



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:15 PM GMT

PDB ID : 1O66
Title : Crystal structure of 3-methyl-2-oxobutanoate hydroxymethyltransferase
Authors : Structural GenomiX
Deposited on : 2003-10-23
Resolution : 1.75 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

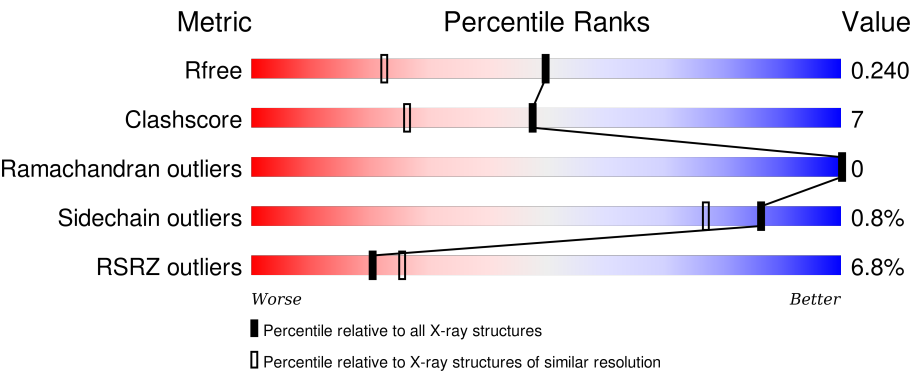
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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}	91344	1609 (1.76-1.76)
Clashscore	102246	1730 (1.76-1.76)
Ramachandran outliers	100387	1711 (1.76-1.76)
Sidechain outliers	100360	1711 (1.76-1.76)
RSRZ outliers	91569	1610 (1.76-1.76)

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 on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. 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	275	<div><div>4%</div><div><div></div><div>75%</div><div>15%</div><div>9%</div></div></div>
1	B	275	<div><div>5%</div><div><div></div><div>81%</div><div>8%</div><div>10%</div></div></div>
1	C	275	<div><div>3%</div><div><div></div><div>75%</div><div>14%</div><div>11%</div></div></div>
1	D	275	<div><div>5%</div><div><div></div><div>83%</div><div>8%</div><div>9%</div></div></div>
1	E	275	<div><div>11%</div><div><div></div><div>67%</div><div>19%</div><div>14%</div></div></div>

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
2	GOL	C	274	-	-	-	X
2	GOL	D	274	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9550 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 3-methyl-2-oxobutanoate hydroxymethyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	Se	0	3	0
			1818	1147	306	342	6	17			
1	B	248	Total	C	N	O	S	Se	0	4	0
			1793	1130	299	341	7	16			
1	C	246	Total	C	N	O	S	Se	0	5	0
			1810	1145	300	342	6	17			
1	D	249	Total	C	N	O	S	Se	0	4	0
			1816	1146	306	342	6	16			
1	E	237	Total	C	N	O	S	Se	0	2	0
			1713	1081	288	322	6	16			

There are 140 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	cloning artifact	UNP Q9JZW6
A	0	SER	-	cloning artifact	UNP Q9JZW6
A	1	LEU	-	cloning artifact	UNP Q9JZW6
A	10	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	19	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	31	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	38	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	47	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	63	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	78	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	103	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	109	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	118	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	127	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	178	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	213	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	216	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	232	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
A	264	GLU	-	cloning artifact	UNP Q9JZW6

