



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 3ZVH / pdb_00003zvh
Title : Methylaspartate ammonia lyase from Clostridium tetanomorphum mutant Q73A
Authors : Raj, H.; Szymanski, W.; de Villiers, J.; Rozeboom, H.J.; Veetil, V.P.; Reis, C.R.; de Villiers, M.; de Wildeman, S.; Dekker, F.J.; Quax, W.J.; Thunnissen, A.M.W.H.; Feringa, B.L.; Janssen, D.B.; Poelarends, G.J.
Deposited on : 2011-07-25
Resolution : 1.99 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

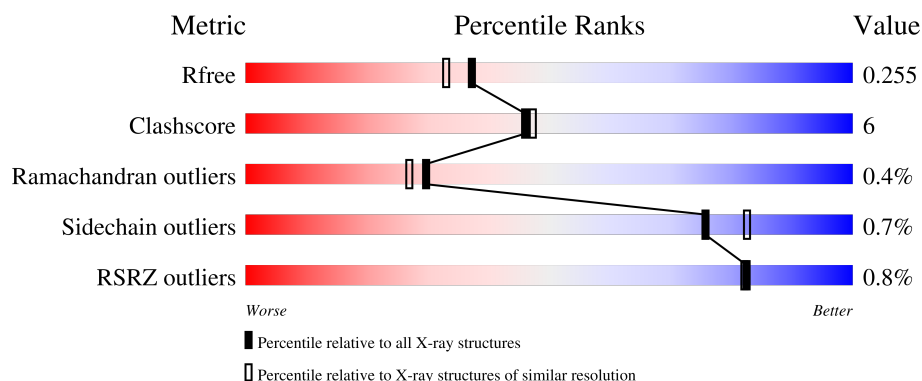
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	438	 82% 12% 5%
2	B	438	 79% 15% 5%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7001 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 METHYLASPARTATE AMMONIA-LYASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	415	3217	2020	560	611	26	0	2	0

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	414	LYS	-	expression tag	UNP Q05514
A	415	LEU	-	expression tag	UNP Q05514
A	416	GLY	-	expression tag	UNP Q05514
A	417	PRO	-	expression tag	UNP Q05514
A	418	GLU	-	expression tag	UNP Q05514
A	419	GLN	-	expression tag	UNP Q05514
A	420	LYS	-	expression tag	UNP Q05514
A	421	LEU	-	expression tag	UNP Q05514
A	422	ILE	-	expression tag	UNP Q05514
A	423	SER	-	expression tag	UNP Q05514
A	424	GLU	-	expression tag	UNP Q05514
A	425	GLU	-	expression tag	UNP Q05514
A	426	ASP	-	expression tag	UNP Q05514
A	427	LEU	-	expression tag	UNP Q05514
A	428	ASN	-	expression tag	UNP Q05514
A	429	SER	-	expression tag	UNP Q05514
A	430	ALA	-	expression tag	UNP Q05514
A	431	VAL	-	expression tag	UNP Q05514
A	432	ASP	-	expression tag	UNP Q05514
A	433	HIS	-	expression tag	UNP Q05514
A	434	HIS	-	expression tag	UNP Q05514
A	435	HIS	-	expression tag	UNP Q05514
A	436	HIS	-	expression tag	UNP Q05514
A	437	HIS	-	expression tag	UNP Q05514
A	438	HIS	-	expression tag	UNP Q05514
A	73	ALA	GLN	engineered mutation	UNP Q05514

