



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:43 AM UTC

PDB ID : 3PK7 / pdb_00003pk7
Title : Crystal structure of D-mannonate dehydratase from *Chromohalobacter salexigens* with MG and Glycerol bound in the active site
Authors : Fedorov, A.A.; Fedorov, E.V.; Wichelecki, D.; Gerlt, J.A.; Almo, S.C.
Deposited on : 2010-11-11
Resolution : 1.64 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

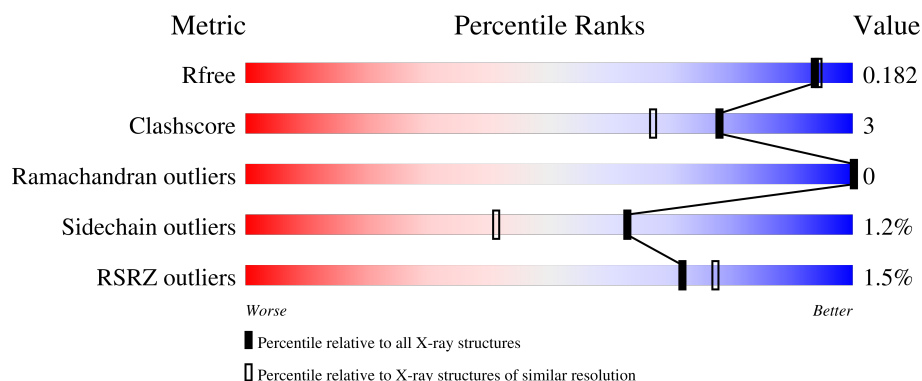
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.64 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	1141 (1.64-1.64)
Clashscore	190562	1171 (1.64-1.64)
Ramachandran outliers	187476	1151 (1.64-1.64)
Sidechain outliers	187428	1150 (1.64-1.64)
RSRZ outliers	180081	1141 (1.64-1.64)

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	405	90% 7% ..
1	B	405	89% 5% 5%
1	C	405	89% 8% .
1	D	405	91% . .
1	E	405	91% 6% .

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Mol	Chain	Length	Quality of chain
1	F	405	% 91% 5%
1	G	405	% 86% 9%
1	H	405	3% 90% 6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 27832 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 Mandelate racemase/muconate lactonizing enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	396	Total	C	N	O	S	0	2	0
			3160	2002	565	579	14			
1	B	384	Total	C	N	O	S	0	1	0
			3052	1937	546	555	14			
1	C	395	Total	C	N	O	S	0	1	0
			3133	1986	557	576	14			
1	D	387	Total	C	N	O	S	0	2	0
			3088	1955	556	563	14			
1	E	396	Total	C	N	O	S	0	1	0
			3141	1990	559	578	14			
1	F	386	Total	C	N	O	S	0	2	0
			3075	1949	549	563	14			
1	G	388	Total	C	N	O	S	0	3	0
			3107	1970	555	568	14			
1	H	394	Total	C	N	O	S	0	1	0
			3130	1982	559	575	14			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q1QT89
A	2	SER	-	expression tag	UNP Q1QT89
A	3	LEU	-	expression tag	UNP Q1QT89
B	1	MET	-	expression tag	UNP Q1QT89
B	2	SER	-	expression tag	UNP Q1QT89
B	3	LEU	-	expression tag	UNP Q1QT89
C	1	MET	-	expression tag	UNP Q1QT89
C	2	SER	-	expression tag	UNP Q1QT89
C	3	LEU	-	expression tag	UNP Q1QT89
D	1	MET	-	expression tag	UNP Q1QT89
D	2	SER	-	expression tag	UNP Q1QT89
D	3	LEU	-	expression tag	UNP Q1QT89
E	1	MET	-	expression tag	UNP Q1QT89

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Chain	Residue	Modelled	Actual	Comment	Reference
E	2	SER	-	expression tag	UNP Q1QT89
E	3	LEU	-	expression tag	UNP Q1QT89
F	1	MET	-	expression tag	UNP Q1QT89
F	2	SER	-	expression tag	UNP Q1QT89
F	3	LEU	-	expression tag	UNP Q1QT89
G	1	MET	-	expression tag	UNP Q1QT89
G	2	SER	-	expression tag	UNP Q1QT89
G	3	LEU	-	expression tag	UNP Q1QT89
H	1	MET	-	expression tag	UNP Q1QT89
H	2	SER	-	expression tag	UNP Q1QT89
H	3	LEU	-	expression tag	UNP Q1QT89

