WORLDWIDE wwPDB PROTEIN DATA BANK

Full wwPDB X-ray Structure Validation Report ⓘ

Jan 4, 2024 – 12:00 am GMT

PDB ID	:	5J78
Title	:	Crystal structure of an Acetylating Aldehyde Dehydrogenase from Geobacillus thermoglucosidasius
Authors	:	Crennell, S.J.; Extance, J.P.; Danson, M.J.
Deposited on	:	2016-04-06
Resolution	:	2.10 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	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.10 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	Percentile Ranks	Value
Rfree		0.199
Clashscore		2
Ramachandran outliers		0
Sidechain outliers		1.3%
RSRZ outliers		0.4%

■ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution
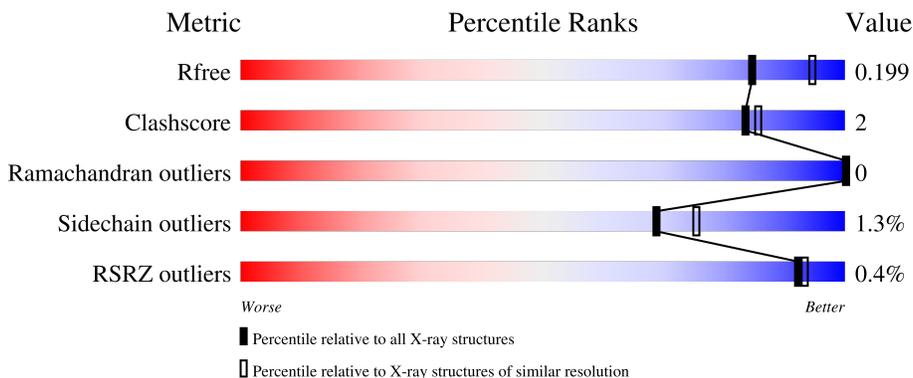
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	488	89% · 7%
1	B	488	.% 87% 6% 7%
1	C	488	88% 5% 7%
1	D	488	89% · 7%
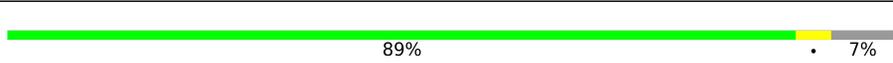
The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

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	ACT	B	1005	-	-	X	-
2	ACT	C	1005	-	-	X	-
2	ACT	D	1003	-	-	X	-

2 Entry composition ⓘ

There are 5 unique types of molecules in this entry. The entry contains 30202 atoms, of which 14701 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 Acetaldehyde dehydrogenase (Acetylating).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	453	Total	C	H	N	O	S	0	22	0
			7161	2222	3639	603	683	14			
1	B	453	Total	C	H	N	O	S	0	18	0
			7078	2202	3587	592	683	14			
1	C	453	Total	C	H	N	O	S	0	13	0
			7052	2193	3577	593	675	14			
1	D	452	Total	C	H	N	O	S	0	20	0
			7092	2209	3594	593	682	14			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP A0A0M1QQ83
A	2	GLY	-	expression tag	UNP A0A0M1QQ83
A	3	SER	-	expression tag	UNP A0A0M1QQ83
A	4	SER	-	expression tag	UNP A0A0M1QQ83
A	5	HIS	-	expression tag	UNP A0A0M1QQ83
A	6	HIS	-	expression tag	UNP A0A0M1QQ83
A	7	HIS	-	expression tag	UNP A0A0M1QQ83
A	8	HIS	-	expression tag	UNP A0A0M1QQ83
A	9	HIS	-	expression tag	UNP A0A0M1QQ83
A	10	HIS	-	expression tag	UNP A0A0M1QQ83
A	11	SER	-	expression tag	UNP A0A0M1QQ83
A	12	SER	-	expression tag	UNP A0A0M1QQ83
A	13	GLY	-	expression tag	UNP A0A0M1QQ83
A	14	LEU	-	expression tag	UNP A0A0M1QQ83
A	15	VAL	-	expression tag	UNP A0A0M1QQ83
A	16	PRO	-	expression tag	UNP A0A0M1QQ83
A	17	ARG	-	expression tag	UNP A0A0M1QQ83
A	18	GLY	-	expression tag	UNP A0A0M1QQ83
A	19	SER	-	expression tag	UNP A0A0M1QQ83
A	20	HIS	-	expression tag	UNP A0A0M1QQ83
A	21	MET	-	expression tag	UNP A0A0M1QQ83

