Full wwPDB X-ray Structure Validation Report (i)

Mar 9, 2026 – 09:48 PM UTC

PDB ID	:	6BM9 / pdb_00006bm9
Title	:	Directed evolutionary changes in MBL super family - VIM-2 Round 10
Authors	:	Hong, N.-S.; Jackson, C.J.; Carr, P.D.
Deposited on	:	2017-11-13
Resolution	:	2.19 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 (1)) were used in the production of this report:

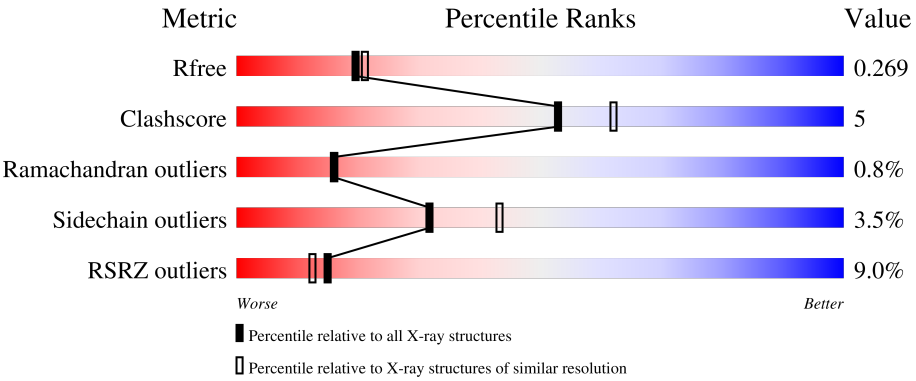
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (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 2.19 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	243	7% 79% 12% • 6%
1	B	243	6% 82% 10% • 7%
1	C	243	9% 79% 12% • 8%
1	D	243	12% 79% 14% 7%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 7053 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 Metallo-beta-lactamase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	0	2	0
			1737	1097	303	336	1			
1	B	225	Total	C	N	O	S	0	1	0
			1705	1080	296	328	1			
1	C	224	Total	C	N	O	S	0	2	0
			1698	1072	295	330	1			
1	D	225	Total	C	N	O	S	0	1	0
			1702	1075	296	330	1			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	GLY	-	expression tag	UNP Q5U7L7
A	25	HIS	-	expression tag	UNP Q5U7L7
A	26	MET	-	expression tag	UNP Q5U7L7
A	31	ARG	GLY	conflict	UNP Q5U7L7
A	36	ALA	VAL	conflict	UNP Q5U7L7
A	41	ASP	VAL	conflict	UNP Q5U7L7
A	59	ALA	THR	conflict	UNP Q5U7L7
A	61	PRO	SER	conflict	UNP Q5U7L7
A	62	LEU	PHE	conflict	UNP Q5U7L7
A	66	ALA	VAL	conflict	UNP Q5U7L7
A	144	LYS	GLU	conflict	UNP Q5U7L7
A	148	THR	ASN	conflict	UNP Q5U7L7
A	181	PRO	THR	conflict	UNP Q5U7L7
A	185	ILE	VAL	conflict	UNP Q5U7L7
A	192	ARG	SER	conflict	UNP Q5U7L7
A	213	ALA	ASP	conflict	UNP Q5U7L7
A	253	SER	THR	conflict	UNP Q5U7L7
A	254	ASP	ASN	conflict	UNP Q5U7L7
B	24	GLY	-	expression tag	UNP Q5U7L7
B	25	HIS	-	expression tag	UNP Q5U7L7
B	26	MET	-	expression tag	UNP Q5U7L7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	31	ARG	GLY	conflict	UNP Q5U7L7
B	36	ALA	VAL	conflict	UNP Q5U7L7
B	41	ASP	VAL	conflict	UNP Q5U7L7
B	59	ALA	THR	conflict	UNP Q5U7L7
B	61	PRO	SER	conflict	UNP Q5U7L7
B	62	LEU	PHE	conflict	UNP Q5U7L7
B	66	ALA	VAL	conflict	UNP Q5U7L7
B	144	LYS	GLU	conflict	UNP Q5U7L7
B	148	THR	ASN	conflict	UNP Q5U7L7
B	181	PRO	THR	conflict	UNP Q5U7L7
B	185	ILE	VAL	conflict	UNP Q5U7L7
B	192	ARG	SER	conflict	UNP Q5U7L7
B	213	ALA	ASP	conflict	UNP Q5U7L7
B	253	SER	THR	conflict	UNP Q5U7L7
B	254	ASP	ASN	conflict	UNP Q5U7L7
C	24	GLY	-	expression tag	UNP Q5U7L7
C	25	HIS	-	expression tag	UNP Q5U7L7
C	26	MET	-	expression tag	UNP Q5U7L7
C	31	ARG	GLY	conflict	UNP Q5U7L7
C	36	ALA	VAL	conflict	UNP Q5U7L7
C	41	ASP	VAL	conflict	UNP Q5U7L7
C	59	ALA	THR	conflict	UNP Q5U7L7
C	61	PRO	SER	conflict	UNP Q5U7L7
C	62	LEU	PHE	conflict	UNP Q5U7L7
C	66	ALA	VAL	conflict	UNP Q5U7L7
C	144	LYS	GLU	conflict	UNP Q5U7L7
C	148	THR	ASN	conflict	UNP Q5U7L7
C	181	PRO	THR	conflict	UNP Q5U7L7
C	185	ILE	VAL	conflict	UNP Q5U7L7
C	192	ARG	SER	conflict	UNP Q5U7L7
C	213	ALA	ASP	conflict	UNP Q5U7L7
C	253	SER	THR	conflict	UNP Q5U7L7
C	254	ASP	ASN	conflict	UNP Q5U7L7
D	24	GLY	-	expression tag	UNP Q5U7L7
D	25	HIS	-	expression tag	UNP Q5U7L7
D	26	MET	-	expression tag	UNP Q5U7L7
D	31	ARG	GLY	conflict	UNP Q5U7L7
D	36	ALA	VAL	conflict	UNP Q5U7L7
D	41	ASP	VAL	conflict	UNP Q5U7L7
D	59	ALA	THR	conflict	UNP Q5U7L7
D	61	PRO	SER	conflict	UNP Q5U7L7
D	62	LEU	PHE	conflict	UNP Q5U7L7

