



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:12 PM UTC

PDB ID : 6O41 / pdb_00006o41
Title : Crystal structure of the unbound PGZL1 germline Fab fragment (PGZL1_gVmDmJ)
Authors : Irimia, A.; Wilson, I.A.
Deposited on : 2019-02-27
Resolution : 2.46 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:

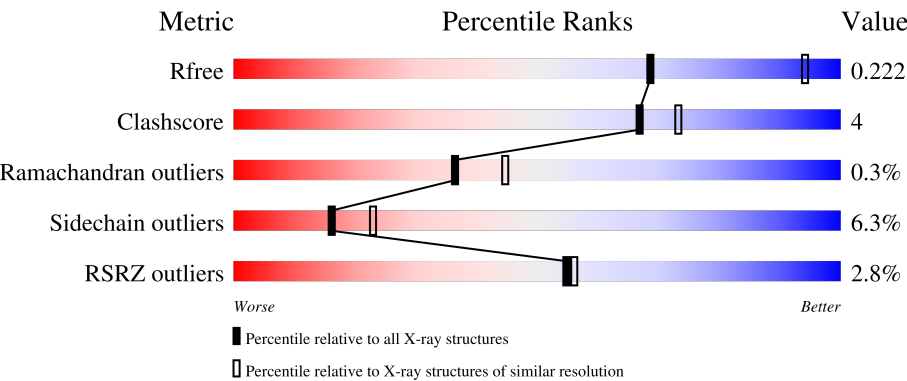
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry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

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.46 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	215	89%8%..
1	C	215	90%8%..
1	L	215	88%10%. .
2	B	227	4% 88%9%.
2	D	227	5% 82%12%. .

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Mol	Chain	Length	Quality of chain
2	H	227	
3	M	66	
3	N	66	
3	O	66	

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
4	GOL	L	305	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11744 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 germline PGZL1_gVmDmJ light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	213	Total	C	N	O	S	0	4	0
			1668	1037	291	335	5			
1	A	213	Total	C	N	O	S	0	1	0
			1635	1019	279	332	5			
1	C	213	Total	C	N	O	S	0	1	0
			1635	1019	279	332	5			

- Molecule 2 is a protein called germline PGZL1_gVmDmJ heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	223	Total	C	N	O	S	0	4	0
			1652	1044	275	324	9			
2	B	222	Total	C	N	O	S	0	3	0
			1643	1039	274	321	9			
2	D	222	Total	C	N	O	S	0	3	0
			1639	1036	273	321	9			

- Molecule 3 is a protein called Immunoglobulin G-binding protein G (DOMAIN III).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	M	60	Total	C	N	O	0	0	0
			465	292	76	97			
3	N	60	Total	C	N	O	0	1	0
			473	296	78	99			
3	O	59	Total	C	N	O	0	0	0
			460	289	75	96			

There are 18 discrepancies between the modelled and reference sequences:

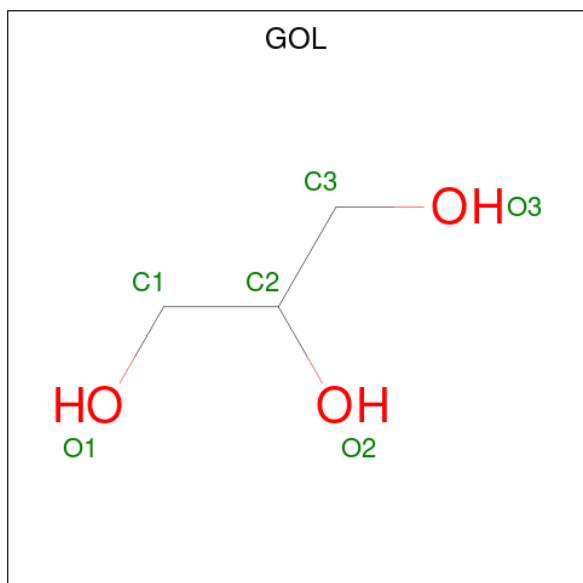
Chain	Residue	Modelled	Actual	Comment	Reference
M	1	MET	-	initiating methionine	UNP P19909
M	62	HIS	-	expression tag	UNP P19909

