



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:09 AM GMT

PDB ID : 3O8R
Title : Visualizing ATP-dependent RNA Translocation by the NS3 Helicase from HCV
Authors : Appleby, T.C.; Somoza, J.R.
Deposited on : 2010-08-03
Resolution : 2.30 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

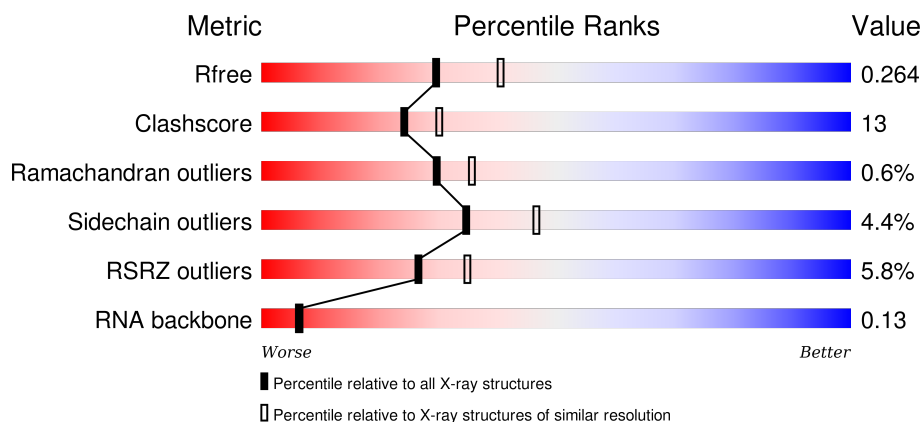
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)
RNA backbone	2183	1011 (2.84-1.76)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	666	<div> <div>3%</div> <div>79%</div> <div>17%</div> <div>..</div> </div>
1	B	666	<div> <div>8%</div> <div>68%</div> <div>25%</div> <div>..</div> </div>
2	C	6	<div> <div>17%</div> <div>17%</div> <div>33%</div> <div>50%</div> </div>

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 10085 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 HCV NS3 protease/helicase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	645	Total	C	N	O	S	0	0	0
			4807	3026	834	917	30			
1	B	645	Total	C	N	O	S	0	0	0
			4807	3026	834	917	30			

