Full wwPDB X-ray Structure Validation Report (i)

Jun 16, 2024 – 04:17 PM EDT

PDB ID	:	4ZE4
Title	:	Structure of Gan1D, a putative 6-phospho-beta-galactosidase from Geobacillus stearothermophilus
Authors	:	Lansky, S.; Zehavi, A.; Dvir, H.; Shoham, Y.; Shoham, G.
Deposited on	:	2015-04-20
Resolution	:	1.92 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:

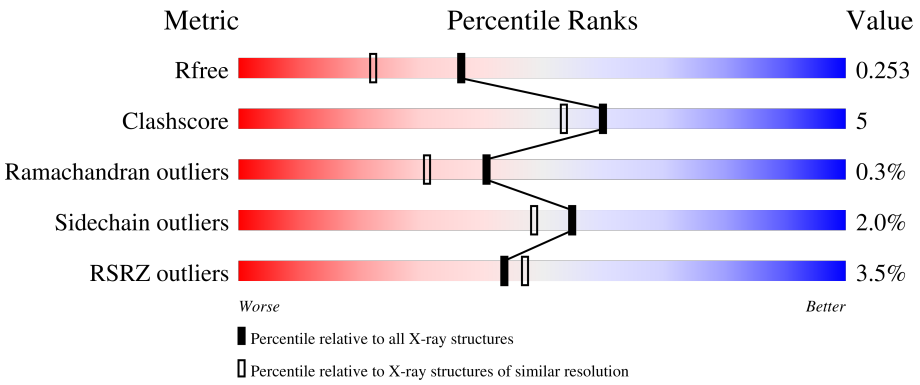
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

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

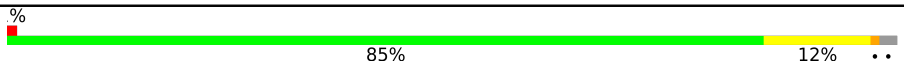
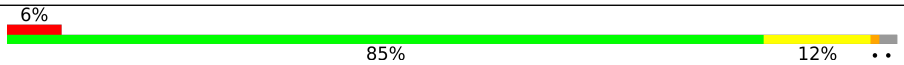
The reported resolution of this entry is 1.92 Å.

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	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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% 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	
1	B	485	

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 8358 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.

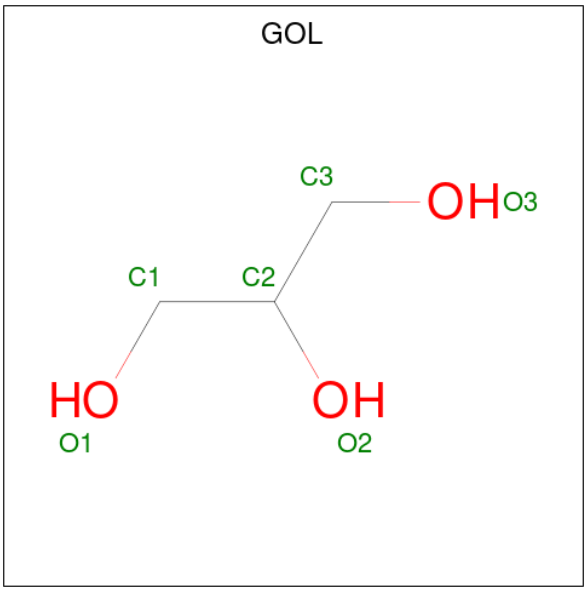
- Molecule 1 is a protein called Putative 6-phospho-beta-galactobiosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	474	3920	2518	674	717	11	0	8	0
1	B	474	3880	2490	664	715	11	0	1	0

There are 16 discrepancies between the modelled and reference sequences:

