



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 1OGB / pdb_00001ogb
Title : Chitinase b from Serratia marcescens mutant D142N
Authors : Vaaje-Kolstad, G.; Houston, D.R.; Rao, F.V.; Peter, M.G.; Synstad, B.; van Aalten, D.M.F.; Eijsink, V.G.H.
Deposited on : 2003-04-29
Resolution : 1.85 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

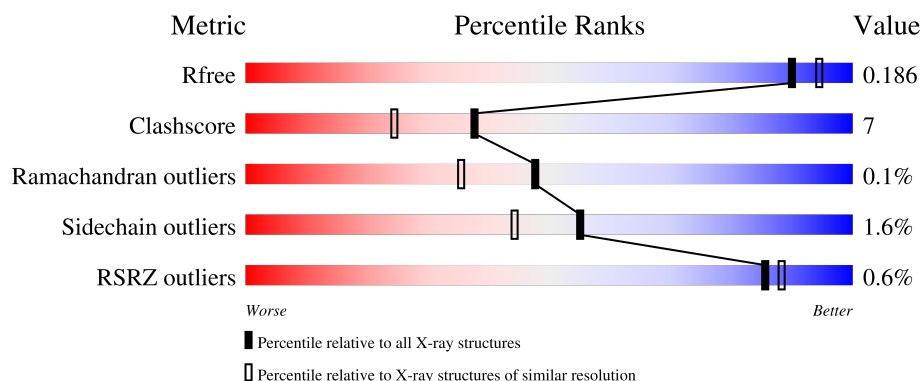
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	499	% 88% 12%
1	B	499	% 87% 12%

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	A	1505	-	-	X	-
2	GOL	A	1506	-	-	X	-
2	GOL	B	1503	-	-	X	-
2	GOL	B	1506	-	-	X	-
2	GOL	B	1507	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9344 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.

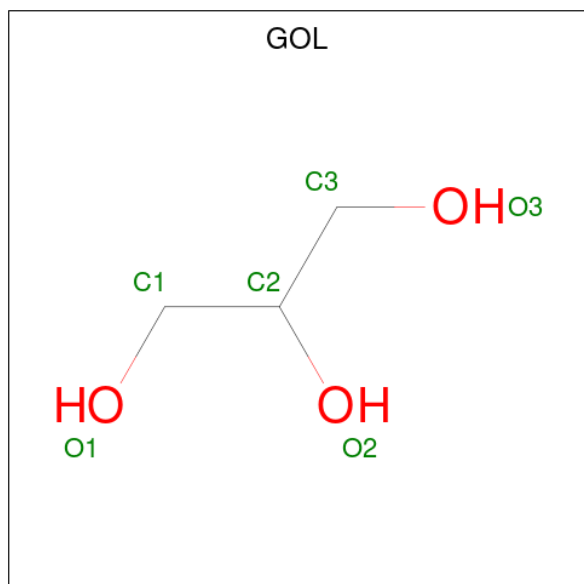
- Molecule 1 is a protein called CHITINASE B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	498	Total	C	N	O	S	12	4	1
			3922	2506	662	740	14			
1	B	498	Total	C	N	O	S	5	9	0
			3962	2530	675	743	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	142	ASN	ASP	engineered mutation	UNP P11797
B	142	ASN	ASP	engineered mutation	UNP P11797

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

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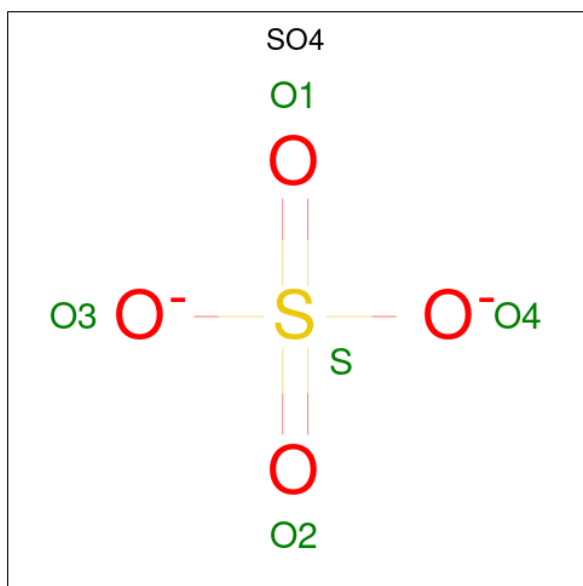
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total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