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Chain	Residue	Modelled	Actual	Comment	Reference
A	265	GLY	-	cloning artifact	UNP Q9JZW6
A	266	GLY	-	cloning artifact	UNP Q9JZW6
A	267	SER	-	cloning artifact	UNP Q9JZW6
A	268	HIS	-	cloning artifact	UNP Q9JZW6
A	269	HIS	-	cloning artifact	UNP Q9JZW6
A	270	HIS	-	cloning artifact	UNP Q9JZW6
A	271	HIS	-	cloning artifact	UNP Q9JZW6
A	272	HIS	-	cloning artifact	UNP Q9JZW6
A	273	HIS	-	cloning artifact	UNP Q9JZW6
B	-1	MSE	-	cloning artifact	UNP Q9JZW6
B	0	SER	-	cloning artifact	UNP Q9JZW6
B	1	LEU	-	cloning artifact	UNP Q9JZW6
B	10	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	19	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	31	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	38	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	47	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	63	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	78	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	103	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	109	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	118	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	127	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	178	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	213	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	216	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	232	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
B	264	GLU	-	cloning artifact	UNP Q9JZW6
B	265	GLY	-	cloning artifact	UNP Q9JZW6
B	266	GLY	-	cloning artifact	UNP Q9JZW6
B	267	SER	-	cloning artifact	UNP Q9JZW6
B	268	HIS	-	cloning artifact	UNP Q9JZW6
B	269	HIS	-	cloning artifact	UNP Q9JZW6
B	270	HIS	-	cloning artifact	UNP Q9JZW6
B	271	HIS	-	cloning artifact	UNP Q9JZW6
B	272	HIS	-	cloning artifact	UNP Q9JZW6
B	273	HIS	-	cloning artifact	UNP Q9JZW6
C	-1	MSE	-	cloning artifact	UNP Q9JZW6
C	0	SER	-	cloning artifact	UNP Q9JZW6
C	1	LEU	-	cloning artifact	UNP Q9JZW6
C	10	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	19	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6

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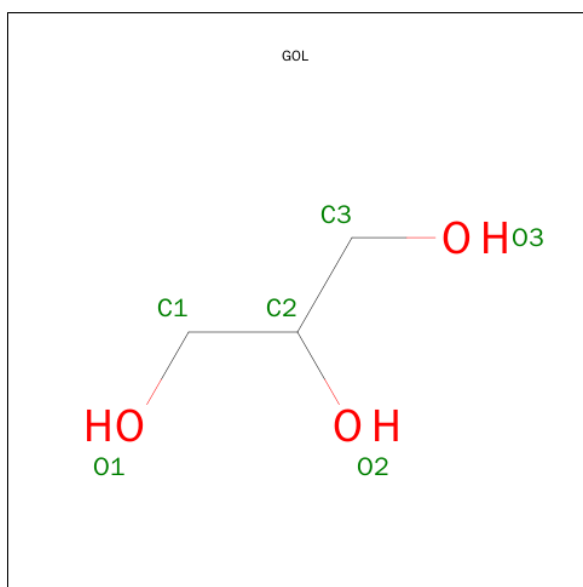
Chain	Residue	Modelled	Actual	Comment	Reference
C	31	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	38	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	47	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	63	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	78	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	103	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	109	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	118	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	127	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	178	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	213	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	216	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	232	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
C	264	GLU	-	cloning artifact	UNP Q9JZW6
C	265	GLY	-	cloning artifact	UNP Q9JZW6
C	266	GLY	-	cloning artifact	UNP Q9JZW6
C	267	SER	-	cloning artifact	UNP Q9JZW6
C	268	HIS	-	cloning artifact	UNP Q9JZW6
C	269	HIS	-	cloning artifact	UNP Q9JZW6
C	270	HIS	-	cloning artifact	UNP Q9JZW6
C	271	HIS	-	cloning artifact	UNP Q9JZW6
C	272	HIS	-	cloning artifact	UNP Q9JZW6
C	273	HIS	-	cloning artifact	UNP Q9JZW6
D	-1	MSE	-	cloning artifact	UNP Q9JZW6
D	0	SER	-	cloning artifact	UNP Q9JZW6
D	1	LEU	-	cloning artifact	UNP Q9JZW6
D	10	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	19	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	31	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	38	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	47	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	63	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	78	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	103	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	109	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	118	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	127	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	178	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	213	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	216	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	232	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
D	264	GLU	-	cloning artifact	UNP Q9JZW6