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	0	0
2	B	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0
2	D	2	Total Mg 2 2	0	0
2	E	1	Total Mg 1 1	0	0
2	F	1	Total Mg 1 1	0	0
2	G	1	Total Mg 1 1	0	0
2	H	1	Total Mg 1 1	0	0

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	F	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	354	Total	O	0	0
			354	354		
4	B	364	Total	O	0	0
			364	364		
4	C	385	Total	O	0	0
			385	385		
4	D	345	Total	O	0	0
			345	345		

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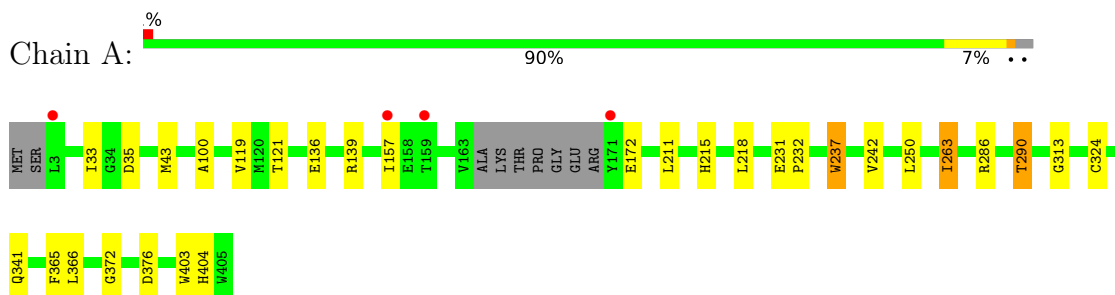
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	372	Total 372	O 372	0	0
4	F	378	Total 378	O 378	0	0
4	G	353	Total 353	O 353	0	0
4	H	338	Total 338	O 338	0	0

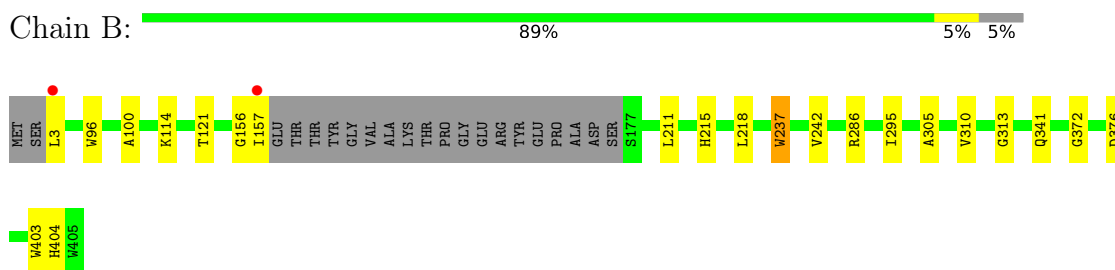
3 Residue-property plots [i](#)

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.

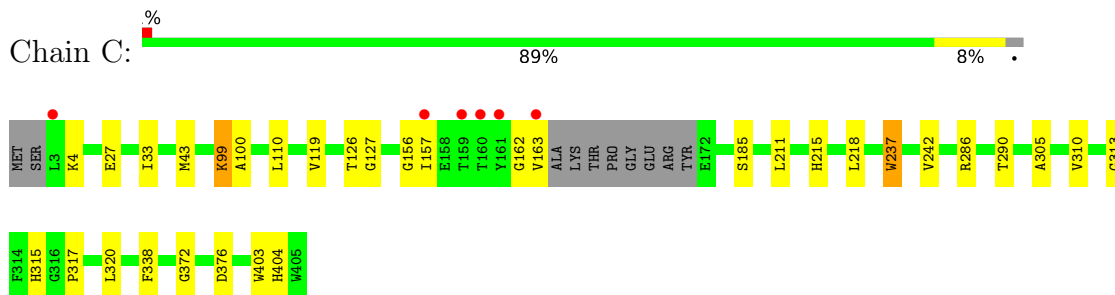
- Molecule 1: Mandelate racemase/muconate lactonizing enzyme