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

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	593	Total	O	0	0
			593	593		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	694	Total 694	O 694	0	0

i

- Molecule 1: CHITINASE B



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.95Å 103.98Å 186.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.97 – 1.85 14.97 – 1.85	Depositor EDS
% Data completeness (in resolution range)	96.7 (14.97-1.85) 96.5 (14.97-1.85)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.44 (at 1.84Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.159 , 0.193 0.153 , 0.186	Depositor DCC
R_{free} test set	1323 reflections (1.46%)	wwPDB-VP
Wilson B-factor (Å ²)	18.5	Xtriage
Anisotropy	0.260	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 58.3	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	0.96	EDS
Total number of atoms	9344	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% 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: GOL, 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	0.81	1/4049 (0.0%)	1.04	13/5519 (0.2%)
1	B	0.82	1/4109 (0.0%)	1.05	16/5594 (0.3%)
All	All	0.81	2/8158 (0.0%)	1.05	29/11113 (0.3%)

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	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	131	MET	SD-CE	6.81	1.96	1.79
1	B	228	ALA	CA-CB	5.02	1.58	1.53

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	115	LYS	N-CA-C	10.19	121.97	111.07
1	A	485	ALA	N-CA-C	9.69	123.34	110.36
1	A	115	LYS	N-CA-C	8.87	120.56	111.07
1	B	309	SER	N-CA-C	-7.57	100.54	110.53
1	A	45	THR	N-CA-C	-7.29	104.92	113.88
1	A	100	SER	N-CA-C	7.26	123.03	113.30
1	A	63	ASP	N-CA-C	-6.93	101.05	109.57
1	B	319	PRO	N-CA-C	6.67	122.24	114.03
1	B	146	PRO	CA-C-N	-6.65	111.12	123.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	146	PRO	C-N-CA	-6.65	111.12	123.03
1	B	63	ASP	N-CA-C	-6.65	99.91	109.48
1	B	100	SER	N-CA-C	6.64	122.20	113.30
1	B	423	ASN	N-CA-C	6.22	122.33	114.31
1	A	309	SER	N-CA-C	-5.86	102.64	110.55
1	A	151	VAL	N-CA-C	5.83	116.56	110.62
1	B	45	THR	N-CA-C	-5.80	106.75	113.88
1	A	423	ASN	N-CA-C	5.76	122.13	114.12
1	B	189	ALA	N-CA-C	5.72	118.29	111.71
1	A	413	ASP	N-CA-C	5.67	117.47	111.28
1	A	324	TRP	N-CA-C	5.50	119.98	113.16
1	B	184	ALA	N-CA-C	-5.40	100.44	109.24
1	B	218	GLY	CA-C-N	5.39	125.06	119.56
1	B	218	GLY	C-N-CA	5.39	125.06	119.56
1	A	184	ALA	N-CA-C	-5.29	100.56	108.96
1	A	483	THR	N-CA-C	5.09	121.25	113.61
1	B	324	TRP	N-CA-C	5.08	119.46	113.16
1	B	318	TYR	CA-C-N	5.05	125.12	119.87
1	B	318	TYR	C-N-CA	5.05	125.12	119.87
1	A	107	HIS	N-CA-C	5.02	117.41	111.33

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	110	TYR	Sidechain

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	3922	0	3749	43	0
1	B	3962	0	3798	63	0
2	A	66	0	88	17	0
2	B	72	0	96	25	0
3	A	25	0	0	0	0
3	B	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	593	0	0	5	0
4	B	694	0	0	5	0
All	All	9344	0	7731	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 7.