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Chain	Residue	Modelled	Actual	Comment	Reference
A	22	ALA	-	expression tag	UNP A0A0M1QQ83
A	23	SER	-	expression tag	UNP A0A0M1QQ83
A	372	ILE	THR	conflict	UNP A0A0M1QQ83
B	1	MET	-	initiating methionine	UNP A0A0M1QQ83
B	2	GLY	-	expression tag	UNP A0A0M1QQ83
B	3	SER	-	expression tag	UNP A0A0M1QQ83
B	4	SER	-	expression tag	UNP A0A0M1QQ83
B	5	HIS	-	expression tag	UNP A0A0M1QQ83
B	6	HIS	-	expression tag	UNP A0A0M1QQ83
B	7	HIS	-	expression tag	UNP A0A0M1QQ83
B	8	HIS	-	expression tag	UNP A0A0M1QQ83
B	9	HIS	-	expression tag	UNP A0A0M1QQ83
B	10	HIS	-	expression tag	UNP A0A0M1QQ83
B	11	SER	-	expression tag	UNP A0A0M1QQ83
B	12	SER	-	expression tag	UNP A0A0M1QQ83
B	13	GLY	-	expression tag	UNP A0A0M1QQ83
B	14	LEU	-	expression tag	UNP A0A0M1QQ83
B	15	VAL	-	expression tag	UNP A0A0M1QQ83
B	16	PRO	-	expression tag	UNP A0A0M1QQ83
B	17	ARG	-	expression tag	UNP A0A0M1QQ83
B	18	GLY	-	expression tag	UNP A0A0M1QQ83
B	19	SER	-	expression tag	UNP A0A0M1QQ83
B	20	HIS	-	expression tag	UNP A0A0M1QQ83
B	21	MET	-	expression tag	UNP A0A0M1QQ83
B	22	ALA	-	expression tag	UNP A0A0M1QQ83
B	23	SER	-	expression tag	UNP A0A0M1QQ83
B	372	ILE	THR	conflict	UNP A0A0M1QQ83
C	1	MET	-	initiating methionine	UNP A0A0M1QQ83
C	2	GLY	-	expression tag	UNP A0A0M1QQ83
C	3	SER	-	expression tag	UNP A0A0M1QQ83
C	4	SER	-	expression tag	UNP A0A0M1QQ83
C	5	HIS	-	expression tag	UNP A0A0M1QQ83
C	6	HIS	-	expression tag	UNP A0A0M1QQ83
C	7	HIS	-	expression tag	UNP A0A0M1QQ83
C	8	HIS	-	expression tag	UNP A0A0M1QQ83
C	9	HIS	-	expression tag	UNP A0A0M1QQ83
C	10	HIS	-	expression tag	UNP A0A0M1QQ83
C	11	SER	-	expression tag	UNP A0A0M1QQ83
C	12	SER	-	expression tag	UNP A0A0M1QQ83
C	13	GLY	-	expression tag	UNP A0A0M1QQ83
C	14	LEU	-	expression tag	UNP A0A0M1QQ83
C	15	VAL	-	expression tag	UNP A0A0M1QQ83

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Chain	Residue	Modelled	Actual	Comment	Reference
C	16	PRO	-	expression tag	UNP A0A0M1QQ83
C	17	ARG	-	expression tag	UNP A0A0M1QQ83
C	18	GLY	-	expression tag	UNP A0A0M1QQ83
C	19	SER	-	expression tag	UNP A0A0M1QQ83
C	20	HIS	-	expression tag	UNP A0A0M1QQ83
C	21	MET	-	expression tag	UNP A0A0M1QQ83
C	22	ALA	-	expression tag	UNP A0A0M1QQ83
C	23	SER	-	expression tag	UNP A0A0M1QQ83
C	372	ILE	THR	conflict	UNP A0A0M1QQ83
D	1	MET	-	initiating methionine	UNP A0A0M1QQ83
D	2	GLY	-	expression tag	UNP A0A0M1QQ83
D	3	SER	-	expression tag	UNP A0A0M1QQ83
D	4	SER	-	expression tag	UNP A0A0M1QQ83
D	5	HIS	-	expression tag	UNP A0A0M1QQ83
D	6	HIS	-	expression tag	UNP A0A0M1QQ83
D	7	HIS	-	expression tag	UNP A0A0M1QQ83
D	8	HIS	-	expression tag	UNP A0A0M1QQ83
D	9	HIS	-	expression tag	UNP A0A0M1QQ83
D	10	HIS	-	expression tag	UNP A0A0M1QQ83
D	11	SER	-	expression tag	UNP A0A0M1QQ83
D	12	SER	-	expression tag	UNP A0A0M1QQ83
D	13	GLY	-	expression tag	UNP A0A0M1QQ83
D	14	LEU	-	expression tag	UNP A0A0M1QQ83
D	15	VAL	-	expression tag	UNP A0A0M1QQ83
D	16	PRO	-	expression tag	UNP A0A0M1QQ83
D	17	ARG	-	expression tag	UNP A0A0M1QQ83
D	18	GLY	-	expression tag	UNP A0A0M1QQ83
D	19	SER	-	expression tag	UNP A0A0M1QQ83
D	20	HIS	-	expression tag	UNP A0A0M1QQ83
D	21	MET	-	expression tag	UNP A0A0M1QQ83
D	22	ALA	-	expression tag	UNP A0A0M1QQ83
D	23	SER	-	expression tag	UNP A0A0M1QQ83
D	372	ILE	THR	conflict	UNP A0A0M1QQ83

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).

ACT

OXT

O$^-$

C

CH3

O

O

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 7	C 2	H 3	O 2	0	0
2	A	1	Total 7	C 2	H 3	O 2	0	0
2	B	1	Total 7	C 2	H 3	O 2	0	0
2	B	1	Total 7	C 2	H 3	O 2	0	0
2	C	1	Total 7	C 2	H 3	O 2	0	0
2	C	1	Total 7	C 2	H 3	O 2	0	0
2	D	1	Total 7	C 2	H 3	O 2	0	0
2	D	1	Total 7	C 2	H 3	O 2	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).

