



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 11:50 am BST

PDB ID : 4B6E
Title : Discovery of an allosteric mechanism for the regulation of HCV NS3 protein function
Authors : Saalau-Bethell, S.M.; Woodhead, A.J.; Chessari, G.; Carr, M.G.; Coyle, J.; Graham, B.; Hiscock, S.D.; Murray, C.W.; Pathuri, P.; Rich, S.J.; Richardson, C.J.; Williams, P.A.; Jhoti, H.
Deposited on : 2012-08-09
Resolution : 2.46 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

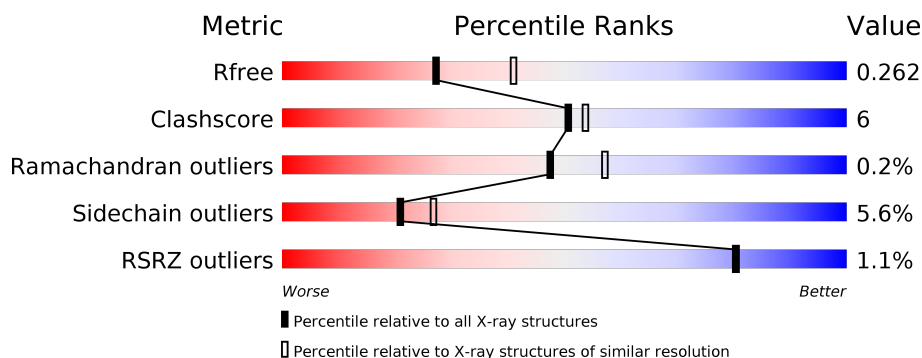
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.46 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}	130704	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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 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	683	<div> <div>2%</div> <div> <div></div> <div>81%</div> <div>12%</div> <div>• 5%</div> </div> </div>
1	B	683	<div> <div>80%</div> <div>13%</div> <div>• 6%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9975 atoms, of which 7 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 NON-STRUCTURAL PROTEIN 4A, SERINE PROTEASE NS3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	646	Total	C	N	O	S	0	0	1
			4808	3026	835	917	30			
1	B	643	Total	C	N	O	S	0	0	1
			4794	3019	832	913	30			

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-34	MET	-	expression tag	UNP P26663
A	-33	GLY	-	expression tag	UNP P26663
A	-32	SER	-	expression tag	UNP P26663
A	-31	SER	-	expression tag	UNP P26663
A	-30	HIS	-	expression tag	UNP P26663
A	-29	HIS	-	expression tag	UNP P26663
A	-28	HIS	-	expression tag	UNP P26663
A	-27	HIS	-	expression tag	UNP P26663
A	-26	HIS	-	expression tag	UNP P26663
A	-25	HIS	-	expression tag	UNP P26663
A	-24	SER	-	expression tag	UNP P26663
A	-23	SER	-	expression tag	UNP P26663
A	-22	GLY	-	expression tag	UNP P26663
A	-21	LEU	-	expression tag	UNP P26663
A	-20	VAL	-	expression tag	UNP P26663
A	-19	PRO	-	expression tag	UNP P26663
A	-18	ARG	-	expression tag	UNP P26663
A	-17	GLY	-	expression tag	UNP P26663
A	-16	SER	-	expression tag	UNP P26663
A	-15	HIS	-	expression tag	UNP P26663
A	-14	MET	-	expression tag	UNP P26663
A	-1	GLY	-	linker	UNP P26663
A	0	SER	-	linker	UNP P26663
A	1	GLY	-	linker	UNP P26663