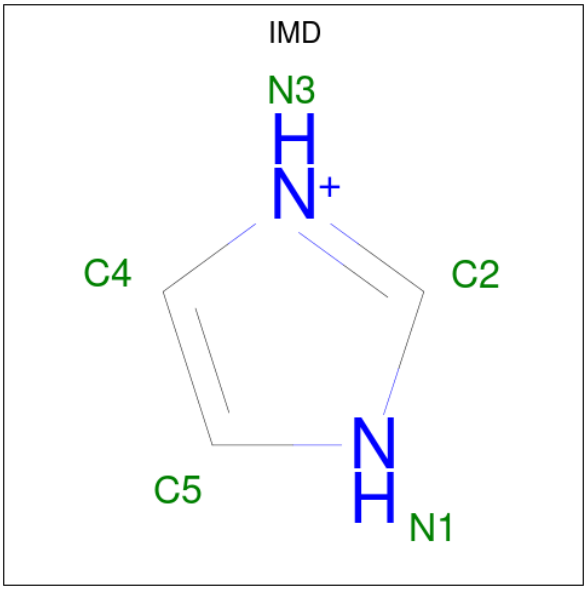
Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	-	initiating methionine	UNP W8QF82
A	-5	ILE	-	expression tag	UNP W8QF82
A	-4	HIS	-	expression tag	UNP W8QF82
A	-3	HIS	-	expression tag	UNP W8QF82
A	-2	HIS	-	expression tag	UNP W8QF82
A	-1	HIS	-	expression tag	UNP W8QF82
A	0	HIS	-	expression tag	UNP W8QF82
A	1	HIS	-	expression tag	UNP W8QF82
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

- 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
			Total	C	O		
2	B	1	6	3	3	0	0

- Molecule 3 is IMIDAZOLE (three-letter code: IMD) (formula: $C_3H_5N_2$).



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

- Molecule 4 is water.

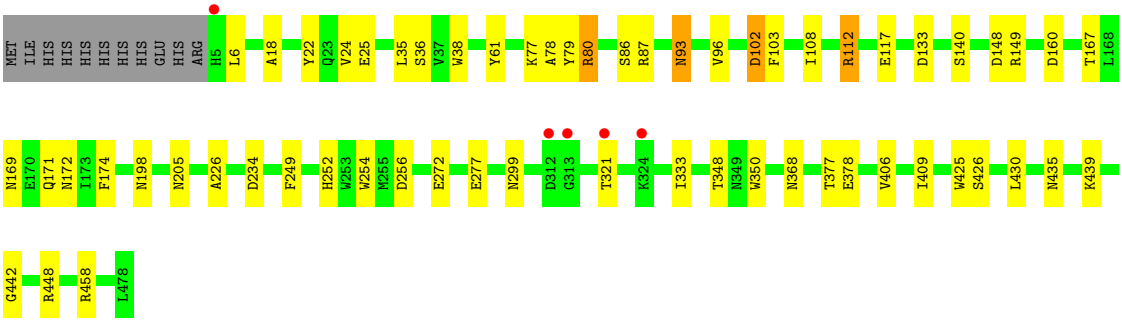
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	365	365	365	0	0
			Total	O		
4	B	171	171	171	0	0

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 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.

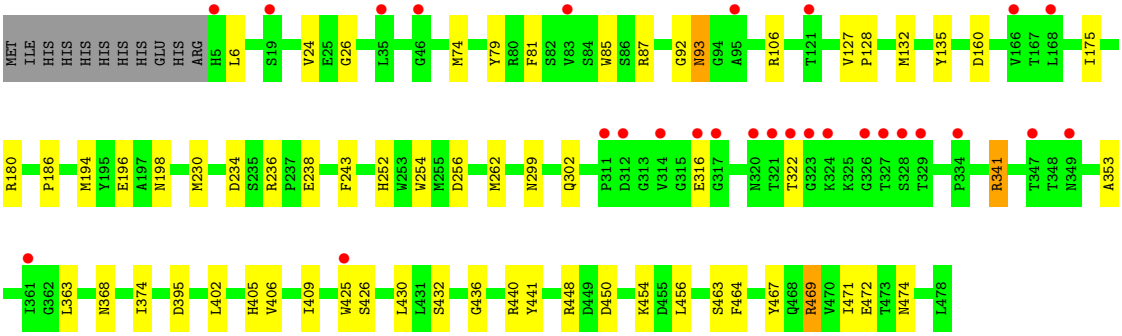
• Molecule 1: Putative 6-phospho-beta-galactobiosidase

Chain A:



• Molecule 1: Putative 6-phospho-beta-galactobiosidase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	108.01Å 68.88Å 152.88Å 90.00° 100.74° 90.00°	Depositor
Resolution (Å)	31.90 – 1.92 31.88 – 1.92	Depositor EDS
% Data completeness (in resolution range)	99.6 (31.90-1.92) 99.7 (31.88-1.92)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.17 (at 1.92Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.194 , 0.245 0.202 , 0.253	Depositor DCC
R_{free} test set	4183 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	25.5	Xtriage
Anisotropy	0.631	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 44.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	8358	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.21% 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: 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	1.19	17/4065 (0.4%)	1.06	17/5528 (0.3%)
1	B	0.76	0/4004	0.82	4/5449 (0.1%)
All	All	1.00	17/8069 (0.2%)	0.95	21/10977 (0.2%)