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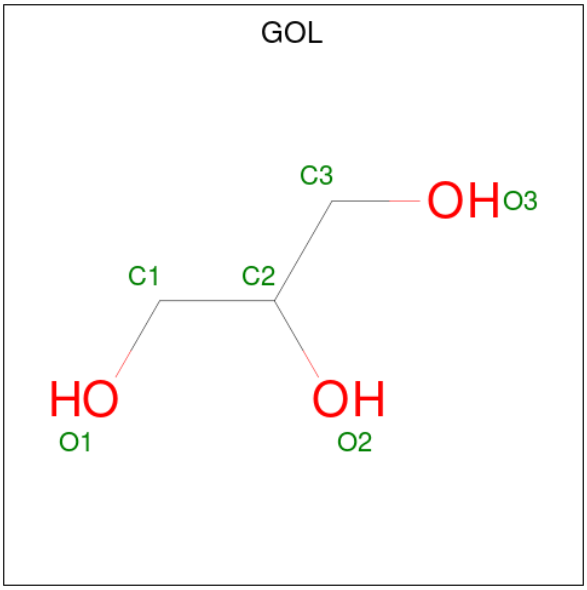
Chain	Residue	Modelled	Actual	Comment	Reference
D	66	ALA	VAL	conflict	UNP Q5U7L7
D	144	LYS	GLU	conflict	UNP Q5U7L7
D	148	THR	ASN	conflict	UNP Q5U7L7
D	181	PRO	THR	conflict	UNP Q5U7L7
D	185	ILE	VAL	conflict	UNP Q5U7L7
D	192	ARG	SER	conflict	UNP Q5U7L7
D	213	ALA	ASP	conflict	UNP Q5U7L7
D	253	SER	THR	conflict	UNP Q5U7L7
D	254	ASP	ASN	conflict	UNP Q5U7L7

