



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 14, 2020 – 01:45 am BST

PDB ID : 1H80
Title : 1,3-ALPHA-1,4-BETA-D-GALACTOSE-4-SULFATE- 3,6-ANHYDRO-D-GALACTOSE-2-SULFATE 4 GALACTOXYDROLASE
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Deposited on : 2001-01-22
Resolution : 1.60 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

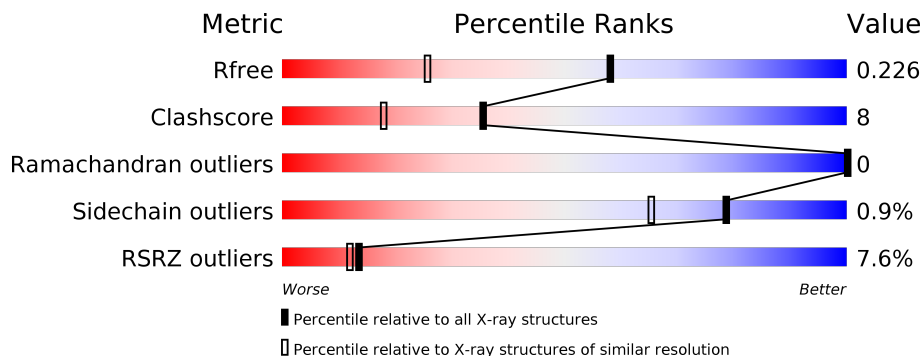
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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	464	 7% 79% 13% • 7%
1	B	464	 8% 79% 13% • 7%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7992 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 IOTA-CARRAGEENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	433	3431	2143	612	657	19	55	11	0
1	B	430	3401	2125	605	652	19	85	9	0

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total	Ca	0	0
			3	3		
2	A	3	Total	Ca	0	0
			3	3		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	3	Total	Na	0	0
			3	3		
4	A	3	Total	Na	0	0
			3	3		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

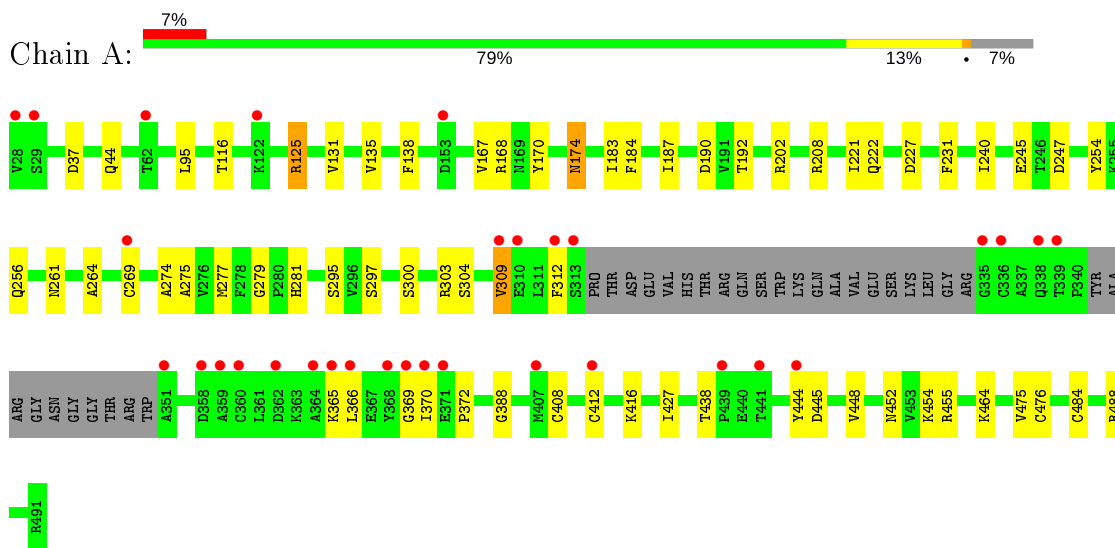
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	531	Total O 531 531	0	0
6	B	579	Total O 579 579	0	0