W O R L D W I D E
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GOL
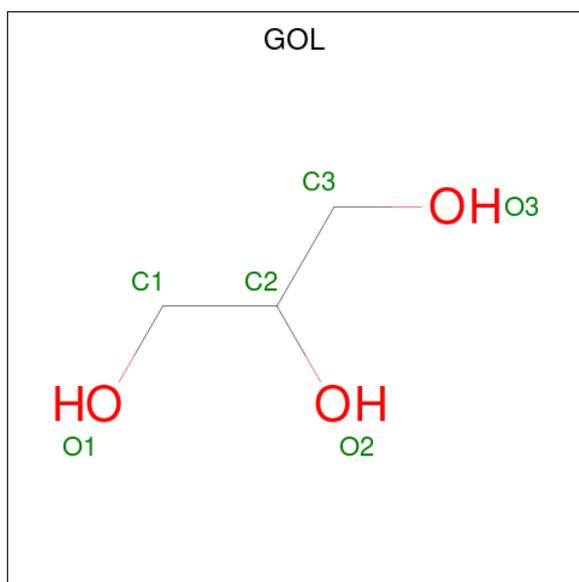
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total 14	C 3	H 8	O 3	0	0
3	B	1	Total 14	C 3	H 8	O 3	0	0
3	B	1	Total 14	C 3	H 8	O 3	0	0
3	C	1	Total 14	C 3	H 8	O 3	0	0
3	C	1	Total 14	C 3	H 8	O 3	0	0
3	C	1	Total 14	C 3	H 8	O 3	0	0
3	C	1	Total 14	C 3	H 8	O 3	0	0
3	C	1	Total 14	C 3	H 8	O 3	0	0
3	C	1	Total 14	C 3	H 8	O 3	0	0
3	C	1	Total 14	C 3	H 8	O 3	0	0
3	C	1	Total 14	C 3	H 8	O 3	0	0
3	C	1	Total 14	C 3	H 8	O 3	0	0
3	C	1	Total 14	C 3	H 8	O 3	0	0
3	C	1	Total 14	C 3	H 8	O 3	0	0
3	D	1	Total 14	C 3	H 8	O 3	0	0
3	D	1	Total 14	C 3	H 8	O 3	0	0
3	D	1	Total 14	C 3	H 8	O 3	0	0
3	D	1	Total 14	C 3	H 8	O 3	0	0
3	D	1	Total 14	C 3	H 8	O 3	0	0
3	D	1	Total 14	C 3	H 8	O 3	0	0
3	D	1	Total 14	C 3	H 8	O 3	0	0

• Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Mg 1	0	0
4	B	1	Total 1	Mg 1	0	0
4	C	1	Total 1	Mg 1	0	0
4	D	1	Total 1	Mg 1	0	0

• Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	321	Total 321	O 321	0	0
5	B	328	Total 328	O 328	0	0
5	C	307	Total 307	O 307	0	0
5	D	313	Total 313	O 313	0	0

3 Residue-property plots ⓘ

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: Acetaldehyde dehydrogenase (Acetylating)
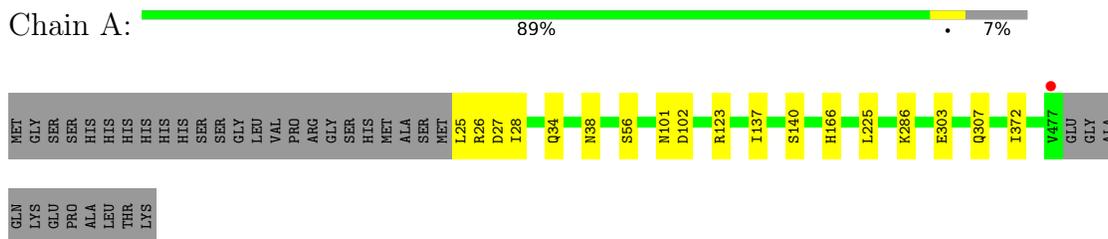
Chain A:

• Molecule 1: Acetaldehyde dehydrogenase (Acetylating)

Chain B:

• Molecule 1: Acetaldehyde dehydrogenase (Acetylating)

Chain C:

• Molecule 1: Acetaldehyde dehydrogenase (Acetylating)

