



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:22 PM UTC

PDB ID : 8IC6 / pdb_00008ic6
Title : exo-beta-D-arabinanase ExoMA2 from Microbacterium arabinogalactanolyticum in complex with Tris
Authors : Fukushima, R.; Kashima, T.; Ishiwata, A.; Fujita, K.; Fushinobu, S.
Deposited on : 2023-02-10
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

<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

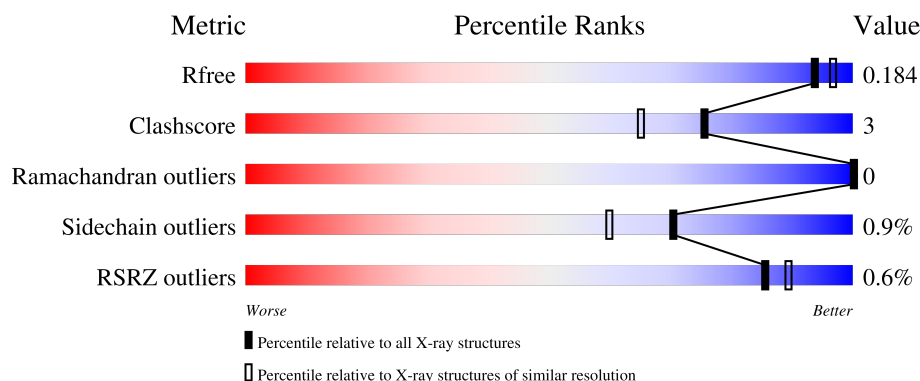
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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}	180053	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (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 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	867	
1	B	867	

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 15006 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.

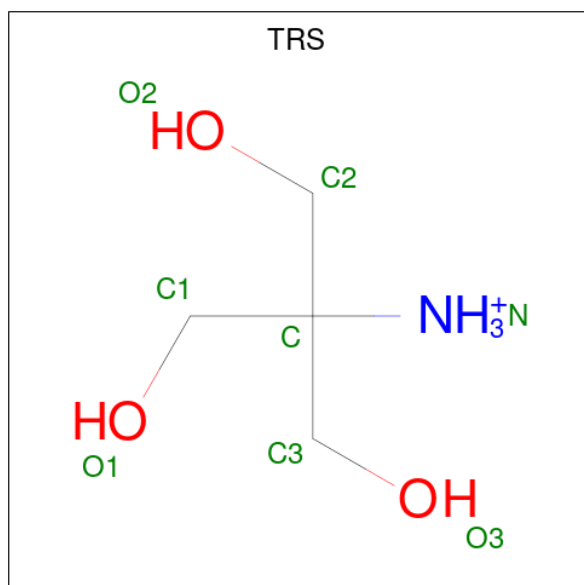
- Molecule 1 is a protein called exo-beta-D-arabinanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	855	Total	C	N	O	S	0	3	0
			6642	4179	1204	1248	11			
1	A	853	Total	C	N	O	S	0	3	0
			6632	4171	1205	1245	11			

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

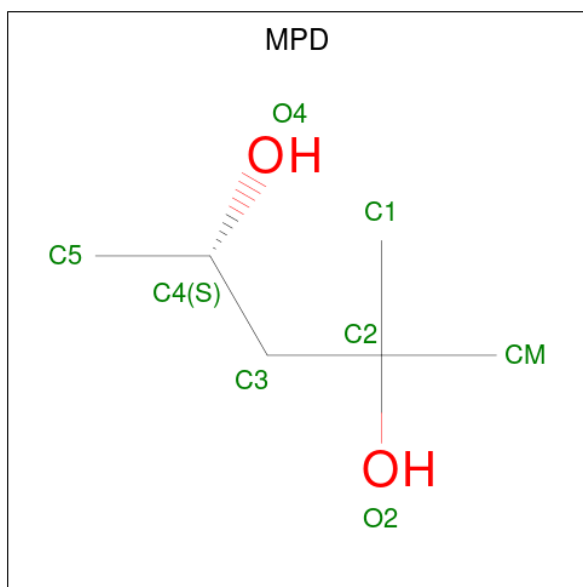
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	8	Total	Mg	0	0
			8	8		
2	A	7	Total	Mg	0	0
			7	7		

- Molecule 3 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: C₄H₁₂NO₃).



