



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 07:53 pm GMT

PDB ID : 4IA6
Title : Hydratase from lactobacillus acidophilus in a ligand bound form (LA LAH)
Authors : Khoshnevis, S.; Neumann, P.; Ficner, R.
Deposited on : 2012-12-06
Resolution : 1.80 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

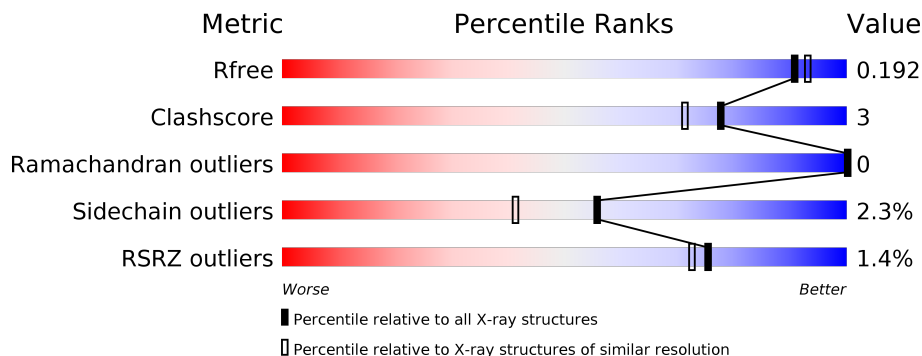
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	100719	4827 (1.80-1.80)
Clashscore	112137	5742 (1.80-1.80)
Ramachandran outliers	110173	5676 (1.80-1.80)
Sidechain outliers	110143	5675 (1.80-1.80)
RSRZ outliers	101464	4906 (1.80-1.80)

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 $\leq 5\%$. 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	591	<div> <div style="width: 92%;"></div> <div style="width: 6%;"></div> <div style="width: 2%;"></div> <div style="width: 2%;"></div> <div style="width: 2%;"></div> </div>
1	B	591	<div> <div style="width: 87%;"></div> <div style="width: 7%;"></div> <div style="width: 6%;"></div> <div style="width: 2%;"></div> <div style="width: 2%;"></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
11	NO3	B	610	-	-	-	X
2	MRD	A	601[A]	-	-	X	X
3	MPD	A	602[B]	-	-	-	X
3	MPD	A	603	-	-	-	X
3	MPD	A	611	-	-	-	X
3	MPD	B	601	-	-	-	X
3	MPD	B	611[B]	-	-	-	X
4	GOL	A	604	-	-	-	X
4	GOL	A	605	-	-	-	X
4	GOL	A	606[A]	-	-	-	X
4	GOL	A	606[B]	-	-	-	X
4	GOL	A	607[A]	-	-	-	X
4	GOL	A	607[B]	-	-	-	X
4	GOL	B	604	-	-	-	X
4	GOL	B	605	-	-	-	X
4	GOL	B	609	-	-	-	X
5	EIC	A	608	-	-	-	X
7	PO4	A	610	-	-	-	X
7	PO4	A	616	-	-	-	X
7	PO4	B	608	-	-	-	X
8	MES	A	615	-	-	-	X
8	MES	B	602	-	-	-	X
8	MES	B	603	-	-	-	X

2 Entry composition [i](#)

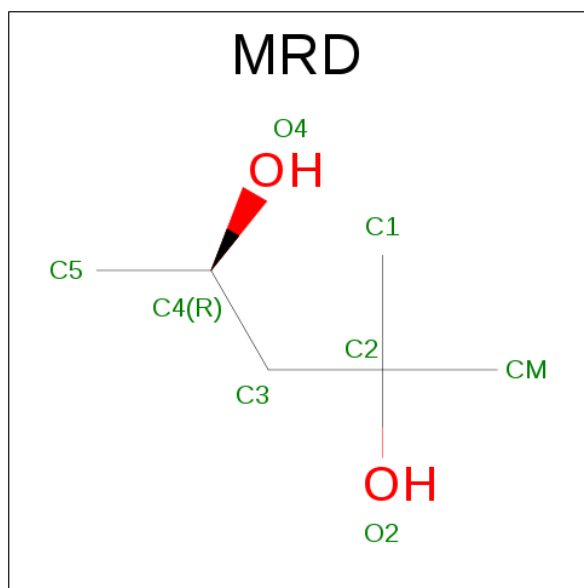
There are 12 unique types of molecules in this entry. The entry contains 10974 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Myosin-crossreactive antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	579	Total	C	N	O	S	0	8	0
			4750	3031	782	912	25			
1	B	558	Total	C	N	O	S	0	6	0
			4581	2928	753	877	23			

- Molecule 2 is (4R)-2-METHYLPENTANE-2,4-DIOL (three-letter code: MRD) (formula: C₆H₁₄O₂).