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0

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



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

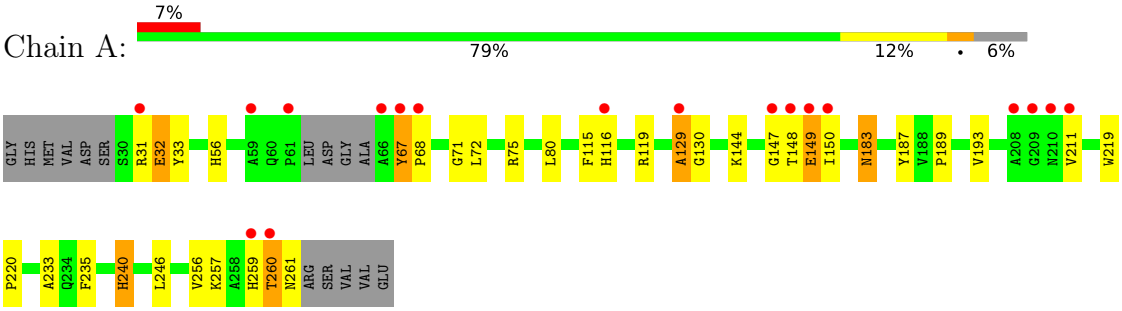
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	58	Total O 58 58	0	0
4	B	66	Total O 66 66	0	0
4	C	53	Total O 53 53	0	0
4	D	18	Total O 18 18	0	0

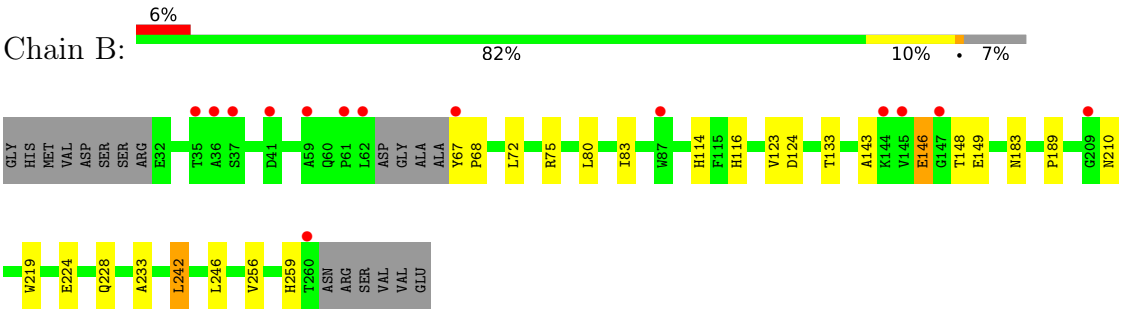
3 Residue-property plots (i)

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.

