



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:05 PM UTC

PDB ID : 5OKA / pdb_00005oka
Title : Non-conservatively refined structure of Gan1D-E170Q, a catalytic mutant of a 6-phospho-beta-galactosidase from *Geobacillus stearothermophilus*
Authors : Lansky, S.; Zehavi, A.; Shoham, Y.; Shoham, G.
Deposited on : 2017-07-25
Resolution : 1.25 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

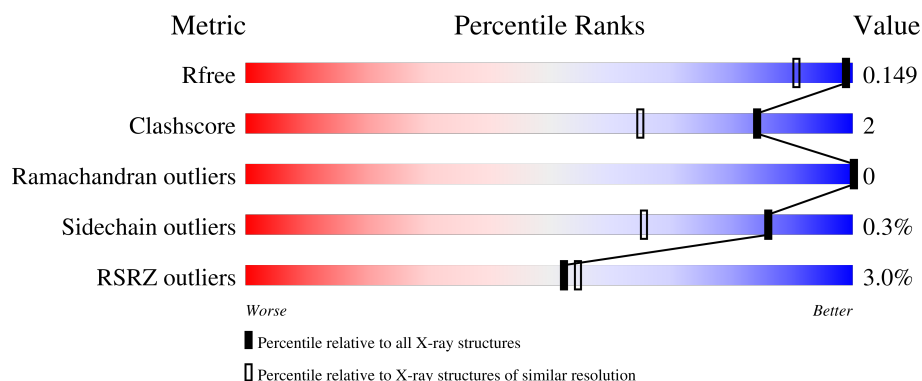
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.25 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	1693 (1.28-1.24)
Clashscore	190562	1730 (1.28-1.24)
Ramachandran outliers	187476	1695 (1.28-1.24)
Sidechain outliers	187428	1694 (1.28-1.24)
RSRZ outliers	180081	1693 (1.28-1.24)

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	485	3% 91% 5%
1	B	485	3% 90% 5% 5%
1	C	485	2% 90% 6% 5%
1	D	485	3% 90% 5% 5%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 18165 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 Putative 6-phospho-beta-galactobiosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	462	Total	C	N	O	S	0	26	0
			3957	2556	674	712	15			
1	B	462	Total	C	N	O	S	0	25	0
			3959	2552	672	721	14			
1	C	462	Total	C	N	O	S	0	23	0
			3939	2542	674	709	14			
1	D	462	Total	C	N	O	S	0	20	0
			3908	2524	670	701	13			

There are 36 discrepancies between the modelled and reference sequences:

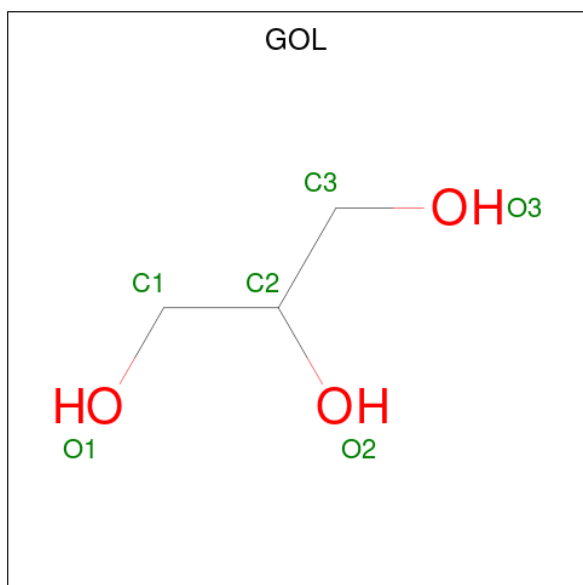
Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	-	initiating methionine	UNP W8QF82
A	-5	ILE	-	expression tag	UNP W8QF82
A	-4	HIS	-	expression tag	UNP W8QF82
A	-3	HIS	-	expression tag	UNP W8QF82
A	-2	HIS	-	expression tag	UNP W8QF82
A	-1	HIS	-	expression tag	UNP W8QF82
A	0	HIS	-	expression tag	UNP W8QF82
A	1	HIS	-	expression tag	UNP W8QF82
A	170	GLN	GLU	engineered mutation	UNP W8QF82
B	-6	MET	-	initiating methionine	UNP W8QF82
B	-5	ILE	-	expression tag	UNP W8QF82
B	-4	HIS	-	expression tag	UNP W8QF82
B	-3	HIS	-	expression tag	UNP W8QF82
B	-2	HIS	-	expression tag	UNP W8QF82
B	-1	HIS	-	expression tag	UNP W8QF82
B	0	HIS	-	expression tag	UNP W8QF82
B	1	HIS	-	expression tag	UNP W8QF82
B	170	GLN	GLU	engineered mutation	UNP W8QF82
C	-6	MET	-	initiating methionine	UNP W8QF82
C	-5	ILE	-	expression tag	UNP W8QF82
C	-4	HIS	-	expression tag	UNP W8QF82