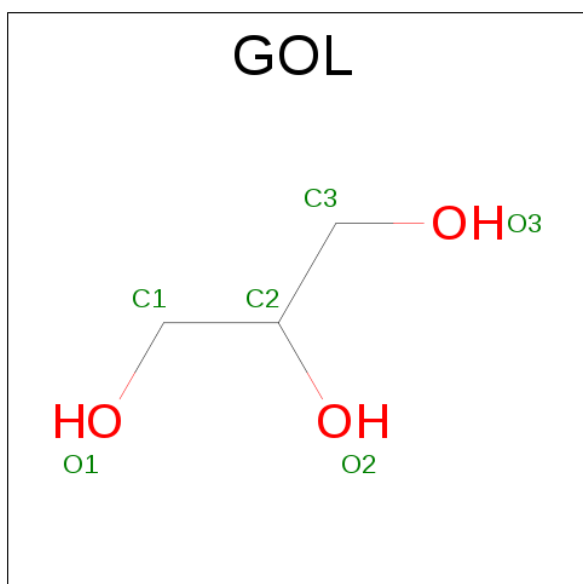
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	1
			8	6	2		

- Molecule 3 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C₆H₁₄O₂).



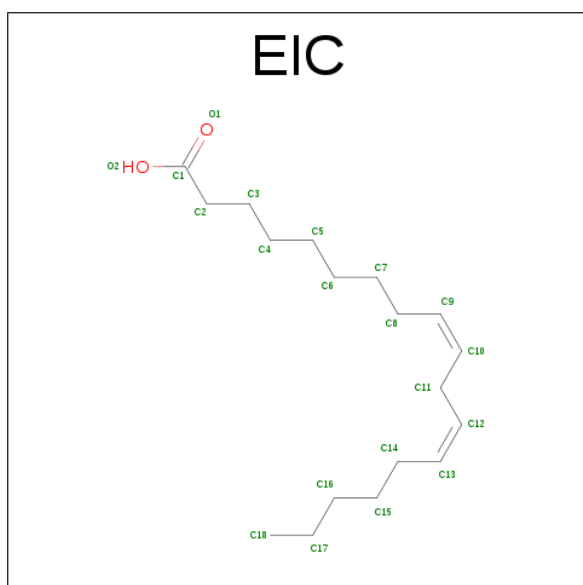
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	1
			8	6	2		
3	A	1	Total	C	O	0	0
			8	6	2		
3	A	1	Total	C	O	0	0
			8	6	2		
3	B	1	Total	C	O	0	0
			8	6	2		
3	B	1	Total	C	O	0	1
			8	6	2		

- Molecule 4 is GLYCEROL (three-letter code: 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 12 6 6	0	1
4	A	1	Total C O 12 6 6	0	1
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
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 LINOLEIC ACID (three-letter code: EIC) (formula: $C_{18}H_{32}O_2$).

