



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:31 AM UTC

PDB ID : 3VGE / pdb_00003vge
Title : Crystal structure of glycosyltrehalose trehalohydrolase (D252S)
Authors : Okazaki, N.; Tamada, T.; Feese, M.D.; Kato, M.; Miura, Y.; Komeda, T.; Kobayashi, K.; Kondo, K.; Kuroki, R.
Deposited on : 2011-08-09
Resolution : 2.70 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

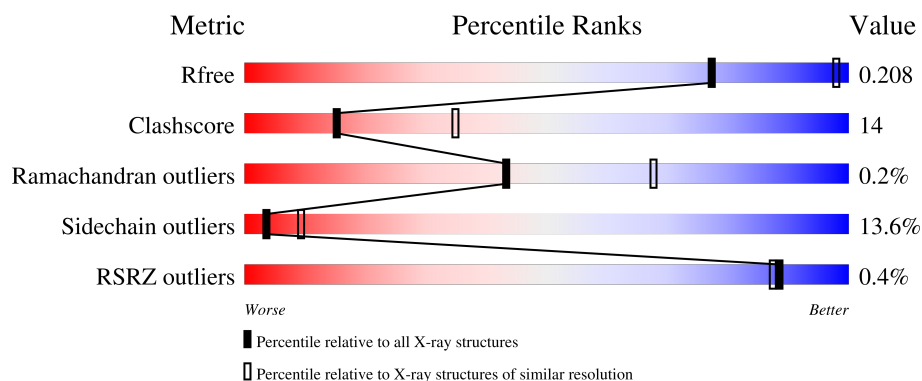
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	558	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4772 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.

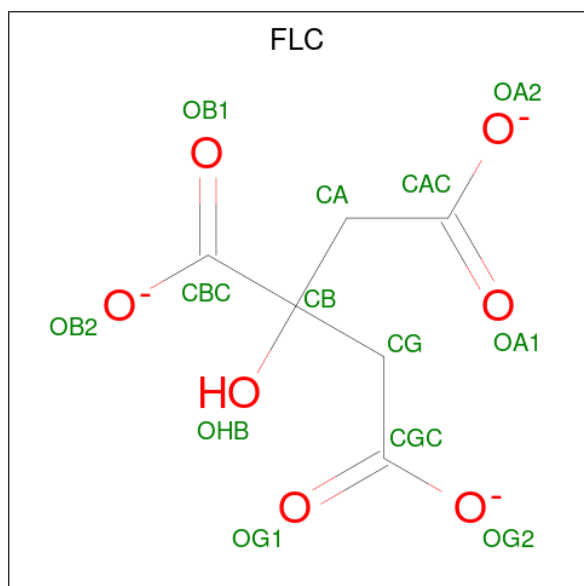
- Molecule 1 is a protein called Malto-oligosyltrehalose trehalohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	555	Total	C	N	O	S	0	0	0
			4552	2934	747	862	9			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	252	SER	ASP	engineered mutation	UNP Q55088

- Molecule 2 is CITRATE ANION (CCD ID: FLC) (formula: $C_6H_5O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	6	7		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