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	87	ARG	CZ-NH2	13.12	1.50	1.33
1	A	36	SER	CB-OG	8.92	1.53	1.42
1	A	25	GLU	CG-CD	8.59	1.64	1.51
1	A	79	TYR	C-N	-6.58	1.19	1.34
1	A	38	TRP	C-O	6.14	1.35	1.23
1	A	38	TRP	N-CA	5.75	1.57	1.46
1	A	149	ARG	CZ-NH1	5.70	1.40	1.33
1	A	93	ASN	C-O	5.64	1.34	1.23
1	A	140	SER	CB-OG	5.50	1.49	1.42
1	A	378	GLU	CD-OE1	5.41	1.31	1.25
1	A	87	ARG	CD-NE	5.31	1.55	1.46
1	A	35	LEU	C-O	5.27	1.33	1.23
1	A	22	TYR	CD1-CE1	5.22	1.47	1.39
1	A	103	PHE	N-CA	5.20	1.56	1.46
1	A	87	ARG	C-N	-5.19	1.22	1.34
1	A	93	ASN	N-CA	5.03	1.56	1.46
1	A	61	TYR	N-CA	5.03	1.56	1.46

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	80	ARG	O-C-N	9.38	137.71	122.70
1	A	80	ARG	NE-CZ-NH2	-8.46	116.07	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	87	ARG	NE-CZ-NH1	8.39	124.49	120.30
1	A	87	ARG	NE-CZ-NH2	-8.08	116.26	120.30
1	A	80	ARG	CA-C-N	-7.93	99.75	117.20
1	A	133	ASP	CB-CG-OD1	-7.79	111.28	118.30
1	A	79	TYR	C-N-CA	7.00	139.19	121.70
1	A	79	TYR	O-C-N	-6.56	112.20	122.70
1	A	22	TYR	CB-CG-CD2	-6.44	117.13	121.00
1	A	102	ASP	CB-CG-OD1	5.95	123.66	118.30
1	A	86	SER	C-N-CA	-5.93	106.88	121.70
1	A	458	ARG	NE-CZ-NH1	5.91	123.25	120.30
1	A	22	TYR	CB-CG-CD1	5.90	124.54	121.00
1	B	256	ASP	CB-CG-OD1	5.67	123.40	118.30
1	A	448	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	B	440	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	A	148	ASP	CB-CG-OD2	-5.45	113.40	118.30
1	B	106	ARG	NE-CZ-NH2	5.38	122.99	120.30
1	A	256	ASP	CB-CG-OD2	-5.32	113.51	118.30
1	B	230	MET	CG-SD-CE	-5.25	91.81	100.20
1	A	448	ARG	NE-CZ-NH1	5.06	122.83	120.30

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	3920	0	3744	41	0
1	B	3880	0	3674	34	0
2	A	6	0	8	0	0
2	B	6	0	8	0	0
3	A	10	0	10	0	0
4	A	365	0	0	10	0
4	B	171	0	0	6	0
All	All	8358	0	7444	70	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 5.