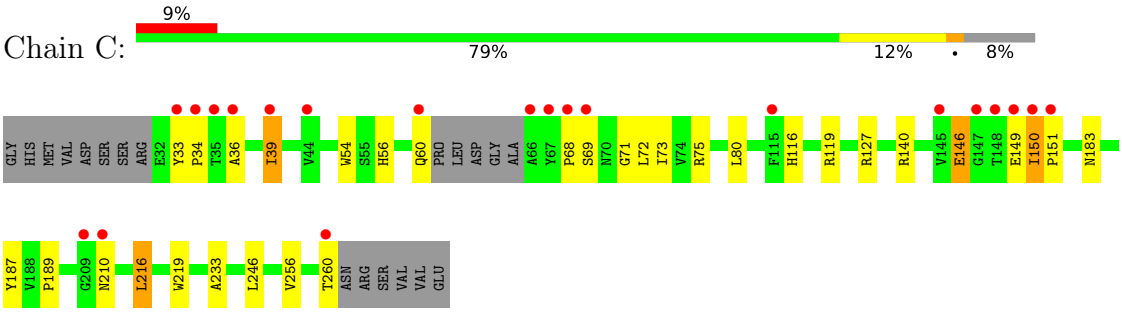
• Molecule 1: Metallo-beta-lactamase



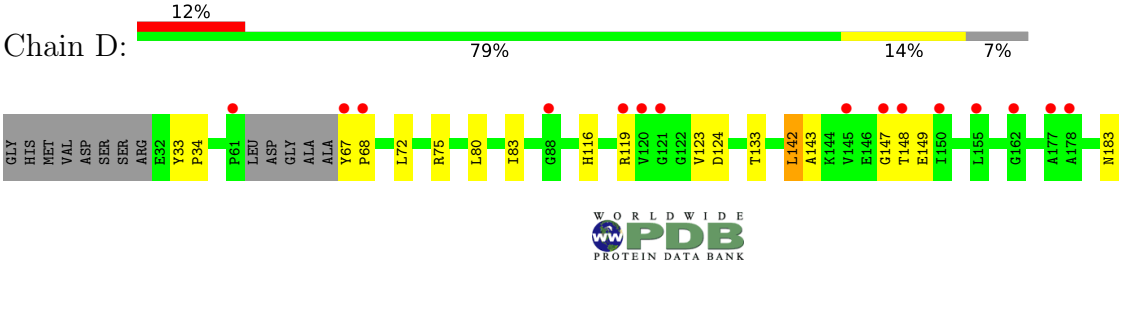
• Molecule 1: Metallo-beta-lactamase

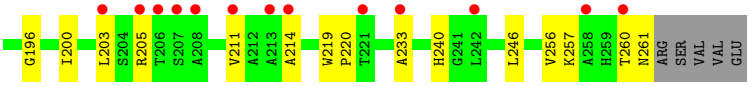


• Molecule 1: Metallo-beta-lactamase



• Molecule 1: Metallo-beta-lactamase




G196
T200
L203
S204
R205
T206
S207
A208
V211
A212
A213
A214
W219
P220
T221
A233
H240
G241
L242
L246
V256
K257
A258
H259
T260
N261
ARG
SER
VAL
VAL
GLU

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	128.60Å 41.67Å 156.76Å 90.00° 99.41° 90.00°	Depositor
Resolution (Å)	37.89 – 2.19 37.89 – 2.19	Depositor EDS
% Data completeness (in resolution range)	97.8 (37.89-2.19) 97.8 (37.89-2.19)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.92 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.221 , 0.267 0.223 , 0.269	Depositor DCC
R_{free} test set	2114 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å^2)	32.4	Xtriage
Anisotropy	0.623	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 56.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.40$, $< L^2 > = 0.23$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	7053	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.64% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, 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	0.31	0/1779	0.74	1/2431 (0.0%)
1	B	0.28	0/1746	0.72	0/2386
1	C	0.30	0/1738	0.71	0/2375
1	D	0.29	0/1743	0.74	0/2383
All	All	0.29	0/7006	0.73	1/9575 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	129	ALA	N-CA-C	5.62	117.48	111.36

There are no chirality outliers.

There are no planarity outliers.

