Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 01:27 AM EDT

PDB ID	:	1PZ3
Title	:	Crystal structure of a family 51 (GH51) alpha-L-arabinofuranosidase from Geobacillus stearothermophilus T6
Authors	:	Hoevel, K.; Shallom, D.; Niefind, K.; Belakhov, V.; Shoham, G.; Baasov, T.; Shoham, Y.; Schomburg, D.
Deposited on	:	2003-07-09
Resolution	:	1.75 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 (i) 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 (i)) were used in the production of this report:

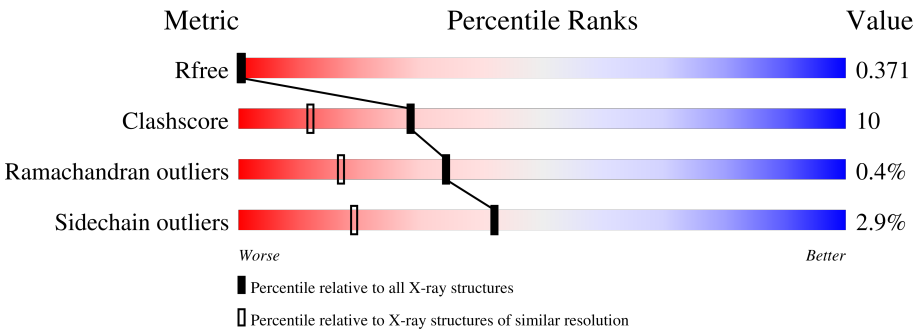
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)

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 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 <=5%

Mol	Chain	Length	Quality of chain
1	A	502	81% 16% • •
1	B	502	83% 15% • •

2 Entry composition (i)

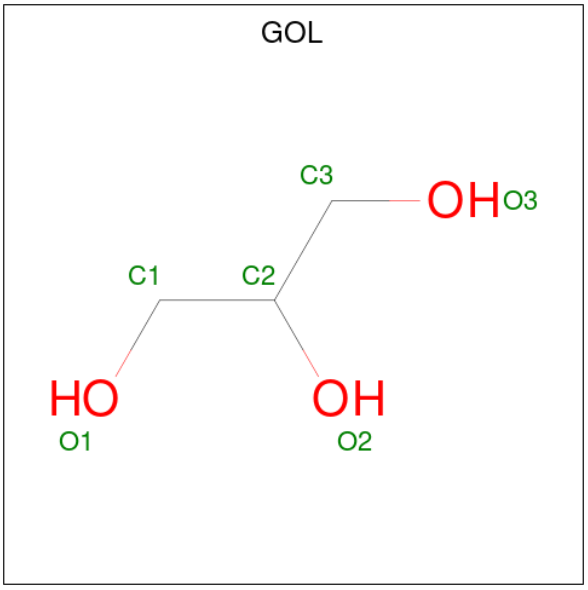
There are 3 unique types of molecules in this entry. The entry contains 9136 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Alpha-L-arabinofuranosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	497	3990	2542	680	748	20	0	0	0
1	B	497	3990	2542	680	748	20	0	0	0

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 3 is water.

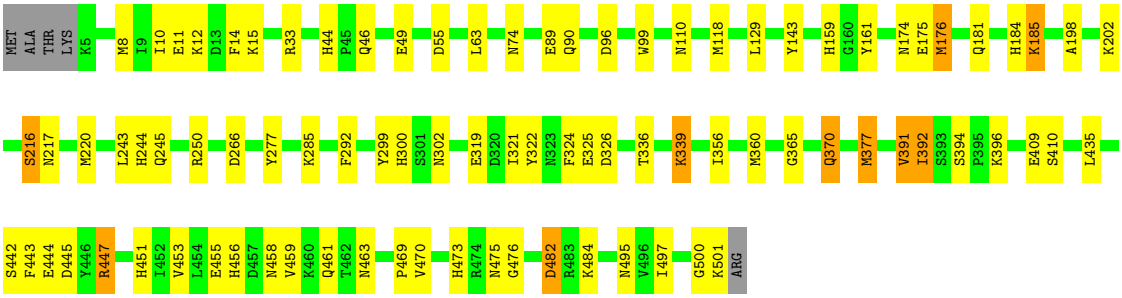
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	583	583	583	0	0
			Total	O		
3	B	561	561	561	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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. 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. 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.