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Chain	Residue	Modelled	Actual	Comment	Reference
D	265	GLY	-	cloning artifact	UNP Q9JZW6
D	266	GLY	-	cloning artifact	UNP Q9JZW6
D	267	SER	-	cloning artifact	UNP Q9JZW6
D	268	HIS	-	cloning artifact	UNP Q9JZW6
D	269	HIS	-	cloning artifact	UNP Q9JZW6
D	270	HIS	-	cloning artifact	UNP Q9JZW6
D	271	HIS	-	cloning artifact	UNP Q9JZW6
D	272	HIS	-	cloning artifact	UNP Q9JZW6
D	273	HIS	-	cloning artifact	UNP Q9JZW6
E	-1	MSE	-	cloning artifact	UNP Q9JZW6
E	0	SER	-	cloning artifact	UNP Q9JZW6
E	1	LEU	-	cloning artifact	UNP Q9JZW6
E	10	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
E	19	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
E	31	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
E	38	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
E	47	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
E	63	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
E	78	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
E	103	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
E	109	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
E	118	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
E	127	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
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E	213	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
E	216	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
E	232	MSE	MET	MODIFIED RESIDUE	UNP Q9JZW6
E	264	GLU	-	cloning artifact	UNP Q9JZW6
E	265	GLY	-	cloning artifact	UNP Q9JZW6
E	266	GLY	-	cloning artifact	UNP Q9JZW6
E	267	SER	-	cloning artifact	UNP Q9JZW6
E	268	HIS	-	cloning artifact	UNP Q9JZW6
E	269	HIS	-	cloning artifact	UNP Q9JZW6
E	270	HIS	-	cloning artifact	UNP Q9JZW6
E	271	HIS	-	cloning artifact	UNP Q9JZW6
E	272	HIS	-	cloning artifact	UNP Q9JZW6
E	273	HIS	-	cloning artifact	UNP Q9JZW6

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).

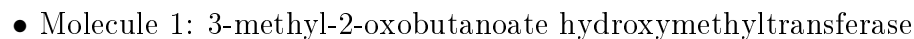
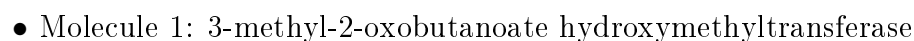


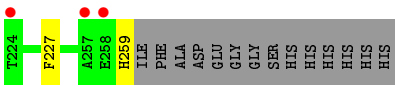
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	E	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is water.

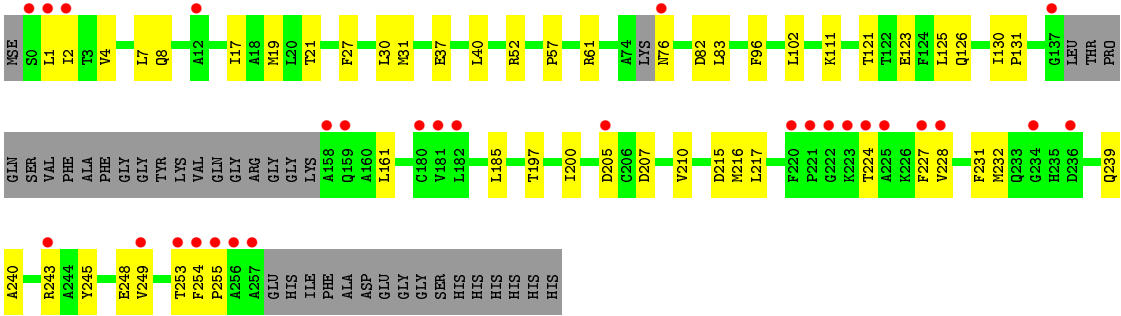
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	148	Total	O	0	0
			148	148		
3	B	98	Total	O	0	0
			98	98		
3	C	148	Total	O	0	0
			148	148		
3	D	103	Total	O	0	0
			103	103		
3	E	73	Total	O	0	0
			73	73		