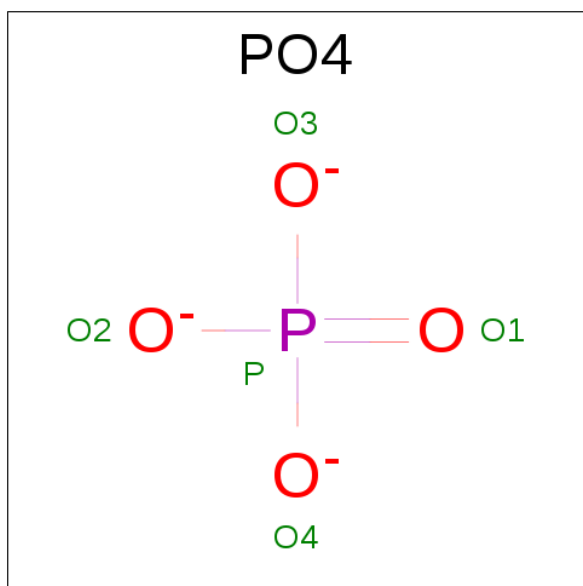


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 20 18 2	0	0

- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

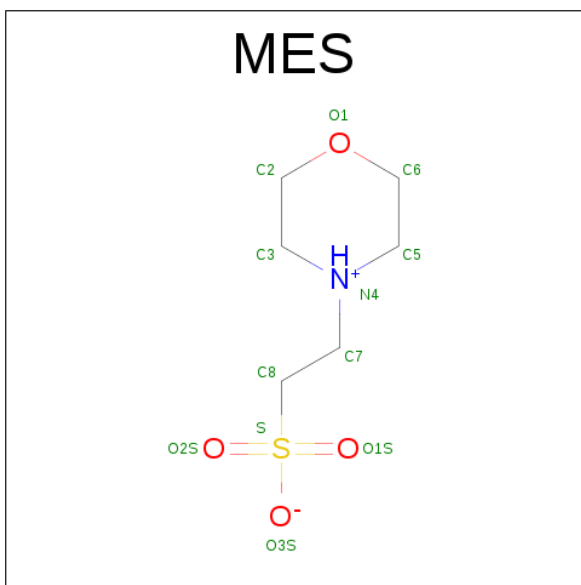
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	2	Total	Na	0	0
			2	2		

- Molecule 7 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	P	0	0
			5	4	1		
7	A	1	Total	O	P	0	0
			5	4	1		
7	B	1	Total	O	P	0	0
			5	4	1		
7	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 8 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).

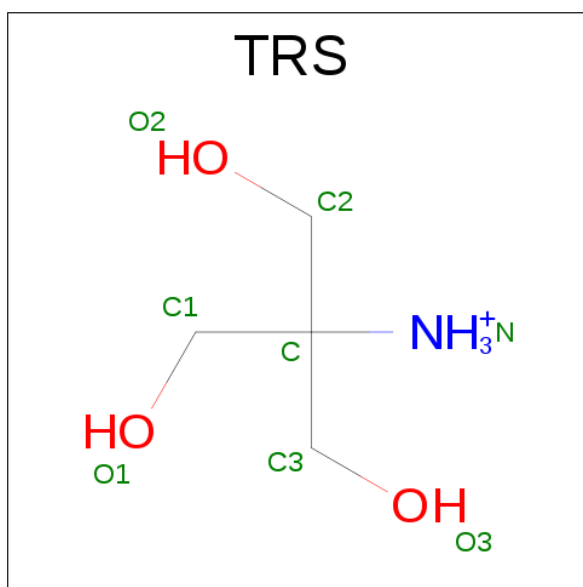


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
8	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
8	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 9 is POTASSIUM ION (three-letter code: K) (formula: K).

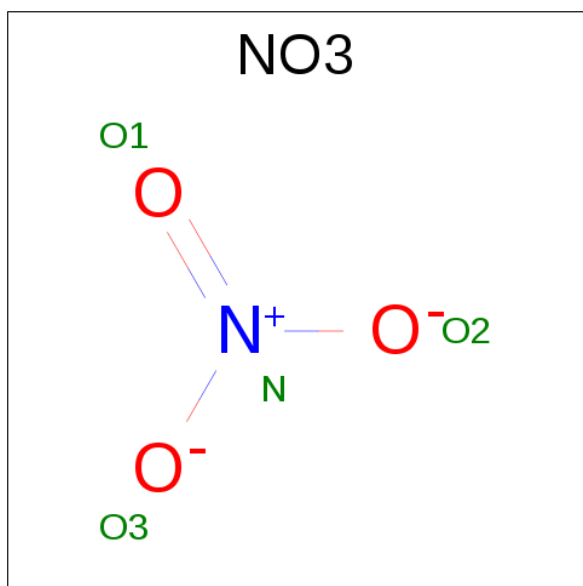
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	B	1	Total	K	0	0
			1	1		
9	A	1	Total	K	0	0
			1	1		

- Molecule 10 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	B	1	Total	C	N	O	
			8	4	1	3	
							0
							0

- Molecule 11 is NITRATE ION (three-letter code: NO3) (formula: NO₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	B	1	Total	N	O		
			4	1	3		
						0	
							0