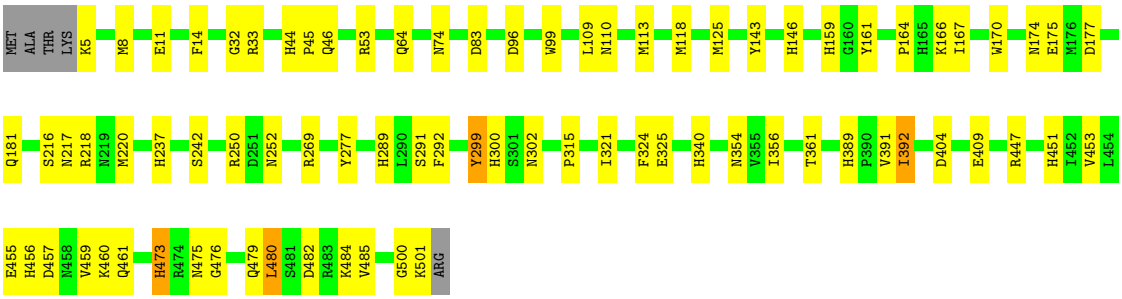
• Molecule 1: Alpha-L-arabinofuranosidase

Chain A:



• Molecule 1: Alpha-L-arabinofuranosidase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α, β, γ	179.43Å 179.43Å 100.23Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 1.75 19.96 – 1.75	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-1.75) 98.3 (19.96-1.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.10 (at 1.74Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.171 , 0.204 0.364 , 0.371	Depositor DCC
R_{free} test set	6001 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	21.5	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.47 , 66.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.010 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.84	EDS
Total number of atoms	9136	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.02% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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

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	0.66	0/4091	0.84	8/5558 (0.1%)
1	B	0.63	0/4091	0.83	7/5558 (0.1%)
All	All	0.64	0/8182	0.84	15/11116 (0.1%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	33	ARG	NE-CZ-NH1	8.11	124.36	120.30
1	B	33	ARG	NE-CZ-NH1	7.64	124.12	120.30
1	B	33	ARG	NE-CZ-NH2	-7.38	116.61	120.30
1	B	457	ASP	CB-CG-OD2	7.26	124.83	118.30
1	B	177	ASP	CB-CG-OD2	6.98	124.58	118.30
1	A	33	ARG	NE-CZ-NH2	-6.24	117.18	120.30
1	A	96	ASP	CB-CG-OD2	6.23	123.91	118.30
1	B	83	ASP	CB-CG-OD1	5.75	123.48	118.30
1	A	266	ASP	CB-CG-OD2	5.54	123.28	118.30
1	B	404	ASP	CB-CG-OD2	5.42	123.18	118.30
1	A	377	MET	CG-SD-CE	-5.15	91.96	100.20
1	A	185	LYS	CD-CE-NZ	-5.14	99.87	111.70
1	A	55	ASP	CB-CG-OD1	5.10	122.89	118.30
1	B	96	ASP	CB-CG-OD2	5.09	122.88	118.30
1	A	482	ASP	CB-CG-OD2	5.09	122.88	118.30

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	3990	0	3891	86	0
1	B	3990	0	3891	71	0
2	A	6	0	8	3	0
2	B	6	0	8	2	0
3	A	583	0	0	28	4
3	B	561	0	0	26	3
All	All	9136	0	7798	155	4

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 10.