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Chain	Residue	Modelled	Actual	Comment	Reference
M	63	HIS	-	expression tag	UNP P19909
M	64	HIS	-	expression tag	UNP P19909
M	65	HIS	-	expression tag	UNP P19909
M	66	HIS	-	expression tag	UNP P19909
N	1	MET	-	initiating methionine	UNP P19909
N	62	HIS	-	expression tag	UNP P19909
N	63	HIS	-	expression tag	UNP P19909
N	64	HIS	-	expression tag	UNP P19909
N	65	HIS	-	expression tag	UNP P19909
N	66	HIS	-	expression tag	UNP P19909
O	1	MET	-	initiating methionine	UNP P19909
O	62	HIS	-	expression tag	UNP P19909
O	63	HIS	-	expression tag	UNP P19909
O	64	HIS	-	expression tag	UNP P19909
O	65	HIS	-	expression tag	UNP P19909
O	66	HIS	-	expression tag	UNP P19909

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	L	1	Total	C	O	0	0
			6	3	3		
4	L	1	Total	C	O	0	0
			6	3	3		
4	H	1	Total	C	O	0	0
			6	3	3		
4	H	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		
4	O	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is water.

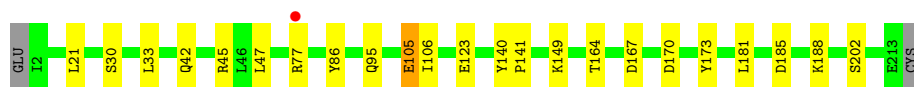
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	79	Total	O	0	0
			79	79		
5	H	58	Total	O	0	0
			58	58		
5	A	59	Total	O	0	0
			59	59		
5	B	47	Total	O	0	0
			47	47		
5	C	72	Total	O	0	0
			72	72		
5	D	42	Total	O	0	0
			42	42		
5	M	16	Total	O	0	0
			16	16		
5	N	13	Total	O	0	0
			13	13		
5	O	10	Total	O	0	0
			10	10		

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.

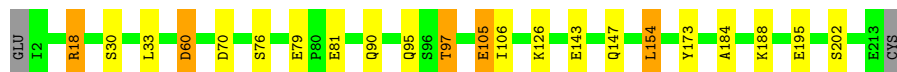
- Molecule 1: germline PGZL1_gVmDmJ light chain

Chain L: 



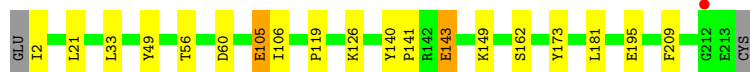
- Molecule 1: germline PGZL1_gVmDmJ light chain

Chain A: 



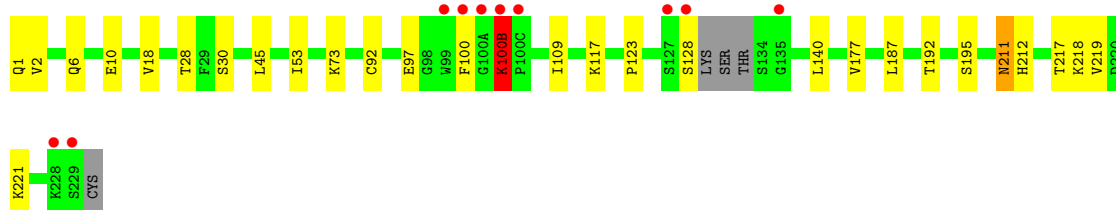
- Molecule 1: germline PGZL1_gVmDmJ light chain

Chain C: 



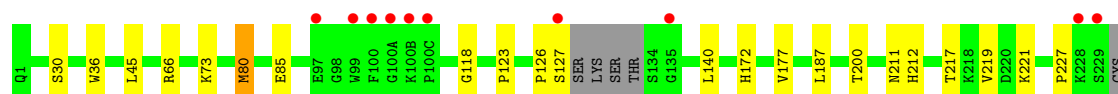
- Molecule 2: germline PGZL1_gVmDmJ heavy chain

Chain H: 

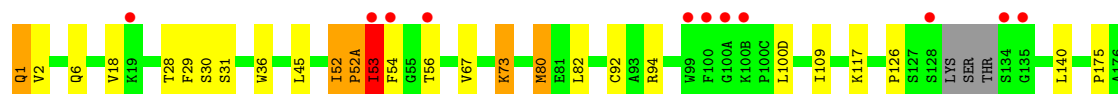
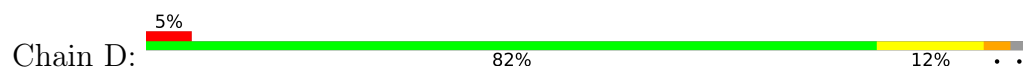