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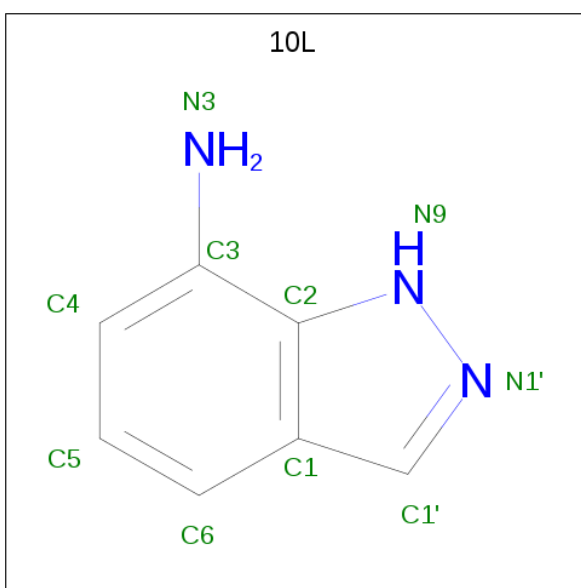
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	SER	-	linker	UNP P26663
B	-34	MET	-	expression tag	UNP P26663
B	-33	GLY	-	expression tag	UNP P26663
B	-32	SER	-	expression tag	UNP P26663
B	-31	SER	-	expression tag	UNP P26663
B	-30	HIS	-	expression tag	UNP P26663
B	-29	HIS	-	expression tag	UNP P26663
B	-28	HIS	-	expression tag	UNP P26663
B	-27	HIS	-	expression tag	UNP P26663
B	-26	HIS	-	expression tag	UNP P26663
B	-25	HIS	-	expression tag	UNP P26663
B	-24	SER	-	expression tag	UNP P26663
B	-23	SER	-	expression tag	UNP P26663
B	-22	GLY	-	expression tag	UNP P26663
B	-21	LEU	-	expression tag	UNP P26663
B	-20	VAL	-	expression tag	UNP P26663
B	-19	PRO	-	expression tag	UNP P26663
B	-18	ARG	-	expression tag	UNP P26663
B	-17	GLY	-	expression tag	UNP P26663
B	-16	SER	-	expression tag	UNP P26663
B	-15	HIS	-	expression tag	UNP P26663
B	-14	MET	-	expression tag	UNP P26663
B	-1	GLY	-	linker	UNP P26663
B	0	SER	-	linker	UNP P26663
B	1	GLY	-	linker	UNP P26663
B	2	SER	-	linker	UNP P26663

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 3 is 1H-indazol-7-amine (three-letter code: 10L) (formula: C₇H₇N₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	H	N	0	0
			17	7	7	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	186	Total 186	O 186	0	0
4	B	160	Total 160	O 160	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	91.40Å 111.31Å 141.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	87.57 – 2.46 87.57 – 2.47	Depositor EDS
% Data completeness (in resolution range)	98.0 (87.57-2.46) 98.0 (87.57-2.47)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.39 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.2.0019I	Depositor
R, R_{free}	0.187 , 0.257 0.193 , 0.262	Depositor DCC
R_{free} test set	2645 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	36.8	Xtriage
Anisotropy	0.009	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.1	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	9975	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, 10L

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	0.58	2/4916 (0.0%)	0.70	2/6713 (0.0%)
1	B	0.59	1/4901 (0.0%)	0.69	0/6692
All	All	0.59	3/9817 (0.0%)	0.69	2/13405 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	159	CYS	CB-SG	-6.15	1.71	1.82
1	A	47	CYS	CB-SG	-5.98	1.72	1.81
1	A	159	CYS	CB-SG	-5.70	1.72	1.81

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	109	ARG	NE-CZ-NH1	5.80	123.20	120.30
1	A	109	ARG	NE-CZ-NH2	-5.19	117.70	120.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	403	ILE	CA