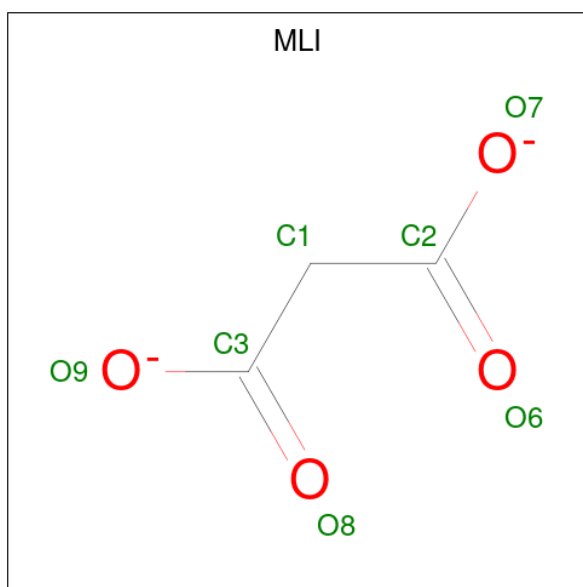
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			8	4	1	3		
3	A	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



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

- Molecule 5 is MALONATE ION (CCD ID: MLI) (formula: C₃H₂O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			7	3	4		

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Cl	0	0
			1	1		

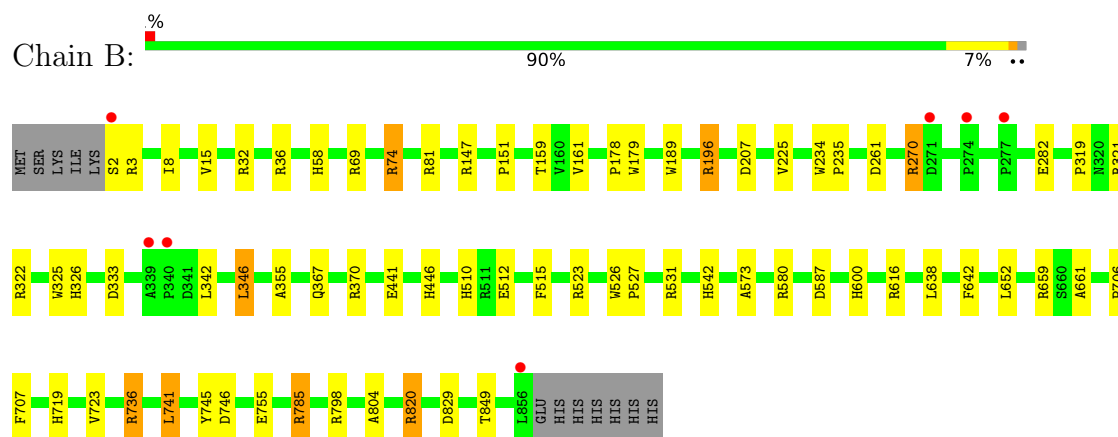
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	814	Total	O	0	0
			814	814		
7	A	847	Total	O	0	0
			847	847		