All (155) 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:319:GLU:CD	3:A:1038:HOH:O	1.67	1.24
1:B:392:ILE:HG21	3:B:1027:HOH:O	1.52	1.09
1:B:118:MET:CG	3:B:1026:HOH:O	2.06	1.03
1:B:118:MET:SD	3:B:1026:HOH:O	2.15	1.01
1:B:118:MET:HG3	3:B:1026:HOH:O	1.59	1.00
1:A:244:HIS:HE1	3:A:1069:HOH:O	1.45	1.00
1:B:175:GLU:OE1	3:B:959:HOH:O	1.83	0.96
1:A:451:HIS:CD2	1:A:497:ILE:HG12	2.01	0.94
1:A:319:GLU:OE1	3:A:1038:HOH:O	1.75	0.93
1:B:354:ASN:HD21	1:B:361:THR:H	1.22	0.88
1:A:244:HIS:CE1	3:A:1069:HOH:O	2.21	0.86
1:A:220:MET:SD	3:A:997:HOH:O	2.35	0.85
1:A:451:HIS:HD2	1:A:497:ILE:HG12	1.40	0.83
1:A:14:PHE:CZ	1:B:391:VAL:HG21	2.15	0.80
1:A:44:HIS:HD2	1:A:46:GLN:H	1.29	0.80
1:A:451:HIS:ND1	1:A:476:GLY:HA3	1.96	0.79
1:B:175:GLU:HG3	1:B:216:SER:HB3	1.63	0.78
1:A:451:HIS:CE1	1:A:495:ASN:HD22	2.03	0.77
1:A:463:ASN:HD21	1:A:470:VAL:H	1.34	0.75
1:A:220:MET:HE1	3:A:1079:HOH:O	1.85	0.74
1:A:175:GLU:CG	1:A:216:SER:HB3	2.18	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:216:SER:OG	3:B:1017:HOH:O	2.06	0.73
1:B:480:LEU:HD22	1:B:485:VAL:HG22	1.70	0.73
1:A:220:MET:SD	3:A:1079:HOH:O	2.46	0.73
1:B:166:LYS:HG2	3:B:789:HOH:O	1.89	0.71
1:B:11:GLU:HG2	1:B:14:PHE:HD1	1.54	0.70
1:A:220:MET:CE	3:A:1079:HOH:O	2.37	0.70
1:A:250:ARG:HH21	1:A:302:ASN:HD21	1.36	0.70
1:A:143:TYR:OH	1:A:159:HIS:HD2	1.75	0.70
1:B:175:GLU:CG	1:B:216:SER:HB3	2.22	0.69
1:B:456:HIS:CE1	1:B:461:GLN:HG2	2.29	0.68
1:B:220:MET:SD	3:B:1017:HOH:O	2.52	0.68
1:B:300:HIS:HD2	1:B:321:ILE:O	1.77	0.67
1:A:175:GLU:OE1	3:A:793:HOH:O	2.12	0.67
2:B:503:GOL:H2	3:B:505:HOH:O	1.94	0.66
1:A:451:HIS:HE1	1:A:495:ASN:HD22	1.41	0.66
1:A:14:PHE:CE2	1:B:391:VAL:HG21	2.30	0.65
1:B:456:HIS:HE1	1:B:461:GLN:HG2	1.63	0.64
1:A:184:HIS:HE1	3:A:676:HOH:O	1.81	0.64
1:B:250:ARG:HH21	1:B:302:ASN:HD21	1.45	0.63
1:A:360:MET:CE	3:A:669:HOH:O	2.47	0.63
1:A:44:HIS:CD2	1:A:46:GLN:H	2.16	0.63
1:B:11:GLU:CG	1:B:14:PHE:HD1	2.12	0.63
2:A:503:GOL:H2	3:A:546:HOH:O	1.98	0.62
1:A:300:HIS:HE1	1:A:326:ASP:OD2	1.82	0.62
1:B:299:TYR:CE1	1:B:300:HIS:CE1	2.88	0.61
1:B:217:ASN:CB	3:B:1017:HOH:O	2.48	0.61
1:A:360:MET:HE2	3:A:669:HOH:O	2.00	0.60
1:B:389:HIS:HD2	3:B:803:HOH:O	1.83	0.59
1:B:44:HIS:ND1	1:B:45:PRO:HD2	2.18	0.59
1:A:391:VAL:HB	1:B:14:PHE:CZ	2.37	0.59
1:A:473:HIS:HD2	1:A:475:ASN:H	1.50	0.58
1:B:174:ASN:HD22	1:B:181:GLN:HE22	1.51	0.58