- Molecule 1: 3-methyl-2-oxobutanoate hydroxymethyltransferase





● Molecule 1: 3-methyl-2-oxobutanoate hydroxymethyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	165.19Å 111.49Å 99.08Å 90.00° 122.65° 90.00°	Depositor
Resolution (Å)	43.44 – 1.75 43.50 – 1.75	Depositor EDS
% Data completeness (in resolution range)	(Not available) (43.44-1.75) 99.8 (43.50-1.75)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 1.75Å)	Xtriage
Refinement program	REFMAC 4	Depositor
R, R_{free}	0.231 , 0.250 0.220 , 0.240	Depositor DCC
R_{free} test set	15130 reflections (11.09%)	DCC
Wilson B-factor (Å ²)	17.0	Xtriage
Anisotropy	0.337	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	2 of 151511 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9550	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 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	0.29	0/1850	0.56	0/2476
1	B	0.29	0/1823	0.56	0/2440
1	C	0.30	0/1851	0.58	0/2476
1	D	0.30	0/1850	0.57	0/2475
1	E	0.27	0/1731	0.52	0/2313
All	All	0.29	0/9105	0.56	0/12180

There are no bond length outliers.

There are no bond angle outliers.

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	1818	0	1821	32	0
1	B	1793	0	1773	15	0
1	C	1810	0	1807	34	0
1	D	1816	0	1806	18	0
1	E	1713	0	1708	39	1
2	A	6	0	8	0	0
2	B	6	0	8	0	0
2	C	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	6	0	8	0	0
2	E	6	0	8	0	0
3	A	148	0	0	4	0
3	B	98	0	0	1	0
3	C	148	0	0	2	0
3	D	103	0	0	0	0
3	E	73	0	0	1	0
All	All	9550	0	8955	133	1

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 7.