Chain D:
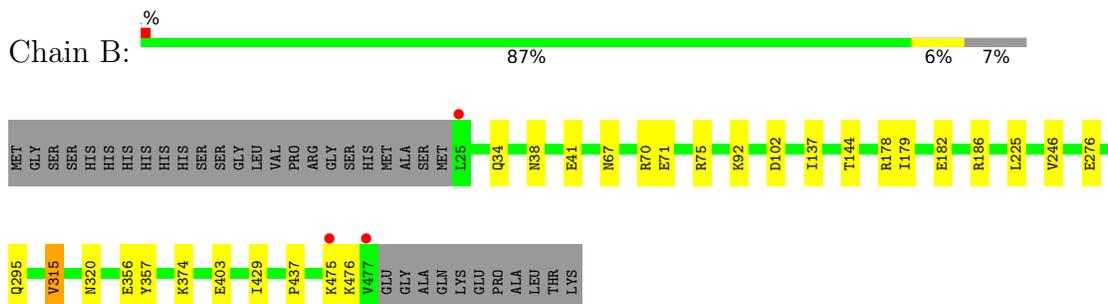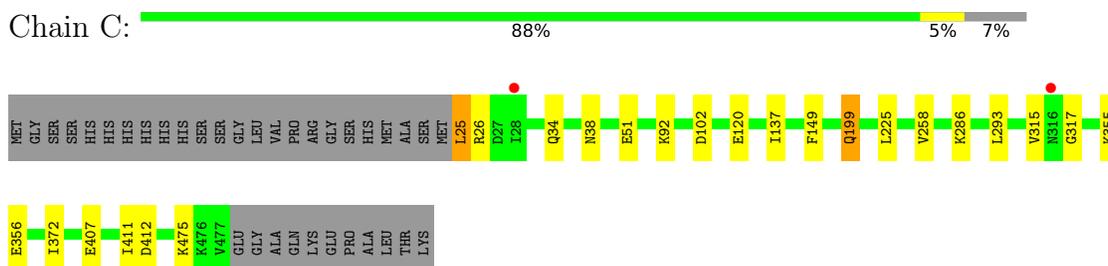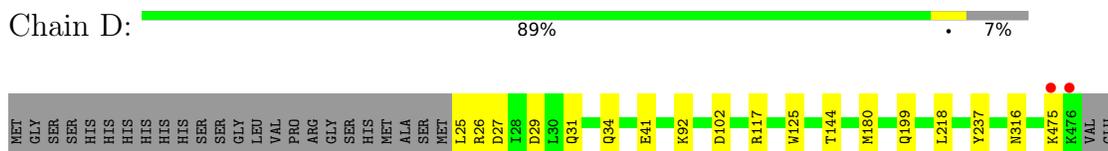
GLY ALA GLN LYS GLU PRO ALA LEU THR LYS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	98.58Å 109.12Å 196.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.29 – 2.10 49.29 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.29-2.10) 99.9 (49.29-2.10)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.29 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.143 , 0.196 0.147 , 0.199	Depositor DCC
R_{free} test set	6236 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å^2)	22.8	Xtriage
Anisotropy	0.616	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.41 , 62.2	EDS
L-test for twinning[2]	$< |L| > = 0.47, < L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	30202	wwPDB-VP
Average B, all atoms (Å^2)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.39% of the height of the origin peak. No significant pseudotranslation is detected.*

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[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: ACT, MG, CSD, 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	0.47	0/3651	0.59	0/4934
1	B	0.48	0/3595	0.58	0/4861
1	C	0.46	0/3556	0.56	0/4809
1	D	0.46	0/3611	0.58	1/4884 (0.0%)
All	All	0.47	0/14413	0.58	1/19488 (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	D	29	ASP	CB-CG-OD1	5.12	122.91	118.30

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	3522	3639	3542	19	0
1	B	3491	3587	3510	24	0
1	C	3475	3577	3533	26	0
1	D	3498	3594	3514	20	0
2	A	8	6	6	0	0
2	B	8	6	6	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	8	6	6	2	0
2	D	8	6	6	3	0
3	A	48	64	64	3	0
3	B	54	72	72	1	0
3	C	66	88	88	1	0
3	D	42	56	56	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	321	0	0	1	0
5	B	328	0	0	0	0
5	C	307	0	0	1	0
5	D	313	0	0	3	0
All	All	15501	14701	14403	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 2.