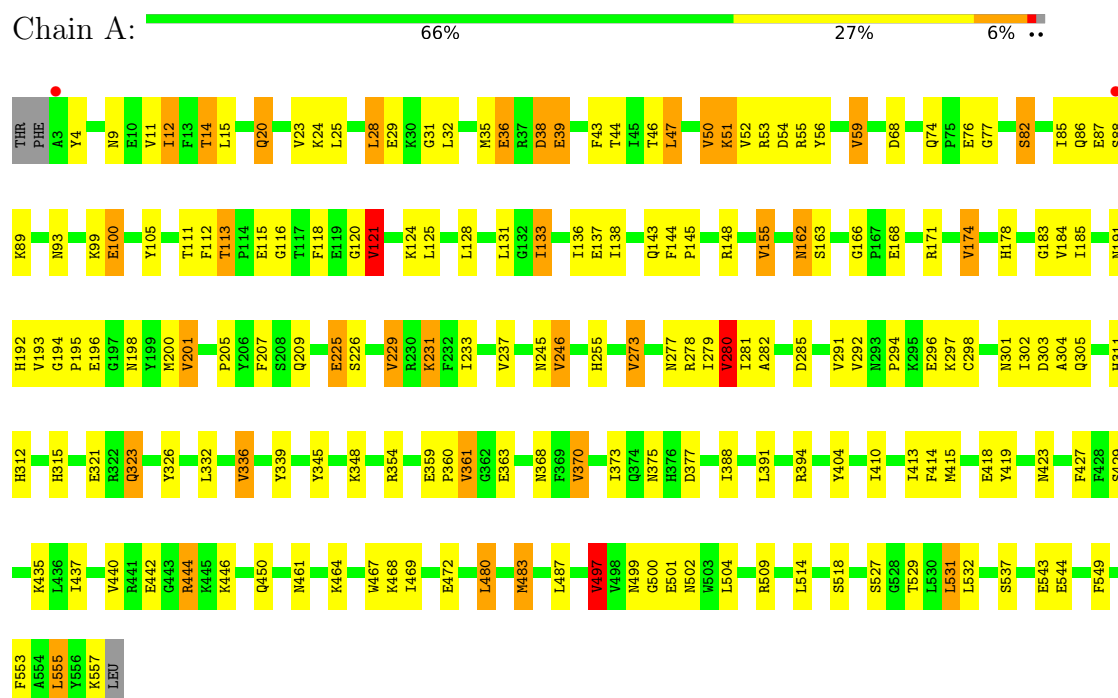
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	165	Total	O	0	0
			165	165		

3 Residue-property plots

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 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: Malto-oligosyltrehalose trehalohydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	78.51Å 78.51Å 282.44Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	61.50 – 2.70 61.50 – 2.71	Depositor EDS
% Data completeness (in resolution range)	95.8 (61.50-2.70) 95.8 (61.50-2.71)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.22 (at 2.69Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.159 , 0.212 0.160 , 0.208	Depositor DCC
R_{free} test set	1384 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	39.3	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 54.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.040 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4772	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 4.15% 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: FLC, 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	1.14	2/4669 (0.0%)	1.22	23/6308 (0.4%)

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	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	195	PRO	N-CA	-5.68	1.43	1.47
1	A	360	PRO	CA-C	-5.20	1.46	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	361	VAL	N-CA-C	-7.87	105.53	113.47
1	A	201	VAL	CB-CA-C	-7.08	101.71	112.05
1	A	361	VAL	N-CA-CB	-6.96	104.05	112.33
1	A	194	GLY	CA-C-N	-6.78	113.21	120.94
1	A	194	GLY	C-N-CA	-6.78	113.21	120.94
1	A	359	GLU	N-CA-C	-6.66	99.52	108.11
1	A	183	GLY	CA-C-N	-6.61	113.94	123.06
1	A	183	GLY	C-N-CA	-6.61	113.94	123.06
1	A	549	PHE	N-CA-C	6.36	120.41	110.17
1	A	166	GLY	CA-C-N	6.31	126.51	119.32
1	A	166	GLY	C-N-CA	6.31	126.51	119.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	497	VAL	N-CA-C	5.84	116.28	107.75
1	A	280	VAL	CB-CA-C	-5.76	101.63	110.50
1	A	500	GLY	N-CA-C	-5.62	104.56	112.60
1	A	423	ASN	CA-C-N	-5.58	114.85	120.21
1	A	423	ASN	C-N-CA	-5.58	114.85	120.21
1	A	121	VAL	CB-CA-C	5.50	119.81	112.22
1	A	277	ASN	N-CA-C	5.47	118.75	111.30
1	A	86	GLN	N-CA-C	5.45	122.41	110.80
1	A	31	GLY	N-CA-C	5.34	118.37	110.63
1	A	444	ARG	CB-CG-CD	-5.32	99.07	111.30
1	A	557	LYS	CG-CD-CE	-5.11	99.55	111.30
1	A	246	VAL	N-CA-CB	-5.09	102.83	111.23

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	112	PHE	Peptide

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	4552	0	4412	120	0
2	A	13	0	5	1	0
3	A	42	0	56	9	0
4	A	165	0	0	2	0
All	All	4772	0	4473	123	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 14.