- Molecule 2 is a protein called METHYLASPARTATE AMMONIA-LYASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	415	Total	C	N	O	S	0	2	0
			3219	2020	560	613	26			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	414	LYS	-	expression tag	UNP Q05514
B	415	LEU	-	expression tag	UNP Q05514
B	416	GLY	-	expression tag	UNP Q05514
B	417	PRO	-	expression tag	UNP Q05514
B	418	GLU	-	expression tag	UNP Q05514
B	419	GLN	-	expression tag	UNP Q05514
B	420	LYS	-	expression tag	UNP Q05514
B	421	LEU	-	expression tag	UNP Q05514
B	422	ILE	-	expression tag	UNP Q05514
B	423	SER	-	expression tag	UNP Q05514
B	424	GLU	-	expression tag	UNP Q05514
B	425	GLU	-	expression tag	UNP Q05514
B	426	ASP	-	expression tag	UNP Q05514
B	427	LEU	-	expression tag	UNP Q05514
B	428	ASN	-	expression tag	UNP Q05514
B	429	SER	-	expression tag	UNP Q05514
B	430	ALA	-	expression tag	UNP Q05514
B	431	VAL	-	expression tag	UNP Q05514
B	432	ASP	-	expression tag	UNP Q05514
B	433	HIS	-	expression tag	UNP Q05514
B	434	HIS	-	expression tag	UNP Q05514
B	435	HIS	-	expression tag	UNP Q05514
B	436	HIS	-	expression tag	UNP Q05514
B	437	HIS	-	expression tag	UNP Q05514
B	438	HIS	-	expression tag	UNP Q05514
B	73	ALA	GLN	engineered mutation	UNP Q05514

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		

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



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

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Cl 1 1	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	261	Total O 261 261	0	0
6	B	271	Total O 271 271	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.79Å 109.70Å 110.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	77.71 – 1.99 77.71 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.7 (77.71-1.99) 97.6 (77.71-2.00)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.195 , 0.249 0.222 , 0.255	Depositor DCC
R_{free} test set	2744 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	14.5	Xtriage
Anisotropy	0.699	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 29.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.038 for -h,l,k	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	7001	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.95% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CL, OCS, CSO, 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	0.62	0/3258	0.92	0/4390
2	B	0.65	0/3258	0.88	2/4390 (0.0%)
All	All	0.63	0/6516	0.90	2/8780 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	279	GLU	N-CA-C	5.59	120.10	113.28
2	B	398	VAL	N-CA-C	5.41	115.62	110.53

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	3217	0	3228	32	0
2	B	3219	0	3227	46	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	18	0	24	0	0
4	B	12	0	16	3	0
5	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	261	0	0	1	0
6	B	271	0	0	5	0
All	All	7001	0	6495	75	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 6.