- Molecule 2: germline PGZL1_gVmDmJ heavy chain

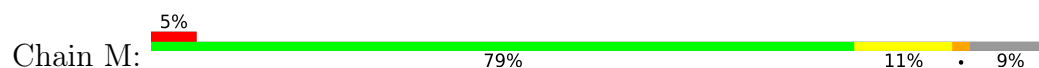
Chain B: 



- Molecule 2: germline PGZL1_gVmDmJ heavy chain



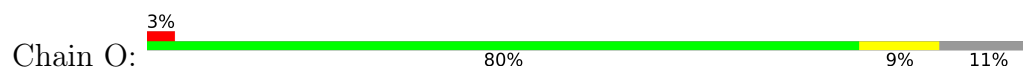
- Molecule 3: Immunoglobulin G-binding protein G (DOMAIN III)



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.84Å 107.49Å 131.96Å 90.00° 94.61° 90.00°	Depositor
Resolution (Å)	39.67 – 2.46 39.67 – 2.46	Depositor EDS
% Data completeness (in resolution range)	97.8 (39.67-2.46) 97.7 (39.67-2.46)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.48Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.169 , 0.219 0.173 , 0.222	Depositor DCC
R_{free} test set	3872 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	50.0	Xtriage
Anisotropy	0.211	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11744	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.12% 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

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	0.24	0/1672	0.47	0/2269
1	C	0.23	0/1672	0.47	0/2269
1	L	0.24	0/1705	0.47	0/2311
2	B	0.23	0/1689	0.44	0/2295
2	D	0.20	0/1685	0.43	1/2291 (0.0%)
2	H	0.25	0/1701	0.46	0/2311
3	M	0.23	0/473	0.40	0/643
3	N	0.19	0/481	0.40	0/654
3	O	0.20	0/468	0.36	0/636
All	All	0.23	0/11546	0.45	1/15679 (0.0%)

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
2	H	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	53	ILE	N-CA-C	-5.85	107.20	112.12

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	100(B)	LYS	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	1635	0	1587	10	0
1	C	1635	0	1587	10	0
1	L	1668	0	1623	16	0
2	B	1643	0	1625	10	0
2	D	1639	0	1614	20	0
2	H	1652	0	1635	9	0
3	M	465	0	444	2	0
3	N	473	0	449	7	0
3	O	460	0	442	1	0
4	A	6	0	8	0	0
4	B	6	0	8	1	0
4	C	12	0	16	0	0
4	D	6	0	8	2	0
4	H	12	0	16	1	0
4	L	30	0	40	7	0
4	O	6	0	8	2	0
5	A	59	0	0	1	0
5	B	47	0	0	0	0
5	C	72	0	0	0	0
5	D	42	0	0	3	0
5	H	58	0	0	1	0
5	L	79	0	0	5	0
5	M	16	0	0	0	0
5	N	13	0	0	0	0
5	O	10	0	0	0	0
All	All	11744	0	11110	82	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 4.