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	4808	0	4783	52	0
1	B	4794	0	4771	59	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	B	10	7	7	1	0
4	A	186	0	0	3	0
4	B	160	0	0	2	0
All	All	9968	7	9561	110	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 (110) 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:97:CYS:SG	1:A:151:VAL:HG23	2.02	0.99
1:A:365:LEU:HD13	1:A:408:VAL:HG23	1.54	0.87
1:A:365:LEU:CD1	1:A:408:VAL:HG23	2.08	0.84
1:A:596:THR:HG22	1:A:607:ASN:HD22	1.44	0.83
1:A:365:LEU:HD13	1:A:408:VAL:CG2	2.10	0.80
1:B:347:ILE:HD13	1:B:381:LEU:HD21	1.68	0.75
1:A:596:THR:CG2	1:A:607:ASN:HD22	2.02	0.73
1:B:280:SER:HB2	1:B:283:ALA:HB2	1.75	0.68
1:A:347:ILE:HD13	1:A:381:LEU:HD21	1.76	0.67
1:B:214:VAL:HG22	1:B:215:PRO:HD3	1.75	0.67
1:A:565:ALA:HB1	1:A:585:LEU:HD21	1.79	0.63
1:A:185:THR:HG22	1:A:186:ASP:H	1.66	0.60
1:B:104:LEU:HD12	1:B:104:LEU:N	2.17	0.60
1:B:347:ILE:HD13	1:B:381:LEU:CD2	2.32	0.58
1:B:712:ARG:HD2	1:B:714:ILE:HD11	1.86	0.58
1:B:209:GLY:O	1:B:214:VAL:HG13	2.04	0.57
1:B:263:ALA:HB1	1:B:264:PRO:HD2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:229:ASN:HD22	1:B:231:SER:H	1.51	0.56
1:A:542:ILE:HD11	1:A:562:ALA:HB3	1.88	0.56
1:A:263:ALA:HB1	1:A:264:PRO:HD2	1.87	0.55
1:A:364:HIS:ND1	1:A:424:SER:OG	2.30	0.55
1:A:565:ALA:CB	1:A:585:LEU:HD21	2.36	0.55
1:A:132:VAL:HG22	1:A:159:CYS:SG	2.47	0.55
1:B:74:MET:HE2	1:B:75:TYR:CE2	2.42	0.55
1:B:116:VAL:HG22	1:B:126:LEU:HD12	1.88	0.54
1:A:587:ARG:HB3	1:A:588:LEU:HD12	1.89	0.54
1:A:356:ILE:HD11	1:A:384:LEU:HD23	1.90	0.54
1:A:596:THR:HG22	1:A:607:ASN:ND2	2.17	0.53
1:A:14:LEU:HD22	1:B:21:LEU:CD2	2.38	0.53
1:B:341:LEU:HD13	1:B:474:ARG:HB3	1.91	0.53
1:B:53:TRP:NE1	1:B:175:MET:HE3	2.23	0.53
1:A:239:GLY:HA3	1:A:252:ILE:HD11	1.91	0.52
1:B:53:TRP:CE2	1:B:175:MET:HE3	2.45	0.52
1:B:53:TRP:CD1	1:B:175:MET:CE	2.93	0.51
1:A:596:THR:CG2	1:A:607:ASN:ND2	2.70	0.51
1:A:477:THR:HG22	1:A:478:PRO:O	2.11	0.51
1:A:185:THR:HG22	1:A:186:ASP:N	2.26	0.50
1:B:251:ASN:HD22	1:B:261:THR:H	1.59	0.50
1:A:365:LEU:HD12	1:A:408:VAL:HG23	1.90	0.49
1:B:557:PHE:O	1:B:561:VAL:HG23	2.12	0.49
1:B:239:GLY:HA3	1:B:252:ILE:HD11	1.93	0.49
1:A:132:VAL:CG2	1:A:159:CYS:SG	3.00	0.49
1:B:251:ASN:ND2	1:B:261:THR:H	2.10	0.49
1:B:588:LEU:O	1:B:592:LEU:HD13	2.11	0.49
1:B:619:ILE:O	1:B:622:CYS:HB2	2.13	0.49
1:B:583:LYS:C	1:B:585:LEU:H	2.15	0.49
1:B:616:THR:O	1:B:620:MET:HG3	2.13	0.48
1:A:210:LYS:HA	1:A:214:VAL:HG13	1.94	0.48
1:B:712:ARG:CD	1:B:714:ILE:HD11	2.43	0.48
1:B:53:TRP:CD1	1:B:175:MET:HE1	2.48	0.48
1:A:396:ASP:O	1:A:399:VAL:HG22	2.13	0.48
1:B:307:LEU:HD13	1:B:519:THR:OG1	2.15	0.47
1:A:214:VAL:HG22	1:A:215:PRO:HD3	1.97	0.47
1:A:263:ALA:HB1	1:A:264:PRO:CD	2.44	0.47