There are 74 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-36	MET	-	EXPRESSION TAG	UNP Q99AU2
A	-35	GLY	-	EXPRESSION TAG	UNP Q99AU2
A	-34	SER	-	EXPRESSION TAG	UNP Q99AU2
A	-33	SER	-	EXPRESSION TAG	UNP Q99AU2
A	-32	HIS	-	EXPRESSION TAG	UNP Q99AU2
A	-31	HIS	-	EXPRESSION TAG	UNP Q99AU2
A	-30	HIS	-	EXPRESSION TAG	UNP Q99AU2
A	-29	HIS	-	EXPRESSION TAG	UNP Q99AU2
A	-28	HIS	-	EXPRESSION TAG	UNP Q99AU2
A	-27	HIS	-	EXPRESSION TAG	UNP Q99AU2
A	-26	SER	-	EXPRESSION TAG	UNP Q99AU2
A	-25	SER	-	EXPRESSION TAG	UNP Q99AU2
A	-24	GLY	-	EXPRESSION TAG	UNP Q99AU2
A	-23	LEU	-	EXPRESSION TAG	UNP Q99AU2
A	-22	VAL	-	EXPRESSION TAG	UNP Q99AU2
A	-21	PRO	-	EXPRESSION TAG	UNP Q99AU2
A	-20	ARG	-	EXPRESSION TAG	UNP Q99AU2
A	-19	GLY	-	EXPRESSION TAG	UNP Q99AU2
A	-18	SER	-	EXPRESSION TAG	UNP Q99AU2
A	-17	HIS	-	EXPRESSION TAG	UNP Q99AU2
A	-16	MET	-	EXPRESSION TAG	UNP Q99AU2
A	-15	GLY	-	EXPRESSION TAG	UNP Q99AU2
A	-14	SER	-	EXPRESSION TAG	UNP Q99AU2
A	-13	VAL	-	EXPRESSION TAG	UNP Q99AU2
A	-12	VAL	-	EXPRESSION TAG	UNP Q99AU2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	ILE	-	EXPRESSION TAG	UNP Q99AU2
A	-10	VAL	-	EXPRESSION TAG	UNP Q99AU2
A	-9	GLY	-	EXPRESSION TAG	UNP Q99AU2
A	-8	ARG	-	EXPRESSION TAG	UNP Q99AU2
A	-7	ILE	-	EXPRESSION TAG	UNP Q99AU2
A	-6	ILE	-	EXPRESSION TAG	UNP Q99AU2
A	-5	LEU	-	EXPRESSION TAG	UNP Q99AU2
A	-4	SER	-	EXPRESSION TAG	UNP Q99AU2
A	-3	GLY	-	EXPRESSION TAG	UNP Q99AU2
A	-2	SER	-	EXPRESSION TAG	UNP Q99AU2
A	-1	GLY	-	EXPRESSION TAG	UNP Q99AU2
A	0	SER	-	EXPRESSION TAG	UNP Q99AU2
B	-36	MET	-	EXPRESSION TAG	UNP Q99AU2
B	-35	GLY	-	EXPRESSION TAG	UNP Q99AU2
B	-34	SER	-	EXPRESSION TAG	UNP Q99AU2
B	-33	SER	-	EXPRESSION TAG	UNP Q99AU2
B	-32	HIS	-	EXPRESSION TAG	UNP Q99AU2
B	-31	HIS	-	EXPRESSION TAG	UNP Q99AU2
B	-30	HIS	-	EXPRESSION TAG	UNP Q99AU2
B	-29	HIS	-	EXPRESSION TAG	UNP Q99AU2
B	-28	HIS	-	EXPRESSION TAG	UNP Q99AU2
B	-27	HIS	-	EXPRESSION TAG	UNP Q99AU2
B	-26	SER	-	EXPRESSION TAG	UNP Q99AU2
B	-25	SER	-	EXPRESSION TAG	UNP Q99AU2
B	-24	GLY	-	EXPRESSION TAG	UNP Q99AU2
B	-23	LEU	-	EXPRESSION TAG	UNP Q99AU2
B	-22	VAL	-	EXPRESSION TAG	UNP Q99AU2
B	-21	PRO	-	EXPRESSION TAG	UNP Q99AU2
B	-20	ARG	-	EXPRESSION TAG	UNP Q99AU2
B	-19	GLY	-	EXPRESSION TAG	UNP Q99AU2
B	-18	SER	-	EXPRESSION TAG	UNP Q99AU2
B	-17	HIS	-	EXPRESSION TAG	UNP Q99AU2
B	-16	MET	-	EXPRESSION TAG	UNP Q99AU2
B	-15	GLY	-	EXPRESSION TAG	UNP Q99AU2
B	-14	SER	-	EXPRESSION TAG	UNP Q99AU2
B	-13	VAL	-	EXPRESSION TAG	UNP Q99AU2
B	-12	VAL	-	EXPRESSION TAG	UNP Q99AU2
B	-11	ILE	-	EXPRESSION TAG	UNP Q99AU2
B	-10	VAL	-	EXPRESSION TAG	UNP Q99AU2
B	-9	GLY	-	EXPRESSION TAG	UNP Q99AU2
B	-8	ARG	-	EXPRESSION TAG	UNP Q99AU2
B	-7	ILE	-	EXPRESSION TAG	UNP Q99AU2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	ILE	-	EXPRESSION TAG	UNP Q99AU2
B	-5	LEU	-	EXPRESSION TAG	UNP Q99AU2
B	-4	SER	-	EXPRESSION TAG	UNP Q99AU2
B	-3	GLY	-	EXPRESSION TAG	UNP Q99AU2
B	-2	SER	-	EXPRESSION TAG	UNP Q99AU2
B	-1	GLY	-	EXPRESSION TAG	UNP Q99AU2
B	0	SER	-	EXPRESSION TAG	UNP Q99AU2

- Molecule 2 is a RNA chain called RNA (5'-R(P*UP*(5BU)P*UP*UP*UP*U)-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	C	6	Total	Br	C	N	O	P	0	0	0
			122	1	54	12	49	6			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

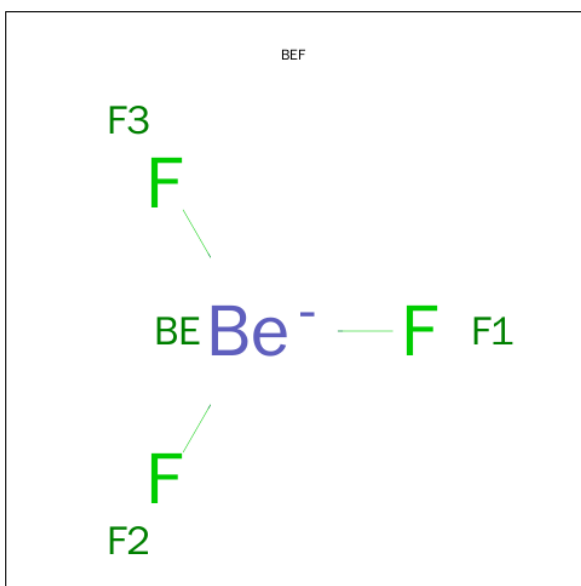
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		

- Molecule 5 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 6 is BERYLLIUM TRIFLUORIDE ION (three-letter code: BEF) (formula: BeF₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	Be	F	0	0
			4	1	3		

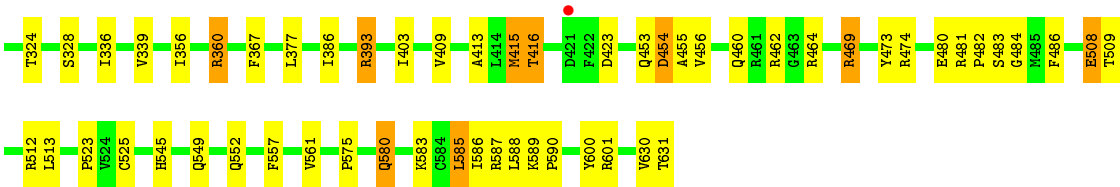
- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	181	Total	O	0	0
			181	181		
8	C	3	Total	O	0	0
			3	3		
8	B	121	Total	O	0	0
			121	121		



