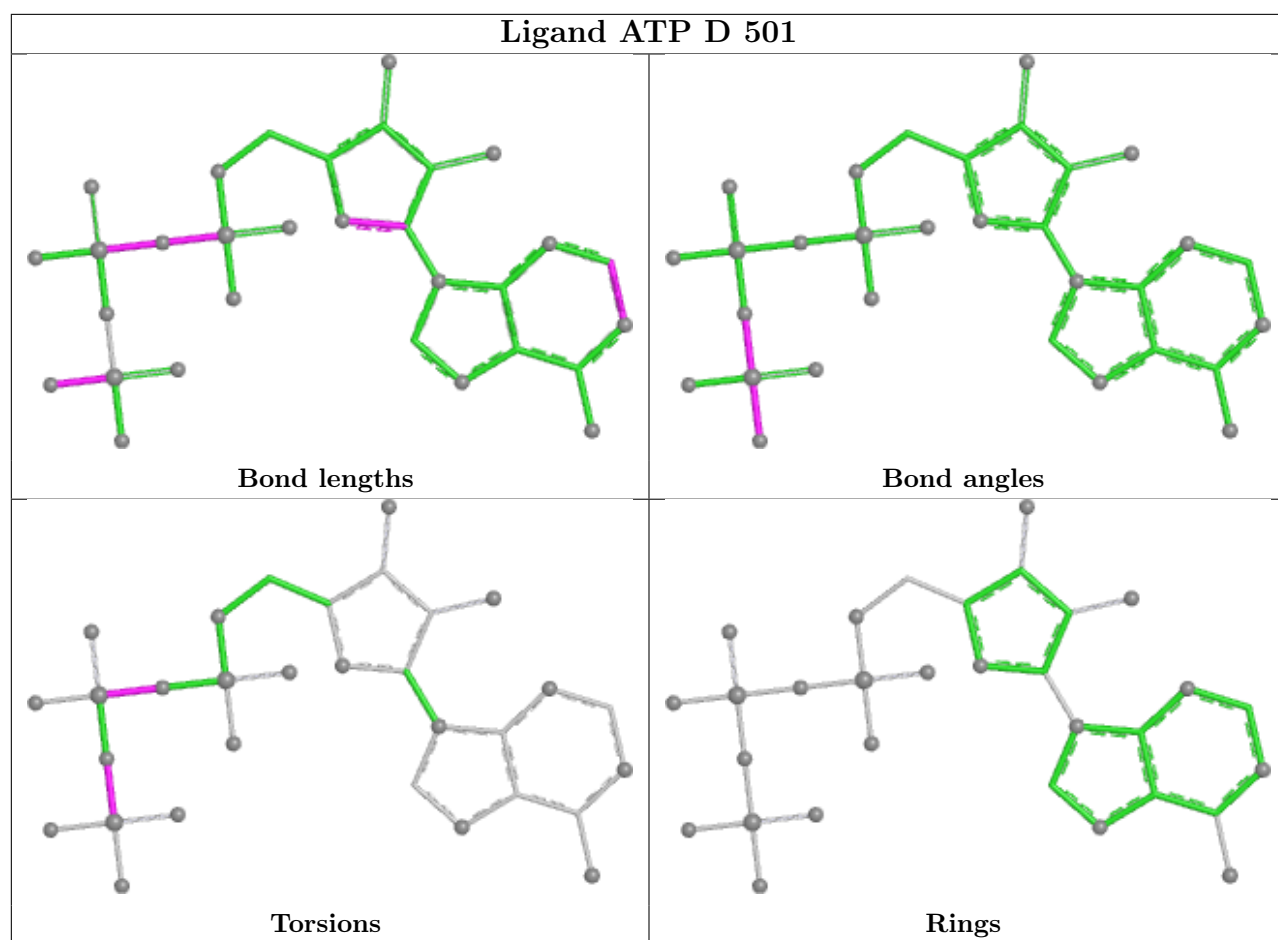






Ligand ATP F 501





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	311/367 (84%)	-0.50	1 (0%) 90 79	69, 97, 169, 214	1 (0%)
1	B	312/367 (85%)	-0.44	2 (0%) 85 69	74, 113, 167, 211	1 (0%)
1	C	313/367 (85%)	-0.43	0 100 100	71, 104, 169, 205	1 (0%)
2	D	300/378 (79%)	-0.48	0 100 100	74, 106, 168, 202	0
2	E	305/378 (80%)	-0.48	1 (0%) 90 79	79, 108, 167, 196	0
2	F	301/378 (79%)	-0.43	0 100 100	66, 97, 160, 188	0
All	All	1842/2235 (82%)	-0.46	4 (0%) 91 83	66, 105, 168, 214	3 (0%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	194	VAL	2.6
1	A	194	VAL	2.4
2	E	146	ILE	2.4
1	B	154	LEU	2.4

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,