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	HIS	-	expression tag	UNP W8QF82
C	-2	HIS	-	expression tag	UNP W8QF82
C	-1	HIS	-	expression tag	UNP W8QF82
C	0	HIS	-	expression tag	UNP W8QF82
C	1	HIS	-	expression tag	UNP W8QF82
C	170	GLN	GLU	engineered mutation	UNP W8QF82
D	-6	MET	-	initiating methionine	UNP W8QF82
D	-5	ILE	-	expression tag	UNP W8QF82
D	-4	HIS	-	expression tag	UNP W8QF82
D	-3	HIS	-	expression tag	UNP W8QF82
D	-2	HIS	-	expression tag	UNP W8QF82
D	-1	HIS	-	expression tag	UNP W8QF82
D	0	HIS	-	expression tag	UNP W8QF82
D	1	HIS	-	expression tag	UNP W8QF82
D	170	GLN	GLU	engineered mutation	UNP W8QF82

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



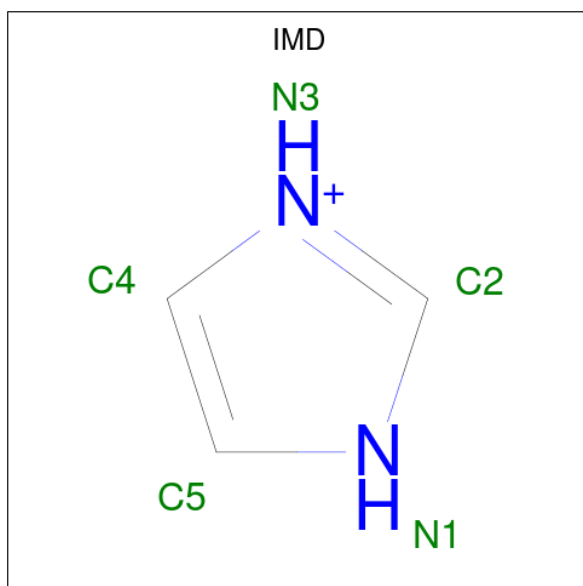
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		

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

- Molecule 3 is IMIDAZOLE (CCD ID: IMD) (formula: $C_3H_5N_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			5	3	2		
3	C	1	Total	C	N	0	0
			5	3	2		
3	D	1	Total	C	N	0	0
			5	3	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	562	Total 562	O 562	0	0
4	B	583	Total 583	O 583	0	0
4	C	624	Total 624	O 624	0	0
4	D	546	Total 546	O 546	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	101.55Å 97.43Å 105.25Å 90.00° 97.82° 90.00°	Depositor
Resolution (Å)	43.91 – 1.25 43.91 – 1.25	Depositor EDS
% Data completeness (in resolution range)	80.2 (43.91-1.25) 80.2 (43.91-1.25)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 1.25Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.127 , 0.149 0.128 , 0.149	Depositor DCC
R_{free} test set	4521 reflections (0.81%)	wwPDB-VP
Wilson B-factor (Å ²)	10.1	Xtriage
Anisotropy	0.421	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 49.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	18165	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 71.14 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.7679e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: IMD, 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	0.28	0/4153	0.53	1/5643 (0.0%)
1	B	0.27	0/4143	0.52	0/5635
1	C	0.28	0/4135	0.52	0/5623
1	D	0.27	0/4091	0.52	0/5562
All	All	0.27	0/16522	0.52	1/22463 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	189	LYS	CB-CG-CD	-5.07	99.64	111.30