● Molecule 2: RNA (5'-R(P*UP*(5BU)P*UP*UP*UP*U)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	91.03 Å 110.83 Å 141.98 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.68 – 2.30 43.68 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.3 (43.68-2.30) 95.6 (43.68-2.30)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.79 (at 2.32 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.215 , 0.263 0.213 , 0.264	Depositor DCC
R_{free} test set	2000 reflections (3.37%)	DCC
Wilson B-factor (Å ²)	37.4	Xtriage
Anisotropy	0.106	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 41.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 64039 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10085	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 5BU, MG, ADP, BEF, ZN, SO4

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	0.45	0/4916	0.59	1/6714 (0.0%)
1	B	0.40	0/4916	0.58	0/6714
2	C	1.19	1/110 (0.9%)	0.99	0/166
All	All	0.44	1/9942 (0.0%)	0.59	1/13594 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	3	U	OP3-P	-10.24	1.48	1.61

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	316	ARG	NE-CZ-NH2	-5.08	117.76	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	4807	0	4783	84	0
1	B	4807	0	4784	167	0
2	C	122	0	60	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
5	A	27	0	12	0	0
6	A	4	0	0	0	0
7	B	10	0	0	0	0
8	A	181	0	0	3	0
8	B	121	0	0	5	0
8	C	3	0	0	0	0
All	All	10085	0	9639	252	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 13.