1:B:143:TYR:OH	1:B:159:HIS:HD2	1.87	0.57
1:B:8:MET:HG3	1:B:392:ILE:HG22	1.87	0.57
1:B:500:GLY:O	1:B:501:LYS:HB2	2.04	0.57
1:B:300:HIS:CD2	1:B:321:ILE:O	2.57	0.57
1:B:64:GLN:NE2	3:B:600:HOH:O	2.38	0.57
1:B:11:GLU:HG2	1:B:14:PHE:CD1	2.39	0.56
1:B:146:HIS:HD2	3:B:611:HOH:O	1.88	0.56
1:B:299:TYR:CZ	1:B:300:HIS:CE1	2.94	0.56
1:A:10:ILE:O	1:A:442:SER:OG	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:GLU:CD	1:A:216:SER:HB3	2.26	0.55
1:A:300:HIS:HD2	3:A:656:HOH:O	1.89	0.55
1:A:49:GLU:HG2	3:A:768:HOH:O	2.05	0.55
1:A:175:GLU:HG2	1:A:216:SER:HB3	1.87	0.55
1:B:217:ASN:HB3	3:B:1017:HOH:O	2.06	0.54
1:A:12:LYS:O	1:A:15:LYS:NZ	2.30	0.54
1:B:174:ASN:HD22	1:B:181:GLN:NE2	2.05	0.54
1:B:167:ILE:HB	1:B:170:TRP:CZ2	2.43	0.54
1:A:325:GLU:HB3	1:A:459:VAL:HB	1.90	0.54
1:B:480:LEU:HD22	1:B:485:VAL:CG2	2.36	0.53
1:A:216:SER:HB2	3:A:997:HOH:O	2.07	0.53
1:B:164:PRO:HG2	1:B:166:LYS:HE2	1.88	0.53
1:A:74:ASN:HA	1:A:181:GLN:HE22	1.73	0.53
1:A:174:ASN:HD22	1:A:181:GLN:NE2	2.07	0.53
1:A:129:LEU:O	1:A:185:LYS:HE3	2.09	0.52
1:A:410:SER:HB3	3:A:1086:HOH:O	2.09	0.52
1:A:174:ASN:HD22	1:A:181:GLN:HE22	1.56	0.52
1:B:74:ASN:HA	1:B:181:GLN:HE22	1.75	0.52
1:B:456:HIS:HD2	3:B:921:HOH:O	1.92	0.52
1:A:14:PHE:CE2	1:B:391:VAL:CG2	2.93	0.52
1:A:198:ALA:O	1:A:202:LYS:HE3	2.10	0.51
1:B:354:ASN:ND2	1:B:361:THR:H	2.00	0.51
1:A:99:TRP:CH2	2:A:503:GOL:H31	2.45	0.51
1:A:217:ASN:ND2	3:A:997:HOH:O	2.43	0.51
1:A:435:LEU:HD22	1:A:435:LEU:N	2.25	0.51
1:A:89:GLU:HG2	1:A:90:GLN:NE2	2.25	0.51
1:A:456:HIS:CE1	1:A:461:GLN:HG2	2.46	0.51
1:B:409:GLU:OE1	3:B:1025:HOH:O	2.19	0.50
1:A:11:GLU:CG	1:A:14:PHE:HD1	2.24	0.50
1:A:216:SER:CB	3:A:997:HOH:O	2.59	0.50
1:A:391:VAL:HG12	1:A:391:VAL:O	2.12	0.50
1:B:453:VAL:HB	1:B:473:HIS:CD2	2.47	0.50
1:A:159:HIS:HE1	3:A:574:HOH:O	1.94	0.50
1:B:389:HIS:CD2	3:B:803:HOH:O	2.61	0.49
1:B:44:HIS:CD2	1:B:53:ARG:CZ	2.95	0.49
1:A:322:TYR:H	1:A:370:GLN:HE22	1.61	0.49
1:A:220:MET:CG	3:A:997:HOH:O	2.57	0.49
1:A:322:TYR:H	1:A:370:GLN:NE2	2.11	0.49
1:A:455:GLU:OE2	1:A:473:HIS:HE1	1.96	0.48
1:B:269:ARG:HG3	1:B:340:HIS:HE1	1.78	0.48
1:B:218:ARG:O	3:B:1018:HOH:O	2.20	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:451:HIS:CE1	1:A:453:VAL:HG22	2.49	0.48
1:A:44:HIS:HE1	1:A:365:GLY:O	1.95	0.48
1:B:109:LEU:O	1:B:113:MET:HG2	2.13	0.48