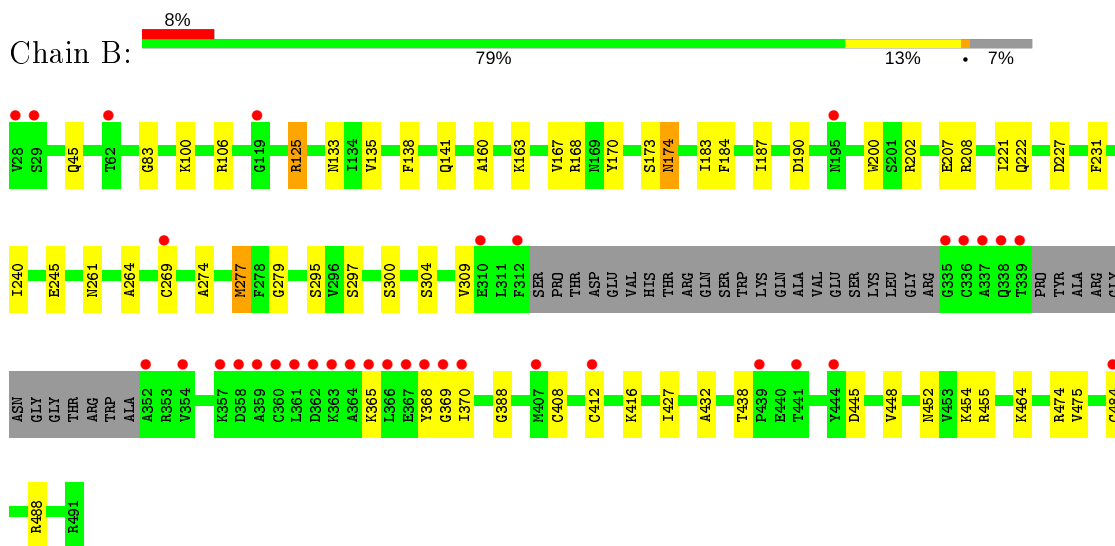
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: IOTA-CARRAGEENASE



- Molecule 1: IOTA-CARRAGEENASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.87Å 90.07Å 124.12Å 90.00° 93.53° 90.00°	Depositor
Resolution (Å)	20.00 – 1.60 8.00 – 1.60	Depositor EDS
% Data completeness (in resolution range)	97.4 (20.00-1.60) 97.6 (8.00-1.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.35 (at 1.60Å)	Xtrriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.207 , 0.223 0.210 , 0.226	Depositor DCC
R_{free} test set	7885 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	15.9	Xtrriage
Anisotropy	0.204	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.46 , 55.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7992	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.51% 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, CA, CL, NA

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	0.32	0/3554	0.65	1/4796 (0.0%)
1	B	0.33	0/3515	0.65	1/4743 (0.0%)
All	All	0.33	0/7069	0.65	2/9539 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	174	ASN	N-CA-C	5.71	126.41	111.00
1	B	174	ASN	N-CA-C	5.50	125.86	111.00