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	1737	0	1687	19	0
1	B	1705	0	1670	15	0
1	C	1698	0	1653	20	0
1	D	1702	0	1657	20	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	B	6	0	8	2	0
3	C	6	0	8	1	0
4	A	58	0	0	0	0
4	B	66	0	0	0	0
4	C	53	0	0	1	0
4	D	18	0	0	0	0
All	All	7053	0	6683	68	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 (68) 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:127:ARG:HD3	1:C:150:ILE:HD13	1.74	0.69
1:A:260:THR:OG1	1:A:261:ASN:N	2.26	0.68
1:B:189:PRO:HG2	3:B:302:GOL:H31	1.74	0.68
1:A:129:ALA:N	1:A:130:GLY:HA2	2.10	0.66
1:D:196:GLY:HA3	1:D:200:ILE:HD13	1.78	0.66
1:C:60:GLN:NE2	1:C:69:SER:OG	2.29	0.61
1:C:119:ARG:HD2	1:D:119:ARG:HD3	1.81	0.60
1:C:233:ALA:HB3	1:C:246:LEU:HD21	1.85	0.58
1:D:148:THR:OG1	1:D:149:GLU:N	2.36	0.58
1:D:260:THR:OG1	1:D:261:ASN:N	2.37	0.57
1:A:233:ALA:HB3	1:A:246:LEU:HD21	1.86	0.56
1:A:115:PHE:CE1	1:B:148:THR:HG21	2.40	0.56
1:A:119:ARG:HH12	1:A:147:GLY:HA3	1.70	0.56
1:C:149:GLU:O	1:C:150:ILE:HG13	2.06	0.56
1:C:140:ARG:NH1	4:C:405:HOH:O	2.39	0.56
1:D:123:VAL:HG13	1:D:133:THR:HG21	1.90	0.53
1:D:72:LEU:HB2	1:D:83:ILE:HB	1.91	0.53
1:A:220:PRO:HB3	1:A:257:LYS:HE3	1.91	0.52
1:D:67:TYR:N	1:D:68:PRO:HD3	2.24	0.52
1:B:189:PRO:HG2	3:B:302:GOL:H11	1.92	0.52
1:B:123:VAL:HG13	1:B:133:THR:HG21	1.92	0.51
1:A:144:LYS:HA	1:A:148:THR:HG22	1.91	0.51
1:B:72:LEU:HB2	1:B:83:ILE:HB	1.93	0.51
1:B:224:GLU:O	1:B:228:GLN:HG2	2.11	0.51
1:D:119:ARG:NH1	1:D:124:ASP:OD2	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:56:HIS:CE1	1:C:71:GLY:HA3	2.47	0.49
1:A:219:TRP:HB3	1:A:256:VAL:HG11	1.94	0.49
1:C:210:ASN:OD1	1:C:210:ASN:N	2.44	0.49
1:A:148:THR:OG1	1:A:149:GLU:N	2.42	0.49
1:C:36:ALA:HA	1:C:39:ILE:HG13	1.94	0.49
1:D:143:ALA:O	1:D:148:THR:HG22	2.13	0.49
1:C:150:ILE:HB	1:C:151:PRO:HD2	1.95	0.48
1:D:147:GLY:HA3	1:D:148:THR:HB	1.96	0.47
1:A:67:TYR:N	1:A:68:PRO:HD3	2.30	0.47
1:A:240:HIS:CD2	1:B:68:PRO:HA	2.49	0.47
1:D:219:TRP:HB3	1:D:256:VAL:HG11	1.96	0.46
1:D:75:ARG:HA	1:D:80:LEU:HD23	1.98	0.46
1:D:220:PRO:HB3	1:D:257:LYS:HE3	1.97	0.46
1:A:183:ASN:HA	1:B:114:HIS:HB3	1.97	0.46
1:A:75:ARG:HA	1:A:80:LEU:HD23	1.97	0.45
1:B:233:ALA:HB3	1:B:246:LEU:HD21	1.97	0.45
1:C:33:TYR:HA	1:C:34:PRO:HD3	1.84	0.45
1:C:75:ARG:HA	1:C:80:LEU:HD23	1.97	0.45
1:D:205:ARG:NH2	1:D:240:HIS:HB2	2.32	0.45
1:B:75:ARG:HA	1:B:80:LEU:HD23	1.98	0.44
1:D:211:VAL:HG13	1:D:214:ALA:HB3	1.99	0.44
1:D:142:LEU:HD13	1:D:142:LEU:HA	1.88	0.44
1:C:56:HIS:NE2	1:C:71:GLY:HA3	2.33	0.43
1:D:233:ALA:HB3	1:D:246:LEU:HD21	2.01	0.43
1:A:56:HIS:CE1	1:A:71:GLY:HA3	2.53	0.43
1:C:68:PRO:O	1:D:240:HIS:HA	2.18	0.43
1:C:219:TRP:HB3	1:C:256:VAL:HG11	2.00	0.43
1:A:193:VAL:HG22	1:A:235[A]:PHE:HB2	2.01	0.43
1:B:219:TRP:HB3	1:B:256:VAL:HG11	2.01	0.42
1:A:193:VAL:HG22	1:A:235[B]:PHE:HB3	2.01	0.42
1:B:149:GLU:H	1:B:149:GLU:HG2	1.64	0.42
1:C:216:LEU:HB3	1:C:260:THR:HG22	2.01	0.42
1:B:148:THR:OG1	1:B:149:GLU:N	2.52	0.42
1:D:33:TYR:HA	1:D:34:PRO:HD3	1.90	0.42
1:D:119:ARG:HH12	1:D:149:GLU:HB2	1.85	0.42
1:A:31:ARG:HB3	1:A:32:GLU:H	1.66	0.42
1:C:187:TYR:O	1:C:189:PRO:HD3	2.20	0.42
1:A:33:TYR:HB3	1:B:242:LEU:HD22	2.02	0.41
1:A:187:TYR:O	1:A:189:PRO:HD3	2.21	0.41
1:C:54:TRP:HB2	1:C:73:ILE:HB	2.03	0.41
1:C:149:GLU:C	1:C:150:ILE:HG13	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:189:PRO:HG2	3:C:302:GOL:H31	2.03	0.41
1:B:143:ALA:O	1:B:148:THR:HG22	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	226/243 (93%)	214 (95%)	9 (4%)	3 (1%)	9 8
1	B	222/243 (91%)	214 (96%)	6 (3%)	2 (1%)	14 14
1	C	222/243 (91%)	212 (96%)	8 (4%)	2 (1%)	14 14
1	D	222/243 (91%)	214 (96%)	8 (4%)	0	100 100
All	All	892/972 (92%)	854 (96%)	31 (4%)	7 (1%)	16 16