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	3957	0	3825	14	0
1	B	3959	0	3789	20	0
1	C	3939	0	3789	19	0
1	D	3908	0	3754	19	0
2	A	24	0	32	3	0
2	B	30	0	40	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	12	0	16	0	0
2	D	6	0	8	0	0
3	A	5	0	5	0	0
3	C	5	0	5	0	0
3	D	5	0	5	1	0
4	A	562	0	0	4	0
4	B	583	0	0	4	0
4	C	624	0	0	3	0
4	D	546	0	0	7	0
All	All	18165	0	15268	71	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 2.

All (71) 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:C:214[A]:HIS:HE1	1:D:383:GLU:HG2	1.59	0.68
1:A:97[B]:ASN:ND2	4:A:603:HOH:O	2.26	0.66
1:A:438:GLN:NE2	4:A:601:HOH:O	2.23	0.60
1:D:149:ARG:NH1	4:D:603:HOH:O	2.35	0.59
1:B:472[A]:GLU:OE1	4:B:601:HOH:O	2.17	0.59
1:D:249[A]:PHE:CE1	1:D:333:ILE:HD11	2.42	0.55
1:B:348[B]:THR:HB	1:B:354:ILE:HD11	1.88	0.55
1:B:341:ARG:NH2	1:B:351[A]:ASP:OD2	2.37	0.55
1:C:473:THR:HG21	1:C:477[B]:GLU:HG3	1.92	0.52
1:A:149:ARG:HD3	4:A:825:HOH:O	2.09	0.52
1:D:149:ARG:NH2	4:D:612:HOH:O	2.44	0.50
1:D:245:ASN:ND2	4:D:613:HOH:O	2.44	0.50
1:D:97[B]:ASN:OD1	1:D:99:LYS:HB2	2.12	0.50
1:A:266:ALA:HB1	2:A:504:GOL:H32	1.95	0.48
1:B:182:GLY:HA3	1:B:189:LYS:HE2	1.95	0.48
1:A:336:LEU:CD2	2:A:504:GOL:H2	2.43	0.48
1:B:378:GLU:HG3	1:B:425:TRP:HB2	1.96	0.48
2:B:504:GOL:O1	4:B:602:HOH:O	2.20	0.48
1:A:378:GLU:HG3	1:A:425:TRP:HB2	1.95	0.47
1:D:312:ASP:OD1	1:D:312:ASP:N	2.46	0.47
1:C:153:THR:O	1:C:157[B]:ARG:HG2	2.15	0.47
1:C:312:ASP:OD1	1:C:312:ASP:N	2.46	0.46
1:A:112:ARG:HE	1:A:112:ARG:HB2	1.65	0.46
1:D:378:GLU:HG3	1:D:425:TRP:HB2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:180[B]:ARG:HD3	1:B:194[B]:MET:HE1	1.98	0.45
1:D:292:LYS:HB3	4:D:702:HOH:O	2.16	0.45
1:B:341:ARG:HH22	1:B:351[A]:ASP:CG	2.24	0.45
1:B:312:ASP:OD1	1:B:312:ASP:N	2.50	0.45
1:C:249:PHE:CE2	1:C:333:ILE:HD11	2.51	0.45
1:B:425:TRP:HA	1:B:426[B]:SER:HA	1.79	0.45
1:D:92:GLY:HA2	1:D:128:PRO:HG2	1.99	0.45
1:A:112:ARG:NH1	1:A:161:ARG:O	2.50	0.44
1:B:341:ARG:NH1	1:B:351[A]:ASP:OD1	2.37	0.44
1:D:157[A]:ARG:HD2	3:D:502:IMD:C4	2.47	0.44
1:D:262:MET:HE3	4:D:1072:HOH:O	2.18	0.44
1:C:378:GLU:HG3	1:C:425:TRP:HB2	1.99	0.43