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	3431	0	3336	60	0
1	B	3401	0	3298	50	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	1	0	0	1	0
3	B	1	0	0	1	0
4	A	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	3	0	0	0	0
5	A	18	0	24	3	0
5	B	18	0	24	3	0
6	A	531	0	0	4	0
6	B	579	0	0	1	0
All	All	7992	0	6682	108	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 (108) 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:309:VAL:HG12	1:A:370:ILE:HD12	1.24	1.16
1:A:408[B]:CYS:SG	1:A:475:VAL:HG23	2.03	0.98
1:A:309:VAL:CG1	1:A:370:ILE:HD12	1.97	0.94
1:B:269[B]:CYS:SG	1:B:295:SER:HB2	2.14	0.87
1:A:269[B]:CYS:SG	1:A:295:SER:HB2	2.15	0.86
1:B:408[B]:CYS:SG	1:B:475:VAL:HG23	2.14	0.86
1:A:309:VAL:HG12	1:A:370:ILE:CD1	2.06	0.84
1:A:269[B]:CYS:SG	1:A:295:SER:CB	2.72	0.78
1:A:202:ARG:HD2	5:A:1501:GOL:O3	1.83	0.78
1:B:269[B]:CYS:SG	1:B:295:SER:CB	2.72	0.77
1:A:412:CYS:SG	1:A:427:ILE:HB	2.28	0.72
1:A:309:VAL:CG1	1:A:370:ILE:CD1	2.65	0.72
1:B:125:ARG:HG3	1:B:160:ALA:HB3	1.70	0.72
1:B:412:CYS:SG	1:B:427:ILE:HB	2.31	0.71
1:A:44[B]:GLN:HG3	1:B:208:ARG:NH1	2.06	0.70
1:A:131:VAL:HA	1:A:192:THR:HG21	1.73	0.69
1:B:200:TRP:CE2	5:B:1501:GOL:H2	2.27	0.69
1:B:484[B]:CYS:SG	1:B:488:ARG:HB3	2.34	0.68
1:A:309:VAL:HG11	1:A:370:ILE:HG13	1.74	0.67
1:B:304[B]:SER:OG	1:B:448:VAL:HG21	1.98	0.64
1:B:412:CYS:SG	1:B:484[B]:CYS:HB2	2.40	0.60
1:B:455:ARG:HD3	1:B:464:LYS:HA	1.84	0.60
1:A:416:LYS:H	1:A:416:LYS:CD	2.16	0.58
1:A:484[B]:CYS:SG	1:A:488:ARG:HB3	2.44	0.57
1:B:202:ARG:NH2	5:B:1501:GOL:O3	2.37	0.57
1:A:416:LYS:HD3	1:A:416:LYS:N	2.19	0.57
1:B:106:ARG:NH1	1:B:141[B]:GLN:HG3	2.20	0.57
1:B:438:THR:HB	1:B:445:ASP:OD1	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:ARG:HG2	1:B:163:LYS:HB2	1.87	0.56
1:B:416:LYS:H	1:B:416:LYS:CE	2.19	0.55
1:B:416:LYS:CD	1:B:416:LYS:H	2.18	0.55
1:B:416:LYS:HD3	1:B:416:LYS:N	2.22	0.55
1:A:309:VAL:HG11	1:A:370:ILE:CG1	2.37	0.54
1:A:416:LYS:H	1:A:416:LYS:CE	2.20	0.54
1:A:240:ILE:CD1	1:A:277[B]:MET:SD	2.96	0.54
1:A:44[B]:GLN:HG3	1:B:208:ARG:HH12	1.73	0.53
1:A:412:CYS:SG	1:A:484[B]:CYS:HB2	2.47	0.53
1:A:416:LYS:HD3	1:A:416:LYS:H	1.72	0.53
1:B:240:ILE:CD1	1:B:277[B]:MET:SD	2.97	0.53
1:B:200:TRP:NE1	5:B:1501:GOL:H2	2.26	0.51
1:A:372:PRO:HG2	1:A:444:TYR:HB3	1.92	0.51
1:A:452:ASN:OD1	1:A:454:LYS:HE2	2.10	0.51
1:B:416:LYS:HD3	1:B:416:LYS:H	1.76	0.51
1:A:408[B]:CYS:SG	1:A:475:VAL:C	2.89	0.50
1:A:183:ILE:O	1:A:184:PHE:HB2	2.12	0.49
1:A:416:LYS:N	1:A:416:LYS:CD	2.75	0.49
1:A:455:ARG:HD3	1:A:464:LYS:HA	1.95	0.49
1:A:261:ASN:HB3	6:A:2298:HOH:O	2.13	0.48
1:B:183:ILE:O	1:B:184:PHE:HB2	2.13	0.48
1:B:416:LYS:N	1:B:416:LYS:CD	2.77	0.48
1:A:231:PHE:O	1:A:264:ALA:HA	2.14	0.47