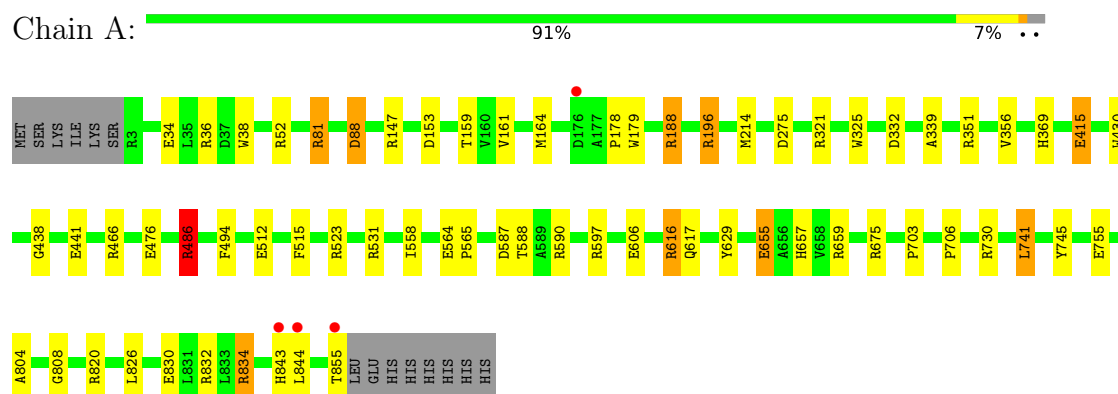
3 Residue-property plots

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: exo-beta-D-arabinanase



- Molecule 1: exo-beta-D-arabinanase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.55Å 97.51Å 139.98Å 90.00° 100.55° 90.00°	Depositor
Resolution (Å)	48.80 – 1.75 48.80 – 1.75	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.80-1.75) 100.0 (48.80-1.75)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.8.0403	Depositor
R, R_{free}	0.138 , 0.174 0.151 , 0.184	Depositor DCC
R_{free} test set	8854 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	18.9	Xtriage
Anisotropy	0.113	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	15006	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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: MLI, MG, TRS, CL, MPD

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	0.95	3/6823 (0.0%)	1.20	12/9301 (0.1%)
1	B	0.94	2/6833 (0.0%)	1.19	15/9316 (0.2%)
All	All	0.95	5/13656 (0.0%)	1.20	27/18617 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	13
1	B	0	13
All	All	0	26

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	415	GLU	CD-OE1	8.18	1.40	1.25
1	A	88	ASP	CG-OD1	7.01	1.38	1.25
1	A	88	ASP	CG-OD2	6.30	1.37	1.25
1	B	32	ARG	CD-NE	-5.26	1.38	1.46
1	B	58	HIS	CE1-NE2	5.18	1.37	1.32

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	188	ARG	NE-CZ-NH2	-9.71	110.47	119.20
1	A	855	THR	CA-CB-OG1	-7.50	98.36	109.60
1	A	675	ARG	CB-CG-CD	-7.29	94.53	111.30
1	A	588	THR	CA-CB-OG1	-7.23	98.75	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	32	ARG	NE-CZ-NH1	6.56	128.06	121.50
1	B	446	HIS	CA-CB-CG	6.42	120.22	113.80
1	B	32	ARG	NE-CZ-NH2	-6.42	113.42	119.20
1	B	326	HIS	CA-CB-CG	-6.05	107.75	113.80
1	B	587	ASP	CA-CB-CG	5.92	118.52	112.60
1	A	369	HIS	CA-CB-CG	5.89	119.69	113.80
1	A	486	ARG	CG-CD-NE	5.87	124.91	112.00
1	B	600	HIS	CA-CB-CG	-5.72	108.08	113.80
1	A	332	ASP	CA-CB-CG	5.58	118.18	112.60
1	A	597	ARG	NE-CZ-NH1	-5.47	116.03	121.50
1	B	207	ASP	CA-CB-CG	5.39	117.99	112.60
1	A	415	GLU	CB-CG-CD	5.39	121.76	112.60
1	B	346	LEU	CA-C-N	-5.38	113.43	121.87
1	B	346	LEU	C-N-CA	-5.38	113.43	121.87
1	A	34	GLU	CB-CG-CD	5.32	121.65	112.60
1	B	746	ASP	CA-CB-CG	5.32	117.92	112.60
1	B	282	GLU	CB-CA-C	5.30	119.87	110.85
1	B	333	ASP	CA-CB-CG	5.29	117.89	112.60
1	A	153	ASP	CA-CB-CG	5.19	117.79	112.60
1	B	74	ARG	NE-CZ-NH2	-5.12	114.60	119.20
1	B	642	PHE	CA-CB-CG	5.11	118.91	113.80
1	A	655	GLU	CB-CA-C	5.02	119.22	110.68
1	B	261	ASP	CB-CA-C	5.00	117.81	109.56

There are no chirality outliers.