All (70) 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:112[A]:ARG:HH11	1:A:112[A]:ARG:HG3	1.09	1.16
1:B:316:GLU:HG2	4:B:602:HOH:O	1.53	1.08
1:A:112[A]:ARG:HH11	1:A:112[A]:ARG:CG	1.66	1.05
1:A:112[A]:ARG:HG3	1:A:112[A]:ARG:NH1	1.80	0.90
1:A:102:ASP:HB2	4:A:601:HOH:O	1.72	0.90
1:B:198:ASN:HD21	1:B:254:TRP:HE1	1.31	0.78
1:A:112[A]:ARG:HH11	1:A:112[A]:ARG:CB	1.99	0.76
1:B:469:ARG:NH1	1:B:472:GLU:OE1	2.24	0.71
1:A:198:ASN:HD21	1:A:254:TRP:HE1	1.40	0.70
1:A:252:HIS:HD2	1:B:368:ASN:HD21	1.47	0.63
1:A:252:HIS:CD2	1:B:368:ASN:HD21	2.17	0.62
1:A:112[A]:ARG:NH1	1:A:112[A]:ARG:CB	2.66	0.56
1:B:175:ILE:HG22	1:B:194:MET:HG3	1.87	0.55
1:B:180:ARG:NH1	4:B:602:HOH:O	2.22	0.55
1:B:132:MET:CE	1:B:186:PRO:HB3	2.36	0.55
1:B:448:ARG:NH1	4:B:613:HOH:O	2.38	0.55
1:A:18:ALA:HA	1:A:80:ARG:O	2.07	0.54
1:A:102:ASP:CB	4:A:601:HOH:O	2.42	0.54
1:B:402:LEU:O	1:B:406:VAL:HG23	2.08	0.54
1:A:112[B]:ARG:NH2	4:A:605:HOH:O	2.37	0.53
1:B:236:ARG:HD3	1:B:238:GLU:OE1	2.09	0.53
1:A:77:LYS:HE2	4:A:918:HOH:O	2.09	0.53
1:B:467:TYR:O	1:B:471:ILE:HG13	2.10	0.52
1:A:112[B]:ARG:NE	4:A:605:HOH:O	2.39	0.51
1:A:160:ASP:HB2	4:A:904:HOH:O	2.10	0.51
1:A:350:TRP:CE2	1:A:439[A]:LYS:HD2	2.45	0.51
1:A:272:GLU:HA	1:A:277:GLU:OE2	2.11	0.50
1:B:132:MET:HE3	1:B:186:PRO:HB3	1.94	0.49
1:A:368:ASN:HD21	1:B:252:HIS:CD2	2.31	0.48
1:A:77:LYS:HE3	4:A:674:HOH:O	2.12	0.48
1:A:80:ARG:HH22	1:A:169:ASN:HD22	1.61	0.48
1:B:405:HIS:O	1:B:409:ILE:HG13	2.14	0.47
1:B:93:ASN:C	1:B:93:ASN:HD22	2.18	0.47
1:B:363:LEU:CD2	1:B:374:ILE:HG21	2.44	0.47
1:B:302:GLN:HB2	1:B:353:ALA:HB3	1.96	0.47
1:B:432:SER:HG	1:B:436:GLY:N	2.11	0.47
1:A:96:VAL:CG1	4:A:951:HOH:O	2.63	0.46
1:B:26:GLY:O	1:B:87:ARG:NH2	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112[A]:ARG:NH1	1:A:112[A]:ARG:HB3	2.31	0.46
1:A:426:SER:O	1:A:442:GLY:HA2	2.16	0.46
1:A:368:ASN:HD21	1:B:252:HIS:HD2	1.63	0.45
1:A:108:ILE:O	1:A:112[A]:ARG:HG2	2.16	0.45
1:B:395:ASP:OD2	1:B:463:SER:OG	2.23	0.45
1:B:198:ASN:ND2	1:B:254:TRP:HE1	2.05	0.45
1:A:425:TRP:HA	1:A:426:SER:HA	1.78	0.45
1:B:425:TRP:HD1	1:B:441:TYR:CD2	2.36	0.45
1:A:167:THR:OG1	1:A:205:ASN:ND2	2.51	0.44
1:B:450:ASP:OD2	4:B:601:HOH:O	2.21	0.44
1:A:112[A]:ARG:CG	1:A:112[A]:ARG:NH1	2.38	0.44
1:A:169:ASN:HD21	1:A:299:ASN:HD21	1.66	0.43
1:A:234:ASP:HA	1:B:234:ASP:HA	2.00	0.43
1:B:448:ARG:HB3	1:B:456:LEU:HD23	2.00	0.43
1:A:160:ASP:HB2	4:A:827:HOH:O	2.18	0.43
1:A:321:THR:HG22	1:A:435:ASN:HD21	1.83	0.43
1:A:171:GLN:HA	1:A:174:PHE:CE2	2.54	0.43
1:A:321:THR:CG2	1:A:435:ASN:HD21	2.32	0.43
1:A:348:THR:HB	4:A:619:HOH:O	2.19	0.42
1:B:85:TRP:CZ3	1:B:127:VAL:HG23	2.54	0.42
1:A:172:ASN:HD21	1:A:226:ALA:HB3	1.84	0.42
1:B:74:MET:HB2	1:B:464:PHE:CZ	2.55	0.41
1:B:79:TYR:CE2	1:B:81:PHE:HB3	2.55	0.41
1:B:196:GLU:HG2	4:B:705:HOH:O	2.21	0.41
1:B:341:ARG:HG2	4:B:735:HOH:O	2.20	0.41
1:A:78:ALA:HA	1:A:117:GLU:O	2.21	0.41
1:A:249:PHE:CE2	1:A:333:ILE:CD1	3.04	0.41
1:B:92:GLY:HA2	1:B:128:PRO:HG2	2.03	0.41
1:A:80:ARG:HH22	1:A:169:ASN:ND2	2.18	0.40
1:A:406:VAL:HA	1:A:409:ILE:HD12	2.03	0.40
1:A:299:ASN:OD1	1:A:377:THR:OG1	2.23	0.40
1:B:425:TRP:HA	1:B:426:SER:HA	1.80	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	480/485 (99%)	462 (96%)	18 (4%)	0	100	100
1	B	473/485 (98%)	446 (94%)	24 (5%)	3 (1%)	25	14
All	All	953/970 (98%)	908 (95%)	42 (4%)	3 (0%)	41	31