All (133) 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:127[B]:MSE:HE3	1:A:128:ARG:HH12	1.10	1.11
1:C:127[A]:MSE:HE3	1:C:128:ARG:HH12	1.32	0.90
1:A:127[B]:MSE:HE3	1:A:128:ARG:NH1	1.93	0.83
1:C:161:LEU:HD23	1:C:185:LEU:HD22	1.64	0.80
1:D:5:ASN:HD21	1:D:9:LYS:HE2	1.50	0.76
1:A:123:GLU:HG2	1:A:127[A]:MSE:HE2	1.67	0.75
1:A:121:THR:O	1:A:125[B]:LEU:HD23	1.87	0.74
1:D:82:ASP:OD1	1:D:111:LYS:HE3	1.88	0.73
1:C:96:PHE:HA	1:C:125[A]:LEU:HD11	1.70	0.72
1:D:5:ASN:ND2	1:D:9:LYS:HE2	2.05	0.72
1:D:117[B]:TRP:CZ3	1:D:118:MSE:HE3	2.25	0.71
1:E:52:ARG:HD3	1:E:57:PRO:HG2	1.70	0.71
1:D:117[B]:TRP:CE3	1:D:118:MSE:HE3	2.26	0.71
1:E:96:PHE:HA	1:E:125[A]:LEU:HD11	1.73	0.70
1:E:2:ILE:HD11	1:E:7:LEU:HD21	1.74	0.69
1:C:117[B]:TRP:CZ3	1:C:118:MSE:HE3	2.28	0.69
1:B:96:PHE:HA	1:B:125[B]:LEU:HD11	1.75	0.69
1:E:2:ILE:HG22	1:E:76:ASN:O	1.93	0.68
1:C:127[A]:MSE:HE3	1:C:128:ARG:NH1	2.07	0.68
1:A:96:PHE:HA	1:A:125[B]:LEU:HD11	1.76	0.66
1:A:94:GLN:HB3	3:A:411:HOH:O	1.96	0.64
1:C:117[B]:TRP:CE3	1:C:118:MSE:HE3	2.33	0.63
1:D:162:LEU:HD23	1:D:189:VAL:HG22	1.79	0.63
1:E:245:TYR:O	1:E:249:VAL:HG23	1.99	0.63
1:A:161:LEU:HD23	1:A:185:LEU:HD22	1.80	0.62
1:A:95:ALA:N	3:A:411:HOH:O	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:161:LEU:HD23	1:E:185:LEU:HD22	1.82	0.61
1:B:92:LYS:HE3	3:B:363:HOH:O	2.01	0.61
1:B:27:PHE:HB3	1:B:31:MSE:HE2	1.84	0.60
1:A:140:PRO:HB2	3:A:394:HOH:O	2.03	0.59
1:A:123:GLU:O	1:A:127[A]:MSE:HG3	2.03	0.58
1:B:161:LEU:HD23	1:B:185:LEU:HD13	1.84	0.58
1:C:184:GLU:HG3	1:C:185:LEU:N	2.20	0.57
1:E:52:ARG:HD3	1:E:57:PRO:CG	2.34	0.56
1:E:82:ASP:OD1	1:E:111:LYS:HE2	2.05	0.56
1:C:83:LEU:HD23	1:C:102:LEU:HD12	1.88	0.55
1:A:5:ASN:HD21	1:A:9:LYS:HE3	1.72	0.55
1:D:27:PHE:HB3	1:D:31:MSE:HE2	1.87	0.55
1:A:1:LEU:HD22	1:A:1:LEU:N	2.22	0.55
1:D:161:LEU:HD23	1:D:185:LEU:HD22	1.88	0.55
1:E:4:VAL:O	1:E:8:GLN:HG3	2.07	0.55
1:B:4:VAL:O	1:B:8:GLN:HG3	2.07	0.55
1:E:231:PHE:CE2	1:E:245:TYR:HA	2.42	0.55
1:E:248:GLU:HB3	1:E:253:THR:HB	1.90	0.54
1:E:161:LEU:HD23	1:E:185:LEU:CD2	2.38	0.54
1:C:246:VAL:O	1:C:250:LYS:HG2	2.08	0.54
1:E:30:LEU:C	1:E:30:LEU:HD23	2.29	0.53
1:E:19:MSE:O	1:E:210:VAL:HA	2.09	0.53
1:E:121:THR:O	1:E:125[A]:LEU:HD23	2.09	0.52
1:D:118:MSE:HE2	1:D:118:MSE:HA	1.92	0.52
1:E:83:LEU:HD23	1:E:102:LEU:HD12	1.91	0.52
1:C:188:LYS:O	1:C:192:THR:HG23	2.10	0.52
1:A:162:LEU:O	1:A:166:LYS:HG3	2.11	0.51
1:A:88:TYR:HA	3:A:411:HOH:O	2.11	0.51
1:A:42:GLY:HA2	1:A:82:ASP:OD2	2.10	0.51
1:A:161:LEU:HD23	1:A:185:LEU:CD2	2.40	0.51
1:A:27:PHE:HB3	1:A:31:MSE:HE2	1.93	0.50
1:B:92:LYS:HD2	1:B:124:PHE:CD1	2.47	0.50
1:C:158:ALA:HA	1:C:185:LEU:HD11	1.94	0.50
1:C:125[A]:LEU:HD12	1:C:130:ILE:HD12	1.95	0.49
1:E:83:LEU:HD23	1:E:102:LEU:CD1	2.42	0.49
1:E:200:ILE:O	1:E:200:ILE:HG23	2.13	0.49
1:E:185:LEU:O	1:E:185:LEU:HD23	2.13	0.49
1:C:4:VAL:O	1:C:8:GLN:HG3	2.14	0.48
1:C:64:CYS:O	1:C:68:GLU:HG3	2.14	0.48
1:B:64:CYS:O	1:B:68:GLU:HG3	2.13	0.48