1:B:149:ARG:HD3	4:B:753:HOH:O	2.17	0.43
1:C:94:GLY:O	1:C:149:ARG:NH2	2.27	0.43
1:C:132[A]:MET:HE1	4:C:694:HOH:O	2.19	0.43
1:B:262[A]:MET:HE3	1:B:262[A]:MET:HB2	1.81	0.43
1:D:249[B]:PHE:HZ	1:D:270:TYR:HB3	1.84	0.43
1:A:394:ASN:ND2	4:A:622:HOH:O	2.52	0.43
1:C:112:ARG:HD2	1:C:161:ARG:HB3	2.01	0.42
1:B:153:THR:O	1:B:157[B]:ARG:HG2	2.19	0.42
1:C:157[A]:ARG:HG3	1:C:158:PHE:CE1	2.53	0.42
1:D:310:PRO:HD2	4:D:682:HOH:O	2.19	0.42
1:A:16:GLY:HA2	1:A:78:ALA:O	2.19	0.42
1:A:47:ARG:HD3	1:A:47:ARG:HA	1.87	0.42
1:A:350[B]:TRP:HZ2	1:A:384:PHE:HZ	1.68	0.42
1:B:348[A]:THR:OG1	1:B:352[A]:TRP:HB2	2.20	0.42
1:C:191[B]:MET:HE2	1:C:191[B]:MET:HB3	1.98	0.42
1:C:171:GLN:HA	1:C:174:PHE:CE2	2.55	0.42
1:C:292:LYS:HB3	4:C:745:HOH:O	2.20	0.42
1:D:175:ILE:HG21	1:D:198:ASN:HB2	2.02	0.42
1:C:214[A]:HIS:CE1	1:D:383:GLU:HG2	2.45	0.42
1:C:32:GLY:HA3	1:C:97[B]:ASN:HD21	1.84	0.41
1:B:16:GLY:HA2	1:B:78:ALA:O	2.20	0.41
1:B:180[B]:ARG:HB2	1:B:194[B]:MET:SD	2.59	0.41
1:A:180[A]:ARG:HD3	1:A:191[A]:MET:HG2	2.02	0.41
1:C:112:ARG:CD	1:C:161:ARG:HB3	2.50	0.41
1:A:336:LEU:HD21	2:A:504:GOL:H2	2.02	0.41
1:D:31:ASP:HB3	1:D:99:LYS:HB3	2.03	0.41
1:D:132[A]:MET:HE1	4:D:635:HOH:O	2.21	0.41
1:B:171:GLN:HA	1:B:174:PHE:CE2	2.56	0.41
1:C:425:TRP:HA	1:C:426:SER:HA	1.80	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:192:LYS:NZ	4:B:622:HOH:O	2.54	0.41
1:D:171:GLN:HA	1:D:174:PHE:CE2	2.55	0.40
1:B:194[B]:MET:HE3	1:B:195:TYR:CE2	2.56	0.40
1:B:425:TRP:HA	1:B:426[A]:SER:HA	1.80	0.40
1:C:16:GLY:HA2	1:C:78:ALA:O	2.21	0.40
1:C:192:LYS:HD2	4:C:795:HOH: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	485/485 (100%)	471 (97%)	14 (3%)	0	100	100
1	B	484/485 (100%)	469 (97%)	15 (3%)	0	100	100
1	C	482/485 (99%)	467 (97%)	15 (3%)	0	100	100
1	D	478/485 (99%)	465 (97%)	13 (3%)	0	100	100
All	All	1929/1940 (99%)	1872 (97%)	57 (3%)	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	418/412 (102%)	416 (100%)	2 (0%)	81	56
1	B	418/412 (102%)	416 (100%)	2 (0%)	81	56
1	C	416/412 (101%)	414 (100%)	2 (0%)	81	56
1	D	410/412 (100%)	407 (99%)	3 (1%)	76	47
All	All	1662/1648 (101%)	1653 (100%)	9 (0%)	86	56