All (82) 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)
2:D:30:SER:HB2	2:D:53:ILE:HG12	1.63	0.80
1:L:45:ARG:HD3	4:L:305:GOL:H11	1.66	0.76
2:B:126:PRO:HG3	2:B:140:LEU:HB3	1.71	0.73
1:A:60:ASP:OD1	1:A:60:ASP:N	2.25	0.69
2:B:118:GLY:H	4:O:101:GOL:H31	1.57	0.69
2:D:126:PRO:HG3	2:D:140:LEU:HB3	1.78	0.65
1:L:77[A]:ARG:NH1	5:L:403:HOH:O	2.30	0.65
1:A:18:ARG:HD2	1:A:76:SER:HA	1.79	0.64
2:H:100:PHE:O	2:H:100(B):LYS:NZ	2.30	0.64
1:L:185:ASP:HA	1:L:188:LYS:HE2	1.82	0.61
1:C:105:GLU:HG2	1:C:106:ILE:N	2.16	0.60
2:D:1:GLN:HG2	2:D:2:VAL:H	1.69	0.58
3:N:14:GLY:O	3:N:18:LYS:NZ	2.36	0.58
1:C:126:LYS:HD2	3:N:43:GLY:HA2	1.86	0.57
1:L:77[B]:ARG:NH1	2:B:200:THR:O	2.37	0.57
2:D:94:ARG:NH2	5:D:401:HOH:O	2.37	0.56
1:L:45:ARG:NH2	5:L:405:HOH:O	2.37	0.56
1:L:42:GLN:OE1	4:L:305:GOL:H31	2.06	0.55
1:A:105:GLU:HG2	1:A:106:ILE:N	2.23	0.54
1:A:147:GLN:OE1	1:A:154:LEU:HD13	2.09	0.53
2:D:36:TRP:CE2	2:D:80:MET:HB3	2.43	0.53
2:B:36:TRP:CE2	2:B:80:MET:HB3	2.44	0.52
1:L:95:GLN:NE2	5:L:402:HOH:O	2.30	0.52
2:D:30:SER:HA	2:D:52(A):PRO:HB2	1.92	0.52
2:B:172:HIS:ND1	4:B:301:GOL:H31	2.24	0.51
2:H:6:GLN:HG2	2:H:92[A]:CYS:SG	2.51	0.51
1:L:105:GLU:HG2	1:L:106:ILE:N	2.24	0.51
3:N:7:THR:OG1	3:N:24:LYS:NZ	2.43	0.50
1:C:60:ASP:OD1	1:C:60:ASP:N	2.42	0.50
4:D:301:GOL:H31	3:N:38:TYR:CD1	2.47	0.50
2:H:117:LYS:NZ	5:H:402:HOH:O	2.44	0.50
2:D:18:VAL:HG21	2:D:109:ILE:HD12	1.93	0.49
2:D:222:LYS:NZ	2:D:226:GLU:OE2	2.45	0.49
1:L:105:GLU:HG3	1:L:173:TYR:OH	2.12	0.49
2:D:6:GLN:HG2	2:D:92[A]:CYS:SG	2.53	0.49
2:H:187:LEU:HD22	4:H:302:GOL:H12	1.94	0.48
1:A:90:GLN:HE21	1:A:97:THR:HG23	1.79	0.48
1:L:170:ASP:HA	4:L:304:GOL:H31	1.95	0.48
1:C:162:SER:HB2	2:D:175:PRO:HG2	1.95	0.48
2:H:123:PRO:HD3	2:H:221:LYS:HE2	1.96	0.47
3:M:25:ALA:HB1	3:M:30:THR:HB	1.96	0.47
2:D:211:ASN:ND2	5:D:404:HOH:O	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:52:ILE:HG13	2:D:54:PHE:HD1	1.80	0.47
1:A:105:GLU:HG3	1:A:173:TYR:OH	2.16	0.46
2:B:118:GLY:N	4:O:101:GOL:H31	2.25	0.46
2:H:212:HIS:HB3	2:H:217:THR:HB	1.98	0.46
1:A:81:GLU:N	1:A:81:GLU:OE1	2.49	0.46
1:C:119:PRO:HB3	1:C:209:PHE:CE2	2.51	0.46
2:H:97:GLU:OE1	2:H:97:GLU:N	2.49	0.45
3:M:6:THR:HG23	3:M:25:ALA:O	2.16	0.45
2:H:211:ASN:OD1	2:H:218:LYS:HG2	2.17	0.45
2:D:73:LYS:HE3	2:D:73:LYS:HB2	1.78	0.45
2:D:218:LYS:HB2	2:D:218:LYS:HE3	1.64	0.45
1:A:147:GLN:HB3	1:A:195:GLU:HB3	1.99	0.44
2:D:1:GLN:HE21	2:D:1:GLN:N	2.16	0.44
2:D:211:ASN:OD1	2:D:218:LYS:HG3	2.17	0.44
4:D:301:GOL:H12	3:N:22:THR:O	2.17	0.44
1:C:49:TYR:CG	2:D:100(D):LEU:HD21	2.53	0.44
1:L:164:THR:HB	4:L:302:GOL:H31	1.98	0.44
1:A:95:GLN:OE1	5:A:401:HOH:O	2.21	0.44
2:D:67:VAL:HG22	2:D:82:LEU:HD13	1.99	0.44
2:D:117:LYS:NZ	5:D:405:HOH:O	2.51	0.44
2:B:212:HIS:HB3	2:B:217:THR:HB	2.00	0.44
2:B:123:PRO:HD3	2:B:221:LYS:HE2	2.01	0.43
2:H:18:VAL:HG21	2:H:109:ILE:HD12	2.01	0.43
3:O:44:VAL:HG13	3:O:61:GLU:HG3	2.00	0.43
3:N:11:VAL:HG22	3:N:58:THR:HG23	2.01	0.43
1:C:143:GLU:CD	1:C:143:GLU:H	2.27	0.43
1:C:149:LYS:NZ	1:C:195:GLU:OE1	2.33	0.42
2:B:126:PRO:HD2	2:B:227:PRO:HA	2.00	0.42
1:L:149:LYS:NZ	5:L:411:HOH:O	2.52	0.42
1:L:45:ARG:HB2	4:L:305:GOL:H2	2.00	0.42
2:D:1:GLN:HE21	2:D:1:GLN:H1	1.65	0.41
1:C:105:GLU:HG3	1:C:173:TYR:OH	2.19	0.41
1:L:140:TYR:CG	1:L:141:PRO:HA	2.55	0.41
2:B:73:LYS:H	2:B:73:LYS:HD2	1.85	0.41
1:L:47:LEU:HD11	1:L:86:TYR:HE2	1.86	0.41
4:L:305:GOL:O1	5:L:401:HOH:O	2.22	0.41
1:C:140:TYR:CG	1:C:141:PRO:HA	2.56	0.41
3:N:10:LEU:HD11	3:N:59:VAL:HG12	2.03	0.41
1:A:184:ALA:O	1:A:188:LYS:HG3	2.21	0.40
1:L:167:ASP:HA	4:L:302:GOL:H12	2.03	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	212/215 (99%)	208 (98%)	4 (2%)	0	100	100
1	C	212/215 (99%)	207 (98%)	5 (2%)	0	100	100
1	L	215/215 (100%)	210 (98%)	5 (2%)	0	100	100
2	B	221/227 (97%)	213 (96%)	8 (4%)	0	100	100
2	D	221/227 (97%)	206 (93%)	14 (6%)	1 (0%)	24	32
2	H	223/227 (98%)	215 (96%)	6 (3%)	2 (1%)	14	17
3	M	58/66 (88%)	56 (97%)	1 (2%)	1 (2%)	7	6
3	N	59/66 (89%)	56 (95%)	3 (5%)	0	100	100
3	O	57/66 (86%)	56 (98%)	1 (2%)	0	100	100
All	All	1478/1524 (97%)	1427 (96%)	47 (3%)	4 (0%)	36	45