1:D:64:CYS:O	1:D:68:GLU:HG3	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:LYS:HG3	1:B:158:ALA:N	2.29	0.48
1:C:179:GLU:HG3	1:C:200:ILE:HG23	1.95	0.48
1:B:30:LEU:C	1:B:30:LEU:HD23	2.34	0.48
1:A:30:LEU:C	1:A:30:LEU:HD23	2.34	0.47
1:E:2:ILE:HD11	1:E:7:LEU:CD2	2.44	0.47
1:D:4:VAL:O	1:D:8:GLN:HG3	2.15	0.47
1:E:254:PHE:CD1	1:E:255:PRO:HA	2.50	0.47
1:E:7:LEU:HD22	1:E:17:ILE:HD13	1.98	0.46
1:C:161:LEU:HD23	1:C:185:LEU:CD2	2.39	0.46
1:E:27:PHE:HB3	1:E:31:MSE:HE2	1.96	0.46
1:B:19:MSE:HG2	1:B:210:VAL:HG23	1.98	0.46
1:A:1:LEU:H1	1:A:1:LEU:HD22	1.81	0.46
1:C:128:ARG:HG2	1:D:117[B]:TRP:CZ3	2.51	0.45
1:C:162:LEU:HD23	1:C:189:VAL:HG22	1.98	0.45
1:C:30:LEU:HD23	1:C:30:LEU:C	2.37	0.45
1:B:121:THR:O	1:B:125[B]:LEU:HD23	2.16	0.45
1:A:113:GLU:OE2	1:A:139:THR:HB	2.16	0.45
1:E:19:MSE:HG2	1:E:210:VAL:HG23	1.99	0.45
1:A:200:ILE:HG23	1:A:200:ILE:O	2.17	0.45
1:A:123:GLU:O	1:A:126:GLN:HG3	2.16	0.44
1:E:205:ASP:HB2	3:E:339:HOH:O	2.16	0.44
1:A:128:ARG:HG2	1:C:117[B]:TRP:CZ3	2.52	0.44
1:E:240:ALA:HA	1:E:243:ARG:HE	1.82	0.44
1:B:158:ALA:HA	1:B:185:LEU:HD11	2.00	0.44
1:E:227:PHE:CD1	1:E:228:VAL:HG23	2.53	0.44
1:E:1:LEU:HD23	1:E:76:ASN:HA	1.99	0.44
1:E:2:ILE:CD1	1:E:37:GLU:HB3	2.48	0.44
1:C:0:SER:C	1:C:76:ASN:HD22	2.21	0.44
1:C:215:ASP:HB3	1:C:228:VAL:HG11	2.00	0.44
1:E:217:LEU:O	1:E:232:MSE:HG3	2.18	0.44
1:A:83:LEU:HD12	1:A:112:LEU:HB3	1.99	0.43
1:A:4:VAL:O	1:A:8:GLN:HG3	2.19	0.43
1:C:27:PHE:HB3	1:C:31:MSE:HE2	2.01	0.43
1:E:123:GLU:O	1:E:126:GLN:HG3	2.18	0.43
1:D:19:MSE:HG2	1:D:210:VAL:HG23	2.01	0.43
1:C:200:ILE:HG23	1:C:200:ILE:O	2.18	0.43
1:E:224:THR:HG22	1:E:228:VAL:HB	2.01	0.43
1:C:61:ARG:NH1	3:C:318:HOH:O	2.51	0.43
1:C:138:LEU:C	1:C:138:LEU:HD13	2.40	0.43
1:A:231:PHE:O	1:A:235:HIS:HD2	2.01	0.42
1:A:113:GLU:OE2	1:A:141:GLN:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:ARG:HG2	1:D:117[B]:TRP:CH2	2.55	0.42
1:C:125[A]:LEU:CD1	1:C:130:ILE:HD12	2.49	0.42
1:E:2:ILE:HD12	1:E:37:GLU:HB3	2.01	0.42
1:A:7:LEU:HD22	1:A:17:ILE:HD13	2.01	0.42
1:A:19:MSE:O	1:A:210:VAL:HA	2.20	0.41
1:E:216:MSE:SE	1:E:217:LEU:HG	2.70	0.41
1:C:121:THR:O	1:C:125[A]:LEU:HD23	2.20	0.41
1:C:131:PRO:HG3	1:D:145:ALA:HA	2.03	0.41
1:E:21:THR:HA	1:E:40:LEU:O	2.21	0.41
1:C:177:LEU:HD23	1:C:177:LEU:C	2.41	0.41
1:E:125[A]:LEU:CD1	1:E:130:ILE:HD12	2.50	0.41
1:C:140:PRO:HB2	3:C:354:HOH:O	2.21	0.41
1:D:30:LEU:C	1:D:30:LEU:HD23	2.41	0.41
1:A:145:ALA:HA	1:E:131:PRO:HG3	2.02	0.41
1:E:197:THR:O	1:E:207:ASP:HB2	2.21	0.41
1:A:64:CYS:O	1:A:68:GLU:HG3	2.21	0.41
1:C:83:LEU:HD23	1:C:102:LEU:CD1	2.50	0.41
1:E:215:ASP:HB3	1:E:228:VAL:HG11	2.03	0.41
1:B:215:ASP:HB3	1:B:228:VAL:HG11	2.01	0.41
1:B:52:ARG:HD3	1:B:57:PRO:HB2	2.02	0.41
1:B:123:GLU:O	1:B:127[B]:MSE:HG3	2.21	0.40
1:D:227:PHE:HA	1:D:259:HIS:HB3	2.02	0.40
1:C:123:GLU:O	1:C:127[B]:MSE:HG3	2.21	0.40
1:D:19:MSE:O	1:D:210:VAL:HA	2.21	0.40
1:A:224:THR:HG22	1:A:228:VAL:HB	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:61:ARG:NH1	1:E:61:ARG:NH1[2_656]	2.12	0.08