All (75) 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:104:ARG:HH22	1:A:115:GLU:CD	1.84	0.84
2:B:117:ASP:O	2:B:131:ARG:HD3	1.77	0.84
1:A:30:ASP:O	1:A:33:THR:HG22	1.83	0.79
2:B:263:GLU:HA	2:B:266:LYS:HD2	1.68	0.76
1:A:104:ARG:NH2	1:A:115:GLU:OE1	2.20	0.75
2:B:213:GLU:HG2	6:B:2165:HOH:O	1.88	0.73
2:B:330:ILE:HD13	2:B:345:ILE:HG12	1.69	0.73
2:B:259:GLN:O	2:B:263:GLU:HG2	1.90	0.72
1:A:2:LYS:N	1:A:60:GLU:OE1	2.15	0.70
2:B:201:GLU:HB2	4:B:1417:GOL:O1	1.90	0.70
2:B:199:ASN:HD21	4:B:1417:GOL:H32	1.58	0.69
2:B:117:ASP:OD1	2:B:131:ARG:HD2	1.94	0.68
2:B:80:ARG:HG2	2:B:80:ARG:HH11	1.57	0.68
2:B:330:ILE:HD11	2:B:355:ALA:HB1	1.75	0.67
2:B:204:LEU:HG	2:B:208:GLY:HA2	1.78	0.65
2:B:199:ASN:HD21	4:B:1417:GOL:C3	2.09	0.65
1:A:332:THR:OG1	1:A:333:PRO:HD3	2.00	0.62
2:B:80:ARG:HG2	2:B:80:ARG:NH1	2.14	0.62
1:A:100:LYS:HD3	1:A:119:MET:HE3	1.80	0.62
2:B:228:ARG:NH1	2:B:230:ASP:OD1	2.33	0.61
2:B:330:ILE:HD13	2:B:345:ILE:CG1	2.31	0.61
2:B:55:VAL:HG11	2:B:134:ILE:HD13	1.83	0.60
2:B:167:VAL:HG11	2:B:383:VAL:HG13	1.89	0.54
2:B:81:ASP:OD2	2:B:128:THR:HG23	2.08	0.54
1:A:118:LYS:HA	1:A:125:ARG:HH21	1.74	0.53
2:B:357:CYS:HB3	2:B:383:VAL:HG12	1.89	0.53
2:B:250:ASP:HB3	2:B:253:ALA:HB3	1.90	0.52
2:B:275:PRO:HG3	2:B:304:LEU:HD22	1.91	0.52
1:A:81:ASP:OD2	1:A:128:THR:HG23	2.10	0.52
2:B:127:HIS:HB3	2:B:130:ILE:HD12	1.92	0.52
1:A:271:ARG:HG2	1:A:303:GLU:HB2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:LEU:HD21	2:B:404:ARG:HD2	1.94	0.50
2:B:44:THR:HG23	6:B:2031:HOH:O	2.11	0.50
2:B:330:ILE:CD1	2:B:345:ILE:HG12	2.41	0.49
1:A:327:MET:HE3	1:A:356:TYR:HB2	1.94	0.49
1:A:1:MET:SD	1:A:60:GLU:HG2	2.54	0.48
2:B:86:ALA:O	2:B:90:ILE:HG13	2.13	0.48
2:B:144:LYS:HA	6:B:2108:HOH:O	2.13	0.48
1:A:333:PRO:HD3	1:A:359:GLY:HA2	1.95	0.47
1:A:211:LEU:HD11	1:A:237:ILE:HD13	1.96	0.47
2:B:269:HIS:HB2	6:B:2078:HOH:O	2.14	0.47
2:B:57:LEU:HB2	2:B:65:ALA:HB3	1.95	0.47
6:A:2260:HOH:O	2:B:64:VAL:HB	2.15	0.47
2:B:14:GLY:O	2:B:388:GLY:HA2	2.14	0.47
2:B:200:VAL:O	2:B:205:GLY:N	2.46	0.47
1:A:333:PRO:HG3	1:A:360:THR:HG22	1.96	0.46
1:A:43:PHE:CE2	1:A:83:LEU:HD13	2.51	0.46
1:A:357:CYS:HB3	1:A:383:VAL:HG12	1.97	0.46
2:B:405:VAL:O	2:B:409:VAL:HG23	2.16	0.46
1:A:75:SER:HB3	1:A:83:LEU:HD12	1.97	0.45
1:A:314:GLU:HG2	1:A:315:ASP:N	2.31	0.45
1:A:56:LEU:HA	1:A:65:ALA:O	2.17	0.45
2:B:368:ALA:HA	2:B:386:LYS:HE3	2.00	0.44
1:A:98:ALA:HB3	1:A:99:PRO:HD3	1.99	0.43
1:A:274:GLY:O	1:A:275:PRO:C	2.61	0.43
1:A:187:LYS:HE3	2:B:16:TYR:CE2	2.53	0.43
2:B:59:LEU:HD22	2:B:106:ILE:HD11	2.00	0.43
2:B:98:ALA:HB3	2:B:99:PRO:HD3	2.00	0.43
1:A:70:ALA:HB1	1:A:363:GLU:HA	2.01	0.42
1:A:175:ASP:O	1:A:177:ARG:N	2.52	0.42
2:B:327:MET:SD	2:B:356:TYR:HB2	2.59	0.42
2:B:139:LEU:HG	2:B:150:MET:HG2	2.02	0.41
2:B:395[B]:MET:HE2	2:B:395[B]:MET:HB3	1.95	0.41
2:B:192:LEU:HD22	2:B:233:PRO:HB2	2.02	0.41
2:B:331:LYS:HA	6:B:2243:HOH:O	2.20	0.41
1:A:386:LYS:HB2	1:A:387:PRO:HA	2.03	0.41
1:A:395[B]:MET:HE3	1:A:395[B]:MET:HB3	1.85	0.41
1:A:44:THR:N	1:A:48:GLN:OE1	2.42	0.41
1:A:69:CYS:SG	1:A:130:ILE:HA	2.61	0.41
1:A:407:ALA:HB2	2:B:9:THR:CG2	2.50	0.41
2:B:1:MET:HE2	2:B:1:MET:HB2	1.94	0.41
2:B:89:PHE:HB3	2:B:130:ILE:HD11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:263:GLU:CA	2:B:266:LYS:HD2	2.46	0.41
1:A:109:PHE:CE1	1:A:139:LEU:HB2	2.55	0.40
1:A:255:ALA:CB	1:A:294:GLU:HG2	2.51	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	414/438 (94%)	394 (95%)	18 (4%)	2 (0%)	24	21
2	B	414/438 (94%)	389 (94%)	24 (6%)	1 (0%)	43	42
All	All	828/876 (94%)	783 (95%)	42 (5%)	3 (0%)	30	27