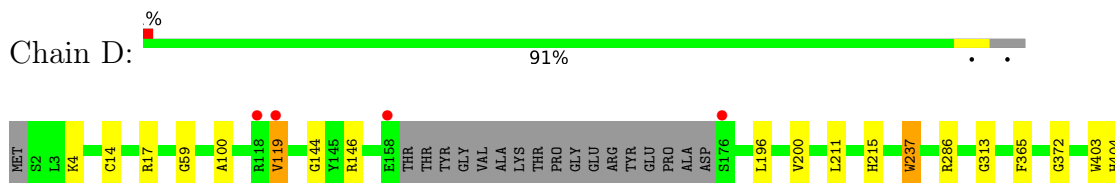
- Molecule 1: Mandelate racemase/muconate lactonizing enzyme



- Molecule 1: Mandelate racemase/muconate lactonizing enzyme

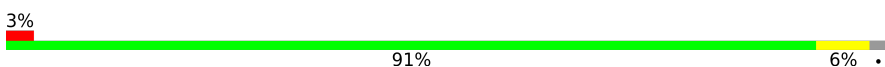


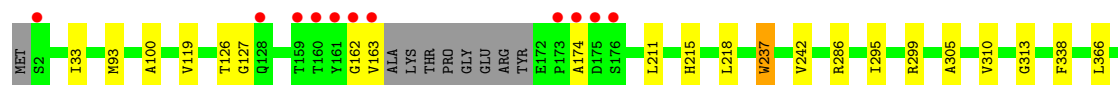
- Molecule 1: Mandelate racemase/muconate lactonizing enzyme



W405

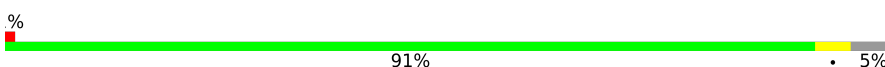
- Molecule 1: Mandelate racemase/muconate lactonizing enzyme

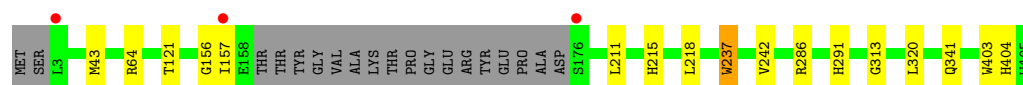
Chain E:  3% 91% 6%



G372 D376 W403 W404 W405

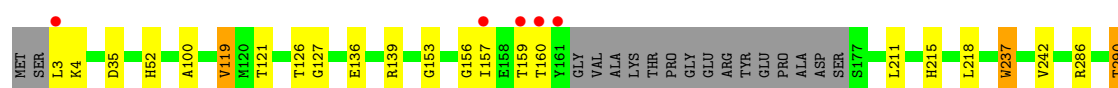
- Molecule 1: Mandelate racemase/muconate lactonizing enzyme

Chain F:  0% 91% 5%



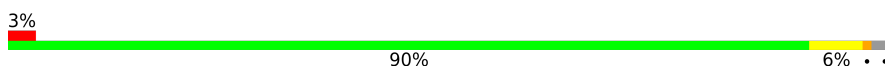
- Molecule 1: Mandelate racemase/muconate lactonizing enzyme

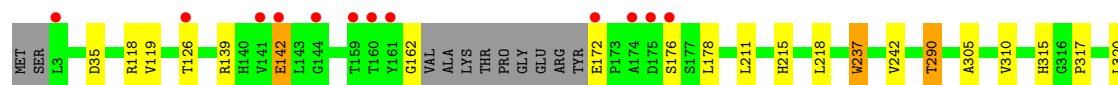
Chain G:  0% 86% 9%



A305 V310 G313 T318 D319 L320 C324 Q341 M344 R359 F360 E361 F365 G372 K384 W403 W404 W405

- Molecule 1: Mandelate racemase/muconate lactonizing enzyme

Chain H:  3% 90% 6%



C324 F338 E349 F365 L366 L381 K384 W403 W404 W405

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	195.27Å 85.76Å 195.10Å 90.00° 110.31° 90.00°	Depositor
Resolution (Å)	40.05 – 1.64 40.05 – 1.64	Depositor EDS
% Data completeness (in resolution range)	99.8 (40.05-1.64) 99.8 (40.05-1.64)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.82 (at 1.64Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.158 , 0.184 0.156 , 0.182	Depositor DCC
R_{free} test set	18479 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	17.9	Xtriage
Anisotropy	0.263	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 33.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	27832	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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, MG