All (123) 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:273:VAL:HG21	1:A:280:VAL:HG22	1.42	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:394:ARG:NH2	1:A:472:GLU:OE1	2.08	0.86
1:A:85:ILE:HD12	3:A:1106:GOL:H32	1.56	0.86
1:A:321:GLU:HB3	1:A:323:GLN:HE22	1.44	0.83
1:A:143:GLN:HE22	1:A:162:ASN:H	1.26	0.81
1:A:323:GLN:HE21	1:A:323:GLN:H	1.30	0.80
1:A:51:LYS:HE2	1:A:52:VAL:HG13	1.62	0.79
1:A:311:HIS:HD2	1:A:312:HIS:ND1	1.81	0.78
3:A:1103:GOL:H2	4:A:2145:HOH:O	1.84	0.77
1:A:113:THR:HG22	1:A:116:GLY:H	1.51	0.73
1:A:133:ILE:HD11	1:A:413:ILE:HD12	1.70	0.72
1:A:323:GLN:H	1:A:323:GLN:NE2	1.88	0.71
1:A:273:VAL:CG2	1:A:280:VAL:HG22	2.19	0.71
1:A:191:ASN:OD1	1:A:192:HIS:HD2	1.74	0.71
1:A:100:GLU:H	1:A:100:GLU:CD	2.00	0.70
1:A:162:ASN:HD22	1:A:163:SER:H	1.41	0.68
1:A:377:ASP:OD2	3:A:1107:GOL:H31	1.95	0.67
1:A:105:TYR:HB2	1:A:133:ILE:HG12	1.77	0.66
1:A:537:SER:O	3:A:1102:GOL:H32	1.96	0.66
1:A:444:ARG:HD2	1:A:450:GLN:NE2	2.11	0.66
1:A:444:ARG:HH11	1:A:450:GLN:HE21	1.45	0.64
1:A:255:HIS:CD2	1:A:285:ASP:H	2.16	0.63
1:A:461:ASN:O	1:A:464:LYS:HG3	1.98	0.63
1:A:113:THR:HG21	1:A:120:GLY:HA3	1.83	0.61
1:A:174:VAL:HG23	1:A:178:HIS:CE1	2.35	0.61
1:A:85:ILE:HD12	3:A:1106:GOL:C3	2.28	0.60
1:A:444:ARG:HH11	1:A:450:GLN:NE2	2.00	0.59
1:A:113:THR:CG2	1:A:116:GLY:H	2.15	0.59
1:A:144:PHE:HB2	1:A:145:PRO:CD	2.33	0.58
1:A:191:ASN:OD1	1:A:192:HIS:CD2	2.57	0.57
1:A:200:MET:CE	1:A:207:PHE:CZ	2.87	0.57
1:A:368:ASN:HD21	3:A:1103:GOL:H12	1.70	0.57
1:A:144:PHE:HB2	1:A:145:PRO:HD2	1.87	0.56
1:A:162:ASN:ND2	1:A:163:SER:H	2.02	0.56
1:A:278:ARG:O	1:A:279:ILE:HD13	2.05	0.56
1:A:226:SER:HA	1:A:229:VAL:HG13	1.86	0.56
1:A:23:VAL:HG12	1:A:35:MET:CE	2.36	0.56
1:A:162:ASN:HD22	1:A:163:SER:N	2.06	0.54
1:A:444:ARG:HD2	1:A:450:GLN:HE21	1.74	0.53
1:A:143:GLN:NE2	1:A:162:ASN:H	2.00	0.52
1:A:461:ASN:HA	1:A:464:LYS:HE3	1.90	0.52
1:A:39:GLU:H	1:A:39:GLU:CD	2.18	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:388:ILE:HA	1:A:391:LEU:HD12	1.91	0.52
1:A:345:TYR:CE2	1:A:348:LYS:HE2	2.46	0.51
1:A:14:THR:HG23	1:A:44:THR:CG2	2.40	0.51
1:A:255:HIS:CD2	1:A:255:HIS:H	2.29	0.51
1:A:282:ALA:O	1:A:305:GLN:HA	2.11	0.51
1:A:315:HIS:CD2	1:A:326:TYR:CZ	2.99	0.51
1:A:162:ASN:ND2	1:A:163:SER:N	2.59	0.51
1:A:28:LEU:H	1:A:28:LEU:CD2	2.24	0.50
1:A:38:ASP:C	1:A:38:ASP:OD1	2.54	0.50
1:A:14:THR:HG23	1:A:44:THR:HG22	1.92	0.50
1:A:87:GLU:O	1:A:89:LYS:HG2	2.12	0.50
1:A:118:PHE:HA	1:A:121:VAL:HG13	1.93	0.50
1:A:440:VAL:O	1:A:444:ARG:HB2	2.11	0.50
1:A:200:MET:HE3	1:A:207:PHE:CZ	2.47	0.50
1:A:51:LYS:HD3	1:A:52:VAL:H	1.76	0.50
1:A:483:MET:HE2	1:A:531:LEU:HG	1.92	0.49
1:A:113:THR:HG23	1:A:115:GLU:H	1.77	0.49
1:A:280:VAL:O	1:A:303:ASP:HB2	2.11	0.49
1:A:427:PHE:HB3	1:A:450:GLN:HE22	1.77	0.49
1:A:131:LEU:CD2	1:A:418:GLU:HG3	2.43	0.49