1:B:125:MET:HB3	1:B:170:TRP:CE3	2.49	0.48
1:B:325:GLU:HB3	1:B:459:VAL:HB	1.96	0.48
1:A:447:ARG:HD2	3:A:1042:HOH:O	2.14	0.48
1:B:217:ASN:HB2	3:B:1017:HOH:O	2.13	0.48
1:A:482:ASP:OD1	3:A:955:HOH:O	2.20	0.47
1:A:110:ASN:HB3	1:A:161:TYR:CZ	2.50	0.47
1:B:32:GLY:HA2	1:B:315:PRO:O	2.15	0.47
1:B:125:MET:HG2	1:B:170:TRP:CZ3	2.50	0.47
1:A:324:PHE:CZ	1:A:455:GLU:HA	2.50	0.46
1:A:339:LYS:HE3	1:A:409:GLU:OE2	2.15	0.46
1:A:245:GLN:NE2	3:A:685:HOH:O	2.48	0.46
1:A:11:GLU:HG2	1:A:14:PHE:HD1	1.81	0.46
1:B:299:TYR:CE1	1:B:300:HIS:HE1	2.33	0.46
1:B:242:SER:HA	1:B:291:SER:O	2.17	0.45
1:A:336:THR:HA	1:A:339:LYS:HG2	1.98	0.45
1:B:453:VAL:HB	1:B:473:HIS:NE2	2.31	0.45
1:A:339:LYS:HD2	3:A:993:HOH:O	2.17	0.45
1:B:159:HIS:HE1	3:B:612:HOH:O	2.00	0.45
1:B:99:TRP:CH2	2:B:503:GOL:H31	2.52	0.44
1:B:252:ASN:HD21	1:B:460:LYS:HE3	1.81	0.43
1:A:463:ASN:ND2	1:A:470:VAL:H	2.10	0.43
1:A:456:HIS:HE1	1:A:469:PRO:HB2	1.83	0.43
1:A:11:GLU:HG2	1:A:14:PHE:CD1	2.54	0.43
1:A:484:LYS:HB2	1:A:484:LYS:NZ	2.33	0.43
1:A:321:ILE:HA	1:A:370:GLN:HE22	1.84	0.43
1:B:220:MET:CG	3:B:1017:HOH:O	2.66	0.42
1:A:63:LEU:HD23	1:A:377:MET:HG3	2.02	0.42
1:B:146:HIS:HE1	3:B:687:HOH:O	2.02	0.42
1:B:473:HIS:CE1	3:B:988:HOH:O	2.72	0.42
1:A:392:ILE:CD1	1:A:394:SER:HB3	2.49	0.42
1:A:10:ILE:HG22	1:A:443:PHE:CE1	2.55	0.42
1:A:500:GLY:O	1:A:501:LYS:HB2	2.20	0.42
1:B:110:ASN:HB3	1:B:161:TYR:CZ	2.55	0.42
1:A:243:LEU:HD12	1:A:243:LEU:N	2.35	0.42
1:B:289:HIS:HE1	3:B:839:HOH:O	2.03	0.42
1:B:237:HIS:HD2	3:B:867:HOH:O	2.03	0.41
1:B:479:GLN:NE2	3:B:820:HOH:O	2.53	0.41
1:A:456:HIS:HD2	1:A:458:ASN:H	1.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:473:HIS:CE1	1:B:475:ASN:HB2	2.56	0.41
1:A:99:TRP:CZ2	2:A:503:GOL:H31	2.56	0.41
1:A:220:MET:HB3	1:A:220:MET:HE2	1.85	0.41
1:A:447:ARG:NE	3:A:1018:HOH:O	2.53	0.41
1:A:175:GLU:HB2	3:A:1069:HOH:O	2.20	0.41
1:A:10:ILE:HG22	1:A:443:PHE:HE1	1.86	0.41
1:A:176:MET:O	1:A:184:HIS:HD2	2.04	0.41
1:B:324:PHE:CZ	1:B:455:GLU:HA	2.55	0.40
1:B:451:HIS:CG	1:B:476:GLY:HA3	2.57	0.40
1:A:10:ILE:HA	3:A:834:HOH:O	2.22	0.40
1:A:11:GLU:HA	1:A:442:SER:HG	1.87	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:763:HOH:O	3:B:511:HOH:O[6_554]	2.07	0.13
3:A:1063:HOH:O	3:B:992:HOH:O[1_554]	2.13	0.07
3:A:1041:HOH:O	3:B:688:HOH:O[1_554]	2.16	0.04
3:A:507:HOH:O	3:A:564:HOH:O[2_655]	2.17	0.03