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	0.41	0/3248	0.73	2/4423 (0.0%)
1	B	0.42	0/3137	0.71	0/4273
1	C	0.41	0/3220	0.71	0/4388
1	D	0.40	0/3173	0.72	0/4318
1	E	0.41	0/3228	0.73	0/4396
1	F	0.43	0/3160	0.71	0/4304
1	G	0.41	0/3193	0.73	0/4349
1	H	0.41	0/3217	0.72	2/4381 (0.0%)
All	All	0.41	0/25576	0.72	4/34832 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	172	GLU	CA-C-N	5.53	124.99	119.24
1	H	172	GLU	C-N-CA	5.53	124.99	119.24
1	A	172	GLU	CA-C-N	5.07	126.17	119.84
1	A	172	GLU	C-N-CA	5.07	126.17	119.84

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	3160	0	3043	18	0
1	B	3052	0	2951	10	0
1	C	3133	0	3019	22	0
1	D	3088	0	2982	10	0
1	E	3141	0	3027	20	0
1	F	3075	0	2965	12	0
1	G	3107	0	2997	27	0
1	H	3130	0	3013	20	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	2	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
3	A	6	0	8	0	0
3	B	6	0	8	0	0
3	C	6	0	8	0	0
3	D	6	0	8	0	0
3	E	6	0	8	0	0
3	F	6	0	8	0	0
3	G	6	0	8	0	0
3	H	6	0	8	0	0
4	A	354	0	0	3	0
4	B	364	0	0	0	0
4	C	385	0	0	3	0
4	D	345	0	0	0	0
4	E	372	0	0	4	0
4	F	378	0	0	3	0
4	G	353	0	0	3	0
4	H	338	0	0	2	0
All	All	27832	0	24061	136	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 3.