All (252) 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:B:19:THR:HA	1:B:22:THR:HG23	1.43	0.98
1:B:213:LYS:HZ1	1:B:217:ALA:HB2	1.27	0.96
1:B:254:THR:HG22	1:B:256:VAL:H	1.28	0.96
1:B:14:LEU:HD12	1:B:15:GLY:H	1.27	0.96
1:A:24:ARG:HH11	1:A:24:ARG:HB3	1.35	0.89
1:A:118:ARG:HH12	1:A:121:ASP:HA	1.37	0.88
1:B:14:LEU:CD1	1:B:15:GLY:H	1.86	0.87
1:B:13:LEU:CD1	1:B:17:ILE:HD13	2.06	0.85
1:B:11:ARG:HH22	1:B:27:ASN:HB2	1.42	0.84
1:A:470:ARG:HH11	1:A:470:ARG:HG2	1.42	0.84
1:B:356:ILE:HD11	1:B:386:ILE:HD11	1.62	0.82
1:B:74:MET:HE3	1:B:86:GLN:H	1.45	0.81
1:B:231:SER:HB2	1:B:416:THR:HA	1.65	0.79
1:A:413:ALA:HB2	2:C:5:U:H5'	1.65	0.78
1:B:14:LEU:HD12	1:B:15:GLY:N	1.98	0.78
1:A:24:ARG:NH1	1:A:24:ARG:HB3	2.01	0.76
1:A:181:SER:HA	8:A:677:HOH:O	1.86	0.75
1:A:425:VAL:HG23	1:A:465:THR:HB	1.69	0.75
1:B:213:LYS:NZ	1:B:217:ALA:HB2	2.02	0.74
1:A:230:PRO:O	2:C:6:U:H5'	1.89	0.73
1:A:32:GLU:HG3	1:A:94:LEU:HD21	1.71	0.72
1:B:305:THR:OG1	1:B:512:ARG:HD3	1.90	0.72
1:B:356:ILE:CD1	1:B:386:ILE:HD11	2.20	0.72
1:B:11:ARG:NH2	1:B:27:ASN:HB2	2.04	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:LYS:HD3	1:B:631:THR:HG22	1.72	0.70
1:B:-8:ARG:HH11	1:B:8:GLN:NE2	1.90	0.70
1:B:74:MET:CE	1:B:86:GLN:H	2.05	0.69
1:B:73:GLN:OE1	1:B:76:THR:HG22	1.93	0.69
1:B:11:ARG:NH2	1:B:25:ASP:OD2	2.26	0.69
1:A:371:LYS:HG2	2:C:4:5BU:OP2	1.92	0.68
1:B:213:LYS:HZ1	1:B:217:ALA:CB	2.03	0.68
1:B:13:LEU:HD12	1:B:17:ILE:HD13	1.75	0.68
1:A:29:VAL:HG21	1:A:88:PRO:HB2	1.76	0.67
1:B:175:MET:O	1:B:179:MET:HG3	1.95	0.67
1:B:75:TYR:CE1	1:B:179:MET:HA	2.30	0.67
1:B:296:ASP:CG	1:B:299:THR:HG22	2.17	0.66
1:A:431:CYS:SG	8:A:741:HOH:O	2.54	0.66
1:B:232:VAL:HG12	8:B:693:HOH:O	1.96	0.65
1:B:23:GLY:HA3	1:B:70:PRO:HG3	1.80	0.64
1:B:13:LEU:HA	1:B:16:CYS:HB3	1.80	0.63
1:A:232:VAL:HG12	2:C:6:U:OP1	1.99	0.63
1:B:73:GLN:OE1	1:B:76:THR:CG2	2.48	0.62
1:B:19:THR:HA	1:B:22:THR:CG2	2.24	0.62
1:A:114:ILE:HD11	1:A:134:TYR:HE1	1.64	0.61
1:B:293:HIS:CE1	1:B:460:GLN:NE2	2.68	0.61
1:A:470:ARG:HH11	1:A:470:ARG:CG	2.11	0.61
1:B:263:ALA:HB1	1:B:264:PRO:HD2	1.83	0.61
1:B:77:ASN:ND2	1:B:80:GLN:HG3	2.16	0.61
1:A:258:THR:HG22	1:A:258:THR:O	2.00	0.60
1:B:12:GLY:C	1:B:14:LEU:H	2.03	0.60
1:B:226:LEU:HA	1:B:266:THR:HG23	1.84	0.59
1:B:136:LYS:NZ	1:B:630:VAL:O	2.34	0.59
1:A:371:LYS:HE3	2:C:4:5BU:OP1	2.01	0.59
1:A:23:GLY:HA3	1:A:70:PRO:HG3	1.85	0.59
1:B:213:LYS:HA	1:B:246:HIS:CE1	2.38	0.58
1:A:102:SER:HB2	1:A:117:ARG:NH1	2.19	0.58
1:A:360:ARG:HA	1:A:386:ILE:HD11	1.84	0.58
1:B:293:HIS:HE1	1:B:460:GLN:NE2	2.00	0.57
1:B:393:ARG:CZ	1:B:413:ALA:HB2	2.35	0.57
1:B:29:VAL:HG11	1:B:88:PRO:HG2	1.85	0.57
1:B:294:SER:OG	1:B:299:THR:HG21	2.04	0.57
1:A:445:THR:HG22	1:A:447:GLU:HG2	1.87	0.57
1:B:-15:GLY:O	1:B:62:LYS:HD3	2.04	0.57
1:B:232:VAL:CG2	1:B:254:THR:HG21	2.35	0.57
1:B:197:PHE:HA	1:B:316:ARG:O	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:74:MET:HE3	1:B:86:GLN:CB	2.35	0.57
1:A:392:TYR:HB2	2:C:5:U:OP1	2.04	0.57
1:B:588:LEU:C	1:B:590:PRO:HD2	2.24	0.57
1:B:229:ASN:OD1	1:B:230:PRO:HD2	2.05	0.57
1:B:-8:ARG:HD2	1:B:8:GLN:NE2	2.20	0.56
1:B:109:ARG:HD2	1:B:137:GLY:O	2.06	0.56
1:A:42:SER:OG	1:A:109:ARG:NH2	2.38	0.56
1:B:193:VAL:HG11	1:B:223:TYR:CE1	2.41	0.56
1:B:423:ASP:OD2	1:B:469:ARG:NH2	2.38	0.56
1:A:116:VAL:HG22	1:A:126:LEU:HD23	1.88	0.55
1:A:336:ILE:HD13	1:A:462:ARG:HG2	1.87	0.55
1:B:194:PRO:HG3	1:B:198:GLN:HB2	1.89	0.55
1:B:25:ASP:N	1:B:66:GLY:O	2.35	0.55
1:B:455:ALA:HB3	1:B:483:SER:HB3	1.88	0.55
1:A:505:THR:OG1	1:A:508:GLU:HG3	2.06	0.55
1:A:194:PRO:HG3	1:A:198:GLN:HB3	1.89	0.55
1:B:130:ARG:HD3	1:B:134:TYR:CD1	2.42	0.55
1:B:455:ALA:HA	1:B:480:GLU:OE2	2.07	0.55
1:B:480:GLU:HG2	1:B:481:ARG:H	1.71	0.54