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:A:26[B]:ARG:HH21	1:C:34[B]:GLN:NE2	1.78	0.80
1:A:26[B]:ARG:HH22	1:C:38:ASN:ND2	1.80	0.78
1:A:34[B]:GLN:NE2	5:A:1101:HOH:O	2.17	0.76
1:A:26[A]:ARG:N	1:C:34[A]:GLN:OE1	2.18	0.74
1:C:92:LYS:HB3	2:C:1005:ACT:H1	1.70	0.73
1:D:117:ARG:NH1	5:D:1102:HOH:O	2.22	0.72
1:C:26[A]:ARG:O	5:C:1101:HOH:O	2.10	0.69
1:D:34[B]:GLN:NE2	5:D:1103:HOH:O	2.26	0.68
1:D:41[A]:GLU:OE1	5:D:1101:HOH:O	2.11	0.68
1:A:26[B]:ARG:HH22	1:C:38:ASN:HD21	1.43	0.66
1:A:26[B]:ARG:HH21	1:C:34[B]:GLN:HE22	1.43	0.65
1:B:34[A]:GLN:NE2	1:D:27[A]:ASP:OD2	2.34	0.60
1:C:51:GLU:OE1	3:C:1011:GOL:O3	2.20	0.58
1:A:26[B]:ARG:NH2	1:C:34[B]:GLN:NE2	2.51	0.58
1:B:178:ARG:NH1	1:B:182:GLU:OE1	2.37	0.57
1:C:92:LYS:CB	2:C:1005:ACT:H1	2.34	0.57
1:A:27[B]:ASP:CG	1:C:34[B]:GLN:HE22	2.09	0.56
1:B:67:ASN:OD1	1:B:70:ARG:NH1	2.37	0.56
1:A:303[B]:GLU:OE2	1:A:307:GLN:NE2	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27[B]:ASP:OD2	1:C:34[B]:GLN:NE2	2.40	0.54
1:B:34[A]:GLN:HE22	1:D:26[A]:ARG:HH21	1.56	0.53
1:B:144:THR:OG1	2:B:1005:ACT:H1	2.08	0.53
1:C:286:LYS:HG3	1:C:372:ILE:HD12	1.92	0.52
1:D:92:LYS:HB3	2:D:1003:ACT:H1	1.91	0.52
1:B:34[B]:GLN:OE1	1:D:26[B]:ARG:N	2.34	0.52
1:B:41:GLU:OE1	1:D:26[A]:ARG:NH1	2.42	0.52
1:D:92:LYS:HD3	2:D:1003:ACT:H1	1.93	0.51
1:C:317:GLY:O	1:C:355:LYS:NZ	2.44	0.51
1:A:101:ASN:HD22	3:A:1008:GOL:C1	2.24	0.50
1:A:38:ASN:HD21	1:C:26[B]:ARG:NH1	2.10	0.49
1:D:180:MET:HA	1:D:180:MET:HE2	1.93	0.49
1:B:34[A]:GLN:NE2	1:D:26[A]:ARG:HH21	2.10	0.49
1:B:92:LYS:HD3	2:B:1005:ACT:H2	1.95	0.49
1:C:25[A]:LEU:HD12	1:C:25[A]:LEU:O	2.12	0.49
1:A:101:ASN:HD22	3:A:1008:GOL:H12	1.78	0.49
1:A:123:ARG:HH21	3:A:1011:GOL:C1	2.26	0.48
1:C:258:VAL:HG13	1:C:293:LEU:HD23	1.97	0.47
1:B:137:ILE:HG21	1:B:225:LEU:HD21	1.95	0.47
1:D:144:THR:OG1	2:D:1003:ACT:H2	2.15	0.46
1:B:71:GLU:O	3:B:1010:GOL:O2	2.30	0.46
1:A:137:ILE:HG21	1:A:225:LEU:HD21	1.98	0.46
1:B:75:ARG:NH2	1:B:179:ILE:HD11	2.32	0.45
1:B:34[A]:GLN:NE2	1:D:26[A]:ARG:NH2	2.64	0.45
1:B:356:GLU:HG2	1:B:357:TYR:CD2	2.52	0.44
1:C:407:GLU:HG2	1:D:125[A]:TRP:CH2	2.52	0.44
1:C:34[B]:GLN:NE2	1:C:34[B]:GLN:HA	2.34	0.43
1:A:34[A]:GLN:NE2	1:C:26[A]:ARG:NH2	2.67	0.43
1:B:246:VAL:HG21	1:B:276:GLU:HG3	1.99	0.43
1:C:137:ILE:HG21	1:C:225:LEU:CD2	2.48	0.43
1:C:356:GLU:OE1	1:C:356:GLU:N	2.44	0.43
1:B:34[A]:GLN:NE2	1:B:34[A]:GLN:HA	2.33	0.43
1:B:92:LYS:HD3	2:B:1005:ACT:CH3	2.49	0.43
1:A:25[A]:LEU:HA	1:C:34[A]:GLN:OE1	2.20	0.42
1:A:140:SER:HB3	1:A:166:HIS:CG	2.55	0.42
1:D:218:LEU:HD23	1:D:237:TYR:HB2	2.02	0.42
1:B:429:ILE:O	1:B:437:PRO:HA	2.20	0.42
1:B:41:GLU:CD	1:D:26[A]:ARG:HH11	2.23	0.42
1:B:137:ILE:HG21	1:B:225:LEU:CD2	2.50	0.42
1:B:34[B]:GLN:OE1	1:D:25[B]:LEU:HA	2.20	0.41
1:A:286:LYS:HG3	1:A:372:ILE:HD12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:182:GLU:HG2	1:B:186:ARG:NH1	2.36	0.41
1:C:411:ILE:HG23	1:C:412:ASP:N	2.35	0.41
1:B:38:ASN:HD21	1:D:26[B]:ARG:NH1	2.18	0.40
1:C:120[B]:GLU:CD	1:C:120[B]:GLU:H	2.24	0.40
1:D:26[B]:ARG:O	1:D:31:GLN:NE2	2.54	0.40
1:B:315:VAL:HG13	1:B:320:ASN:HB2	2.04	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	473/488 (97%)	460 (97%)	13 (3%)	0	100	100
1	B	467/488 (96%)	458 (98%)	9 (2%)	0	100	100
1	C	462/488 (95%)	453 (98%)	9 (2%)	0	100	100
1	D	468/488 (96%)	456 (97%)	12 (3%)	0	100	100
All	All	1870/1952 (96%)	1827 (98%)	43 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	385/389 (99%)	381 (99%)	4 (1%)	76	82
1	B	379/389 (97%)	372 (98%)	7 (2%)	59	65
1	C	374/389 (96%)	367 (98%)	7 (2%)	57	63
1	D	379/389 (97%)	376 (99%)	3 (1%)	81	86
All	All	1517/1556 (98%)	1496 (99%)	21 (1%)	69	73