All (136) 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:E:126:THR:HG21	1:E:163:VAL:HA	1.62	0.80
1:E:126:THR:HG21	1:E:162:GLY:O	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:126:THR:HG21	1:C:163:VAL:HA	1.68	0.75
1:A:157:ILE:HG13	4:A:2991:HOH:O	1.87	0.74
1:F:218:LEU:O	1:F:242[A]:VAL:HG12	1.90	0.72
1:H:290:THR:HG22	1:H:324:CYS:SG	2.34	0.67
1:C:126:THR:HG21	1:C:162:GLY:O	1.95	0.67
1:F:43:MET:H	1:G:52:HIS:HE1	1.43	0.67
1:C:4:LYS:HB3	1:C:27:GLU:HG3	1.76	0.67
1:G:218:LEU:O	1:G:242[A]:VAL:HG12	1.96	0.66
1:A:136:GLU:OE2	1:A:139:ARG:HD3	1.96	0.66
1:A:290:THR:HG22	1:A:324:CYS:SG	2.35	0.65
1:A:286:ARG:HD2	1:A:313:GLY:O	1.95	0.65
1:G:159:THR:HG22	4:G:3103:HOH:O	1.98	0.63
1:G:35:ASP:OD2	1:G:290:THR:HG21	1.99	0.62
1:C:286:ARG:HD2	1:C:313:GLY:O	2.01	0.60
1:G:3:LEU:HD21	4:G:2357:HOH:O	2.01	0.60
1:D:286:ARG:HD2	1:D:313:GLY:O	2.01	0.60
1:C:126:THR:HG23	1:C:163:VAL:HG22	1.82	0.60
1:C:157:ILE:HG13	4:C:2189:HOH:O	2.02	0.60
1:C:156:GLY:O	1:C:157:ILE:HD13	2.02	0.59
1:E:126:THR:CG2	1:E:163:VAL:HA	2.33	0.59
1:H:139:ARG:O	1:H:142:GLU:HB3	2.02	0.59
1:A:35:ASP:OD2	1:A:290:THR:HG21	2.04	0.58
1:G:290:THR:HG22	1:G:324:CYS:SG	2.44	0.58
1:B:156:GLY:O	1:B:157:ILE:HD13	2.03	0.57
1:F:43:MET:H	1:G:52:HIS:CE1	2.22	0.57
1:H:290:THR:CG2	1:H:324:CYS:SG	2.92	0.57
4:F:1291:HOH:O	1:G:52:HIS:HD2	1.88	0.57
1:G:286:ARG:HD2	1:G:313:GLY:O	2.05	0.56
1:B:114:LYS:HE3	1:F:64:ARG:NH1	2.21	0.56
1:B:218:LEU:O	1:B:242[A]:VAL:HG12	2.06	0.55
1:E:295:ILE:O	1:E:299:ARG:HG3	2.07	0.55
1:D:119:VAL:HG22	1:D:365:PHE:HB2	1.88	0.54
1:H:35:ASP:OD2	1:H:290:THR:HG21	2.06	0.54
1:D:4:LYS:HE3	1:D:59:GLY:O	2.08	0.54
1:E:174:ALA:CB	4:E:3055:HOH:O	2.55	0.54
1:C:99:LYS:HG2	1:C:110:LEU:HD21	1.89	0.54
1:E:126:THR:CG2	1:E:162:GLY:O	2.56	0.53
1:C:126:THR:CG2	1:C:163:VAL:HA	2.37	0.53
1:G:153:GLY:HA2	1:G:160:THR:HG21	1.90	0.53
1:C:157:ILE:HD11	1:C:185:SER:HB3	1.91	0.52
1:H:366:LEU:HD22	1:H:366:LEU:N	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:218:LEU:O	1:H:242:VAL:HG12	2.09	0.52
1:F:403:TRP:CG	1:F:404:HIS:H	2.28	0.52
1:E:286:ARG:HD2	1:E:313:GLY:O	2.09	0.51
1:G:290:THR:HG23	1:G:320:LEU:HD22	1.91	0.51
1:D:403:TRP:CG	1:D:404:HIS:H	2.28	0.51
1:G:136:GLU:OE2	1:G:139:ARG:HD3	2.10	0.51
1:H:176:SER:HB2	1:H:178:LEU:O	2.10	0.51
1:F:156:GLY:O	1:F:157:ILE:HD13	2.10	0.51
1:A:290:THR:CG2	1:A:324:CYS:SG	2.98	0.51
1:B:403:TRP:CG	1:B:404:HIS:H	2.29	0.51
1:E:403:TRP:CG	1:E:404:HIS:H	2.30	0.50
1:E:174:ALA:HB2	4:E:3055:HOH:O	2.11	0.50
1:C:403:TRP:CG	1:C:404:HIS:H	2.28	0.49
1:E:126:THR:HG23	1:E:163:VAL:HG22	1.94	0.49
1:G:403:TRP:CG	1:G:404:HIS:H	2.31	0.49
1:E:100:ALA:HB3	1:E:372:GLY:HA2	1.95	0.49
1:C:33:ILE:HG23	4:C:2459:HOH:O	2.12	0.49
1:H:403:TRP:CG	1:H:404:HIS:H	2.31	0.48
1:A:403:TRP:CG	1:A:404:HIS:H	2.31	0.48
1:E:211:LEU:HD23	1:E:237:TRP:CE2	2.49	0.48
1:F:211:LEU:HD23	1:F:237:TRP:CE2	2.49	0.48
1:E:305:ALA:HB1	1:E:310:VAL:HB	1.96	0.48
1:D:100:ALA:HB3	1:D:372:GLY:HA2	1.95	0.48
1:H:365:PHE:C	1:H:366:LEU:HD22	2.39	0.47
1:C:126:THR:CG2	1:C:162:GLY:O	2.60	0.47
1:H:118:ARG:NH1	4:H:2077:HOH:O	2.46	0.47
1:B:211:LEU:HD23	1:B:237:TRP:CE2	2.49	0.47
1:E:126:THR:HG22	1:E:127:GLY:N	2.28	0.47
1:E:218:LEU:O	1:E:242:VAL:HG12	2.14	0.47
1:C:218:LEU:O	1:C:242[A]:VAL:HG12	2.14	0.47
1:A:218:LEU:O	1:A:242:VAL:HG12	2.16	0.46
1:C:211:LEU:HD23	1:C:237:TRP:CE2	2.51	0.46
1:G:100:ALA:HB3	1:G:372:GLY:HA2	1.97	0.46
1:G:153:GLY:HA2	1:G:160:THR:CG2	2.46	0.46
1:A:100:ALA:HB3	1:A:372:GLY:HA2	1.98	0.45