1:B:226:LEU:HA	1:B:266:THR:CG2	2.38	0.54
1:B:225:VAL:HB	1:B:265:VAL:HG22	1.88	0.54
1:B:601:ARG:NH2	8:B:679:HOH:O	2.40	0.54
1:A:470:ARG:NH1	1:A:470:ARG:HG2	2.19	0.54
1:A:456:VAL:HG23	1:A:483:SER:HB2	1.90	0.54
1:B:423:ASP:HA	1:B:469:ARG:NH2	2.23	0.54
1:A:115:PRO:HB2	1:A:127:LEU:HD22	1.90	0.53
1:A:136:LYS:NZ	1:A:631:THR:HA	2.23	0.53
1:B:14:LEU:O	1:B:18:ILE:CD1	2.57	0.53
1:B:194:PRO:HD3	1:B:317:LEU:HD13	1.90	0.53
1:B:12:GLY:O	1:B:13:LEU:HB3	2.08	0.53
1:A:434:GLN:HA	1:A:447:GLU:O	2.09	0.52
1:A:97:CYS:HB2	1:A:149:HIS:HB2	1.92	0.52
1:A:583:LYS:HA	1:A:586:ILE:HG13	1.91	0.52
1:B:232:VAL:HG23	1:B:254:THR:HG21	1.91	0.52
1:B:12:GLY:H	1:B:14:LEU:HG	1.74	0.52
1:B:14:LEU:O	1:B:18:ILE:HD13	2.10	0.52
1:B:232:VAL:CG1	8:B:693:HOH:O	2.54	0.52
1:B:32:GLU:HG3	1:B:94:LEU:HD21	1.91	0.51
1:A:413:ALA:HA	2:C:5:U:H4'	1.91	0.51
1:B:145:CYS:HB2	1:B:146:PRO:HD2	1.91	0.51
1:B:393:ARG:NH2	1:B:413:ALA:HB2	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:43:PHE:HA	1:B:137:GLY:O	2.11	0.51
1:A:575:PRO:HG2	1:A:577:SER:O	2.10	0.51
1:A:229:ASN:O	1:A:269:THR:HA	2.11	0.51
1:B:240:ALA:O	1:B:244:LYS:HG3	2.11	0.51
1:B:228:LEU:HA	1:B:268:SER:O	2.11	0.51
1:B:248:ILE:HG12	1:B:249:ASP:N	2.26	0.51
1:A:118:ARG:NH1	1:A:121:ASP:HA	2.15	0.51
1:B:136:LYS:CD	1:B:631:THR:HG22	2.38	0.50
1:B:25:ASP:HB3	1:B:66:GLY:HA2	1.92	0.50
1:A:418:TYR:CE2	1:A:420:GLY:HA3	2.46	0.50
1:B:64:LEU:HB2	1:B:85:TRP:CD1	2.46	0.50
1:B:360:ARG:O	1:B:360:ARG:HG3	2.11	0.50
1:B:589:LYS:N	1:B:590:PRO:HD2	2.27	0.50
1:B:480:GLU:HG2	1:B:481:ARG:N	2.27	0.50
1:B:51:VAL:O	1:B:53:TRP:CD1	2.65	0.49
1:A:32:GLU:HG3	1:A:94:LEU:CD2	2.41	0.49
1:B:282:GLY:HA2	1:B:313:ALA:O	2.12	0.49
1:B:213:LYS:HZ3	1:B:213:LYS:C	2.11	0.49
1:A:371:LYS:HB2	1:A:392:TYR:CD2	2.47	0.49
1:B:583:LYS:O	1:B:586:ILE:HG13	2.12	0.49
1:B:183:VAL:HG13	1:B:184:PHE:H	1.77	0.49
1:B:557:PHE:O	1:B:561:VAL:HG23	2.11	0.49
1:B:296:ASP:OD1	1:B:299:THR:HG22	2.14	0.48
1:A:393:ARG:HH11	1:A:393:ARG:HG3	1.78	0.48
1:B:-6:ILE:O	1:B:5:ALA:HA	2.13	0.48
1:B:60:GLY:HA3	1:B:482:PRO:HG3	1.93	0.48
1:A:17:ILE:HG13	1:B:17:ILE:HG21	1.96	0.48
1:B:293:HIS:CE1	1:B:460:GLN:HE22	2.31	0.48
1:B:423:ASP:HA	1:B:469:ARG:HH22	1.79	0.48
1:B:145:CYS:HB2	1:B:146:PRO:CD	2.44	0.48
1:B:508:GLU:O	1:B:512:ARG:HG3	2.14	0.47
1:B:-11:ILE:O	1:B:11:ARG:NH1	2.46	0.47
1:A:188:SER:HA	1:A:203:HIS:O	2.15	0.47
1:B:254:THR:HG22	1:B:256:VAL:N	2.12	0.47
1:B:12:GLY:C	1:B:14:LEU:N	2.69	0.47
1:A:363:ARG:HG2	1:A:406:VAL:HG12	1.97	0.47
1:A:114:ILE:HD11	1:A:134:TYR:CE1	2.48	0.47
1:B:339:VAL:O	1:B:474:ARG:HA	2.14	0.47
1:B:244:LYS:O	1:B:246:HIS:N	2.47	0.46
1:B:248:ILE:CD1	1:B:265:VAL:HB	2.44	0.46
1:A:256:VAL:HG23	1:A:257:ARG:N	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:232:VAL:HG13	1:B:233:ALA:N	2.30	0.46
1:B:235:THR:HG23	1:B:267:TYR:HB3	1.97	0.46
1:B:266:THR:HG21	1:B:284:TYR:CZ	2.50	0.46
1:A:363:ARG:O	1:A:423:ASP:N	2.44	0.46
1:B:244:LYS:HB3	1:B:244:LYS:HE3	1.68	0.46
2:C:5:U:O2	2:C:5:U:H2'	2.16	0.46
1:B:254:THR:HG22	1:B:255:GLY:N	2.30	0.46
1:B:183:VAL:HG13	1:B:184:PHE:N	2.31	0.46
1:B:324:THR:O	1:B:324:THR:HG23	2.15	0.46
1:A:132:VAL:HG13	1:A:159:CYS:SG	2.56	0.46
1:A:619:ILE:O	1:A:622:CYS:HB2	2.15	0.46
1:B:266:THR:HG21	1:B:284:TYR:OH	2.15	0.46
1:A:413:ALA:CB	2:C:5:U:H5'	2.42	0.46
1:A:390:ALA:HA	1:A:409:VAL:O	2.16	0.46
1:B:-12:VAL:O	1:B:-10:VAL:HG13	2.16	0.46
1:B:197:PHE:HD1	1:B:315:ALA:O	1.99	0.45
1:B:286:ILE:HG23	1:B:317:LEU:HD23	1.98	0.45
1:A:132:VAL:HG22	1:A:162:GLY:O	2.16	0.45
1:A:175:MET:O	1:A:179:MET:HG3	2.15	0.45
1:A:118:ARG:NH1	1:A:120:GLY:O	2.49	0.45
1:B:115:PRO:HB2	1:B:127:LEU:HD22	1.98	0.45
1:A:519:THR:HA	1:A:520:PRO:HD3	1.78	0.45
1:A:347:ILE:HB	1:A:354:ILE:HB	1.98	0.45