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28[A]	ILE
1	A	28[B]	ILE
1	A	56	SER
1	A	102	ASP
1	B	102	ASP
1	B	295	GLN
1	B	315	VAL
1	B	374	LYS
1	B	403	GLU
1	B	475	LYS
1	B	476	LYS
1	C	25[A]	LEU
1	C	25[B]	LEU
1	C	102	ASP
1	C	149	PHE
1	C	199	GLN
1	C	315	VAL
1	C	475	LYS
1	D	102	ASP
1	D	199	GLN
1	D	316	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	38	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues 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	$\#|Z| > 2$
1	CSD	C	273	1	3,7,8	0.70	0	1,8,10	1.94	0
1	CSD	A	273	1	3,7,8	0.85	0	1,8,10	1.18	0
1	CSD	D	273	1	3,7,8	0.77	0	1,8,10	0.37	0
1	CSD	B	273	1	3,7,8	0.76	0	1,8,10	0.28	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
1	CSD	C	273	1	-	0/2/6/8	-
1	CSD	A	273	1	-	1/2/6/8	-
1	CSD	D	273	1	-	1/2/6/8	-
1	CSD	B	273	1	-	0/2/6/8	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	273	CSD	N-CA-CB-SG
1	D	273	CSD	N-CA-CB-SG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 47 ligands modelled in this entry, 4 are monoatomic - leaving 43 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	C	1007	-	5,5,5	0.39	0	5,5,5	0.31	0
3	GOL	B	1007	-	5,5,5	0.35	0	5,5,5	0.27	0
3	GOL	D	1005	-	5,5,5	0.41	0	5,5,5	0.22	0
3	GOL	A	1002	-	5,5,5	0.44	0	5,5,5	0.37	0
3	GOL	B	1012	-	5,5,5	0.38	0	5,5,5	0.31	0
3	GOL	C	1006	-	5,5,5	0.39	0	5,5,5	0.54	0
3	GOL	C	1014	-	5,5,5	0.44	0	5,5,5	0.34	0
3	GOL	D	1008	-	5,5,5	0.35	0	5,5,5	0.54	0
3	GOL	B	1010	-	5,5,5	0.47	0	5,5,5	0.26	0
3	GOL	C	1011	-	5,5,5	0.34	0	5,5,5	0.23	0
3	GOL	B	1006	-	5,5,5	0.35	0	5,5,5	0.61	0
2	ACT	A	1005	-	3,3,3	1.05	0	3,3,3	0.98	0
3	GOL	B	1011	-	5,5,5	0.33	0	5,5,5	0.36	0
2	ACT	C	1001	-	3,3,3	0.64	0	3,3,3	1.44	0
3	GOL	D	1004	-	5,5,5	0.34	0	5,5,5	0.35	0
2	ACT	B	1001	-	3,3,3	0.73	0	3,3,3	1.43	0
3	GOL	A	1009	-	5,5,5	0.34	0	5,5,5	0.30	0
3	GOL	D	1006	-	5,5,5	0.33	0	5,5,5	0.26	0
3	GOL	C	1013	-	5,5,5	0.35	0	5,5,5	0.55	0
3	GOL	D	1010	-	5,5,5	0.39	0	5,5,5	0.53	0
3	GOL	B	1003	-	5,5,5	0.32	0	5,5,5	0.44	0
2	ACT	C	1005	-	3,3,3	0.67	0	3,3,3	1.62	1 (33%)
3	GOL	C	1012	-	5,5,5	0.42	0	5,5,5	0.38	0
2	ACT	D	1001	-	3,3,3	0.75	0	3,3,3	1.65	2 (66%)
3	GOL	A	1008	-	5,5,5	0.40	0	5,5,5	0.51	0
3	GOL	A	1003	-	5,5,5	0.29	0	5,5,5	0.58	0
3	GOL	C	1002	-	5,5,5	0.45	0	5,5,5	0.45	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	ACT	A	1001	-	3,3,3	0.67	0	3,3,3	1.52	0
3	GOL	B	1002	-	5,5,5	0.60	0	5,5,5	0.66	0
3	GOL	B	1009	-	5,5,5	0.33	0	5,5,5	0.24	0
3	GOL	C	1009	-	5,5,5	0.32	0	5,5,5	0.35	0
3	GOL	A	1006	-	5,5,5	0.26	0	5,5,5	0.19	0
3	GOL	A	1010	-	5,5,5	0.42	0	5,5,5	0.26	0
3	GOL	B	1008	-	5,5,5	0.33	0	5,5,5	0.24	0
3	GOL	C	1008	-	5,5,5	0.32	0	5,5,5	0.47	0
2	ACT	B	1005	-	3,3,3	1.00	0	3,3,3	1.09	0
3	GOL	A	1007	-	5,5,5	0.38	0	5,5,5	0.28	0
3	GOL	A	1011	-	5,5,5	0.42	0	5,5,5	0.42	0
3	GOL	C	1003	-	5,5,5	0.29	0	5,5,5	0.34	0
2	ACT	D	1003	-	3,3,3	0.76	0	3,3,3	1.30	0
3	GOL	D	1009	-	5,5,5	0.38	0	5,5,5	0.53	0
3	GOL	D	1007	-	5,5,5	0.30	0	5,5,5	0.32	0
3	GOL	C	1010	-	5,5,5	0.32	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	C	1007	-	-	1/4/4/4	-
3	GOL	B	1007	-	-	2/4/4/4	-
3	GOL	D	1005	-	-	2/4/4/4	-
3	GOL	A	1002	-	-	2/4/4/4	-
3	GOL	B	1012	-	-	1/4/4/4	-
3	GOL	C	1006	-	-	0/4/4/4	-
3	GOL	C	1014	-	-	2/4/4/4	-
3	GOL	D	1008	-	-	2/4/4/4	-
3	GOL	B	1010	-	-	4/4/4/4	-
3	GOL	C	1011	-	-	4/4/4/4	-
3	GOL	B	1006	-	-	0/4/4/4	-
3	GOL	B	1011	-	-	0/4/4/4	-
3	GOL	D	1004	-	-	3/4/4/4	-
3	GOL	A	1009	-	-	0/4/4/4	-
3	GOL	D	1006	-	-	4/4/4/4	-
3	GOL	C	1013	-	-	4/4/4/4	-
3	GOL	D	1010	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	1003	-	-	0/4/4/4	-
3	GOL	C	1012	-	-	1/4/4/4	-
3	GOL	A	1008	-	-	2/4/4/4	-
3	GOL	A	1003	-	-	0/4/4/4	-
3	GOL	C	1002	-	-	4/4/4/4	-
3	GOL	B	1002	-	-	2/4/4/4	-
3	GOL	B	1009	-	-	0/4/4/4	-
3	GOL	C	1009	-	-	1/4/4/4	-
3	GOL	A	1006	-	-	2/4/4/4	-
3	GOL	A	1010	-	-	3/4/4/4	-
3	GOL	B	1008	-	-	2/4/4/4	-
3	GOL	C	1008	-	-	0/4/4/4	-
3	GOL	A	1011	-	-	2/4/4/4	-
3	GOL	A	1007	-	-	2/4/4/4	-
3	GOL	C	1003	-	-	2/4/4/4	-
3	GOL	D	1009	-	-	4/4/4/4	-
3	GOL	D	1007	-	-	2/4/4/4	-
3	GOL	C	1010	-	-	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(^{o})	Ideal(^{o})
2	C	1005	ACT	OXT-C-CH3	2.23	124.39	115.18
2	D	1001	ACT	OXT-C-CH3	2.03	123.57	115.18
2	D	1001	ACT	OXT-C-O	-2.02	114.62	122.05