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	495/502 (99%)	471 (95%)	22 (4%)	2 (0%)	34 17
1	B	495/502 (99%)	477 (96%)	16 (3%)	2 (0%)	34 17
All	All	990/1004 (99%)	948 (96%)	38 (4%)	4 (0%)	34 17

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	299	TYR
1	B	299	TYR
1	A	356	ILE
1	B	356	ILE

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	430/434 (99%)	415 (96%)	15 (4%)	36	13
1	B	430/434 (99%)	420 (98%)	10 (2%)	50	28
All	All	860/868 (99%)	835 (97%)	25 (3%)	42	19

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	MET
1	A	118	MET
1	A	176	MET
1	A	216	SER
1	A	277	TYR
1	A	285	LYS
1	A	292	PHE
1	A	339	LYS
1	A	370	GLN
1	A	391	VAL
1	A	392	ILE
1	A	396	LYS
1	A	444	GLU
1	A	445	ASP
1	A	447	ARG
1	B	5	LYS
1	B	46	GLN
1	B	277	TYR
1	B	292	PHE
1	B	392	ILE

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Mol	Chain	Res	Type
1	B	447	ARG
1	B	473	HIS
1	B	480	LEU
1	B	482	ASP
1	B	484	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	HIS
1	A	64	GLN
1	A	159	HIS
1	A	181	GLN
1	A	184	HIS
1	A	219	ASN
1	A	237	HIS
1	A	245	GLN
1	A	300	HIS
1	A	302	ASN
1	A	370	GLN
1	A	451	HIS
1	A	456	HIS
1	A	463	ASN
1	A	473	HIS
1	A	495	ASN
1	B	64	GLN
1	B	146	HIS
1	B	159	HIS
1	B	181	GLN
1	B	252	ASN
1	B	289	HIS
1	B	300	HIS
1	B	302	ASN
1	B	354	ASN
1	B	461	GLN
1	B	475	ASN
1	B	479	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

2 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	503	-	5,5,5	0.68	0	5,5,5	1.07	1 (20%)
2	GOL	B	503	-	5,5,5	0.67	0	5,5,5	0.90	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	503	-	-	2/4/4/4	-
2	GOL	B	503	-	-	2/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	503	GOL	O2-C2-C1	2.26	118.55	109.18

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	503	GOL	C1-C2-C3-O3
2	B	503	GOL	C1-C2-C3-O3
2	A	503	GOL	O2-C2-C3-O3
2	B	503	GOL	O2-C2-C3-O3

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	503	GOL	3	0
2	B	503	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)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.