1:A:191:PRO:HG2	1:A:317:LEU:HD21	1.95	0.47
1:A:363:ARG:HG2	1:A:406:VAL:HG12	1.97	0.47
1:B:82:LEU:HG	1:B:170:VAL:HG11	1.97	0.47
1:A:432:VAL:HG13	1:A:448:THR:HG23	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:625:ALA:HB2	4:B:2139:HOH:O	2.14	0.46
1:A:356:ILE:CD1	1:A:384:LEU:HD23	2.44	0.46
1:B:449:THR:HG22	4:B:2114:HOH:O	2.15	0.46
1:A:371:LYS:NZ	4:A:2111:HOH:O	2.48	0.46
1:B:63:THR:HG22	1:B:72:THR:HA	1.98	0.46
1:B:576:PRO:HB3	1:B:605:VAL:HG22	1.97	0.46
1:A:538:GLY:HA3	1:A:618:TYR:CE2	2.51	0.46
1:A:571:ALA:HB1	4:A:2160:HOH:O	2.16	0.45
1:B:187:ASN:O	1:B:202:LEU:HD12	2.16	0.45
1:A:126:LEU:HD11	1:A:167:VAL:HG13	1.98	0.45
1:B:55:VAL:HG21	1:B:139:SER:HB3	1.99	0.45
1:B:126:LEU:CD1	1:B:167:VAL:HG13	2.47	0.44
1:B:53:TRP:CD1	1:B:175:MET:HE3	2.53	0.44
1:A:104:LEU:HD21	1:A:118:ARG:HG3	2.00	0.44
1:A:214:VAL:N	1:A:215:PRO:HD2	2.33	0.44
1:B:104:LEU:HD11	1:B:118:ARG:HB2	1.99	0.44
1:A:561:VAL:HG13	1:A:582:TRP:CZ3	2.53	0.44
1:A:137:GLY:H	1:A:632:LEU:N	2.15	0.44
1:B:589:LYS:HB3	1:B:590:PRO:HD3	2.00	0.43
1:B:445:THR:HG22	1:B:447:GLU:HG3	2.01	0.43
1:A:392:TYR:CE1	1:A:395:LEU:HD13	2.53	0.43
1:A:29:VAL:HA	4:A:2014:HOH:O	2.19	0.43
1:B:228:LEU:HD23	1:B:268:SER:HB3	2.00	0.43
1:B:514:ARG:HG3	1:B:518:ASN:ND2	2.34	0.42
1:A:565:ALA:HB1	1:A:585:LEU:CD2	2.45	0.42
1:A:145:CYS:SG	1:A:147:SER:OG	2.68	0.42
1:A:158:VAL:HB	1:A:166:ALA:HB3	2.02	0.42
1:A:446:ILE:HG21	1:A:557:PHE:CE2	2.55	0.42
1:A:180:ARG:O	1:A:181:SER:OG	2.35	0.42
1:B:104:LEU:CD1	1:B:104:LEU:N	2.81	0.42
1:B:606:GLN:HB3	1:B:606:GLN:HE21	1.65	0.42
1:B:293:HIS:CD2	1:B:322:THR:OG1	2.73	0.42
1:B:526:GLN:NE2	1:B:628:GLU:HB3	2.35	0.42
1:B:214:VAL:HG22	1:B:215:PRO:CD	2.47	0.42
1:B:229:ASN:ND2	1:B:231:SER:H	2.15	0.42
1:A:432:VAL:CG1	1:A:448:THR:HG23	2.50	0.42
1:B:481:ARG:HB2	1:B:482:PRO:HD2	2.02	0.42
1:A:82:LEU:HG	1:A:170:VAL:HG11	2.02	0.41
1:B:333:HIS:CE1	1:B:335:ASN:HB2	2.55	0.41
1:B:293:HIS:HD2	1:B:322:THR:OG1	2.03	0.41
1:A:421:ASP:C	1:A:422:PHE:CD1	2.93	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:LEU:HD11	1:B:167:VAL:HG13	2.02	0.41
1:B:485:MET:SD	3:B:1722:10L:H4	2.61	0.41
1:B:104:LEU:CD1	1:B:118:ARG:HB2	2.50	0.41
1:B:390:ALA:HA	1:B:409:VAL:O	2.21	0.41
1:B:51:VAL:CG1	1:B:53:TRP:CE2	3.04	0.41
1:A:251:ASN:HD22	1:A:261:THR:H	1.69	0.41
1:B:535:VAL:O	1:B:539:LEU:HD13	2.21	0.41
1:B:347:ILE:HB	1:B:354:ILE:HB	2.03	0.41
1:B:51:VAL:HG11	1:B:53:TRP:CZ2	2.55	0.40
1:A:396:ASP:OD2	1:A:398:SER:OG	2.29	0.40
1:B:392:TYR:CE1	1:B:395:LEU:HD13	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	642/683 (94%)	614 (96%)	26 (4%)	2 (0%)	41	49
1	B	638/683 (93%)	616 (97%)	21 (3%)	1 (0%)	47	57
All	All	1280/1366 (94%)	1230 (96%)	47 (4%)	3 (0%)	47	57