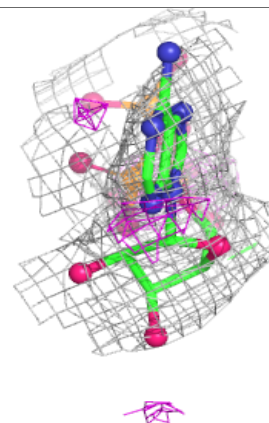
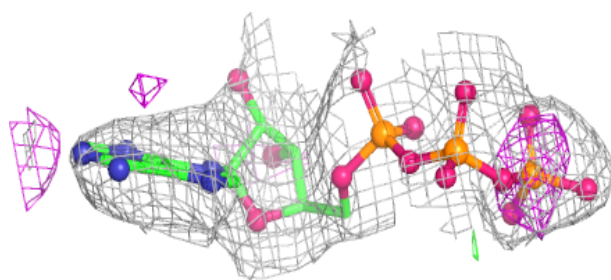
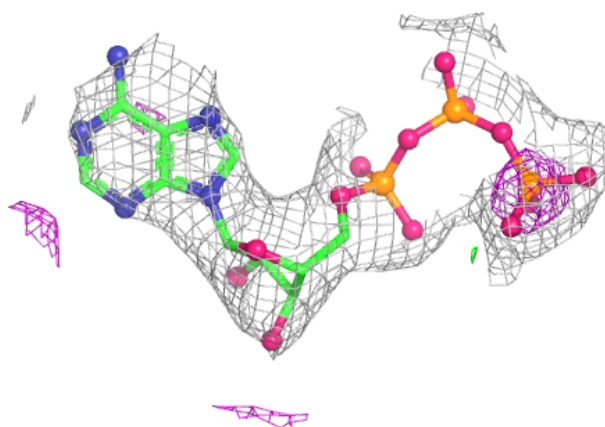
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	ATP	B	501	31/31	0.95	0.09	93,97,133,134	0
3	ATP	C	501	31/31	0.95	0.09	83,88,134,134	0
3	ATP	A	501	31/31	0.97	0.07	77,81,123,124	0
3	ATP	D	501	31/31	0.97	0.06	90,96,120,121	0
3	ATP	E	501	31/31	0.97	0.06	87,92,124,126	0
3	ATP	F	501	31/31	0.97	0.06	77,84,117,119	0

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

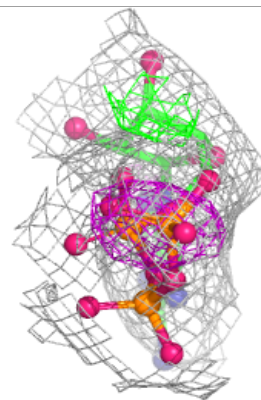
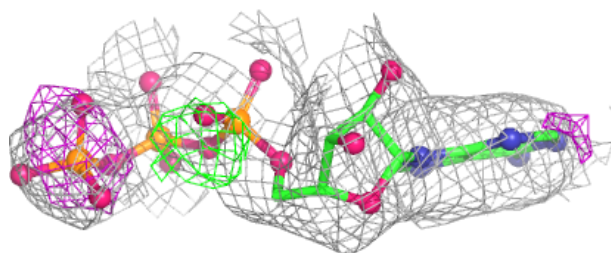
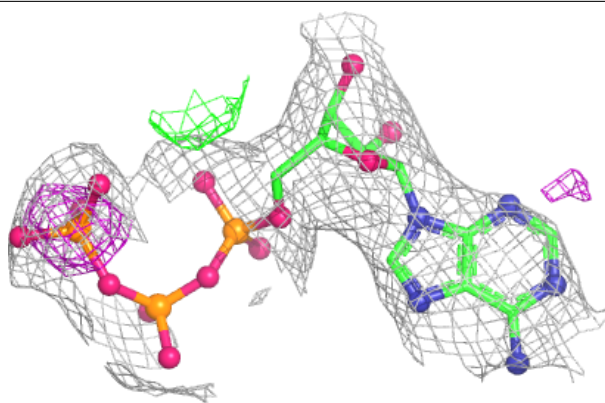
Electron density around ATP B 501:

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

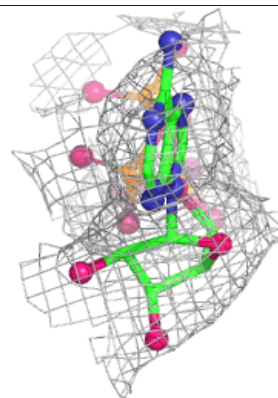
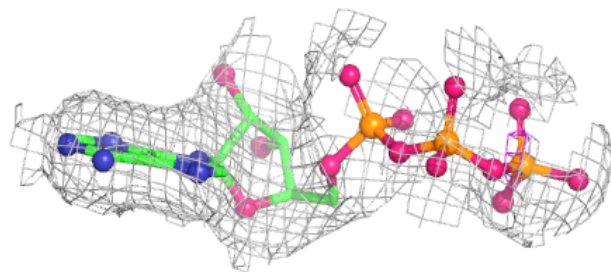
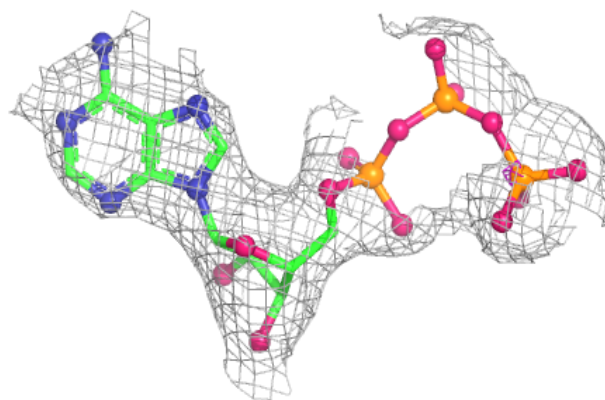


Electron density around ATP C 501:

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

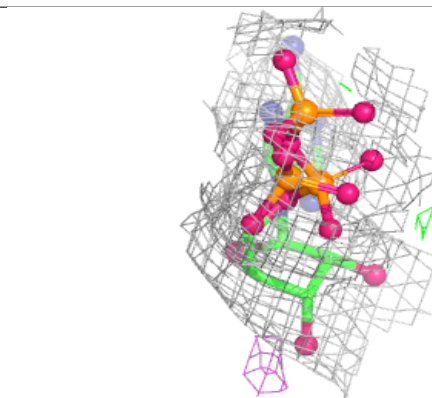
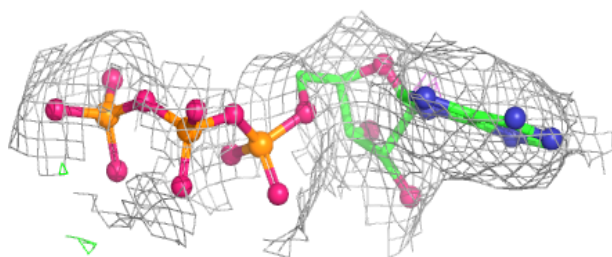
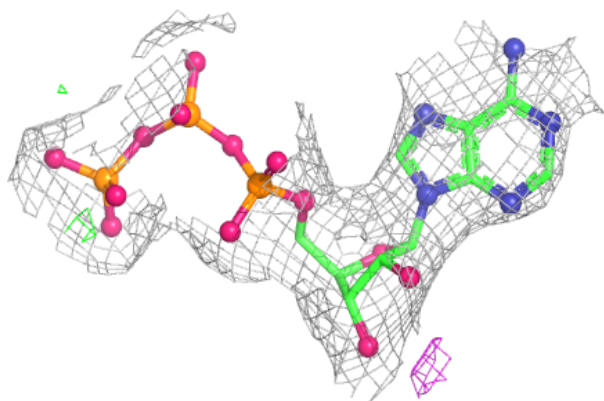
**Electron density around ATP A 501:**