There are no chirality outliers.

All (66) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1002	GOL	O1-C1-C2-O2
3	A	1002	GOL	O1-C1-C2-C3
3	A	1006	GOL	O1-C1-C2-O2
3	A	1006	GOL	O1-C1-C2-C3
3	A	1007	GOL	C1-C2-C3-O3
3	A	1010	GOL	C1-C2-C3-O3
3	B	1010	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	B	1010	GOL	O2-C2-C3-O3
3	C	1002	GOL	O1-C1-C2-O2
3	C	1002	GOL	C1-C2-C3-O3
3	C	1010	GOL	O1-C1-C2-C3
3	C	1011	GOL	C1-C2-C3-O3
3	C	1013	GOL	O1-C1-C2-C3
3	C	1014	GOL	O1-C1-C2-C3
3	D	1004	GOL	O1-C1-C2-C3
3	D	1006	GOL	O1-C1-C2-O2
3	D	1006	GOL	O1-C1-C2-C3
3	D	1006	GOL	C1-C2-C3-O3
3	D	1007	GOL	O1-C1-C2-C3
3	D	1010	GOL	O1-C1-C2-C3
3	D	1010	GOL	C1-C2-C3-O3
3	B	1010	GOL	O1-C1-C2-O2
3	C	1010	GOL	O1-C1-C2-O2
3	C	1014	GOL	O1-C1-C2-O2
3	D	1007	GOL	O1-C1-C2-O2
3	A	1008	GOL	O1-C1-C2-C3
3	A	1011	GOL	O1-C1-C2-C3
3	B	1002	GOL	C1-C2-C3-O3
3	B	1010	GOL	O1-C1-C2-C3
3	B	1012	GOL	O1-C1-C2-C3
3	C	1002	GOL	O1-C1-C2-C3
3	C	1003	GOL	O1-C1-C2-C3
3	C	1007	GOL	C1-C2-C3-O3
3	C	1011	GOL	O1-C1-C2-C3
3	C	1013	GOL	C1-C2-C3-O3
3	D	1008	GOL	O1-C1-C2-C3
3	D	1009	GOL	O1-C1-C2-C3
3	D	1009	GOL	C1-C2-C3-O3
3	A	1010	GOL	O2-C2-C3-O3
3	B	1002	GOL	O2-C2-C3-O3
3	C	1002	GOL	O2-C2-C3-O3
3	C	1003	GOL	O1-C1-C2-O2
3	C	1011	GOL	O1-C1-C2-O2
3	C	1011	GOL	O2-C2-C3-O3
3	D	1004	GOL	O1-C1-C2-O2
3	D	1008	GOL	O1-C1-C2-O2
3	D	1010	GOL	O1-C1-C2-O2
3	D	1010	GOL	O2-C2-C3-O3
3	A	1007	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	C	1013	GOL	O1-C1-C2-O2
3	C	1013	GOL	O2-C2-C3-O3
3	D	1006	GOL	O2-C2-C3-O3
3	A	1008	GOL	O1-C1-C2-O2
3	B	1008	GOL	O2-C2-C3-O3
3	D	1005	GOL	O1-C1-C2-O2
3	B	1007	GOL	C1-C2-C3-O3
3	A	1010	GOL	O1-C1-C2-C3
3	B	1007	GOL	O2-C2-C3-O3
3	B	1008	GOL	C1-C2-C3-O3
3	D	1004	GOL	C1-C2-C3-O3
3	D	1005	GOL	O1-C1-C2-C3
3	A	1011	GOL	O1-C1-C2-O2
3	D	1009	GOL	O1-C1-C2-O2
3	D	1009	GOL	O2-C2-C3-O3
3	C	1012	GOL	O2-C2-C3-O3
3	C	1009	GOL	C1-C2-C3-O3

There are no ring outliers.