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

Mol	Chain	Res	Type
1	A	478[A]	LEU
1	A	478[B]	LEU
1	B	262[A]	MET
1	B	262[B]	MET
1	C	24[A]	VAL
1	C	24[B]	VAL
1	D	430	LEU
1	D	452[A]	SER
1	D	452[B]	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	344	HIS
1	A	379	ASN
1	B	214	HIS
1	B	274	GLN
1	B	379	ASN
1	B	410	GLN
1	C	274	GLN
1	C	379	ASN
1	D	274	GLN
1	D	379	ASN
1	D	410	GLN
1	D	438	GLN

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](#)

15 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	501	-	5,5,5	0.41	0	5,5,5	0.43	0
2	GOL	B	504	-	5,5,5	0.37	0	5,5,5	0.37	0
3	IMD	C	503	-	5,5,5	0.66	0	5,5,5	0.40	0
2	GOL	A	504	-	5,5,5	0.40	0	5,5,5	0.41	0
2	GOL	C	502	-	5,5,5	0.36	0	5,5,5	0.44	0
2	GOL	D	501	-	5,5,5	0.35	0	5,5,5	0.36	0
2	GOL	C	501	-	5,5,5	0.41	0	5,5,5	0.54	0
2	GOL	B	503	-	5,5,5	0.36	0	5,5,5	0.34	0
3	IMD	A	505	-	5,5,5	0.62	0	5,5,5	0.45	0
2	GOL	B	505	-	5,5,5	0.41	0	5,5,5	0.29	0
2	GOL	B	502	-	5,5,5	0.37	0	5,5,5	0.37	0
2	GOL	A	503	-	5,5,5	0.40	0	5,5,5	0.33	0
3	IMD	D	502	-	5,5,5	0.64	0	5,5,5	0.41	0
2	GOL	A	502	-	5,5,5	0.39	0	5,5,5	0.38	0
2	GOL	B	501	-	5,5,5	0.35	0	5,5,5	0.39	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	501	-	-	0/4/4/4	-
2	GOL	B	504	-	-	2/4/4/4	-
3	IMD	C	503	-	-	-	0/1/1/1
2	GOL	A	504	-	-	2/4/4/4	-
2	GOL	C	502	-	-	1/4/4/4	-
2	GOL	D	501	-	-	1/4/4/4	-
2	GOL	C	501	-	-	2/4/4/4	-
2	GOL	B	503	-	-	2/4/4/4	-
3	IMD	A	505	-	-	-	0/1/1/1
2	GOL	B	505	-	-	2/4/4/4	-
2	GOL	B	502	-	-	0/4/4/4	-
2	GOL	A	503	-	-	1/4/4/4	-
3	IMD	D	502	-	-	-	0/1/1/1
2	GOL	A	502	-	-	0/4/4/4	-
2	GOL	B	501	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	504	GOL	O1-C1-C2-C3
2	B	505	GOL	O1-C1-C2-C3
2	C	501	GOL	O1-C1-C2-O2
2	C	501	GOL	O1-C1-C2-C3
2	B	504	GOL	O1-C1-C2-O2
2	B	505	GOL	O1-C1-C2-O2
2	B	503	GOL	O1-C1-C2-O2
2	D	501	GOL	O1-C1-C2-O2
2	B	503	GOL	O1-C1-C2-C3
2	C	502	GOL	O1-C1-C2-O2
2	A	504	GOL	C1-C2-C3-O3
2	A	504	GOL	O2-C2-C3-O3
2	A	503	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	504	GOL	1	0
2	A	504	GOL	3	0
3	D	502	IMD	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	462/485 (95%)	-0.27	16 (3%)	47 49	6, 13, 30, 57	26 (5%)
1	B	462/485 (95%)	-0.31	13 (2%)	55 57	6, 13, 26, 58	25 (5%)
1	C	462/485 (95%)	-0.40	12 (2%)	57 59	5, 12, 24, 72	23 (4%)
1	D	462/485 (95%)	-0.25	15 (3%)	50 52	7, 14, 28, 69	20 (4%)
All	All	1848/1940 (95%)	-0.31	56 (3%)	52 55	5, 13, 27, 72	94 (5%)