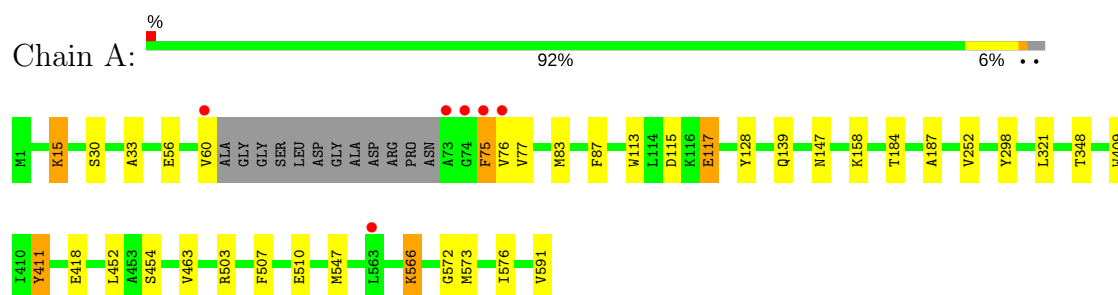
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	700	Total 717	O 717	0	17
12	B	689	Total 714	O 714	0	25

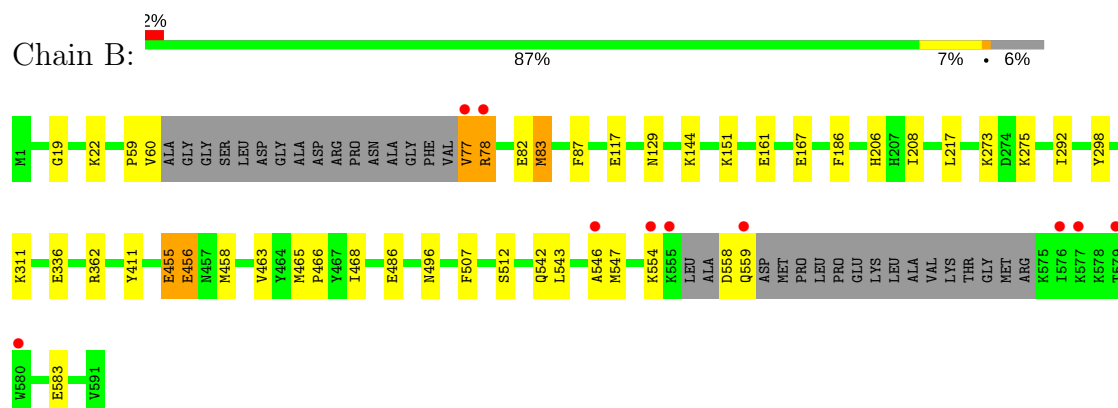
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Myosin-crossreactive antigen



• Molecule 1: Myosin-crossreactive antigen



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	165.10Å 77.74Å 104.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.73 – 1.80 48.72 – 1.78	Depositor EDS
% Data completeness (in resolution range)	97.6 (48.73-1.80) 96.2 (48.72-1.78)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 1.78Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.154 , 0.192 0.154 , 0.192	Depositor DCC
R_{free} test set	6030 reflections (4.93%)	DCC
Wilson B-factor (Å ²)	16.4	Xtriage
Anisotropy	0.320	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 58.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10974	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.87% 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, MPD, NA, K, EIC, MES, MRD, TRS, PO4, NO3

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	0.59	0/4872	0.65	0/6599
1	B	0.60	0/4698	0.66	0/6365
All	All	0.59	0/9570	0.66	0/12964

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

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	4750	0	4593	30	0
1	B	4581	0	4402	31	0
2	A	8	0	14	6	0
3	A	24	0	42	2	0
3	B	16	0	28	2	0
4	A	48	0	64	4	0
4	B	24	0	32	3	0
5	A	20	0	32	1	0
6	A	2	0	0	0	0
7	A	10	0	0	0	0
7	B	10	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	12	0	12	0	0
8	B	24	0	24	0	0
9	A	1	0	0	0	0
9	B	1	0	0	0	0
10	B	8	0	12	0	0
11	B	4	0	0	0	0
12	A	717	0	0	8	0
12	B	714	0	0	4	2
All	All	10974	0	9255	64	2