7 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1010	GOL	1	0
3	C	1011	GOL	1	0
2	C	1005	ACT	2	0
3	A	1008	GOL	2	0
2	B	1005	ACT	3	0
3	A	1011	GOL	1	0
2	D	1003	ACT	3	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(Å²)	Q<0.9
1	A	452/488 (92%)	-0.51	1 (0%)	95	95	15, 23, 46, 74	0
1	B	452/488 (92%)	-0.50	3 (0%)	87	89	14, 21, 41, 77	0
1	C	452/488 (92%)	-0.43	2 (0%)	92	93	14, 24, 48, 77	0
1	D	451/488 (92%)	-0.50	2 (0%)	92	93	14, 22, 45, 79	0
All	All	1807/1952 (92%)	-0.48	8 (0%)	92	93	14, 23, 45, 79	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	476	LYS	3.3
1	D	475	LYS	2.8
1	A	477	VAL	2.7
1	B	25[A]	LEU	2.4
1	B	477	VAL	2.4
1	B	475	LYS	2.3
1	C	316	ASN	2.2
1	C	28[A]	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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(Å²)	Q<0.9
1	CSD	A	273	8/9	0.99	0.12	15,21,33,40	0
1	CSD	B	273	8/9	0.99	0.10	15,19,31,37	0
1	CSD	C	273	8/9	0.99	0.13	13,20,26,30	0
1	CSD	D	273	8/9	0.99	0.09	16,20,31,37	0

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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	MG	B	1004	1/1	0.68	0.11	49,49,49,49	0
3	GOL	A	1009	6/6	0.77	0.25	41,57,75,90	0
4	MG	C	1004	1/1	0.80	0.11	43,43,43,43	0
3	GOL	B	1011	6/6	0.83	0.19	55,66,80,82	0
3	GOL	D	1009	6/6	0.83	0.21	42,53,74,89	0
3	GOL	B	1009	6/6	0.85	0.16	40,59,88,101	0
3	GOL	D	1010	6/6	0.85	0.14	41,55,66,67	0
3	GOL	A	1011	6/6	0.85	0.15	36,55,67,72	0
3	GOL	C	1013	6/6	0.85	0.20	44,56,67,74	0
3	GOL	C	1014	6/6	0.86	0.15	52,62,71,79	0
3	GOL	A	1003	6/6	0.88	0.12	43,64,83,83	0
3	GOL	C	1003	6/6	0.88	0.12	48,65,79,81	0
3	GOL	B	1007	6/6	0.89	0.17	36,51,80,96	0
3	GOL	B	1003	6/6	0.89	0.17	51,63,76,80	0
3	GOL	C	1002	6/6	0.90	0.18	29,53,77,77	0
3	GOL	A	1010	6/6	0.90	0.17	37,50,66,66	0
3	GOL	D	1007	6/6	0.90	0.24	41,60,71,72	0
3	GOL	D	1008	6/6	0.90	0.13	36,52,63,73	0
3	GOL	C	1011	6/6	0.91	0.24	36,52,70,84	0
3	GOL	B	1010	6/6	0.91	0.17	40,53,64,64	0
3	GOL	B	1002	6/6	0.91	0.16	28,46,71,71	0
3	GOL	B	1008	6/6	0.92	0.10	35,50,58,60	0
3	GOL	D	1006	6/6	0.92	0.16	35,52,71,72	0
3	GOL	C	1012	6/6	0.92	0.10	52,63,71,82	0
3	GOL	A	1007	6/6	0.92	0.17	34,58,92,110	0
3	GOL	C	1010	6/6	0.93	0.15	38,56,74,89	0
3	GOL	A	1006	6/6	0.93	0.16	29,44,68,76	0
3	GOL	B	1012	6/6	0.93	0.13	35,51,70,80	0
3	GOL	A	1002	6/6	0.93	0.15	31,49,59,64	0
3	GOL	A	1008	6/6	0.93	0.13	35,43,52,62	0
3	GOL	C	1007	6/6	0.93	0.16	39,61,73,77	0
4	MG	A	1004	1/1	0.94	0.04	55,55,55,55	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	GOL	D	1005	6/6	0.94	0.09	32,42,59,59	0
3	GOL	C	1009	6/6	0.94	0.14	38,53,61,64	0
4	MG	D	1002	1/1	0.94	0.12	49,49,49,49	0
3	GOL	C	1006	6/6	0.95	0.18	31,44,60,72	0
3	GOL	C	1008	6/6	0.95	0.12	35,43,51,61	0
2	ACT	B	1001	4/4	0.96	0.10	30,34,39,40	0
2	ACT	C	1001	4/4	0.96	0.11	24,28,34,36	0
3	GOL	B	1006	6/6	0.96	0.14	32,44,73,88	0
2	ACT	D	1003	4/4	0.97	0.21	18,34,64,64	0
3	GOL	D	1004	6/6	0.97	0.15	30,48,58,67	0
2	ACT	A	1001	4/4	0.98	0.09	29,34,34,35	0
2	ACT	B	1005	4/4	0.98	0.17	12,32,54,54	0
2	ACT	A	1005	4/4	0.98	0.18	11,26,59,59	0
2	ACT	C	1005	4/4	0.98	0.21	14,30,60,60	0
2	ACT	D	1001	4/4	0.98	0.10	25,35,38,41	0

6.5 Other polymers (i)

There are no such residues in this entry.