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:B:444:GLY:N	1:B:447:ASN:HD21	1.50	1.08
1:A:89:ARG:HH22	2:A:1505:GOL:H31	1.21	1.04
1:B:444:GLY:H	1:B:447:ASN:ND2	1.62	0.97
1:B:327:GLY:H	2:B:1501:GOL:H11	1.27	0.94
1:B:445:PRO:HA	1:B:448:LEU:HD21	1.54	0.88
1:A:89:ARG:NH2	2:A:1505:GOL:H31	1.91	0.86
2:A:1505:GOL:H12	4:A:2149:HOH:O	1.78	0.83
1:B:444:GLY:H	1:B:447:ASN:HD21	0.83	0.81
1:B:370:HIS:HD2	1:B:373:ASN:H	1.25	0.79
2:A:1507:GOL:H32	2:A:1509:GOL:O1	1.83	0.78
1:B:31:PHE:HB3	2:B:1507:GOL:H11	1.70	0.72
1:A:479:TRP:HE1	2:B:1506:GOL:H2	1.56	0.70
1:B:241:ASN:O	1:B:244[B]:ARG:HG3	1.90	0.69
1:B:444:GLY:CA	1:B:447:ASN:HD21	2.07	0.68
1:B:390:LYS:HZ3	2:B:1509:GOL:H12	1.57	0.68
2:A:1502:GOL:H32	4:A:2212:HOH:O	1.93	0.67
1:B:221:GLU:OE2	2:B:1503:GOL:H31	1.95	0.67
1:B:359:TRP:CH2	1:B:383[A]:GLU:HG2	2.31	0.66
1:A:89:ARG:HH22	2:A:1505:GOL:C3	2.05	0.66
1:A:261:SER:N	2:A:1506:GOL:H32	2.11	0.66
1:B:263:PHE:HB2	2:B:1500:GOL:H31	1.76	0.65
1:B:444:GLY:N	1:B:447:ASN:ND2	2.33	0.64
1:B:445:PRO:CA	1:B:448:LEU:HD21	2.29	0.62
1:B:31:PHE:O	2:B:1507:GOL:H11	2.01	0.61
1:B:404:HIS:HE1	2:B:1507:GOL:H32	1.67	0.59
1:A:89:ARG:HH12	2:A:1505:GOL:H31	1.68	0.59
1:B:327:GLY:N	2:B:1501:GOL:H11	2.09	0.59
1:A:359:TRP:CH2	1:A:383[A]:GLU:HG2	2.38	0.58
1:A:261:SER:O	2:A:1506:GOL:H11	2.04	0.58
1:B:448:LEU:HD22	1:B:448:LEU:N	2.19	0.57
1:B:478:LYS:HE3	1:B:490:SER:O	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:390:LYS:NZ	2:B:1509:GOL:H12	2.19	0.56
1:B:372:GLN:NE2	4:B:2519:HOH:O	2.37	0.56
1:A:225:ASN:ND2	1:A:226:HIS:H	2.03	0.55
1:A:89:ARG:HH12	2:A:1505:GOL:C3	2.19	0.55
1:B:353:TYR:O	1:B:370:HIS:HE1	1.89	0.55
1:B:31:PHE:CG	1:B:406:GLY:HA2	2.40	0.55
1:A:479:TRP:HE1	2:B:1506:GOL:C2	2.20	0.55
1:A:141:ILE:C	1:A:142[B]:ASN:HD22	2.15	0.54
1:A:484:SER:HB2	1:B:147:GLN:NE2	2.23	0.54
1:B:383[B]:GLU:OE2	2:B:1510:GOL:O3	2.25	0.54
2:A:1507:GOL:H12	2:A:1509:GOL:O2	2.08	0.54
1:B:327:GLY:H	2:B:1501:GOL:C1	2.09	0.54
1:B:370:HIS:CD2	1:B:373:ASN:H	2.16	0.53
1:A:89:ARG:CZ	2:A:1505:GOL:H31	2.39	0.53
1:A:89:ARG:NH1	2:A:1505:GOL:H31	2.23	0.53
1:A:278:GLU:HG3	1:A:445:PRO:HB2	1.90	0.52
1:A:475:TRP:CD2	1:A:486:PRO:HB3	2.44	0.52
1:B:31:PHE:CB	2:B:1507:GOL:H11	2.39	0.51
1:B:10:TYR:HD2	1:B:50[A]:SER:OG	1.93	0.51
1:A:479:TRP:NE1	2:B:1506:GOL:H2	2.24	0.51
1:B:129[A]:ARG:NH2	4:B:2250:HOH:O	2.43	0.50
1:B:444:GLY:CA	1:B:447:ASN:ND2	2.72	0.50
1:A:480:GLY:CA	2:B:1503:GOL:H12	2.42	0.49
1:A:140:ASP:OD2	1:A:142[B]:ASN:ND2	2.44	0.49
1:A:162:ARG:CZ	2:A:1503:GOL:H32	2.42	0.49