2:A:1001:FLC:OA1	3:A:1104:GOL:H11	2.13	0.48
1:A:11:VAL:HG11	1:A:50:VAL:CG1	2.43	0.48
1:A:4:TYR:HA	1:A:12:ILE:O	2.13	0.48
1:A:136:ILE:HD12	1:A:138:ILE:HD11	1.94	0.48
1:A:499:ASN:CB	1:A:504:LEU:HD12	2.44	0.47
1:A:28:LEU:HD23	1:A:56:TYR:HA	1.96	0.47
1:A:200:MET:CE	1:A:207:PHE:HZ	2.26	0.47
1:A:87:GLU:O	1:A:87:GLU:HG3	2.15	0.47
1:A:291:VAL:O	1:A:302:ILE:HB	2.14	0.47
1:A:297:LYS:O	1:A:298:CYS:HB2	2.15	0.47
1:A:336:VAL:HB	1:A:497:VAL:HG21	1.97	0.47
1:A:480:LEU:HD13	1:A:532:LEU:CD2	2.45	0.47
1:A:209:GLN:H	1:A:209:GLN:CD	2.24	0.46
1:A:74:GLN:NE2	1:A:82:SER:HA	2.30	0.46
3:A:1104:GOL:H2	4:A:2057:HOH:O	2.16	0.46
1:A:137:GLU:HG3	1:A:185:ILE:HG22	1.98	0.46
1:A:55:ARG:NH1	1:A:76:GLU:OE2	2.49	0.45
1:A:155:VAL:HG22	1:A:198:ASN:HB2	1.98	0.45
1:A:231:LYS:HA	1:A:231:LYS:HD2	1.57	0.45
1:A:323:GLN:HE21	1:A:323:GLN:N	2.08	0.45
1:A:171:ARG:HG2	3:A:1106:GOL:H11	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:GLN:HG2	1:A:23:VAL:CG2	2.47	0.45
1:A:375:ASN:C	1:A:375:ASN:OD1	2.59	0.45
1:A:28:LEU:CD2	1:A:28:LEU:N	2.80	0.44
1:A:373:ILE:HG12	1:A:404:TYR:CE2	2.52	0.44
1:A:28:LEU:H	1:A:28:LEU:HD23	1.83	0.44
1:A:148:ARG:O	1:A:429:SER:HA	2.18	0.44
1:A:273:VAL:HG21	1:A:280:VAL:CG2	2.29	0.44
1:A:499:ASN:HB3	1:A:504:LEU:HD12	1.99	0.44
1:A:171:ARG:HD2	1:A:171:ARG:HA	1.81	0.43
1:A:444:ARG:NH1	1:A:450:GLN:NE2	2.65	0.43
1:A:205:PRO:HG2	1:A:225:GLU:HB2	2.01	0.43
1:A:502:ASN:HB3	1:A:518:SER:HB3	2.00	0.43
1:A:36:GLU:O	1:A:43:PHE:HA	2.19	0.43
1:A:543:GLU:O	1:A:544:GLU:C	2.60	0.43
1:A:294:PRO:HA	1:A:301:ASN:ND2	2.34	0.43
1:A:437:ILE:O	1:A:440:VAL:HG12	2.18	0.43
1:A:419:TYR:CD1	1:A:419:TYR:C	2.96	0.42
1:A:131:LEU:HD21	1:A:418:GLU:HG3	2.01	0.42
1:A:74:GLN:HB3	1:A:77:GLY:O	2.19	0.42
1:A:124:LYS:HE3	1:A:467:TRP:CE2	2.55	0.42
1:A:444:ARG:HD3	1:A:444:ARG:HA	1.52	0.42
1:A:24:LYS:HB2	1:A:59:VAL:HG13	2.02	0.42
1:A:294:PRO:HA	1:A:301:ASN:HD22	1.85	0.41
1:A:296:GLU:H	1:A:296:GLU:CD	2.29	0.41
1:A:311:HIS:CD2	1:A:311:HIS:C	2.98	0.41
1:A:315:HIS:CD2	1:A:326:TYR:CE1	3.07	0.41
1:A:414:PHE:O	1:A:415:MET:C	2.64	0.41
1:A:553:PHE:CZ	1:A:555:LEU:HB2	2.56	0.41
1:A:105:TYR:HB2	1:A:133:ILE:CG1	2.46	0.41
1:A:480:LEU:HD13	1:A:532:LEU:HD23	2.01	0.41
1:A:469:ILE:H	1:A:469:ILE:HG13	1.70	0.41
1:A:56:TYR:HE1	1:A:68:ASP:HB2	1.85	0.40
1:A:111:THR:HB	1:A:464:LYS:HA	2.03	0.40
1:A:200:MET:HE2	1:A:200:MET:HB2	1.92	0.40
1:A:370:VAL:HA	1:A:410:ILE:O	2.21	0.40
1:A:51:LYS:O	1:A:54:ASP:HB2	2.20	0.40
1:A:124:LYS:HE3	1:A:467:TRP:CZ2	2.56	0.40
1:A:20:GLN:HG2	1:A:23:VAL:HG22	2.03	0.40
1:A:46:THR:O	1:A:47:LEU:HD13	2.22	0.40
1:A:281:ILE:HA	1:A:304:ALA:O	2.21	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	553/558 (99%)	521 (94%)	31 (6%)	1 (0%)	43 68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	88	SER

