



Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2020 – 10:49 pm BST

PDB ID : 3L2R
Title : Crystal structure of the Prototype Foamy Virus (PFV) intasome in complex with magnesium
Authors : Hare, S.; Gupta, S.S.; Cherepanov, P.
Deposited on : 2009-12-15
Resolution : 2.88 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

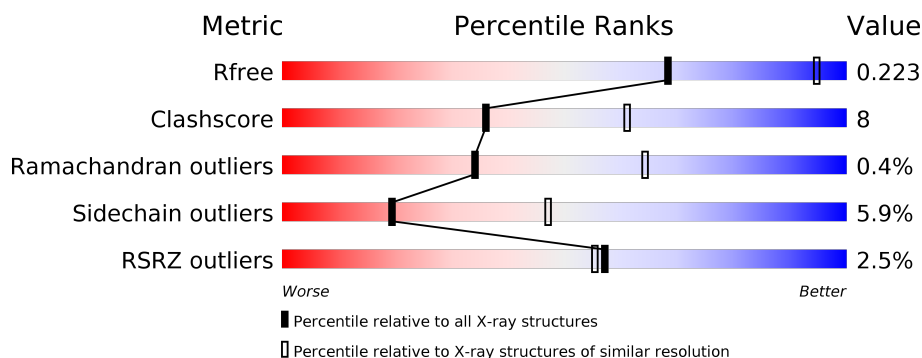
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.88 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	2691 (2.90-2.86)
Clashscore	141614	2947 (2.90-2.86)
Ramachandran outliers	138981	2868 (2.90-2.86)
Sidechain outliers	138945	2871 (2.90-2.86)
RSRZ outliers	127900	2629 (2.90-2.86)

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	395	<div> <div>2%</div> <div> <div></div> <div>72%</div> <div>18%</div> <div>•</div> <div>8%</div> </div> </div>
1	B	395	<div> <div>%</div> <div> <div></div> <div>34%</div> <div>7%</div> <div>•</div> <div>59%</div> </div> </div>
2	C	19	<div> <div></div> <div> <div>26%</div> <div>53%</div> <div>21%</div> </div> </div>
3	D	17	<div> <div>6%</div> <div> <div></div> <div>53%</div> <div>41%</div> <div>6%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	GOL	A	805	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 4928 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 Integrase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	365	Total	C	N	O	S	0	0	0
			2886	1853	506	523	4			
1	B	163	Total	C	N	O	S	2	0	0
			1264	830	201	232	1			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP P14350
A	-1	PRO	-	EXPRESSION TAG	UNP P14350
A	0	GLY	-	EXPRESSION TAG	UNP P14350
A	217	SER	GLY	VARIANT	UNP P14350
A	218	GLY	SER	VARIANT	UNP P14350
B	-2	GLY	-	EXPRESSION TAG	UNP P14350
B	-1	PRO	-	EXPRESSION TAG	UNP P14350
B	0	GLY	-	EXPRESSION TAG	UNP P14350
B	217	SER	GLY	VARIANT	UNP P14350
B	218	GLY	SER	VARIANT	UNP P14350

- Molecule 2 is a DNA chain called 5'-D(*AP*TP*TP*GP*TP*CP*AP*TP*GP*GP*AP*A P*TP*TP*TP*TP*GP*TP*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	19	Total	C	N	O	P	0	0	0
			389	189	66	116	18			

- Molecule 3 is a DNA chain called 5'-D(*TP*AP*CP*AP*AP*AP*AP*TP*TP*CP*CP*AP *TP*GP*AP*CP*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	17	Total	C	N	O	P	0	0	0
			343	166	65	96	16			