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	149	GLU
1	A	211	VAL
1	B	146	GLU
1	C	146	GLU
1	C	150	ILE
1	B	210	ASN
1	A	67	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	181/191 (95%)	173 (96%)	8 (4%)	25	34
1	B	178/191 (93%)	171 (96%)	7 (4%)	28	39
1	C	177/191 (93%)	171 (97%)	6 (3%)	32	44
1	D	178/191 (93%)	174 (98%)	4 (2%)	45	61
All	All	714/764 (94%)	689 (96%)	25 (4%)	32	43

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

Mol	Chain	Res	Type
1	A	32	GLU
1	A	72	LEU
1	A	116	HIS
1	A	150	ILE
1	A	183	ASN
1	A	240	HIS
1	A	259	HIS
1	A	260	THR
1	B	67	TYR
1	B	116	HIS
1	B	124	ASP
1	B	146	GLU
1	B	183	ASN
1	B	242	LEU
1	B	259	HIS
1	C	39	ILE
1	C	72	LEU
1	C	116	HIS
1	C	146	GLU
1	C	183	ASN
1	C	216	LEU
1	D	116	HIS
1	D	142	LEU
1	D	183	ASN
1	D	203	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	210	ASN
1	A	240	HIS
1	B	153	HIS
1	B	228	GLN
1	D	179	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 (i)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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	B	302	-	5,5,5	0.36	0	5,5,5	0.23	0
3	GOL	C	302	-	5,5,5	0.37	0	5,5,5	0.21	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	B	302	-	-	4/4/4/4	-
3	GOL	C	302	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	302	GOL	C1-C2-C3-O3
3	C	302	GOL	C1-C2-C3-O3
3	C	302	GOL	O2-C2-C3-O3
3	B	302	GOL	O1-C1-C2-C3
3	B	302	GOL	O2-C2-C3-O3
3	B	302	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	302	GOL	2	0
3	C	302	GOL	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	228/243 (93%)	0.48	18 (7%)	18	16	15, 36, 76, 101	2 (0%)
1	B	225/243 (92%)	0.32	14 (6%)	26	23	17, 34, 65, 100	4 (1%)
1	C	224/243 (92%)	0.68	21 (9%)	14	11	20, 42, 84, 96	3 (1%)
1	D	225/243 (92%)	0.95	28 (12%)	8	6	19, 52, 83, 97	1 (0%)
All	All	902/972 (92%)	0.61	81 (8%)	15	12	15, 40, 79, 101	10 (1%)