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	491/494 (99%)	424 (86%)	67 (14%)	3 9

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	12	ILE
1	A	14	THR
1	A	15	LEU
1	A	20	GLN
1	A	25	LEU
1	A	28	LEU
1	A	29	GLU
1	A	32	LEU
1	A	36	GLU

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Mol	Chain	Res	Type
1	A	38	ASP
1	A	39	GLU
1	A	47	LEU
1	A	50	VAL
1	A	51	LYS
1	A	53	ARG
1	A	59	VAL
1	A	82	SER
1	A	93	ASN
1	A	99	LYS
1	A	100	GLU
1	A	113	THR
1	A	121	VAL
1	A	125	LEU
1	A	128	LEU
1	A	133	ILE
1	A	155	VAL
1	A	162	ASN
1	A	168	GLU
1	A	174	VAL
1	A	184	VAL
1	A	193	VAL
1	A	196	GLU
1	A	201	VAL
1	A	225	GLU
1	A	229	VAL
1	A	231	LYS
1	A	233	ILE
1	A	237	VAL
1	A	245	ASN
1	A	246	VAL
1	A	273	VAL
1	A	280	VAL
1	A	292	VAL
1	A	323	GLN
1	A	332	LEU
1	A	336	VAL
1	A	339	TYR
1	A	354	ARG
1	A	361	VAL
1	A	363	GLU
1	A	370	VAL

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Mol	Chain	Res	Type
1	A	435	LYS
1	A	442	GLU
1	A	446	LYS
1	A	468	LYS
1	A	480	LEU
1	A	483	MET
1	A	487	LEU
1	A	497	VAL
1	A	501	GLU
1	A	509	ARG
1	A	514	LEU
1	A	527	SER
1	A	529	THR
1	A	531	LEU
1	A	555	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	GLN
1	A	93	ASN
1	A	143	GLN
1	A	162	ASN
1	A	192	HIS
1	A	245	ASN
1	A	255	HIS
1	A	301	ASN
1	A	311	HIS
1	A	323	GLN
1	A	368	ASN
1	A	448	ASN
1	A	450	GLN
1	A	499	ASN
1	A	541	HIS