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 (64) 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:411:TYR:HB2	2:A:601[A]:MRD:H5C2	1.59	0.85
1:A:15:LYS:NZ	1:B:117:GLU:OE2	2.15	0.80
1:A:184:THR:O	2:A:601[A]:MRD:HMC1	1.86	0.76
1:B:208[A]:ILE:HG21	1:B:546:ALA:HA	1.67	0.75
1:B:82:GLU:OE2	3:B:601:MPD:HM3	1.88	0.72
1:B:554:LYS:O	1:B:559:GLN:NE2	2.23	0.70
1:A:75:PHE:HE2	1:A:77:VAL:HG22	1.58	0.69
1:A:147[B]:ASN:ND2	12:A:1348:HOH:O	2.26	0.69
2:A:601[A]:MRD:H5C3	2:A:601[A]:MRD:HMC2	1.76	0.68
7:B:612:PO4:O2	12:B:1085:HOH:O	2.14	0.64
1:A:115:ASP:OD1	12:A:1014:HOH:O	2.15	0.63
1:B:456:GLU:H	1:B:456:GLU:CD	2.03	0.62
2:A:601[A]:MRD:HMC3	12:A:1023:HOH:O	1.99	0.61
1:B:554:LYS:HB2	1:B:559:GLN:HG2	1.83	0.60
1:B:82:GLU:HG3	1:B:217:LEU:HD22	1.83	0.60
1:A:75:PHE:CE2	1:A:77:VAL:HG22	2.36	0.59
1:B:273:LYS:HD2	1:B:275:LYS:HE2	1.84	0.58
1:B:186:PHE:HE1	3:B:601:MPD:HM2	1.67	0.58
1:A:76:VAL:HG22	1:A:348:THR:HG23	1.86	0.58
1:A:33:ALA:H	4:A:607[A]:GOL:C1	2.18	0.57
1:A:418:GLU:OE1	12:A:1187:HOH:O	2.18	0.55
1:B:167:GLU:OE2	4:B:605:GOL:H2	2.06	0.55
1:B:311:LYS:NZ	1:B:336:GLU:OE2	2.40	0.55
1:A:113:TRP:O	1:A:117[A]:GLU:HG2	2.08	0.53
1:A:158:LYS:HD3	1:B:583:GLU:HG2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:59:PRO:O	1:B:60:VAL:HB	2.08	0.51
1:A:30:SER:OG	1:A:56:GLU:OE1	2.21	0.50
1:A:187:ALA:HB2	3:A:602[B]:MPD:H32	1.94	0.50
1:A:187:ALA:HB2	2:A:601[A]:MRD:H3C1	1.95	0.49
1:B:455:GLU:HG2	12:B:1064:HOH:O	2.13	0.48
1:B:496:ASN:HD22	4:B:609:GOL:H12	1.78	0.48
12:A:942:HOH:O	1:B:117:GLU:HG3	2.14	0.48
1:B:77:VAL:HG22	1:B:78:ARG:N	2.29	0.48
1:A:298:TYR:CD1	1:A:463:VAL:HG22	2.49	0.47
1:B:83[B]:MET:HG2	1:B:87:PHE:CD2	2.49	0.47
1:B:206:HIS:HA	1:B:542:GLN:OE1	2.15	0.46
1:B:82:GLU:H	1:B:82:GLU:CD	2.19	0.46
1:A:573:MET:CE	1:A:591:VAL:HG11	2.46	0.46
1:A:411:TYR:HB2	3:A:602[B]:MPD:H11	1.97	0.46
1:A:76:VAL:HB	1:A:463:VAL:HG21	1.98	0.46
1:B:78:ARG:HB2	12:B:1387:HOH:O	2.16	0.46
1:B:455:GLU:HG2	1:B:455:GLU:H	1.56	0.45
1:B:298:TYR:CD1	1:B:463:VAL:HG22	2.52	0.45
1:A:87:PHE:CG	1:A:510:GLU:HB2	2.51	0.45
1:B:512:SER:OG	4:B:609:GOL:H12	2.16	0.45
1:A:573:MET:HE1	1:A:591:VAL:HG11	1.99	0.44
1:B:292:ILE:HD12	1:B:468:ILE:HD12	2.00	0.43
1:A:454[A]:SER:HB2	4:A:612:GOL:H11	1.99	0.43
1:B:19:GLY:O	1:B:22:LYS:HG2	2.19	0.43
1:A:503:ARG:HE	4:A:606[A]:GOL:H2	1.83	0.43
1:A:139:GLN:NE2	12:A:1255[B]:HOH:O	2.52	0.43
1:B:455:GLU:HA	1:B:458:MET:O	2.19	0.42
1:A:566:LYS:HG3	12:A:1248[A]:HOH:O	2.19	0.42
2:A:601[A]:MRD:H5C1	12:A:732:HOH:O	2.20	0.42
1:B:129:ASN:ND2	12:B:1212:HOH:O	2.43	0.42
1:A:452:LEU:HD23	1:A:452:LEU:HA	1.89	0.42
1:A:572:GLY:O	1:A:576:ILE:HG13	2.20	0.42
1:B:83[B]:MET:HG2	1:B:87:PHE:CE2	2.55	0.41
1:B:543:LEU:O	1:B:547:MET:HG3	2.20	0.41
1:A:33:ALA:H	4:A:607[A]:GOL:H11	1.85	0.41
1:A:76:VAL:HG13	1:A:409:TRP:CZ3	2.56	0.40
5:A:608:EIC:H131	5:A:608:EIC:H161	1.63	0.40
1:A:252:VAL:HB	1:A:321:LEU:HD21	2.04	0.40
1:B:465:MET:HA	1:B:466:PRO:HD2	1.99	0.40