1:A:222:GLN:NE2	3:A:1495:CL:CL	2.84	0.47
1:B:100:LYS:HD3	1:B:133:ASN:HD21	1.78	0.47
1:A:125[B]:ARG:NH1	6:A:2192:HOH:O	2.47	0.47
1:B:187:ILE:HB	1:B:221:ILE:HD13	1.97	0.47
1:B:231:PHE:O	1:B:264:ALA:HA	2.15	0.46
1:B:408[B]:CYS:SG	1:B:475:VAL:C	2.94	0.46
1:A:372:PRO:CG	1:A:444:TYR:HB3	2.46	0.46
1:B:222:GLN:NE2	3:B:1495:CL:CL	2.86	0.46
1:A:202:ARG:HA	1:A:227:ASP:O	2.16	0.46
1:A:269[B]:CYS:HB3	1:A:275:ALA:CA	2.45	0.46
1:B:174:ASN:HA	1:B:208:ARG:O	2.15	0.46
1:B:202:ARG:HA	1:B:227:ASP:O	2.17	0.45
1:B:365:LYS:O	1:B:369:GLY:HA2	2.17	0.45
1:A:208[B]:ARG:HG3	5:A:1499:GOL:O3	2.16	0.45
1:A:416:LYS:HE2	6:A:2437:HOH:O	2.16	0.45
1:A:408[B]:CYS:SG	1:A:476:CYS:N	2.89	0.45
1:A:269[B]:CYS:HB3	1:A:275:ALA:HA	1.98	0.45
1:A:135:VAL:O	1:A:167:VAL:HA	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:245:GLU:HA	1:B:279:GLY:O	2.17	0.44
1:A:365:LYS:O	1:A:369:GLY:HA2	2.17	0.44
1:B:135:VAL:O	1:B:167:VAL:HA	2.17	0.44
1:A:125[B]:ARG:NH1	6:A:2191:HOH:O	2.51	0.44
1:A:245:GLU:HA	1:A:279:GLY:O	2.18	0.44
1:A:187:ILE:HB	1:A:221:ILE:HD13	1.99	0.44
1:B:297:SER:HB2	1:B:388:GLY:HA3	1.99	0.43
1:B:261:ASN:HB3	6:B:2352:HOH:O	2.19	0.43
1:B:106:ARG:NH2	1:B:141[B]:GLN:OE1	2.52	0.43
1:B:300:SER:HA	1:B:432:ALA:O	2.19	0.43
1:B:297:SER:CB	1:B:388:GLY:HA3	2.48	0.43
1:A:138:PHE:O	1:A:170:TYR:HA	2.19	0.43
1:B:309:VAL:HG13	1:B:370:ILE:CD1	2.50	0.42
1:A:240:ILE:HD13	1:A:277[B]:MET:SD	2.59	0.42
1:B:168:ARG:HA	1:B:202:ARG:O	2.19	0.42
1:B:274:ALA:HA	1:B:300:SER:O	2.20	0.42
1:A:247:ASP:HB3	1:A:281:HIS:CG	2.54	0.42
1:A:304[A]:SER:OG	1:A:448:VAL:HG21	2.20	0.42
1:B:309:VAL:HG13	1:B:370:ILE:HD12	2.02	0.42
1:B:45:GLN:HB3	1:B:83:GLY:HA2	2.02	0.42
1:A:254:TYR:O	1:A:256:GLN:HG3	2.20	0.41
1:A:438:THR:HB	1:A:445:ASP:OD1	2.20	0.41
1:B:416:LYS:H	1:B:416:LYS:HE2	1.85	0.41
1:B:452:ASN:OD1	1:B:454:LYS:HE3	2.20	0.41
1:A:269[B]:CYS:SG	1:A:295:SER:HB3	2.60	0.41
1:A:95:LEU:HD22	1:A:116:THR:O	2.20	0.41
1:A:168:ARG:HA	1:A:202:ARG:O	2.20	0.41
1:B:173:SER:HA	1:B:207:GLU:O	2.20	0.41
1:B:138:PHE:O	1:B:170:TYR:HA	2.20	0.41
1:A:309:VAL:CG1	1:A:370:ILE:CG1	2.97	0.41
1:A:312:PHE:CD1	1:A:312:PHE:N	2.89	0.41
1:A:274:ALA:HA	1:A:300:SER:O	2.21	0.40
1:A:416:LYS:HE2	1:A:416:LYS:H	1.84	0.40
1:B:368:TYR:HB2	1:B:370:ILE:HG12	2.02	0.40
1:A:297:SER:CB	1:A:388:GLY:HA3	2.52	0.40
1:A:37:ASP:HB3	5:A:1500:GOL:H12	2.03	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	438/464 (94%)	417 (95%)	21 (5%)	0	100	100
1	B	433/464 (93%)	412 (95%)	21 (5%)	0	100	100
All	All	871/928 (94%)	829 (95%)	42 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	385/399 (96%)	380 (99%)	5 (1%)	69	50
1	B	381/399 (96%)	377 (99%)	4 (1%)	76	61
All	All	766/798 (96%)	757 (99%)	9 (1%)	78	54