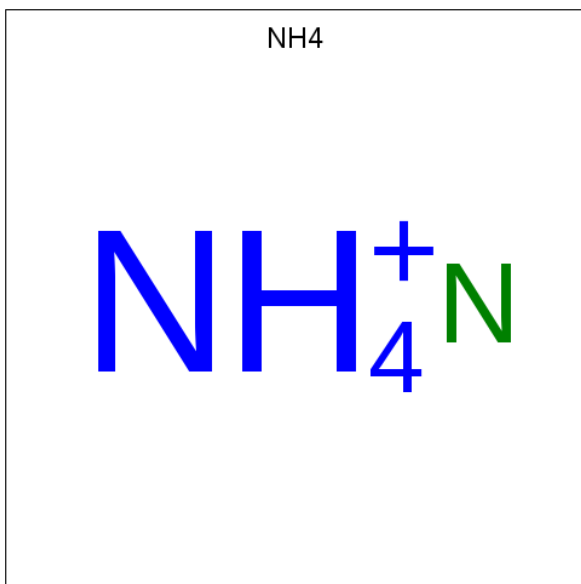
- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Mg	0	0
			1	1		
5	A	1	Total	Mg	0	0
			1	1		

- Molecule 6 is AMMONIUM ION (three-letter code: NH4) (formula: H₄N).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	N	0	0
			1	1		

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	5	Total	O	0	0
			5	5		
8	B	4	Total	O	0	0
			4	4		
8	C	1	Total	O	0	0
			1	1		
8	D	2	Total	O	0	0
			2	2		



● Molecule 3: 5'-D(*TP*AP*CP*AP*AP*AP*AP*TP*TP*CP*CP*AP*TP*GP*AP*CP*A)-3',



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	158.48 Å 158.48 Å 124.19 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.00 – 2.88 39.00 – 2.88	Depositor EDS
% Data completeness (in resolution range)	99.0 (39.00-2.88) 99.0 (39.00-2.88)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 2.90 Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.202 , 0.222 0.206 , 0.223	Depositor DCC
R_{free} test set	1777 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	78.8	Xtriage
Anisotropy	0.282	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 50.9	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	4928	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% 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: GOL, ZN, MG, NH4

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	0.61	0/2965	0.74	2/4050 (0.0%)
1	B	0.63	0/1304	0.66	1/1789 (0.1%)
2	C	1.22	0/435	2.07	18/671 (2.7%)
3	D	1.08	0/385	1.91	11/591 (1.9%)
All	All	0.73	0/5089	1.06	32/7101 (0.5%)

There are no bond length outliers.

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	12	DA	O4'-C1'-N9	-8.31	102.19	108.00
3	D	17	DA	O4'-C1'-N9	7.99	113.59	108.00
2	C	8	DT	O4'-C1'-N1	-7.84	102.51	108.00
2	C	8	DT	O4'-C4'-C3'	-7.68	101.39	106.00
2	C	7	DA	O4'-C1'-N9	-7.59	102.68	108.00
3	D	4	DA	O4'-C1'-N9	-7.33	102.86	108.00
3	D	1	DT	P-O3'-C3'	7.21	128.35	119.70
2	C	8	DT	O4'-C1'-C2'	-6.88	100.40	105.90
2	C	8	DT	P-O3'-C3'	-6.84	111.50	119.70
1	B	235	LEU	CA-CB-CG	6.77	130.87	115.30
2	C	15	DT	N3-C4-O4	6.63	123.88	119.90
2	C	2	DT	P-O3'-C3'	6.43	127.41	119.70
3	D	17	DA	O4'-C4'-C3'	-6.42	101.93	104.50
2	C	2	DT	N3-C2-O2	-6.40	118.46	122.30
2	C	3	DT	P-O3'-C3'	6.30	127.27	119.70
3	D	1	DT	O4'-C1'-N1	6.26	112.38	108.00
2	C	5	DT	N3-C4-O4	6.18	123.61	119.90
3	D	15	DA	O5'-P-OP2	-6.09	100.22	105.70
2	C	14	DT	P-O3'-C3'	6.09	127.01	119.70
2	C	1	DA	P-O3'-C3'	6.09	127.01	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	220	VAL	CB-CA-C	-6.09	99.83	111.40
2	C	11	DA	P-O3'-C3'	5.78	126.64	119.70
3	D	8	DT	N3-C4-O4	5.72	123.33	119.90
2	C	19	DA	N1-C6-N6	5.42	121.85	118.60
2	C	15	DT	C5-C4-O4	-5.37	121.14	124.90
3	D	10	DC	O4'-C1'-N1	-5.36	104.25	108.00
3	D	15	DA	O4'-C4'-C3'	5.30	109.18	106.00
2	C	17	DG	C5-C6-N1	5.22	114.11	111.50
2	C	16	DT	C4-C5-C7	5.20	122.12	119.00
1	A	251	LEU	CA-CB-CG	5.04	126.88	115.30
2	C	7	DA	N1-C6-N6	5.03	121.62	118.60
3	D	4	DA	C2-N3-C4	-5.01	108.10	110.60