1:B:43:PHE:N	1:B:43:PHE:CD1	2.85	0.45
1:A:470:ARG:NH1	1:A:470:ARG:CG	2.71	0.45
1:B:307:LEU:HD21	1:B:320:LEU:HD21	1.97	0.45
1:A:380:LYS:O	1:A:384:LEU:HD23	2.17	0.45
1:B:239:GLY:HA3	1:B:252:ILE:HD11	1.99	0.45
1:B:-4:SER:OG	1:B:4:THR:HG23	2.16	0.45
1:B:213:LYS:HD2	1:B:246:HIS:CE1	2.52	0.44
1:A:97:CYS:HA	1:A:149:HIS:CD2	2.53	0.44
1:A:13:LEU:O	1:A:17:ILE:HG12	2.16	0.44
1:B:213:LYS:HD2	1:B:246:HIS:NE2	2.32	0.44
1:B:74:MET:HE3	1:B:86:GLN:HB3	1.99	0.44
1:B:235:THR:O	1:B:238:PHE:HB2	2.17	0.44
1:B:580:GLN:HG2	1:B:580:GLN:H	1.56	0.44
1:A:414:LEU:HA	1:A:414:LEU:HD12	1.81	0.44
1:B:367:PHE:HZ	1:B:415:MET:HE1	1.82	0.44
1:B:217:ALA:O	1:B:220:ALA:HB3	2.18	0.44
2:C:3:U:H6	2:C:3:U:O5'	2.01	0.44
1:B:77:ASN:HD21	1:B:80:GLN:HG3	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:-6:ILE:O	1:A:5:ALA:HA	2.17	0.43
1:B:180:ARG:HA	1:B:180:ARG:HD3	1.63	0.43
1:A:617:LYS:HA	1:A:620:MET:HE2	1.99	0.43
1:B:336:ILE:HD13	1:B:462:ARG:HG2	1.98	0.43
1:B:44:LEU:O	1:B:140:GLY:HA3	2.18	0.43
1:B:190:PRO:HB3	1:B:214:VAL:HG13	1.99	0.43
1:B:213:LYS:HD2	1:B:213:LYS:HA	1.66	0.43
1:B:-8:ARG:HH11	1:B:8:GLN:HE21	1.66	0.43
1:B:194:PRO:HB2	1:B:196:SER:O	2.19	0.43
1:A:198:GLN:HG3	1:A:199:VAL:N	2.34	0.43
1:A:254:THR:C	1:A:272:LYS:HD3	2.39	0.43
1:A:102:SER:O	1:A:117:ARG:HD2	2.19	0.43
1:B:209:GLY:O	1:B:214:VAL:HG23	2.19	0.43
1:A:393:ARG:CG	2:C:5:U:H5"	2.49	0.42
1:A:393:ARG:HG2	2:C:5:U:OP2	2.18	0.42
1:B:232:VAL:HG21	1:B:254:THR:HG21	2.01	0.42
1:B:303:ILE:O	1:B:307:LEU:HG	2.19	0.42
1:A:431:CYS:O	1:A:450:THR:HA	2.20	0.42
1:B:229:ASN:O	1:B:269:THR:HA	2.19	0.42
1:B:462:ARG:HG3	1:B:473:TYR:CG	2.54	0.42
1:B:259:ILE:O	1:B:259:ILE:HG12	2.11	0.42
1:B:377:LEU:HD23	1:B:409:VAL:CG1	2.50	0.42
1:B:-14:SER:HB2	8:B:654:HOH:O	2.19	0.42
1:B:17:ILE:O	1:B:20:SER:HB3	2.20	0.42
1:A:377:LEU:HD22	1:A:381:LEU:HG	2.01	0.42
1:B:293:HIS:H	1:B:293:HIS:CD2	2.37	0.42
1:A:477:THR:CG2	1:A:478:PRO:HD2	2.50	0.42
1:A:393:ARG:HG2	2:C:5:U:H5"	2.02	0.42
1:B:453:GLN:O	1:B:481:ARG:HB2	2.19	0.42
1:B:464:ARG:HD2	8:B:692:HOH:O	2.20	0.41
1:A:22:THR:O	1:A:24:ARG:HG3	2.20	0.41
1:B:194:PRO:CB	1:B:198:GLN:HG2	2.50	0.41
1:A:538:GLY:HA3	1:A:618:TYR:CE2	2.55	0.41
1:B:484:GLY:O	1:B:523:PRO:HA	2.19	0.41
1:A:446:ILE:HG21	1:A:557:PHE:CE2	2.55	0.41
1:B:589:LYS:N	1:B:590:PRO:CD	2.84	0.41
1:B:291:GLU:HA	1:B:291:GLU:OE1	2.20	0.41
1:B:224:LYS:HD3	1:B:283:ALA:O	2.19	0.41
1:A:293:HIS:HA	8:A:768:HOH:O	2.20	0.41
1:B:171:PRO:HB2	1:B:173:GLU:HG2	2.02	0.41
1:B:213:LYS:NZ	1:B:213:LYS:O	2.27	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:LEU:HD13	1:B:17:ILE:HD13	1.95	0.41
1:B:203:HIS:N	1:B:203:HIS:CD2	2.87	0.41
1:A:104:LEU:HD22	1:A:151:VAL:HG21	2.02	0.41
1:A:337:GLU:HB2	1:A:470:ARG:HH12	1.86	0.41
1:B:213:LYS:HZ2	1:B:213:LYS:HG3	1.80	0.41
1:B:299:THR:HG23	1:B:300:ILE:N	2.36	0.41
1:B:228:LEU:HD22	1:B:273:PHE:HB2	2.02	0.41
1:B:214:VAL:HB	1:B:215:PRO:HD3	2.02	0.41
1:B:377:LEU:HD23	1:B:409:VAL:HG11	2.03	0.41
1:B:509:THR:O	1:B:513:LEU:HG	2.20	0.41
1:A:61:SER:HA	1:A:73:GLN:NE2	2.36	0.41
1:B:403:ILE:HA	1:B:403:ILE:HD12	1.79	0.41
1:B:575:PRO:HB3	1:B:585:LEU:HD23	2.03	0.41
1:B:486:PHE:CZ	1:B:525:CYS:HB2	2.56	0.41
1:A:13:LEU:HD11	1:A:17:ILE:HD11	2.03	0.41
1:B:14:LEU:C	1:B:16:CYS:H	2.25	0.41
1:B:454:ASP:OD2	1:B:456:VAL:HG23	2.20	0.41
1:B:67:PRO:HG2	1:B:68:LYS:H	1.86	0.40
1:A:13:LEU:O	1:A:16:CYS:HB3	2.21	0.40
1:A:224:LYS:NZ	1:A:224:LYS:HB2	2.36	0.40
1:A:24:ARG:HH11	1:A:24:ARG:CB	2.20	0.40
1:B:12:GLY:O	1:B:14:LEU:N	2.54	0.40
1:B:545:HIS:O	1:B:549:GLN:HG3	2.22	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	643/666 (96%)	624 (97%)	18 (3%)	1 (0%)	52 64
1	B	643/666 (96%)	603 (94%)	33 (5%)	7 (1%)	17 18