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	249/275 (90%)	245 (98%)	4 (2%)	0	100	100
1	B	248/275 (90%)	245 (99%)	3 (1%)	0	100	100
1	C	248/275 (90%)	244 (98%)	4 (2%)	0	100	100
1	D	249/275 (90%)	246 (99%)	3 (1%)	0	100	100
1	E	233/275 (85%)	227 (97%)	6 (3%)	0	100	100
All	All	1227/1375 (89%)	1207 (98%)	20 (2%)	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	185/189 (98%)	185 (100%)	0	100	100
1	B	180/189 (95%)	179 (99%)	1 (1%)	90	83
1	C	185/189 (98%)	181 (98%)	4 (2%)	60	35
1	D	184/189 (97%)	183 (100%)	1 (0%)	92	87
1	E	170/189 (90%)	169 (99%)	1 (1%)	90	83
All	All	904/945 (96%)	897 (99%)	7 (1%)	86	77

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	185	LEU
1	C	82	ASP
1	C	126	GLN
1	C	184	GLU
1	C	185	LEU
1	D	185	LEU
1	E	239	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	235	HIS
1	B	233	GLN
1	B	235	HIS
1	C	5	ASN
1	C	233	GLN
1	D	5	ASN
1	E	8	GLN
1	E	76	ASN
1	E	159	GLN
1	E	163	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GOL	A	274	2	5,5,5	0.18	0	5,5,5	0.19	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GOL	B	274	2	5,5,5	0.18	0	5,5,5	0.19	0
2	GOL	C	274	2	5,5,5	0.17	0	5,5,5	0.20	0
2	GOL	D	274	2	5,5,5	0.17	0	5,5,5	0.20	0
2	GOL	E	274	-	5,5,5	0.23	0	5,5,5	0.24	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	274	2	-	0/4/4/4	0/0/0/0
2	GOL	B	274	2	-	0/4/4/4	0/0/0/0
2	GOL	C	274	2	-	0/4/4/4	0/0/0/0
2	GOL	D	274	2	-	0/4/4/4	0/0/0/0
2	GOL	E	274	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	234/275 (85%)	0.40	12 (5%) 32 37	7, 16, 31, 39	0
1	B	233/275 (84%)	0.50	15 (6%) 23 28	8, 17, 30, 42	0
1	C	231/275 (84%)	0.24	8 (3%) 48 53	6, 14, 27, 36	0
1	D	234/275 (85%)	0.34	14 (5%) 25 30	7, 15, 29, 40	0
1	E	222/275 (80%)	1.01	29 (13%) 5 6	9, 26, 41, 47	0
All	All	1154/1375 (83%)	0.49	78 (6%) 20 25	6, 17, 35, 47	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	224	THR	9.1
1	E	223	LYS	6.8
1	D	156	GLY	6.7
1	B	155	GLY	6.2
1	E	221	PRO	6.1
1	E	257	ALA	6.0
1	A	260	ILE	6.0
1	E	256	ALA	5.7
1	A	223	LYS	5.4
1	B	0	SER	5.3
1	E	222	GLY	5.2
1	E	253	THR	5.2
1	D	224	THR	5.1
1	B	156	GLY	5.0
1	B	182	LEU	5.0
1	E	1	LEU	4.6
1	E	0	SER	4.5
1	E	180	CYS	4.5
1	C	0	SER	4.4
1	E	227	PHE	4.1