All (2) 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)
12:B:1325:HOH:O	12:B:1325:HOH:O[2_465]	2.05	0.15
12:B:1267:HOH:O	12:B:1267:HOH:O[2_465]	2.13	0.07

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	583/591 (99%)	571 (98%)	12 (2%)	0	100	100
1	B	556/591 (94%)	547 (98%)	9 (2%)	0	100	100
All	All	1139/1182 (96%)	1118 (98%)	21 (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	519/518 (100%)	507 (98%)	12 (2%)	56	41
1	B	499/518 (96%)	484 (97%)	15 (3%)	46	30
All	All	1018/1036 (98%)	991 (97%)	27 (3%)	56	35

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	LYS
1	A	60	VAL

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Mol	Chain	Res	Type
1	A	75	PHE
1	A	83	MET
1	A	117[A]	GLU
1	A	117[B]	GLU
1	A	128	TYR
1	A	411	TYR
1	A	507	PHE
1	A	547[A]	MET
1	A	547[B]	MET
1	A	566	LYS
1	B	77	VAL
1	B	78	ARG
1	B	83[A]	MET
1	B	83[B]	MET
1	B	144	LYS
1	B	151	LYS
1	B	161	GLU
1	B	362	ARG
1	B	411[A]	TYR
1	B	411[B]	TYR
1	B	455	GLU
1	B	456	GLU
1	B	486	GLU
1	B	507	PHE
1	B	558	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 32 ligands modelled in this entry, 4 are monoatomic - leaving 28 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$
2	MRD	A	601[A]	-	7,7,7	0.49	0	9,10,10	0.66	0
3	MPD	A	602[B]	-	7,7,7	0.26	0	9,10,10	0.17	0
3	MPD	A	603	-	7,7,7	0.21	0	9,10,10	0.79	0
4	GOL	A	604	-	5,5,5	0.41	0	5,5,5	0.34	0
4	GOL	A	605	-	5,5,5	0.34	0	5,5,5	0.46	0
4	GOL	A	606[A]	-	5,5,5	0.30	0	5,5,5	0.41	0
4	GOL	A	606[B]	-	5,5,5	0.46	0	5,5,5	0.64	0
4	GOL	A	607[A]	-	5,5,5	0.43	0	5,5,5	0.76	0
4	GOL	A	607[B]	-	5,5,5	0.27	0	5,5,5	0.51	0
5	EIC	A	608	-	16,19,19	1.17	2 (12%)	15,19,19	0.95	0
7	PO4	A	610	-	4,4,4	0.70	0	6,6,6	0.46	0
3	MPD	A	611	-	7,7,7	0.34	0	9,10,10	0.38	0
4	GOL	A	612	-	5,5,5	0.41	0	5,5,5	0.43	0
4	GOL	A	613	-	5,5,5	0.53	0	5,5,5	0.98	0
8	MES	A	615	-	12,12,12	1.98	1 (8%)	14,16,16	2.54	6 (42%)
7	PO4	A	616	-	4,4,4	0.78	0	6,6,6	0.34	0
3	MPD	B	601	-	7,7,7	0.29	0	9,10,10	0.70	0
8	MES	B	602	-	12,12,12	1.72	1 (8%)	14,16,16	2.47	3 (21%)
8	MES	B	603	-	12,12,12	1.95	1 (8%)	14,16,16	2.52	7 (50%)
4	GOL	B	604	-	5,5,5	0.33	0	5,5,5	0.42	0
4	GOL	B	605	-	5,5,5	0.29	0	5,5,5	0.47	0
4	GOL	B	606	-	5,5,5	0.32	0	5,5,5	0.22	0
10	TRS	B	607	-	7,7,7	0.37	0	9,9,9	0.52	0