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1286/1332 (96%)	1227 (95%)	51 (4%)	8 (1%)	30	36

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	14	LEU
1	B	244	LYS
1	B	245	ALA
1	B	281	GLY
1	B	328	SER
1	A	281	GLY
1	B	67	PRO
1	B	282	GLY

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	526/544 (97%)	506 (96%)	20 (4%)	40	54
1	B	526/544 (97%)	500 (95%)	26 (5%)	31	41
All	All	1052/1088 (97%)	1006 (96%)	46 (4%)	35	46

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	SER
1	A	9	GLN
1	A	28	GLN
1	A	68	LYS
1	A	109	ARG
1	A	132	VAL
1	A	173	GLU
1	A	198	GLN
1	A	244	LYS
1	A	257	ARG

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Mol	Chain	Res	Type
1	A	377	LEU
1	A	414	LEU
1	A	450	THR
1	A	470	ARG
1	A	481	ARG
1	A	517	LEU
1	A	529	LEU
1	A	585	LEU
1	A	600	TYR
1	A	606	GLN
1	B	4	THR
1	B	22	THR
1	B	26	LYS
1	B	32	GLU
1	B	165	LYS
1	B	180	ARG
1	B	184	PHE
1	B	187	ASN
1	B	213	LYS
1	B	231	SER
1	B	248	ILE
1	B	256	VAL
1	B	259	ILE
1	B	298	THR
1	B	360	ARG
1	B	393	ARG
1	B	415	MET
1	B	416	THR
1	B	454	ASP
1	B	469	ARG
1	B	508	GLU
1	B	552	GLN
1	B	580	GLN
1	B	585	LEU
1	B	587	ARG
1	B	600	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	GLN
1	B	8	GLN

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Mol	Chain	Res	Type
1	B	195	GLN
1	B	293	HIS
1	B	460	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	5/6 (83%)	3 (60%)	0

All (3) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	5	U
2	C	6	U
2	C	8	U

There are no RNA pucker outliers to report.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is 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	5BU	C	4	2	13,22,23	3.82	4 (30%)	14,32,35	2.30	2 (14%)

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	5BU	C	4	2	-	0/3/25/26	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	4	5BU	O4-C4	-2.21	1.19	1.24
2	C	4	5BU	C4-N3	3.17	1.39	1.33
2	C	4	5BU	C4-C5	9.17	1.50	1.38
2	C	4	5BU	C6-N1	9.48	1.49	1.35