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	100(B)	LYS
2	D	29	PHE
3	M	63	HIS
2	H	30	SER

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	185/186 (100%)	173 (94%)	12 (6%)	15	21
1	C	185/186 (100%)	178 (96%)	7 (4%)	29	43
1	L	188/186 (101%)	181 (96%)	7 (4%)	30	44
2	B	183/187 (98%)	173 (94%)	10 (6%)	19	28
2	D	182/187 (97%)	167 (92%)	15 (8%)	10	13
2	H	185/187 (99%)	171 (92%)	14 (8%)	12	15
3	M	49/55 (89%)	44 (90%)	5 (10%)	7	6
3	N	50/55 (91%)	45 (90%)	5 (10%)	7	7
3	O	49/55 (89%)	45 (92%)	4 (8%)	10	13
All	All	1256/1284 (98%)	1177 (94%)	79 (6%)	16	23

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	21	LEU
1	L	30	SER
1	L	33	LEU
1	L	105	GLU
1	L	123	GLU
1	L	181	LEU
1	L	202	SER
2	H	1	GLN
2	H	2	VAL
2	H	10	GLU
2	H	28	THR
2	H	45	LEU
2	H	53	ILE
2	H	73	LYS
2	H	128	SER
2	H	140	LEU
2	H	177	VAL
2	H	192	THR
2	H	195	SER
2	H	211	ASN
2	H	219	VAL
1	A	18	ARG
1	A	30	SER
1	A	33	LEU
1	A	60	ASP
1	A	70	ASP