7	PO4	B	608	-	4,4,4	0.75	0	6,6,6	0.46	0
4	GOL	B	609	-	5,5,5	0.52	0	5,5,5	0.76	0
11	NO3	B	610	-	1,3,3	4.58	1 (100%)	0,3,3	0.00	-
3	MPD	B	611[B]	-	7,7,7	0.22	0	9,10,10	0.47	0
7	PO4	B	612	-	4,4,4	0.94	0	6,6,6	0.67	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	MRD	A	601[A]	-	-	0/5/5/5	0/0/0/0
3	MPD	A	602[B]	-	-	0/5/5/5	0/0/0/0
3	MPD	A	603	-	-	0/5/5/5	0/0/0/0
4	GOL	A	604	-	-	0/4/4/4	0/0/0/0
4	GOL	A	605	-	-	0/4/4/4	0/0/0/0
4	GOL	A	606[A]	-	-	0/4/4/4	0/0/0/0
4	GOL	A	606[B]	-	-	0/4/4/4	0/0/0/0
4	GOL	A	607[A]	-	-	0/4/4/4	0/0/0/0
4	GOL	A	607[B]	-	-	0/4/4/4	0/0/0/0
5	EIC	A	608	-	-	0/15/17/17	0/0/0/0
7	PO4	A	610	-	-	0/0/0/0	0/0/0/0
3	MPD	A	611	-	-	0/5/5/5	0/0/0/0
4	GOL	A	612	-	-	0/4/4/4	0/0/0/0
4	GOL	A	613	-	-	0/4/4/4	0/0/0/0
8	MES	A	615	-	-	0/6/14/14	0/1/1/1
7	PO4	A	616	-	-	0/0/0/0	0/0/0/0
3	MPD	B	601	-	-	0/5/5/5	0/0/0/0
8	MES	B	602	-	-	0/6/14/14	0/1/1/1
8	MES	B	603	-	-	0/6/14/14	0/1/1/1
4	GOL	B	604	-	-	0/4/4/4	0/0/0/0
4	GOL	B	605	-	-	0/4/4/4	0/0/0/0
4	GOL	B	606	-	-	0/4/4/4	0/0/0/0
10	TRS	B	607	-	-	0/9/9/9	0/0/0/0
7	PO4	B	608	-	-	0/0/0/0	0/0/0/0
4	GOL	B	609	-	-	0/4/4/4	0/0/0/0
11	NO3	B	610	-	-	0/0/0/0	0/0/0/0
3	MPD	B	611[B]	-	-	0/5/5/5	0/0/0/0
7	PO4	B	612	-	-	0/0/0/0	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	603	MES	C8-S	-6.47	1.67	1.77
8	A	615	MES	C8-S	-6.47	1.67	1.77
8	B	602	MES	C8-S	-5.75	1.68	1.77
5	A	608	EIC	C13-C12	3.06	1.48	1.31
5	A	608	EIC	C10-C9	3.18	1.49	1.31
11	B	610	NO3	O1-N	4.58	1.41	1.23

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	615	MES	C2-C3-N4	-3.25	105.55	110.11
8	B	603	MES	C6-C5-N4	-2.85	106.12	110.11
8	B	603	MES	C7-N4-C5	2.31	117.17	111.26
8	B	603	MES	O3S-S-C8	2.62	109.27	106.06
8	B	603	MES	C7-N4-C3	2.69	118.15	111.26
8	A	615	MES	O2S-S-C8	2.85	109.24	106.79
8	B	603	MES	C5-N4-C3	3.11	115.91	108.87
8	A	615	MES	C7-N4-C5	3.23	119.55	111.26
8	B	603	MES	O2S-S-C8	3.39	109.70	106.79
8	A	615	MES	C7-N4-C3	3.42	120.02	111.26
8	B	602	MES	C7-N4-C5	3.91	121.28	111.26
8	A	615	MES	O1S-S-C8	4.21	110.41	106.79
8	B	602	MES	C5-N4-C3	4.46	118.97	108.87
8	A	615	MES	C5-N4-C3	5.01	120.22	108.87
8	B	603	MES	O1S-S-C8	5.37	111.40	106.79
8	B	602	MES	O1S-S-C8	6.14	112.06	106.79