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	2886	0	2910	60	0
1	B	1264	0	1244	15	0
2	C	389	0	220	4	0
3	D	343	0	193	2	0
4	A	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	1	0	0	0	0
7	A	24	0	32	8	0
7	B	6	0	8	0	0
8	A	5	0	0	0	0
8	B	4	0	0	0	0
8	C	1	0	0	0	0
8	D	2	0	0	0	0
All	All	4928	0	4607	76	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 8.

All (76) 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:202:ARG:HH11	1:B:202:ARG:HG2	1.28	0.98
1:A:357:HIS:HA	1:A:358:LEU:HB2	1.55	0.88
1:A:126:PHE:HB3	1:A:220:VAL:HG22	1.58	0.83
1:B:202:ARG:HH11	1:B:202:ARG:CG	1.92	0.83
2:C:1:DA:H2''	2:C:2:DT:H5'	1.60	0.82
1:A:46:LYS:HE2	1:A:77:ALA:O	1.84	0.78
1:A:228:LYS:HE3	3:D:17:DA:OP2	1.90	0.72
1:A:339:LYS:HG2	7:A:805:GOL:H32	1.73	0.70
1:A:339:LYS:NZ	1:A:358:LEU:HD13	2.07	0.68
1:A:136:SER:HA	7:A:802:GOL:H11	1.78	0.65
1:A:175:SER:HB2	1:B:247:PRO:HA	1.77	0.65
1:B:202:ARG:HG2	1:B:202:ARG:NH1	2.03	0.64
1:A:126:PHE:HB3	1:A:220:VAL:CG2	2.28	0.63
1:A:26:LYS:HG3	1:A:27:GLN:H	1.62	0.63
1:A:339:LYS:HG2	7:A:805:GOL:C3	2.28	0.63
1:B:260:VAL:O	1:B:261:LEU:HB2	1.99	0.63
1:A:209:SER:HB3	1:A:216:SER:HB3	1.81	0.63
1:B:130:ILE:HD12	1:B:144:VAL:HG21	1.82	0.61
1:A:328:ALA:O	1:A:330:PRO:HD3	2.01	0.60
1:A:274:SER:OG	1:B:178:ILE:N	2.35	0.60
1:A:339:LYS:H	7:A:805:GOL:C3	2.15	0.60
1:A:357:HIS:HA	1:A:358:LEU:CB	2.27	0.59
1:A:292:LEU:HD22	1:B:271:GLY:HA3	1.84	0.59
1:A:339:LYS:HZ2	1:A:358:LEU:HD13	1.66	0.58
1:A:113:LEU:O	1:A:114:ARG:HD2	2.03	0.58
1:A:344:LEU:HD11	1:A:355:LEU:HB2	1.86	0.57
1:A:209:SER:HB3	1:A:216:SER:CB	2.35	0.57
1:A:73:LEU:HD22	1:A:86:ARG:CZ	2.34	0.57
1:A:70:GLU:CD	1:A:86:ARG:HH21	2.09	0.56
1:A:339:LYS:H	7:A:805:GOL:H32	1.69	0.55
1:A:356:ASP:CG	1:A:357:HIS:N	2.60	0.55
1:A:230:LEU:HD23	1:A:249:VAL:HG13	1.87	0.54
1:B:127:ILE:HA	1:B:144:VAL:O	2.08	0.54
1:A:241:LYS:HG2	7:A:802:GOL:O1	2.08	0.54
3:D:16:DC:H5''	3:D:17:DA:OP2	2.09	0.52
1:B:127:ILE:HG22	1:B:145:VAL:HG22	1.91	0.52