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	125[A]	ARG
1	A	125[B]	ARG
1	A	190	ASP
1	A	309	VAL
1	A	366	LEU
1	B	125	ARG
1	B	190	ASP
1	B	277[A]	MET

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Mol	Chain	Res	Type
1	B	277[B]	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
1	A	222	GLN
1	B	60	ASN
1	B	195	ASN
1	B	222	GLN
1	B	261	ASN
1	B	422	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](#)

Of 20 ligands modelled in this entry, 14 are monoatomic - leaving 6 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	GOL	B	1501	-	5,5,5	0.80	0	5,5,5	0.24	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	GOL	A	1500	-	5,5,5	0.94	0	5,5,5	0.29	0
5	GOL	B	1500	-	5,5,5	0.94	0	5,5,5	0.28	0
5	GOL	A	1501	-	5,5,5	0.96	0	5,5,5	0.33	0
5	GOL	B	1499	-	5,5,5	0.92	0	5,5,5	0.34	0
5	GOL	A	1499	-	5,5,5	0.90	0	5,5,5	0.33	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	GOL	B	1501	-	-	1/4/4/4	-
5	GOL	A	1500	-	-	2/4/4/4	-
5	GOL	B	1500	-	-	2/4/4/4	-
5	GOL	A	1501	-	-	4/4/4/4	-
5	GOL	B	1499	-	-	4/4/4/4	-
5	GOL	A	1499	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1500	GOL	C1-C2-C3-O3
5	A	1500	GOL	C1-C2-C3-O3
5	A	1501	GOL	O1-C1-C2-C3
5	A	1501	GOL	C1-C2-C3-O3
5	B	1499	GOL	O1-C1-C2-C3
5	B	1499	GOL	C1-C2-C3-O3
5	A	1499	GOL	O1-C1-C2-C3
5	A	1499	GOL	C1-C2-C3-O3
5	A	1501	GOL	O1-C1-C2-O2
5	B	1500	GOL	O2-C2-C3-O3
5	A	1500	GOL	O2-C2-C3-O3
5	A	1501	GOL	O2-C2-C3-O3
5	B	1499	GOL	O1-C1-C2-O2
5	B	1499	GOL	O2-C2-C3-O3
5	A	1499	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
5	A	1499	GOL	O2-C2-C3-O3
5	B	1501	GOL	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1501	GOL	3	0
5	A	1500	GOL	1	0
5	A	1501	GOL	1	0
5	A	1499	GOL	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	433/464 (93%)	0.35	31 (7%) 15 14	10, 18, 40, 48	19 (4%)
1	B	430/464 (92%)	0.40	35 (8%) 12 10	10, 16, 41, 49	36 (8%)
All	All	863/928 (92%)	0.37	66 (7%) 13 12	10, 17, 41, 49	55 (6%)