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601[A]	MRD	6	0
3	A	602[B]	MPD	2	0
4	A	606[A]	GOL	1	0
4	A	607[A]	GOL	2	0
5	A	608	EIC	1	0
4	A	612	GOL	1	0
3	B	601	MPD	2	0
4	B	605	GOL	1	0
4	B	609	GOL	2	0
7	B	612	PO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	579/591 (97%)	-0.50	6 (1%) 82 80	8, 16, 35, 87	0
1	B	558/591 (94%)	-0.45	10 (1%) 69 65	8, 15, 40, 78	2 (0%)
All	All	1137/1182 (96%)	-0.48	16 (1%) 75 72	8, 16, 37, 87	2 (0%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	74	GLY	8.1
1	A	75	PHE	7.4
1	A	60	VAL	7.2
1	A	563	LEU	3.5
1	B	576	ILE	3.4
1	B	546	ALA	3.3
1	B	555	LYS	3.2
1	A	76	VAL	3.2
1	B	559	GLN	2.8
1	B	579	THR	2.7
1	B	77	VAL	2.6
1	B	78	ARG	2.4
1	B	554	LYS	2.4
1	B	577	LYS	2.1
1	A	73	ALA	2.0
1	B	580	TRP	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 ⓘ

There are no carbohydrates 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. 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(Å ²)	Q<0.9
4	GOL	A	607[B]	6/6	0.45	0.78	79.84	38,42,43,43	6
4	GOL	A	607[A]	6/6	0.45	0.78	60.76	28,30,36,37	6
7	PO4	A	610	5/5	0.80	0.20	16.18	68,70,71,72	5
3	MPD	A	603	8/8	0.93	0.16	8.20	9,19,21,24	8
8	MES	B	603	12/12	0.85	0.19	6.96	38,42,53,55	12
4	GOL	A	606[A]	6/6	0.73	0.27	6.90	5,21,23,24	6
4	GOL	A	606[B]	6/6	0.73	0.27	6.74	3,21,23,25	6
2	MRD	A	601[A]	8/8	0.81	0.18	6.31	8,17,18,22	8
7	PO4	B	608	5/5	0.77	0.18	6.25	43,49,50,50	5
3	MPD	A	602[B]	8/8	0.85	0.17	5.64	10,16,17,20	8
4	GOL	A	605	6/6	0.82	0.16	5.57	41,44,46,49	0
11	NO3	B	610	4/4	0.83	0.14	4.99	45,47,48,49	0
4	GOL	B	609	6/6	0.88	0.16	4.56	21,25,31,33	0
8	MES	A	615	12/12	0.82	0.21	4.43	32,35,42,43	12
7	PO4	A	616	5/5	0.90	0.21	4.34	49,51,53,54	5
4	GOL	B	605	6/6	0.63	0.21	3.90	45,48,49,51	0
3	MPD	B	601	8/8	0.90	0.19	3.60	6,15,17,23	8
3	MPD	B	611[B]	8/8	0.91	0.18	3.44	3,11,12,13	8
8	MES	B	602	12/12	0.97	0.10	3.36	9,16,24,26	12
5	EIC	A	608	20/20	0.71	0.28	3.03	35,54,75,77	0
4	GOL	A	604	6/6	0.74	0.25	2.91	45,45,47,51	0
3	MPD	A	611	8/8	0.91	0.14	2.69	18,22,23,24	8
4	GOL	B	604	6/6	0.69	0.18	2.66	42,47,50,54	6
10	TRS	B	607	8/8	0.90	0.23	1.57	44,46,50,50	0
4	GOL	A	612	6/6	0.91	0.14	1.28	20,22,27,28	6
4	GOL	B	606	6/6	0.77	0.19	1.26	47,48,48,48	6
6	NA	A	614	1/1	0.96	0.11	1.12	28,28,28,28	0
7	PO4	B	612	5/5	0.97	0.09	1.00	14,14,26,28	5
9	K	B	613	1/1	1.00	0.07	-0.13	11,11,11,11	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
6	NA	A	609	1/1	0.99	0.07	-0.64	22,22,22,22	0
9	K	A	617	1/1	0.99	0.05	-1.85	15,15,15,15	0
4	GOL	A	613	6/6	0.78	0.24	-	22,23,25,33	6

6.5 Other polymers [i](#)

There are no such residues in this entry.