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	278	VAL
1	A	249	VAL
2	B	249	VAL

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	336/355 (95%)	333 (99%)	3 (1%)	70	78
2	B	336/355 (95%)	334 (99%)	2 (1%)	78	85
All	All	672/710 (95%)	667 (99%)	5 (1%)	76	82

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	THR
1	A	249	VAL
1	A	314	GLU
2	B	119	MET
2	B	192	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	ASN
1	A	199	ASN
1	A	329	GLN
2	B	199	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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	CSO	A	361	1	3,6,7	0.64	0	1,6,8	0.10	0
2	OCS	B	361	2	6,8,9	1.12	0	7,11,13	1.72	2 (28%)

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	CSO	A	361	1	-	0/1/5/7	-
2	OCS	B	361	2	-	0/4/7/9	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	361	OCS	OD2-SG-CB	2.61	111.02	105.97
2	B	361	OCS	OD1-SG-CB	2.30	110.19	106.76

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 3 are monoatomic - leaving 5 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
4	GOL	A	1417	-	5,5,5	0.35	0	5,5,5	0.28	0
4	GOL	A	1418	-	5,5,5	0.41	0	5,5,5	0.25	0
4	GOL	B	1416	-	5,5,5	0.41	0	5,5,5	1.56	1 (20%)
4	GOL	A	1416	-	5,5,5	0.37	0	5,5,5	0.49	0
4	GOL	B	1417	-	5,5,5	0.37	0	5,5,5	0.39	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	A	1417	-	-	2/4/4/4	-
4	GOL	A	1418	-	-	2/4/4/4	-
4	GOL	B	1416	-	-	4/4/4/4	-
4	GOL	A	1416	-	-	2/4/4/4	-
4	GOL	B	1417	-	-	4/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1416	GOL	O2-C2-C1	3.03	121.71	109.18

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1416	GOL	O1-C1-C2-C3
4	A	1418	GOL	O1-C1-C2-O2
4	A	1418	GOL	O1-C1-C2-C3
4	B	1416	GOL	O1-C1-C2-C3
4	B	1417	GOL	C1-C2-C3-O3
4	B	1417	GOL	O2-C2-C3-O3
4	B	1416	GOL	O2-C2-C3-O3
4	A	1417	GOL	O1-C1-C2-C3
4	B	1417	GOL	O1-C1-C2-C3
4	A	1416	GOL	O1-C1-C2-O2
4	B	1416	GOL	O1-C1-C2-O2
4	A	1417	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
4	B	1417	GOL	O1-C1-C2-O2
4	B	1416	GOL	C1-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1417	GOL	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	414/438 (94%)	0.22	6 (1%) 73 73	4, 14, 26, 32	2 (0%)
2	B	414/438 (94%)	0.14	1 (0%) 91 90	4, 13, 23, 31	2 (0%)
All	All	828/876 (94%)	0.18	7 (0%) 82 82	4, 14, 25, 32	4 (0%)

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	65	ALA	2.7
2	B	415	LEU	2.4
1	A	253	ALA	2.4
1	A	25	LYS	2.3
1	A	206	LEU	2.3
1	A	258	ILE	2.1
1	A	199	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	OCS	B	361	9/10	0.92	0.11	9,11,19,19	0
1	CSO	A	361	7/8	0.96	0.07	9,10,13,17	0

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($\text{\AA}^2$)	Q<0.9
4	GOL	B	1417	6/6	0.70	0.19	30,32,33,33	0
4	GOL	A	1418	6/6	0.76	0.15	29,31,31,31	0
4	GOL	A	1417	6/6	0.83	0.16	19,22,23,25	0
4	GOL	A	1416	6/6	0.94	0.07	8,11,13,14	0
4	GOL	B	1416	6/6	0.96	0.06	6,8,11,12	0
3	MG	B	1421	1/1	0.97	0.03	8,8,8,8	0
5	CL	B	1418	1/1	0.97	0.04	17,17,17,17	0
3	MG	A	1421	1/1	0.99	0.02	14,14,14,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.