1:A:73:LEU:O	1:A:73:LEU:HG	2.09	0.51
1:A:122:PHE:O	1:A:179:PRO:HA	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:272:ILE:O	1:A:272:ILE:HG13	2.10	0.51
1:A:126:PHE:CB	1:A:220:VAL:HG22	2.37	0.50
1:A:356:ASP:CG	1:A:357:HIS:H	2.14	0.49
1:A:325:GLU:OE2	1:A:362:ARG:NH2	2.43	0.49
1:A:183:HIS:HA	1:A:207:GLU:O	2.13	0.48
1:A:112:ILE:HA	1:A:350:ARG:CD	2.44	0.48
1:A:111:PRO:HG3	1:A:310:ALA:HA	1.94	0.48
1:A:358:LEU:HA	1:A:359:GLY:HA2	1.59	0.47
1:A:339:LYS:HZ1	1:A:358:LEU:HD13	1.77	0.46
1:A:66:HIS:HB3	1:A:99:CYS:SG	2.56	0.46
1:A:26:LYS:HG3	1:A:27:GLN:N	2.29	0.46
1:B:158:THR:HB	1:B:165:ALA:HB1	1.99	0.45
1:A:180:LYS:HE3	1:B:275:ASN:OD1	2.17	0.45
1:A:70:GLU:OE1	1:A:86:ARG:NH2	2.43	0.44
1:A:256:THR:C	1:A:264:THR:HG22	2.38	0.44
1:A:156:TYR:OH	1:A:173:LEU:HB2	2.18	0.43
1:B:121:PRO:O	1:B:122:PHE:HB2	2.19	0.43
1:A:49:PRO:HA	1:A:50:PRO:HD2	1.86	0.43
2:C:4:DG:H2''	2:C:5:DT:O5'	2.18	0.43
1:A:82:TRP:O	1:A:85:MET:HG3	2.18	0.42
1:A:163:THR:H	7:A:801:GOL:H11	1.84	0.42
1:B:134:PRO:HA	1:B:135:PRO:HD3	1.95	0.42
1:A:179:PRO:HG2	1:A:204:ILE:HD13	2.00	0.42
1:A:122:PHE:CE2	1:A:177:ALA:HB3	2.54	0.42
1:A:161:PRO:O	1:A:189:ALA:HB2	2.19	0.42
1:A:176:ILE:HG13	1:A:177:ALA:N	2.35	0.42
1:A:339:LYS:HG2	7:A:805:GOL:O3	2.19	0.42
1:A:41:ARG:O	1:A:44:GLY:N	2.53	0.42
2:C:10:DG:H2''	2:C:11:DA:C8	2.55	0.41
1:A:222:ARG:O	1:A:222:ARG:HD2	2.20	0.41
1:A:111:PRO:O	1:A:350:ARG:HD2	2.20	0.41
1:A:67:THR:HB	1:A:71:ALA:HB3	2.02	0.41
1:B:223:LYS:O	1:B:227:ILE:HG13	2.22	0.40
1:A:315:TRP:CE2	1:A:371:PRO:HD3	2.56	0.40
1:A:174:THR:HA	1:A:177:ALA:O	2.21	0.40
2:C:1:DA:H3'	2:C:2:DT:H71	2.03	0.40
1:A:112:ILE:HA	1:A:350:ARG:HD2	2.03	0.40
1:A:350:ARG:HH22	1:A:367:ASP:CG	2.25	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	363/395 (92%)	349 (96%)	13 (4%)	1 (0%)	41	70
1	B	161/395 (41%)	153 (95%)	7 (4%)	1 (1%)	25	55
All	All	524/790 (66%)	502 (96%)	20 (4%)	2 (0%)	34	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	328	ALA
1	B	261	LEU