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Mol	Chain	Res	Type	RSRZ
1	E	205	ASP	4.0
1	B	1	LEU	4.0
1	A	224	THR	3.9
1	E	236	ASP	3.9
1	B	138	LEU	3.8
1	B	222	GLY	3.8
1	D	1	LEU	3.7
1	D	257	ALA	3.5
1	D	221	PRO	3.4
1	B	12	ALA	3.3
1	C	256	ALA	3.3
1	B	227	PHE	3.3
1	E	158	ALA	3.1
1	C	1	LEU	3.1
1	A	222	GLY	3.0
1	B	223	LYS	3.0
1	D	117[A]	TRP	3.0
1	E	137	GLY	3.0
1	E	225	ALA	3.0
1	C	76	ASN	2.9
1	B	76	ASN	2.8
1	C	227	PHE	2.8
1	E	76	ASN	2.7
1	C	222	GLY	2.7
1	D	222	GLY	2.6
1	E	254	PHE	2.6
1	E	249	VAL	2.6
1	A	144	PHE	2.5
1	A	162	LEU	2.5
1	A	76	ASN	2.5
1	D	182	LEU	2.5
1	E	181	VAL	2.4
1	D	223	LYS	2.4
1	C	221	PRO	2.4
1	B	184	GLU	2.3
1	E	182	LEU	2.3
1	B	224	THR	2.3
1	E	159	GLN	2.3
1	E	234	GLY	2.3
1	D	258	GLU	2.2
1	A	117	TRP	2.2
1	E	2	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	157	LYS	2.2
1	A	142	SER	2.2
1	A	143	VAL	2.2
1	E	228	VAL	2.2
1	D	162	LEU	2.2
1	E	243	ARG	2.2
1	E	220	PHE	2.1
1	D	138	LEU	2.1
1	E	255	PRO	2.1
1	A	225	ALA	2.1
1	E	12	ALA	2.1
1	C	182	LEU	2.1
1	D	157	LYS	2.0
1	D	159	GLN	2.0
1	B	185	LEU	2.0
1	A	236	ASP	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	GOL	D	274	6/6	0.78	0.20	2.14	17,17,17,17	6
2	GOL	C	274	6/6	0.78	0.20	2.14	17,17,17,17	6
2	GOL	B	274	6/6	0.73	0.20	1.86	21,21,21,21	6
2	GOL	A	274	6/6	0.73	0.20	1.86	21,21,21,21	6
2	GOL	E	274	6/6	0.94	0.13	-	22,23,23,24	6

6.5 Other polymers ⓘ

There are no such residues in this entry.