1:A:472:GLY:C	1:A:498:VAL:HG23	2.36	0.49
1:B:63:ASP:OD1	1:B:64:PRO:HD2	2.13	0.49
1:B:244[A]:ARG:HG3	1:B:259:PHE:HB2	1.95	0.49
1:B:445:PRO:C	1:B:448:LEU:HD21	2.38	0.48
1:B:447:ASN:HD22	1:B:447:ASN:N	2.12	0.48
1:B:148:ALA:HB3	4:B:2283:HOH:O	2.13	0.48
1:A:261:SER:H	2:A:1506:GOL:H32	1.80	0.47
1:A:484:SER:CB	1:B:147:GLN:NE2	2.78	0.47
1:A:484:SER:HB2	1:B:147:GLN:HE22	1.80	0.46
1:A:472:GLY:O	1:A:498:VAL:HG23	2.16	0.46
1:B:444:GLY:O	1:B:447:ASN:ND2	2.48	0.46
1:A:480:GLY:H	2:B:1503:GOL:C1	2.27	0.45
1:B:277:MET:HE1	2:B:1508:GOL:O1	2.16	0.45
1:A:482:ILE:HD12	1:A:491:ALA:CB	2.47	0.45
1:A:466:ALA:O	1:A:476:GLN:HA	2.17	0.45
1:A:467:LEU:HA	1:A:475:TRP:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:448:LEU:HB3	1:B:496:GLY:HA2	1.98	0.45
2:A:1505:GOL:H32	4:A:2576:HOH:O	2.17	0.44
1:B:373:ASN:HD22	1:B:373:ASN:N	2.14	0.44
1:A:260:PRO:HA	2:A:1506:GOL:H31	2.00	0.44
1:A:330:GLU:OE1	4:A:2401:HOH:O	2.21	0.44
1:B:482:ILE:HD12	1:B:491:ALA:HB1	1.99	0.44
1:B:31:PHE:C	2:B:1507:GOL:H11	2.43	0.43
1:B:370:HIS:HB3	1:B:375:LEU:HB2	2.00	0.43
1:B:144:GLU:HA	1:B:145:TYR:CG	2.53	0.43
1:A:144:GLU:HA	1:A:145:TYR:CG	2.54	0.43
1:A:132:LYS:NZ	4:A:2212:HOH:O	2.52	0.43
1:B:31:PHE:CA	2:B:1507:GOL:H11	2.49	0.43
1:B:448:LEU:HD22	1:B:448:LEU:H	1.82	0.43
1:B:353:TYR:O	1:B:370:HIS:CE1	2.71	0.43
1:B:445:PRO:C	1:B:448:LEU:CD2	2.92	0.43
1:B:482:ILE:HD12	1:B:491:ALA:CB	2.49	0.43
1:A:342:TYR:CG	1:A:414:LEU:HD11	2.54	0.42
1:A:50:SER:HA	1:A:51:PHE:HA	1.78	0.42
1:A:480:GLY:N	2:B:1503:GOL:H12	2.35	0.42
2:B:1507:GOL:C3	4:B:2685:HOH:O	2.67	0.42
1:A:281:PRO:HG2	1:A:284:LYS:HG3	2.02	0.42
1:B:244[B]:ARG:HG2	1:B:259:PHE:HB2	2.01	0.42
1:A:480:GLY:H	2:B:1503:GOL:H12	1.84	0.41
1:B:97:TRP:CB	2:B:1511:GOL:H31	2.50	0.41
1:B:244[A]:ARG:HG3	1:B:259:PHE:CB	2.51	0.41
1:B:147:GLN:O	1:B:148:ALA:C	2.64	0.41
1:A:31:PHE:CG	1:A:406:GLY:HA2	2.56	0.41
1:A:347:GLN:HE21	1:A:347:GLN:HB2	1.72	0.41
1:B:213:THR:HB	1:B:267:VAL:HG22	2.03	0.41
1:B:372:GLN:HG2	1:B:373:ASN:HD22	1.86	0.41
1:B:224:THR:HG22	1:B:296:PHE:CD2	2.56	0.40
1:B:473:TYR:CD2	1:B:494:LYS:HD3	2.56	0.40
2:B:1506:GOL:H31	4:B:2682:HOH:O	2.20	0.40
1:A:477:THR:HG23	1:A:492:TRP:CZ3	2.57	0.40
1:B:13:ILE:HA	1:B:14:PRO:HD3	1.92	0.40
1:B:342:TYR:CG	1:B:414:LEU:HD11	2.57	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	500/499 (100%)	489 (98%)	11 (2%)	0	100	100
1	B	505/499 (101%)	497 (98%)	7 (1%)	1 (0%)	43	31
All	All	1005/998 (101%)	986 (98%)	18 (2%)	1 (0%)	48	35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	148	ALA