1:A:136:GLU:HA	1:A:139:ARG:HG2	1.97	0.45
1:G:290:THR:CG2	1:G:324:CYS:SG	3.04	0.45
1:G:384:LYS:NZ	4:G:3098:HOH:O	2.48	0.45
1:A:250:LEU:HD21	1:A:263:ILE:HD13	1.98	0.45
1:A:43:MET:HE3	4:A:3059:HOH:O	2.17	0.45
1:G:211:LEU:HD23	1:G:237:TRP:CE2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:315:HIS:CE1	1:H:317:PRO:HG3	2.52	0.45
1:A:211:LEU:HD23	1:A:237:TRP:CE2	2.52	0.45
1:C:305:ALA:HB1	1:C:310:VAL:HB	1.99	0.45
1:C:290:THR:HB	1:C:320:LEU:HD22	1.99	0.44
1:H:305:ALA:HB1	1:H:310:VAL:HB	1.98	0.44
1:E:119:VAL:HG13	1:E:338:PHE:CZ	2.52	0.44
1:A:365:PHE:C	1:A:366:LEU:HD22	2.42	0.44
1:B:121:THR:HA	1:B:341:GLN:O	2.16	0.44
1:B:286:ARG:HD2	1:B:313:GLY:O	2.18	0.44
1:C:119:VAL:HG13	1:C:338:PHE:CZ	2.53	0.44
1:D:144:GLY:O	1:D:146:ARG:NH1	2.51	0.44
1:F:286:ARG:HD2	1:F:313:GLY:O	2.17	0.44
1:G:318:THR:HG23	1:G:344:MET:HE2	2.00	0.43
1:H:119:VAL:HG13	1:H:338:PHE:CZ	2.52	0.43
1:A:33:ILE:HG23	4:A:2123:HOH:O	2.18	0.43
1:G:359:ARG:NH2	1:G:361:GLU:OE2	2.51	0.43
1:E:174:ALA:HB1	4:E:3055:HOH:O	2.19	0.43
1:E:366:LEU:N	1:E:366:LEU:HD22	2.33	0.43
1:H:211:LEU:HD23	1:H:237:TRP:CE2	2.53	0.43
1:F:43:MET:HE3	4:F:1560:HOH:O	2.19	0.42
1:F:157:ILE:HG13	4:F:1808:HOH:O	2.19	0.42
1:H:126:THR:HG21	1:H:162:GLY:O	2.19	0.42
1:C:315:HIS:CE1	1:C:317:PRO:HG3	2.55	0.42
1:E:93:MET:HE3	1:E:93:MET:HB2	1.89	0.42
1:H:126:THR:CG2	1:H:162:GLY:O	2.67	0.42
1:C:100:ALA:HB3	1:C:372:GLY:HA2	2.02	0.42
1:A:366:LEU:HD22	1:A:366:LEU:N	2.34	0.42
1:A:231:GLU:N	1:A:232:PRO:CD	2.82	0.42
1:B:100:ALA:HB3	1:B:372:GLY:HA2	2.01	0.42
1:G:156:GLY:O	1:G:157:ILE:HD13	2.20	0.42
1:G:119:VAL:HG22	1:G:365:PHE:HB2	2.00	0.42
1:H:290:THR:HG23	1:H:320:LEU:HD22	2.02	0.41
1:H:349:GLU:HG2	4:H:1793:HOH:O	2.20	0.41
1:F:291:HIS:CE1	1:F:320:LEU:HD21	2.56	0.41
1:G:305:ALA:HB1	1:G:310:VAL:HB	2.01	0.41
1:H:290:THR:CG2	1:H:324:CYS:HB2	2.50	0.41
1:C:43:MET:HE3	4:C:1528:HOH:O	2.20	0.41
1:D:14:CYS:HA	1:D:17:ARG:O	2.21	0.41
1:E:33:ILE:HG23	4:E:2723:HOH:O	2.20	0.41
1:B:96:TRP:CD1	1:B:295:ILE:HB	2.55	0.41
1:F:121:THR:HA	1:F:341:GLN:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:237:TRP:CD1	1:D:237:TRP:C	2.99	0.41
1:G:121:THR:HA	1:G:341:GLN:O	2.21	0.41
1:G:126:THR:HG22	1:G:127:GLY:N	2.35	0.41
1:C:126:THR:HG22	1:C:127:GLY:N	2.35	0.41
1:D:196:LEU:O	1:D:200:VAL:HG23	2.20	0.41
1:G:290:THR:CG2	1:G:324:CYS:HB2	2.51	0.41
1:G:4:LYS:HE2	1:G:4:LYS:HB3	1.69	0.40
1:H:381:LEU:HD12	1:H:384:LYS:HD2	2.03	0.40
1:B:305:ALA:HB1	1:B:310:VAL:HB	2.04	0.40
1:D:211:LEU:HD23	1:D:237:TRP:CE2	2.57	0.40
1:A:121:THR:HA	1:A:341:GLN:O	2.22	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	394/405 (97%)	382 (97%)	12 (3%)	0	100	100
1	B	381/405 (94%)	370 (97%)	11 (3%)	0	100	100
1	C	392/405 (97%)	381 (97%)	11 (3%)	0	100	100
1	D	385/405 (95%)	373 (97%)	12 (3%)	0	100	100
1	E	393/405 (97%)	382 (97%)	11 (3%)	0	100	100
1	F	384/405 (95%)	373 (97%)	11 (3%)	0	100	100
1	G	387/405 (96%)	377 (97%)	10 (3%)	0	100	100
1	H	391/405 (96%)	379 (97%)	12 (3%)	0	100	100
All	All	3107/3240 (96%)	3017 (97%)	90 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	327/332 (98%)	321 (98%)	6 (2%)	51	26
1	B	316/332 (95%)	312 (99%)	4 (1%)	61	38
1	C	325/332 (98%)	321 (99%)	4 (1%)	63	41
1	D	320/332 (96%)	317 (99%)	3 (1%)	70	54
1	E	326/332 (98%)	323 (99%)	3 (1%)	70	54
1	F	319/332 (96%)	317 (99%)	2 (1%)	78	66
1	G	322/332 (97%)	318 (99%)	4 (1%)	63	41
1	H	324/332 (98%)	319 (98%)	5 (2%)	57	32
All	All	2579/2656 (97%)	2548 (99%)	31 (1%)	63	41