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	322/354 (91%)	304 (94%)	18 (6%)	21	49
1	B	139/354 (39%)	130 (94%)	9 (6%)	17	42
All	All	461/708 (65%)	434 (94%)	27 (6%)	19	47

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LEU
1	A	32	LEU
1	A	149	MET
1	A	176	ILE

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Mol	Chain	Res	Type
1	A	191	THR
1	A	201	GLU
1	A	204	ILE
1	A	209	SER
1	A	220	VAL
1	A	238	ARG
1	A	251	LEU
1	A	262	LYS
1	A	312	SER
1	A	351	THR
1	A	352	VAL
1	A	358	LEU
1	A	360	ASN
1	A	373	SER
1	B	119	GLN
1	B	164	SER
1	B	176	ILE
1	B	202	ARG
1	B	210	THR
1	B	215	GLN
1	B	232	THR
1	B	272	ILE
1	B	274	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 9 ligands modelled in this entry, 1 is modelled with single atom and 3 are monoatomic - leaving 5 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	$\# Z > 2$
7	GOL	A	805	-	5,5,5	0.41	0	5,5,5	1.01	0
7	GOL	A	804	-	5,5,5	0.47	0	5,5,5	0.48	0
7	GOL	A	801	-	5,5,5	0.47	0	5,5,5	0.54	0
7	GOL	A	802	-	5,5,5	0.45	0	5,5,5	0.24	0
7	GOL	B	803	-	5,5,5	0.36	0	5,5,5	0.24	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
7	GOL	A	805	-	-	2/4/4/4	-
7	GOL	A	804	-	-	2/4/4/4	-
7	GOL	A	801	-	-	4/4/4/4	-
7	GOL	A	802	-	-	1/4/4/4	-
7	GOL	B	803	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	801	GOL	C1-C2-C3-O3
7	A	804	GOL	O1-C1-C2-C3
7	A	804	GOL	O1-C1-C2-O2
7	A	805	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
7	A	801	GOL	O1-C1-C2-C3
7	A	801	GOL	O1-C1-C2-O2
7	A	801	GOL	O2-C2-C3-O3
7	A	802	GOL	C1-C2-C3-O3
7	A	805	GOL	O2-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	805	GOL	5	0
7	A	801	GOL	1	0
7	A	802	GOL	2	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 [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	365/395 (92%)	0.07	9 (2%) 57 55	49, 68, 116, 140	0
1	B	163/395 (41%)	-0.09	4 (2%) 57 55	58, 75, 109, 126	1 (0%)
2	C	19/19 (100%)	-0.14	0 100 100	50, 69, 97, 123	0
3	D	17/17 (100%)	-0.33	1 (5%) 22 18	55, 62, 106, 142	0
All	All	564/826 (68%)	0.00	14 (2%) 57 55	49, 70, 113, 142	1 (0%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	116	ASP	3.6
1	A	11	GLU	3.6
1	B	239	PRO	2.9
1	A	16	LEU	2.8
1	A	32	LEU	2.6
3	D	1	DT	2.6
1	B	235	LEU	2.6
1	A	12	LEU	2.4
1	B	278	PHE	2.3
1	A	355	LEU	2.2
1	A	13	ASP	2.2
1	A	357	HIS	2.1
1	A	31	PHE	2.1
1	A	318	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no 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(Å ²)	Q<0.9
7	GOL	A	801	6/6	0.70	0.28	101,104,106,106	0
6	NH4	A	395	1/1	0.85	0.35	34,34,34,34	0
7	GOL	A	804	6/6	0.90	0.29	80,81,82,82	0
7	GOL	B	803	6/6	0.93	0.60	88,90,91,92	0
7	GOL	A	802	6/6	0.94	0.22	93,94,96,96	0
7	GOL	A	805	6/6	0.94	0.14	67,76,77,79	0
5	MG	B	393	1/1	0.96	0.11	89,89,89,89	0
5	MG	A	394	1/1	0.96	0.10	66,66,66,66	0
4	ZN	A	393	1/1	0.99	0.15	62,62,62,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.