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	246	HIS
1	A	403	ILE
1	B	572	GLN

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	526/557 (94%)	496 (94%)	30 (6%)	20	26
1	B	525/557 (94%)	496 (94%)	29 (6%)	21	27
All	All	1051/1114 (94%)	992 (94%)	59 (6%)	21	27

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	LEU
1	A	68	LYS
1	A	82	LEU
1	A	104	LEU
1	A	109	ARG
1	A	114	ILE
1	A	118	ARG
1	A	132	VAL
1	A	160	THR
1	A	198	GLN
1	A	214	VAL
1	A	224	LYS
1	A	232	VAL
1	A	242	MET
1	A	246	HIS
1	A	249	ASP
1	A	337	GLU
1	A	341	LEU
1	A	360	ARG
1	A	380	LYS
1	A	402	THR
1	A	419	THR
1	A	431	CYS
1	A	453	GLN
1	A	470	ARG
1	A	529	LEU
1	A	585	LEU

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Mol	Chain	Res	Type
1	A	596	THR
1	A	611	LEU
1	A	627	LEU
1	B	21	LEU
1	B	26	LYS
1	B	32	GLU
1	B	64	LEU
1	B	72	THR
1	B	82	LEU
1	B	95	THR
1	B	118	ARG
1	B	143	LEU
1	B	206	THR
1	B	214	VAL
1	B	224	LYS
1	B	280	SER
1	B	316	ARG
1	B	329	VAL
1	B	341	LEU
1	B	384	LEU
1	B	431	CYS
1	B	435	THR
1	B	449	THR
1	B	459	SER
1	B	517	LEU
1	B	529	LEU
1	B	535	VAL
1	B	584	CYS
1	B	592	LEU
1	B	600	TYR
1	B	611	LEU
1	B	627	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	229	ASN
1	A	251	ASN
1	A	369	HIS
1	A	607	ASN
1	B	27	ASN
1	B	221	GLN

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Mol	Chain	Res	Type
1	B	229	ASN
1	B	251	ASN
1	B	293	HIS
1	B	518	ASN
1	B	606	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

3 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	SO4	A	1721	-	4,4,4	0.12	0	6,6,6	0.36	0
3	10L	B	1722	-	11,11,11	0.98	1 (9%)	7,15,15	1.26	1 (14%)
2	SO4	B	1721	-	4,4,4	0.15	0	6,6,6	0.13	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
3	10L	B	1722	-	-	-	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1722	10L	C1-C2	-2.73	1.39	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1722	10L	C2-C3-N3	2.59	123.01	118.07

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1722	10L	1	0

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	646/683 (94%)	-0.02	11 (1%) 70 67	14, 38, 71, 82	0
1	B	643/683 (94%)	-0.12	3 (0%) 91 92	19, 38, 67, 96	0
All	All	1289/1366 (94%)	-0.07	14 (1%) 80 80	14, 38, 68, 96	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	183	VAL	6.0
1	A	182	PRO	4.6
1	A	468	GLY	4.1
1	B	184	PHE	4.0
1	A	99	CYS	4.0
1	A	632	LEU	3.4
1	A	184	PHE	3.3
1	A	183	VAL	3.0
1	B	206	THR	2.4
1	A	98	THR	2.2
1	A	466	GLY	2.1
1	A	631	THR	2.1
1	A	588	LEU	2.1
1	A	181	SER	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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. 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(\AA^2)	Q<0.9
2	SO4	B	1721	5/5	0.44	0.30	142,142,143,143	0
3	10L	B	1722	10/10	0.78	0.32	66,70,71,71	0
2	SO4	A	1721	5/5	0.98	0.14	43,43,44,49	0

6.5 Other polymers [i](#)

There are no such residues in this entry.