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	350[A]	TRP	8.5
1	C	314	VAL	7.5
1	B	350[A]	TRP	7.0
1	D	314	VAL	6.8
1	A	352	TRP	6.6
1	C	352[A]	TRP	5.9
1	D	352	TRP	5.7
1	B	352[A]	TRP	5.6
1	D	313	GLY	5.5
1	C	311	PRO	5.5
1	C	313	GLY	5.2
1	C	315	GLY	5.2
1	B	315	GLY	5.1
1	D	315	GLY	5.1
1	A	478[A]	LEU	4.9
1	D	311	PRO	4.5
1	D	310	PRO	4.3
1	B	314	VAL	4.3
1	A	351	ASP	4.1
1	A	329	THR	3.9
1	A	313	GLY	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	5	HIS	3.9
1	B	329	THR	3.8
1	C	312	ASP	3.7
1	B	5	HIS	3.7
1	A	314	VAL	3.6
1	C	310	PRO	3.6
1	B	312	ASP	3.4
1	D	312	ASP	3.4
1	A	315	GLY	3.3
1	C	329	THR	3.3
1	D	329	THR	3.1
1	D	350	TRP	3.1
1	A	160	ASP	2.9
1	B	313	GLY	2.8
1	A	384	PHE	2.8
1	A	346	ASP	2.7
1	A	312	ASP	2.6
1	C	160	ASP	2.5
1	B	478	LEU	2.5
1	B	311	PRO	2.5
1	D	4	ARG	2.5
1	D	290	ALA	2.4
1	A	6	LEU	2.4
1	D	149	ARG	2.3
1	D	99	LYS	2.3
1	A	348	THR	2.3
1	D	384[A]	PHE	2.2
1	B	160	ASP	2.2
1	C	4	ARG	2.2
1	B	346	ASP	2.1
1	D	160	ASP	2.1
1	A	290	ALA	2.1
1	A	311	PRO	2.1
1	C	6	LEU	2.1
1	B	290	ALA	2.0

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

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
2	GOL	B	504	6/6	0.78	0.21	61,62,63,63	0
2	GOL	B	505	6/6	0.82	0.18	52,53,53,54	0
3	IMD	A	505	5/5	0.83	0.13	45,45,45,45	0
3	IMD	D	502	5/5	0.84	0.21	55,56,56,56	5
2	GOL	A	503	6/6	0.86	0.16	43,43,43,43	0
2	GOL	A	504	6/6	0.86	0.15	47,47,47,48	0
2	GOL	B	503	6/6	0.87	0.14	42,42,42,42	0
2	GOL	D	501	6/6	0.87	0.13	44,45,45,45	0
2	GOL	C	502	6/6	0.88	0.13	40,41,41,41	0
2	GOL	C	501	6/6	0.89	0.13	26,32,36,38	0
3	IMD	C	503	5/5	0.90	0.15	55,55,56,56	0
2	GOL	A	501	6/6	0.95	0.08	17,21,23,24	0
2	GOL	B	502	6/6	0.96	0.07	16,19,22,22	0
2	GOL	B	501	6/6	0.97	0.06	16,21,23,25	0
2	GOL	A	502	6/6	0.98	0.06	17,21,23,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.