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	VAL
1	A	215	HIS
1	A	237	TRP
1	A	263	ILE
1	A	290	THR
1	A	376	ASP
1	B	3	LEU
1	B	215	HIS
1	B	237	TRP
1	B	376	ASP
1	C	99	LYS
1	C	215	HIS
1	C	237	TRP
1	C	376	ASP
1	D	119	VAL
1	D	215	HIS
1	D	237	TRP
1	E	215	HIS
1	E	237	TRP

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Mol	Chain	Res	Type
1	E	376	ASP
1	F	215	HIS
1	F	237	TRP
1	G	119	VAL
1	G	215	HIS
1	G	237	TRP
1	G	290	THR
1	H	142	GLU
1	H	215	HIS
1	H	237	TRP
1	H	290	THR
1	H	366	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	128	GLN
1	C	280	GLN
1	D	192	HIS
1	F	278	GLN
1	G	31	HIS
1	G	52	HIS
1	H	192	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 17 ligands modelled in this entry, 9 are monoatomic - leaving 8 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	F	407	-	5,5,5	0.38	0	5,5,5	0.39	0
3	GOL	D	407	-	5,5,5	0.39	0	5,5,5	0.39	0
3	GOL	H	407	-	5,5,5	0.40	0	5,5,5	0.16	0
3	GOL	G	407	-	5,5,5	0.41	0	5,5,5	0.50	0
3	GOL	A	407	-	5,5,5	0.39	0	5,5,5	0.27	0
3	GOL	E	407	-	5,5,5	0.40	0	5,5,5	0.42	0
3	GOL	B	407	-	5,5,5	0.40	0	5,5,5	0.52	0
3	GOL	C	407	-	5,5,5	0.39	0	5,5,5	0.51	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	F	407	-	-	4/4/4/4	-
3	GOL	D	407	-	-	2/4/4/4	-
3	GOL	H	407	-	-	4/4/4/4	-
3	GOL	G	407	-	-	2/4/4/4	-
3	GOL	A	407	-	-	2/4/4/4	-
3	GOL	E	407	-	-	4/4/4/4	-
3	GOL	B	407	-	-	4/4/4/4	-
3	GOL	C	407	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	407	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	C	407	GOL	O1-C1-C2-C3
3	C	407	GOL	C1-C2-C3-O3
3	E	407	GOL	O1-C1-C2-C3
3	E	407	GOL	C1-C2-C3-O3
3	F	407	GOL	O1-C1-C2-C3
3	G	407	GOL	C1-C2-C3-O3
3	H	407	GOL	C1-C2-C3-O3
3	A	407	GOL	C1-C2-C3-O3
3	B	407	GOL	O1-C1-C2-C3
3	F	407	GOL	C1-C2-C3-O3
3	H	407	GOL	O1-C1-C2-C3
3	B	407	GOL	O2-C2-C3-O3
3	C	407	GOL	O1-C1-C2-O2
3	C	407	GOL	O2-C2-C3-O3
3	E	407	GOL	O2-C2-C3-O3
3	F	407	GOL	O1-C1-C2-O2
3	F	407	GOL	O2-C2-C3-O3
3	G	407	GOL	O2-C2-C3-O3
3	H	407	GOL	O2-C2-C3-O3
3	E	407	GOL	O1-C1-C2-O2
3	H	407	GOL	O1-C1-C2-O2
3	A	407	GOL	O2-C2-C3-O3
3	B	407	GOL	O1-C1-C2-O2
3	D	407	GOL	O2-C2-C3-O3
3	D	407	GOL	C1-C2-C3-O3