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	312	PHE	12.8
1	B	335	GLY	10.6
1	A	359	ALA	8.4
1	B	366	LEU	7.9
1	B	441	THR	7.8
1	B	359	ALA	7.3
1	A	312	PHE	6.7
1	B	370	ILE	6.6
1	B	336	CYS	6.4
1	A	366	LEU	6.3
1	B	444	TYR	5.9
1	B	28	VAL	5.8
1	B	369	GLY	5.3
1	B	365	LYS	5.2
1	A	368	TYR	4.9
1	A	313	SER	4.9
1	A	336	CYS	4.9
1	B	368	TYR	4.9
1	A	441	THR	4.6
1	B	352	ALA	4.4
1	A	360	CYS	4.3
1	A	412	CYS	4.3
1	A	369	GLY	4.2
1	B	339	THR	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	335	GLY	4.1
1	A	351	ALA	4.0
1	A	28	VAL	3.9
1	B	484[A]	CYS	3.8
1	A	370	ILE	3.7
1	B	361	LEU	3.6
1	B	338	GLN	3.5
1	B	29	SER	3.5
1	B	337	ALA	3.5
1	B	269[A]	CYS	3.5
1	A	362	ASP	3.4
1	B	362	ASP	3.4
1	B	412	CYS	3.4
1	B	358	ASP	3.3
1	B	310	GLU	3.1
1	B	364	ALA	3.1
1	A	365	LYS	3.0
1	B	407	MET	3.0
1	B	357	LYS	2.9
1	A	310	GLU	2.8
1	A	444	TYR	2.8
1	B	119	GLY	2.8
1	A	29	SER	2.7
1	B	367	GLU	2.7
1	B	354	VAL	2.6
1	A	407	MET	2.6
1	B	62	THR	2.5
1	B	360	CYS	2.5
1	B	439	PRO	2.4
1	A	358	ASP	2.4
1	A	153	ASP	2.4
1	A	62	THR	2.4
1	A	122	LYS	2.4
1	A	338	GLN	2.4
1	A	364	ALA	2.4
1	A	339	THR	2.4
1	B	195	ASN	2.3
1	B	363	LYS	2.2
1	A	269[A]	CYS	2.2
1	A	439	PRO	2.1
1	A	371	GLU	2.1
1	A	309	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(\AA^2)	Q<0.9
5	GOL	A	1501	6/6	0.36	0.31	51,53,53,53	0
5	GOL	A	1500	6/6	0.51	0.30	51,53,53,54	0
5	GOL	B	1500	6/6	0.61	0.29	51,53,53,54	0
5	GOL	B	1501	6/6	0.62	0.32	51,53,53,53	0
5	GOL	B	1499	6/6	0.76	0.24	52,53,53,53	0
5	GOL	A	1499	6/6	0.76	0.26	52,53,53,53	0
4	NA	A	1498	1/1	0.92	0.12	27,27,27,27	0
2	CA	A	1494	1/1	0.95	0.13	27,27,27,27	0
2	CA	B	1494	1/1	0.96	0.08	26,26,26,26	0
4	NA	B	1498	1/1	0.96	0.18	25,25,25,25	0
4	NA	A	1496	1/1	0.97	0.10	15,15,15,15	0
3	CL	B	1495	1/1	0.98	0.08	27,27,27,27	0
2	CA	B	1493	1/1	0.98	0.04	15,15,15,15	0
4	NA	A	1497	1/1	0.99	0.05	16,16,16,16	0
2	CA	A	1493	1/1	0.99	0.04	16,16,16,16	0
2	CA	A	1492	1/1	0.99	0.04	15,15,15,15	0
4	NA	B	1496	1/1	0.99	0.04	15,15,15,15	0
2	CA	B	1492	1/1	0.99	0.04	15,15,15,15	0
3	CL	A	1495	1/1	0.99	0.13	28,28,28,28	0
4	NA	B	1497	1/1	0.99	0.06	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.