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	62	LEU	5.4
1	A	147	GLY	4.9
1	C	66	ALA	4.4
1	A	66	ALA	4.3
1	D	67	TYR	4.2
1	A	260	THR	4.2
1	D	147	GLY	4.1
1	A	67	TYR	4.0
1	C	148	THR	4.0
1	A	211	VAL	3.8
1	C	36	ALA	3.7
1	D	260	THR	3.7
1	B	145	VAL	3.7
1	A	149	GLU	3.7
1	C	67	TYR	3.7
1	D	214	ALA	3.5
1	B	67	TYR	3.5
1	B	61	PRO	3.5
1	C	260	THR	3.2
1	D	88	GLY	3.2
1	A	210	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	115	PHE	3.2
1	A	68	PRO	3.2
1	D	121	GLY	3.1
1	A	208	ALA	3.0
1	D	208	ALA	3.0
1	C	33	TYR	3.0
1	C	210	ASN	3.0
1	C	150	ILE	2.9
1	A	209	GLY	2.9
1	C	68	PRO	2.9
1	D	206	THR	2.8
1	B	59	ALA	2.8
1	D	177	ALA	2.8
1	C	44	VAL	2.8
1	D	120	VAL	2.8
1	A	61	PRO	2.7
1	B	87	TRP	2.7
1	B	209	GLY	2.7
1	D	205	ARG	2.6
1	A	116	HIS	2.6
1	B	260	THR	2.6
1	C	34	PRO	2.6
1	D	203	LEU	2.6
1	D	148	THR	2.6
1	D	258	ALA	2.6
1	C	147	GLY	2.6
1	B	36	ALA	2.6
1	C	60	GLN	2.5
1	C	39	ILE	2.5
1	B	147	GLY	2.5
1	D	145	VAL	2.5
1	A	31	ARG	2.5
1	B	35	THR	2.5
1	D	61	PRO	2.4
1	B	41	ASP	2.4
1	A	148	THR	2.4
1	C	69	SER	2.4
1	D	68	PRO	2.4
1	D	207	SER	2.4
1	C	151	PRO	2.4
1	D	233	ALA	2.3
1	C	145	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	221	THR	2.3
1	B	37	SER	2.3
1	A	259	HIS	2.3
1	D	211	VAL	2.3
1	A	150	ILE	2.3
1	D	119	ARG	2.2
1	C	209	GLY	2.2
1	A	59	ALA	2.2
1	D	242	LEU	2.2
1	C	35	THR	2.1
1	C	149	GLU	2.1
1	D	162	GLY	2.1
1	D	213	ALA	2.1
1	B	144[A]	LYS	2.1
1	D	150	ILE	2.0
1	A	129	ALA	2.0
1	D	178	ALA	2.0
1	D	155	LEU	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
3	GOL	C	302	6/6	0.82	0.18	59,59,59,60	0
3	GOL	B	302	6/6	0.93	0.10	45,46,47,48	0
2	ZN	D	301	1/1	0.93	0.08	68,68,68,68	0
2	ZN	B	301	1/1	0.95	0.06	66,66,66,66	0
2	ZN	A	301	1/1	0.96	0.09	68,68,68,68	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	ZN	C	301	1/1	0.97	0.06	72,72,72,72	0

6.5 Other polymers (i)

There are no such residues in this entry.