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	4	5BU	C5-C4-N3	-4.05	119.67	124.00
2	C	4	5BU	C4-N3-C2	7.37	121.62	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	4	5BU	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 3 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	ADP	A	875	4,6	22,29,29	1.02	1 (4%)	27,45,45	1.95	5 (18%)
6	BEF	A	885	5	0,3,3	0.00	-	0,3,3	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	SO4	B	890	-	4,4,4	0.26	0	6,6,6	0.21	0
7	SO4	B	895	-	4,4,4	0.10	0	6,6,6	0.12	0

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
5	ADP	A	875	4,6	-	0/12/32/32	0/3/3/3
6	BEF	A	885	5	-	0/0/0/0	0/0/0/0
7	SO4	B	890	-	-	0/0/0/0	0/0/0/0
7	SO4	B	895	-	-	0/0/0/0	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	875	ADP	C5-C4	3.18	1.47	1.40

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	875	ADP	N3-C2-N1	-7.49	123.16	128.89
5	A	875	ADP	PA-O3A-PB	-3.20	121.94	132.67
5	A	875	ADP	C4-C5-N7	-2.23	107.43	109.48
5	A	875	ADP	C2-N1-C6	2.25	122.78	118.77
5	A	875	ADP	O3B-PB-O1B	2.55	118.78	110.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	645/666 (96%)	0.18	18 (2%) 56 66	18, 37, 63, 76	0
1	B	645/666 (96%)	0.50	56 (8%) 13 18	22, 47, 77, 94	0
2	C	5/6 (83%)	1.08	1 (20%) 1 2	53, 63, 73, 82	0
All	All	1295/1338 (96%)	0.34	75 (5%) 26 35	18, 41, 73, 94	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	223	TYR	6.3
1	B	262	GLY	6.3
1	B	242	MET	5.6
1	B	195	GLN	5.5
1	B	292	CYS	5.1
1	B	184	PHE	5.0
1	B	287	ILE	4.9
1	B	314	GLY	4.9
1	B	218	TYR	4.4
1	B	24	ARG	4.3
1	B	196	SER	4.2
1	B	241	TYR	4.2
1	B	280	SER	4.1
1	B	316	ARG	4.0
1	A	245	ALA	3.8
1	B	225	VAL	3.8
1	B	12	GLY	3.8
1	B	21	LEU	3.7
1	A	359	ILE	3.7
1	A	182	PRO	3.6
1	A	385	GLY	3.6
1	B	279	CYS	3.5
1	B	282	GLY	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	312	THR	3.5
1	B	318	VAL	3.3
1	A	14	LEU	3.2
1	B	248	ILE	3.2
1	B	255	GLY	3.1
1	B	13	LEU	3.1
1	A	18	ILE	3.1
1	B	286	ILE	3.1
1	B	183	VAL	3.0
1	B	264	PRO	3.0
1	B	220	ALA	2.9
1	A	406	VAL	2.9
1	B	22	THR	2.9
1	B	-15	GLY	2.8
1	B	246	HIS	2.8
1	B	261	THR	2.8
1	A	381	LEU	2.8
1	B	85	TRP	2.8
1	B	197	PHE	2.8
1	B	222	GLY	2.8
1	A	383	GLY	2.7
1	A	388	ALA	2.7
1	B	15	GLY	2.7
1	A	183	VAL	2.6
1	B	260	THR	2.6
1	A	357	GLU	2.6
1	B	14	LEU	2.6
1	B	313	ALA	2.6
2	C	3	U	2.5
1	B	268	SER	2.5
1	A	356	ILE	2.5
1	A	468	GLY	2.4
1	B	208	SER	2.4
1	B	265	VAL	2.4
1	B	202	LEU	2.3
1	A	403	ILE	2.3
1	B	69	GLY	2.3
1	B	74	MET	2.3
1	A	631	THR	2.3
1	B	281	GLY	2.3
1	B	67	PRO	2.2
1	B	211	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	421	ASP	2.2
1	B	247	GLY	2.2
1	B	278	GLY	2.2
1	B	215	PRO	2.1
1	B	19	THR	2.1
1	B	284	TYR	2.1
1	A	384	LEU	2.1
1	B	68	LYS	2.0
1	B	216	ALA	2.0
1	A	15	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	5BU	C	4	21/22	0.89	0.20	-	59,65,73,99	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
5	ADP	A	875	27/27	0.97	0.14	-0.04	30,46,50,52	0
4	MG	A	850	1/1	0.98	0.13	-0.51	37,37,37,37	0
3	ZN	B	800	1/1	0.98	0.10	-0.95	61,61,61,61	0
6	BEF	A	885	4/4	0.96	0.11	-1.00	34,36,36,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	ZN	A	800	1/1	0.99	0.10	-1.08	61,61,61,61	0
7	SO4	B	895	5/5	0.93	0.14	-1.37	60,65,70,72	0
7	SO4	B	890	5/5	0.96	0.11	-1.83	53,55,66,71	0

6.5 Other polymers [i](#)

There are no such residues in this entry.