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	408/405 (101%)	404 (99%)	4 (1%)	68	60
1	B	413/405 (102%)	404 (98%)	9 (2%)	45	32
All	All	821/810 (101%)	808 (98%)	13 (2%)	55	44

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	213	THR
1	A	349	LEU
1	A	378	THR
1	A	380	ASP
1	B	3	THR

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Mol	Chain	Res	Type
1	B	213	THR
1	B	299	VAL
1	B	319	PRO
1	B	373	ASN
1	B	375	LEU
1	B	378	THR
1	B	380	ASP
1	B	447	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	ASN
1	A	147	GLN
1	A	159	GLN
1	A	180	GLN
1	A	225	ASN
1	A	347	GLN
1	A	350	GLN
1	A	372	GLN
1	A	471	GLN
1	B	17	GLN
1	B	35	ASN
1	B	112	ASN
1	B	142	ASN
1	B	147	GLN
1	B	168	GLN
1	B	180	GLN
1	B	272	GLN
1	B	370	HIS
1	B	372	GLN
1	B	373	ASN
1	B	395	GLN
1	B	407	GLN
1	B	411	ASN
1	B	431	GLN
1	B	447	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

30 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	GOL	A	1508	-	5,5,5	0.38	0	5,5,5	0.34	0
2	GOL	B	1508	-	5,5,5	0.41	0	5,5,5	0.30	0
2	GOL	A	1501	-	5,5,5	0.55	0	5,5,5	0.42	0
2	GOL	B	1502	-	5,5,5	0.42	0	5,5,5	0.19	0
2	GOL	B	1506	-	5,5,5	0.16	0	5,5,5	0.27	0
2	GOL	B	1510	-	5,5,5	0.30	0	5,5,5	0.34	0
2	GOL	A	1509	-	5,5,5	0.43	0	5,5,5	0.14	0
3	SO4	A	1512	-	4,4,4	0.45	0	6,6,6	0.14	0
2	GOL	A	1502	-	5,5,5	0.32	0	5,5,5	0.25	0
2	GOL	A	1499	-	5,5,5	0.38	0	5,5,5	0.41	0
2	GOL	A	1506	-	5,5,5	0.47	0	5,5,5	0.43	0
2	GOL	B	1511	-	5,5,5	0.41	0	5,5,5	0.32	0
3	SO4	A	1510	-	4,4,4	0.39	0	6,6,6	0.13	0
2	GOL	A	1500	-	5,5,5	0.51	0	5,5,5	0.24	0
3	SO4	B	1513	-	4,4,4	0.39	0	6,6,6	0.18	0
2	GOL	B	1507	-	5,5,5	0.41	0	5,5,5	0.20	0
2	GOL	B	1504	-	5,5,5	0.45	0	5,5,5	0.21	0
3	SO4	B	1512	-	4,4,4	0.45	0	6,6,6	0.09	0
2	GOL	A	1503	-	5,5,5	0.46	0	5,5,5	0.17	0
2	GOL	B	1505	-	5,5,5	0.46	0	5,5,5	0.54	0
3	SO4	A	1511	-	4,4,4	0.45	0	6,6,6	0.21	0
2	GOL	A	1504	-	5,5,5	0.41	0	5,5,5	0.36	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GOL	A	1507	-	5,5,5	0.41	0	5,5,5	0.08	0
3	SO4	A	1514	-	4,4,4	0.41	0	6,6,6	0.12	0
2	GOL	B	1500	-	5,5,5	0.42	0	5,5,5	0.29	0
2	GOL	A	1505	-	5,5,5	0.39	0	5,5,5	0.21	0
2	GOL	B	1501	-	5,5,5	0.84	0	5,5,5	0.36	0
3	SO4	A	1513	-	4,4,4	0.37	0	6,6,6	0.22	0
2	GOL	B	1503	-	5,5,5	0.54	0	5,5,5	0.32	0
2	GOL	B	1509	-	5,5,5	0.32	0	5,5,5	0.27	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
2	GOL	A	1508	-	-	0/4/4/4	-
2	GOL	B	1508	-	-	0/4/4/4	-
2	GOL	A	1501	-	-	0/4/4/4	-
2	GOL	B	1502	-	-	0/4/4/4	-
2	GOL	B	1506	-	-	0/4/4/4	-
2	GOL	B	1510	-	-	0/4/4/4	-
2	GOL	A	1509	-	-	0/4/4/4	-
2	GOL	A	1502	-	-	0/4/4/4	-
2	GOL	A	1499	-	-	0/4/4/4	-
2	GOL	A	1506	-	-	0/4/4/4	-
2	GOL	B	1511	-	-	0/4/4/4	-
2	GOL	A	1500	-	-	0/4/4/4	-
2	GOL	B	1507	-	-	0/4/4/4	-
2	GOL	B	1504	-	-	0/4/4/4	-
2	GOL	A	1503	-	-	0/4/4/4	-
2	GOL	B	1505	-	-	0/4/4/4	-
2	GOL	A	1504	-	-	0/4/4/4	-
2	GOL	A	1507	-	-	0/4/4/4	-
2	GOL	B	1500	-	-	0/4/4/4	-
2	GOL	A	1505	-	-	0/4/4/4	-
2	GOL	B	1501	-	-	0/4/4/4	-
2	GOL	B	1503	-	-	0/4/4/4	-
2	GOL	B	1509	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