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](#)

8 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	$\# Z > 2$
3	GOL	A	1103	-	5,5,5	0.49	0	5,5,5	0.80	0
3	GOL	A	1102	-	5,5,5	0.49	0	5,5,5	1.19	0
3	GOL	A	1104	-	5,5,5	0.51	0	5,5,5	0.77	0
3	GOL	A	1107	-	5,5,5	0.27	0	5,5,5	1.18	1 (20%)
3	GOL	A	1106	-	5,5,5	0.27	0	5,5,5	0.67	0
3	GOL	A	1105	-	5,5,5	0.59	0	5,5,5	0.63	0
2	FLC	A	1001	-	12,12,12	0.97	0	17,17,17	2.18	6 (35%)
3	GOL	A	1101	-	5,5,5	0.29	0	5,5,5	0.75	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	GOL	A	1103	-	-	4/4/4/4	-
3	GOL	A	1102	-	-	3/4/4/4	-
3	GOL	A	1104	-	-	0/4/4/4	-
3	GOL	A	1107	-	-	1/4/4/4	-
3	GOL	A	1106	-	-	2/4/4/4	-
3	GOL	A	1105	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FLC	A	1001	-	-	4/16/16/16	-
3	GOL	A	1101	-	-	2/4/4/4	-

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	FLC	OB2-CBC-CB	4.73	122.22	113.14
2	A	1001	FLC	OB1-CBC-CB	-4.64	113.10	122.09
2	A	1001	FLC	OA1-CAC-CA	-2.94	114.62	122.95
2	A	1001	FLC	OA2-CAC-CA	2.71	122.94	114.35
2	A	1001	FLC	OHB-CB-CBC	-2.60	105.27	108.96
3	A	1107	GOL	O2-C2-C3	2.22	118.38	109.18
2	A	1001	FLC	OG2-CGC-CG	2.18	121.24	114.35

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1101	GOL	C1-C2-C3-O3
3	A	1106	GOL	C1-C2-C3-O3
3	A	1102	GOL	O1-C1-C2-C3
3	A	1102	GOL	C1-C2-C3-O3
3	A	1103	GOL	O1-C1-C2-C3
3	A	1103	GOL	C1-C2-C3-O3
3	A	1105	GOL	O1-C1-C2-C3
3	A	1101	GOL	O2-C2-C3-O3
3	A	1103	GOL	O2-C2-C3-O3
3	A	1106	GOL	O2-C2-C3-O3
3	A	1103	GOL	O1-C1-C2-O2
3	A	1107	GOL	O1-C1-C2-O2
3	A	1105	GOL	O1-C1-C2-O2
2	A	1001	FLC	CB-CG-CGC-OG2
3	A	1102	GOL	O2-C2-C3-O3
2	A	1001	FLC	CB-CG-CGC-OG1
2	A	1001	FLC	CB-CA-CAC-OA1
2	A	1001	FLC	CB-CA-CAC-OA2

There are no ring outliers.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1103	GOL	2	0
3	A	1102	GOL	1	0
3	A	1104	GOL	2	0
3	A	1107	GOL	1	0
3	A	1106	GOL	3	0
2	A	1001	FLC	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 [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	555/558 (99%)	-0.64	2 (0%) 88 87	22, 35, 73, 94	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	88	SER	3.0
1	A	3	ALA	2.9

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(Å ²)	Q<0.9
3	GOL	A	1103	6/6	0.88	0.15	75,77,77,77	0
3	GOL	A	1105	6/6	0.91	0.12	53,61,63,64	0
3	GOL	A	1106	6/6	0.91	0.21	83,87,87,87	0
3	GOL	A	1102	6/6	0.92	0.18	62,70,71,72	0
3	GOL	A	1101	6/6	0.92	0.13	59,63,68,70	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	A	1107	6/6	0.93	0.11	44,48,48,49	0
3	GOL	A	1104	6/6	0.95	0.11	50,55,58,59	0
2	FLC	A	1001	13/13	0.95	0.07	47,50,59,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.