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](#)

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	396/405 (97%)	-0.27	4 (1%) 79 84	9, 18, 32, 48	2 (0%)
1	B	384/405 (94%)	-0.46	2 (0%) 87 90	7, 16, 27, 43	1 (0%)
1	C	395/405 (97%)	-0.39	6 (1%) 72 77	7, 17, 28, 49	1 (0%)
1	D	387/405 (95%)	-0.16	4 (1%) 79 84	10, 19, 31, 61	2 (0%)
1	E	396/405 (97%)	-0.26	11 (2%) 55 60	12, 18, 30, 56	1 (0%)
1	F	386/405 (95%)	-0.37	3 (0%) 82 86	7, 16, 28, 58	2 (0%)
1	G	388/405 (95%)	-0.27	5 (1%) 75 80	6, 17, 30, 56	3 (0%)
1	H	394/405 (97%)	-0.05	12 (3%) 52 57	10, 19, 33, 54	1 (0%)
All	All	3126/3240 (96%)	-0.28	47 (1%) 72 77	6, 17, 31, 61	13 (0%)

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	174	ALA	4.4
1	E	175	ASP	4.3
1	G	159	THR	4.2
1	H	142	GLU	4.0
1	E	161	TYR	3.7
1	H	159	THR	3.6
1	D	176	SER	3.5
1	G	161	TYR	3.5
1	G	157	ILE	3.4
1	E	173	PRO	3.3
1	B	157	ILE	3.0
1	E	159	THR	3.0
1	E	176	SER	3.0
1	G	3	LEU	3.0
1	F	157	ILE	2.9
1	H	3	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	F	176	SER	2.8
1	A	3	LEU	2.7
1	H	160	THR	2.7
1	B	3	LEU	2.7
1	H	161	TYR	2.7
1	D	118[A]	ARG	2.6
1	H	176	SER	2.6
1	E	160	THR	2.6
1	F	3	LEU	2.6
1	H	175	ASP	2.5
1	C	159	THR	2.5
1	C	161	TYR	2.5
1	H	141	VAL	2.5
1	G	160	THR	2.5
1	H	126	THR	2.5
1	A	171	TYR	2.5
1	H	144	GLY	2.5
1	D	119	VAL	2.4
1	E	162	GLY	2.3
1	H	172	GLU	2.3
1	C	160	THR	2.3
1	A	159	THR	2.2
1	A	157	ILE	2.1
1	C	157	ILE	2.1
1	C	163	VAL	2.1
1	E	163	VAL	2.1
1	H	174	ALA	2.1
1	E	2	SER	2.1
1	C	3	LEU	2.1
1	D	158	GLU	2.0
1	E	128	GLN	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	H	407	6/6	0.84	0.14	21,28,33,39	0
3	GOL	F	407	6/6	0.88	0.12	20,25,33,37	0
3	GOL	C	407	6/6	0.88	0.11	20,26,30,39	0
2	MG	D	408	1/1	0.90	0.10	40,40,40,40	0
3	GOL	E	407	6/6	0.92	0.10	21,24,33,39	0
3	GOL	B	407	6/6	0.92	0.11	20,23,33,39	0
3	GOL	A	407	6/6	0.92	0.09	21,27,29,30	0
3	GOL	G	407	6/6	0.93	0.09	21,23,28,31	0
3	GOL	D	407	6/6	0.93	0.09	20,24,26,35	0
2	MG	H	406	1/1	0.97	0.04	17,17,17,17	0
2	MG	D	406	1/1	0.98	0.03	17,17,17,17	0
2	MG	B	406	1/1	0.99	0.02	14,14,14,14	0
2	MG	E	406	1/1	0.99	0.02	14,14,14,14	0
2	MG	G	406	1/1	0.99	0.02	14,14,14,14	0
2	MG	A	406	1/1	1.00	0.02	16,16,16,16	0
2	MG	F	406	1/1	1.00	0.03	15,15,15,15	0
2	MG	C	406	1/1	1.00	0.01	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.