15 monomers are involved in 42 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1508	GOL	1	0
2	B	1506	GOL	4	0
2	B	1510	GOL	1	0
2	A	1509	GOL	2	0
2	A	1502	GOL	1	0
2	A	1506	GOL	4	0
2	B	1511	GOL	1	0
2	B	1507	GOL	7	0
2	A	1503	GOL	1	0
2	A	1507	GOL	2	0
2	B	1500	GOL	1	0
2	A	1505	GOL	9	0
2	B	1501	GOL	3	0
2	B	1503	GOL	5	0
2	B	1509	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	498/499 (99%)	-0.57	3 (0%) 85 88	9, 18, 31, 45	8 (1%)
1	B	498/499 (99%)	-0.75	3 (0%) 85 88	8, 15, 27, 40	11 (2%)
All	All	996/998 (99%)	-0.66	6 (0%) 85 88	8, 17, 30, 45	19 (1%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	470	TYR	2.9
1	B	488	SER	2.9
1	B	445	PRO	2.6
1	A	2	SER	2.6
1	A	456	TYR	2.5
1	B	446	GLY	2.2

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GOL	A	1509	6/6	0.64	0.15	53,53,54,56	0
2	GOL	A	1508	6/6	0.66	0.15	51,52,52,53	0
2	GOL	B	1508	6/6	0.68	0.15	43,48,49,49	0
2	GOL	A	1507	6/6	0.69	0.15	49,51,51,51	0
3	SO4	A	1511	5/5	0.77	0.16	66,67,68,69	0
2	GOL	B	1503	6/6	0.78	0.15	26,36,40,45	0
3	SO4	A	1510	5/5	0.80	0.14	81,81,82,82	0
2	GOL	B	1501	6/6	0.80	0.14	19,22,24,37	0
2	GOL	A	1505	6/6	0.82	0.13	52,54,55,55	0
2	GOL	B	1510	6/6	0.82	0.14	46,49,51,53	0
2	GOL	A	1501	6/6	0.82	0.15	38,39,40,41	0
2	GOL	B	1505	6/6	0.82	0.11	35,37,39,40	0
3	SO4	B	1513	5/5	0.82	0.12	70,70,71,71	0
2	GOL	B	1511	6/6	0.83	0.13	46,49,49,54	0
2	GOL	A	1506	6/6	0.83	0.18	30,40,41,45	0
3	SO4	B	1512	5/5	0.84	0.20	56,57,58,59	0
2	GOL	B	1509	6/6	0.84	0.12	47,50,51,53	0
2	GOL	B	1500	6/6	0.86	0.17	17,34,35,38	0
2	GOL	A	1502	6/6	0.86	0.14	44,48,49,49	0
2	GOL	B	1506	6/6	0.87	0.13	41,42,44,44	0
3	SO4	A	1514	5/5	0.87	0.10	67,68,68,69	0
2	GOL	A	1503	6/6	0.90	0.09	37,44,44,47	0
2	GOL	B	1504	6/6	0.91	0.08	26,32,35,35	0
2	GOL	B	1502	6/6	0.91	0.09	17,25,29,34	0
2	GOL	A	1504	6/6	0.92	0.08	34,36,38,41	0
2	GOL	A	1499	6/6	0.92	0.08	19,29,32,33	0
2	GOL	A	1500	6/6	0.93	0.07	26,31,33,33	0
2	GOL	B	1507	6/6	0.93	0.13	21,38,41,42	0
3	SO4	A	1512	5/5	0.93	0.12	62,62,62,63	0
3	SO4	A	1513	5/5	0.94	0.07	41,43,43,44	0

6.5 Other polymers ⓘ

There are no such residues in this entry.