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Mol	Chain	Res	Type
1	A	79	GLU
1	A	97	THR
1	A	105	GLU
1	A	126	LYS
1	A	143	GLU
1	A	154	LEU
1	A	202	SER
2	B	30	SER
2	B	45	LEU
2	B	66	ARG
2	B	80	MET
2	B	85	GLU
2	B	127	SER
2	B	177	VAL
2	B	187	LEU
2	B	211	ASN
2	B	219	VAL
1	C	2	ILE
1	C	21	LEU
1	C	33	LEU
1	C	56	THR
1	C	105	GLU
1	C	143	GLU
1	C	181	LEU
2	D	1	GLN
2	D	28	THR
2	D	31	SER
2	D	45	LEU
2	D	52	ILE
2	D	52(A)	PRO
2	D	53	ILE
2	D	56	THR
2	D	73	LYS
2	D	80	MET
2	D	177	VAL
2	D	192	THR
2	D	211	ASN
2	D	219	VAL
2	D	228	LYS
3	M	5	VAL
3	M	6	THR
3	M	11	VAL

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Mol	Chain	Res	Type
3	M	13	ASN
3	M	26	VAL
3	N	5	VAL
3	N	11	VAL
3	N	13[A]	ASN
3	N	13[B]	ASN
3	N	23	THR
3	O	5	VAL
3	O	6	THR
3	O	11	VAL
3	O	26	VAL

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

Mol	Chain	Res	Type
1	L	37	GLN
1	L	160	GLN
1	A	27	GLN
1	A	160	GLN
2	B	203	GLN
1	C	160	GLN
2	D	1	GLN
2	D	58	ASN
2	D	172	HIS
2	D	203	GLN
3	N	62	HIS
3	O	62	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry

13 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	$\# Z > 2$
4	GOL	L	304	-	5,5,5	0.94	0	5,5,5	1.09	0
4	GOL	O	101	-	5,5,5	0.90	0	5,5,5	0.90	0
4	GOL	H	301	-	5,5,5	1.08	0	5,5,5	1.02	0
4	GOL	C	302	-	5,5,5	1.01	0	5,5,5	1.09	0
4	GOL	C	301	-	5,5,5	0.82	0	5,5,5	1.29	1 (20%)
4	GOL	L	302	-	5,5,5	1.02	0	5,5,5	1.00	0
4	GOL	L	301	-	5,5,5	0.97	0	5,5,5	1.05	0
4	GOL	A	301	-	5,5,5	0.74	0	5,5,5	1.29	1 (20%)
4	GOL	D	301	-	5,5,5	0.84	0	5,5,5	1.29	1 (20%)
4	GOL	B	301	-	5,5,5	1.04	0	5,5,5	0.96	0
4	GOL	H	302	-	5,5,5	1.09	0	5,5,5	1.05	0
4	GOL	L	303	-	5,5,5	0.96	0	5,5,5	1.03	0
4	GOL	L	305	-	5,5,5	1.06	0	5,5,5	0.91	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
4	GOL	L	304	-	-	0/4/4/4	-
4	GOL	O	101	-	-	1/4/4/4	-
4	GOL	H	301	-	-	2/4/4/4	-
4	GOL	C	302	-	-	1/4/4/4	-
4	GOL	C	301	-	-	1/4/4/4	-
4	GOL	L	302	-	-	2/4/4/4	-
4	GOL	L	301	-	-	0/4/4/4	-
4	GOL	A	301	-	-	4/4/4/4	-
4	GOL	D	301	-	-	2/4/4/4	-
4	GOL	B	301	-	-	1/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	H	302	-	-	2/4/4/4	-
4	GOL	L	303	-	-	3/4/4/4	-
4	GOL	L	305	-	-	2/4/4/4	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	301	GOL	C3-C2-C1	-2.37	103.11	111.80
4	C	301	GOL	C3-C2-C1	-2.12	104.01	111.80
4	A	301	GOL	C3-C2-C1	-2.01	104.42	111.80

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	L	303	GOL	O1-C1-C2-O2
4	L	303	GOL	O1-C1-C2-C3
4	H	301	GOL	C1-C2-C3-O3
4	H	302	GOL	O1-C1-C2-C3
4	A	301	GOL	O1-C1-C2-C3
4	A	301	GOL	C1-C2-C3-O3
4	D	301	GOL	C1-C2-C3-O3
4	L	302	GOL	O1-C1-C2-O2
4	A	301	GOL	O2-C2-C3-O3
4	L	302	GOL	O1-C1-C2-C3
4	L	305	GOL	C1-C2-C3-O3
4	L	305	GOL	O2-C2-C3-O3
4	H	302	GOL	O1-C1-C2-O2
4	A	301	GOL	O1-C1-C2-O2
4	D	301	GOL	O2-C2-C3-O3
4	L	303	GOL	O2-C2-C3-O3
4	H	301	GOL	O2-C2-C3-O3
4	C	302	GOL	O2-C2-C3-O3
4	O	101	GOL	O1-C1-C2-O2
4	B	301	GOL	O1-C1-C2-C3
4	C	301	GOL	O1-C1-C2-O2

There are no ring outliers.