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	454	LYS
1	B	474	ASN
1	B	135	TYR

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	409/412 (99%)	403 (98%)	6 (2%)	65	61
1	B	402/412 (98%)	391 (97%)	11 (3%)	44	36
All	All	811/824 (98%)	794 (98%)	17 (2%)	55	46

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	A	24	VAL
1	A	93	ASN
1	A	112[A]	ARG
1	A	112[B]	ARG
1	A	430	LEU
1	B	6	LEU
1	B	24	VAL
1	B	93	ASN

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Mol	Chain	Res	Type
1	B	160	ASP
1	B	243	PHE
1	B	262	MET
1	B	299	ASN
1	B	322	THR
1	B	341	ARG
1	B	430	LEU
1	B	469	ARG

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	93	ASN
1	A	129	GLN
1	A	169	ASN
1	A	171	GLN
1	A	198	ASN
1	A	202	ASN
1	A	205	ASN
1	A	251	ASN
1	A	252	HIS
1	A	265	GLN
1	A	269	ASN
1	A	274	GLN
1	A	349	ASN
1	A	379	ASN
1	A	435	ASN
1	B	93	ASN
1	B	171	GLN
1	B	184	HIS
1	B	198	ASN
1	B	202	ASN
1	B	205	ASN
1	B	252	HIS
1	B	269	ASN
1	B	371	GLN
1	B	394	ASN

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)

4 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	IMD	A	503	-	3,5,5	0.58	0	4,5,5	0.28	0
2	GOL	A	501	-	5,5,5	0.35	0	5,5,5	0.29	0
3	IMD	A	502	-	3,5,5	0.30	0	4,5,5	0.48	0
2	GOL	B	501	-	5,5,5	0.39	0	5,5,5	0.44	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	IMD	A	503	-	-	-	0/1/1/1
2	GOL	A	501	-	-	3/4/4/4	-
3	IMD	A	502	-	-	-	0/1/1/1
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 (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	GOL	C1-C2-C3-O3
2	A	501	GOL	O2-C2-C3-O3
2	A	501	GOL	O1-C1-C2-C3

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	79:TYR	C	80:ARG	N	1.18

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	474/485 (97%)	-0.09	5 (1%)	80	82	11, 24, 48, 84	0
1	B	474/485 (97%)	0.40	28 (5%)	22	25	27, 39, 60, 91	0
All	All	948/970 (97%)	0.16	33 (3%)	44	47	11, 33, 56, 91	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	326	GLY	4.9
1	B	322	THR	4.4
1	B	95	ALA	4.4
1	A	5	HIS	4.3
1	B	321	THR	4.2
1	B	320	ASN	3.6
1	B	46	GLY	3.3
1	B	5	HIS	3.1
1	A	324	LYS	3.0
1	B	327	THR	2.9
1	B	168	LEU	2.9
1	B	324	LYS	2.8
1	A	312	ASP	2.8
1	B	312	ASP	2.8
1	B	121	THR	2.7
1	B	323	GLY	2.7
1	A	321	THR	2.6
1	B	425	TRP	2.5
1	B	317	GLY	2.4
1	B	334	PRO	2.4
1	B	316	GLU	2.4
1	B	349	ASN	2.3
1	B	19	SER	2.3
1	B	328	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	313	GLY	2.2
1	B	329	THR	2.2
1	B	347	THR	2.2
1	B	83	VAL	2.2
1	B	311	PRO	2.1
1	B	314	VAL	2.1
1	B	166	VAL	2.0
1	B	35	LEU	2.0
1	B	361	ILE	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 monosaccharides 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	501	6/6	0.59	0.24	65,71,74,75	0
3	IMD	A	502	5/5	0.89	0.23	32,37,39,42	0
3	IMD	A	503	5/5	0.91	0.17	29,29,34,35	0
2	GOL	A	501	6/6	0.92	0.17	51,54,60,61	0

6.5 Other polymers (i)

There are no such residues in this entry.