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

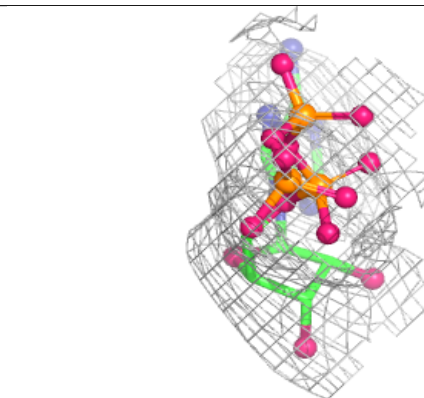
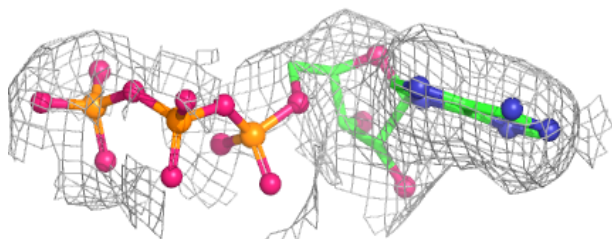
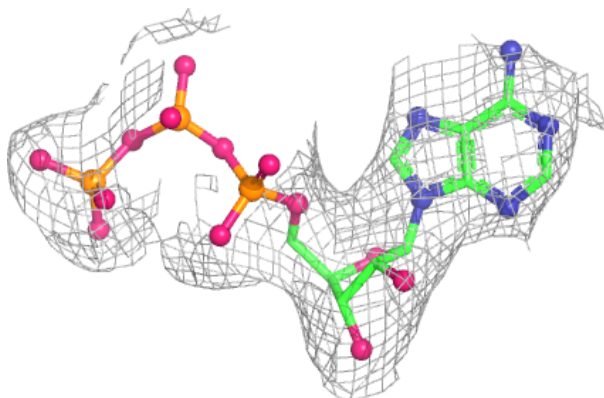


Electron density around ATP D 501:

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

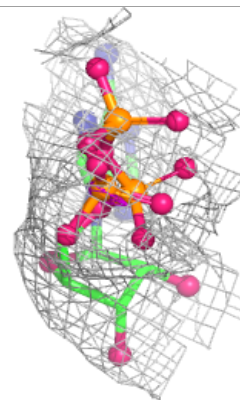
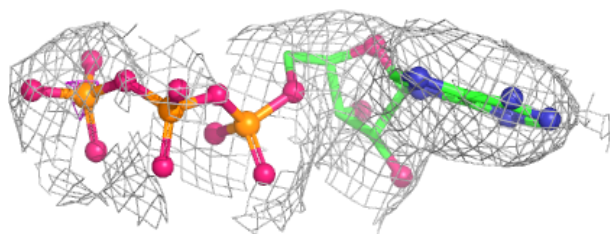
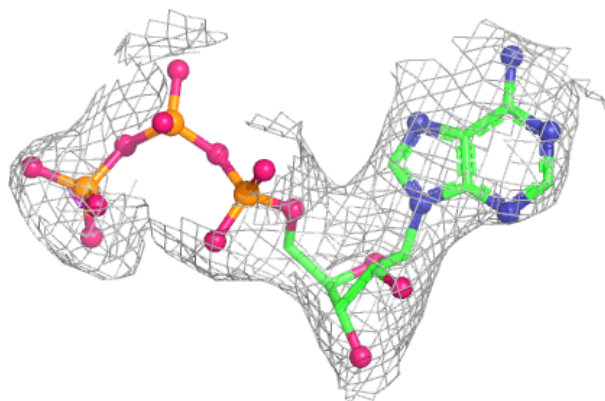
**Electron density around ATP E 501:**

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



Electron density around ATP F 501:

$2mF_o-DF_c$ (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.