7 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	L	304	GOL	1	0
4	O	101	GOL	2	0
4	L	302	GOL	2	0
4	D	301	GOL	2	0
4	B	301	GOL	1	0
4	H	302	GOL	1	0
4	L	305	GOL	4	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	213/215 (99%)	-0.35	0 100 100	31, 48, 80, 116	1 (0%)
1	C	213/215 (99%)	-0.53	1 (0%) 87 88	31, 45, 66, 99	1 (0%)
1	L	213/215 (99%)	-0.53	1 (0%) 87 88	21, 43, 65, 91	4 (1%)
2	B	222/227 (97%)	-0.22	10 (4%) 38 37	27, 47, 87, 123	4 (1%)
2	D	222/227 (97%)	-0.01	12 (5%) 31 29	27, 55, 106, 150	5 (2%)
2	H	223/227 (98%)	-0.23	10 (4%) 38 37	23, 45, 90, 152	6 (2%)
3	M	60/66 (90%)	-0.07	3 (5%) 34 32	33, 52, 80, 114	0
3	N	60/66 (90%)	-0.17	3 (5%) 34 32	24, 53, 80, 110	1 (1%)
3	O	59/66 (89%)	-0.22	2 (3%) 48 48	36, 54, 77, 111	0
All	All	1485/1524 (97%)	-0.29	42 (2%) 55 56	21, 47, 89, 152	22 (1%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	128	SER	6.8
2	H	100	PHE	5.4
2	B	99	TRP	5.2
2	H	99	TRP	5.1
2	D	128	SER	5.0
3	O	5	VAL	4.7
2	D	99	TRP	4.6
3	N	64	HIS	4.5
2	B	100	PHE	4.5
2	D	100	PHE	4.4
2	D	100(B)	LYS	3.9
2	B	229	SER	3.9
2	H	135	GLY	3.8
2	B	100(C)	PRO	3.8
3	M	64	HIS	3.7

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Mol	Chain	Res	Type	RSRZ
2	D	134	SER	3.7
2	D	135	GLY	3.5
3	M	63	HIS	3.5
2	H	100(A)	GLY	3.5
3	O	63	HIS	3.4
3	M	5	VAL	3.4
3	N	63	HIS	3.4
2	B	127	SER	3.2
2	B	100(B)	LYS	3.1
3	N	5	VAL	2.9
2	D	19	LYS	2.9
2	B	100(A)	GLY	2.8
2	B	135	GLY	2.7
2	D	100(A)	GLY	2.6
2	H	100(C)	PRO	2.6
2	D	228	LYS	2.6
2	H	127	SER	2.5
2	H	228	LYS	2.4
2	H	229	SER	2.3
2	D	56	THR	2.2
2	D	53	ILE	2.2
1	L	77[A]	ARG	2.2
2	B	228	LYS	2.2
2	D	54	PHE	2.0
1	C	212	GLY	2.0
2	B	97	GLU	2.0
2	H	100(B)	LYS	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 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
4	GOL	L	302	6/6	0.80	0.20	59,69,70,70	0
4	GOL	L	303	6/6	0.81	0.24	67,69,70,71	0
4	GOL	L	304	6/6	0.85	0.17	88,90,91,91	0
4	GOL	H	302	6/6	0.86	0.23	64,65,70,70	0
4	GOL	B	301	6/6	0.86	0.16	76,78,79,80	0
4	GOL	C	302	6/6	0.86	0.19	81,83,85,85	0
4	GOL	L	305	6/6	0.88	0.23	57,63,73,78	0
4	GOL	H	301	6/6	0.91	0.14	58,64,68,70	0
4	GOL	D	301	6/6	0.91	0.13	68,69,71,73	0
4	GOL	O	101	6/6	0.92	0.14	68,71,75,82	0
4	GOL	C	301	6/6	0.93	0.15	56,71,74,80	0
4	GOL	A	301	6/6	0.93	0.13	58,60,67,73	0
4	GOL	L	301	6/6	0.95	0.11	49,53,60,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.