All (26) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	188	ARG	Sidechain
1	A	196	ARG	Sidechain
1	A	351[A]	ARG	Sidechain
1	A	351[B]	ARG	Sidechain
1	A	36	ARG	Sidechain
1	A	466	ARG	Sidechain
1	A	486	ARG	Sidechain
1	A	523	ARG	Sidechain
1	A	616	ARG	Sidechain
1	A	81[A]	ARG	Sidechain
1	A	81[B]	ARG	Sidechain
1	A	834[A]	ARG	Sidechain
1	A	834[B]	ARG	Sidechain
1	B	147	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	B	196	ARG	Sidechain
1	B	270	ARG	Sidechain
1	B	36	ARG	Sidechain
1	B	531	ARG	Sidechain
1	B	659	ARG	Sidechain
1	B	69[A]	ARG	Sidechain
1	B	69[B]	ARG	Sidechain
1	B	736	ARG	Sidechain
1	B	74	ARG	Sidechain
1	B	785	ARG	Sidechain
1	B	81	ARG	Sidechain
1	B	820	ARG	Sidechain

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	6632	0	6308	42	0
1	B	6642	0	6320	35	0
2	A	7	0	0	0	0
2	B	8	0	0	0	0
3	A	8	0	12	0	0
3	B	8	0	12	0	0
4	A	16	0	28	2	0
4	B	16	0	28	4	0
5	B	7	0	2	0	0
6	A	1	0	0	0	0
7	A	847	0	0	5	0
7	B	814	0	0	10	0
All	All	15006	0	12710	78	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (78) 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)
4:B:910:MPD:HM2	7:B:1055:HOH:O	1.73	0.89
1:A:830:GLU:OE2	1:A:832:ARG:NH1	2.22	0.71
1:B:322:ARG:HG3	1:B:322:ARG:HH11	1.56	0.70
1:A:834[A]:ARG:NH1	7:A:1002:HOH:O	2.30	0.64
1:B:736:ARG:HD2	7:B:1012:HOH:O	1.99	0.62
1:B:321:ARG:NH1	4:B:912:MPD:H4	2.17	0.60
1:A:321:ARG:NH1	4:A:910:MPD:H4	2.17	0.59
1:A:655:GLU:HG2	1:A:659:ARG:NH2	2.19	0.57
1:B:745:TYR:CE2	1:B:755:GLU:HA	2.40	0.56
1:B:523:ARG:NH1	7:B:1010:HOH:O	2.37	0.56
1:A:834[A]:ARG:HD2	7:A:1686:HOH:O	2.09	0.53
1:B:542:HIS:HE1	7:B:1171:HOH:O	1.90	0.53
1:B:325:TRP:CD2	1:B:441:GLU:HB2	2.44	0.53
1:B:322:ARG:HH11	1:B:322:ARG:CG	2.19	0.53
1:B:736:ARG:NE	7:B:1012:HOH:O	2.39	0.51
1:B:515:PHE:O	1:B:804:ALA:HB2	2.10	0.51
1:B:322:ARG:CG	1:B:322:ARG:NH1	2.74	0.51
1:A:515:PHE:O	1:A:804:ALA:HB2	2.11	0.50
1:A:659:ARG:HG2	1:A:730:ARG:NH2	2.27	0.50
1:B:736:ARG:CD	7:B:1012:HOH:O	2.57	0.50
1:A:659:ARG:HD2	1:A:730:ARG:NH1	2.27	0.50
1:A:655:GLU:CG	1:A:659:ARG:NH2	2.75	0.50
1:B:189:TRP:CZ3	1:B:225:VAL:HG11	2.47	0.49
1:A:38:TRP:CH2	1:A:164:MET:HE1	2.47	0.49
1:B:3:ARG:HA	7:B:1476:HOH:O	2.12	0.49
1:B:319:PRO:HB3	1:B:355:ALA:HB1	1.95	0.49
4:B:910:MPD:CM	7:B:1055:HOH:O	2.45	0.49
1:B:319:PRO:HB3	1:B:355:ALA:CB	2.43	0.48
1:A:88:ASP:OD1	7:A:1001:HOH:O	2.20	0.48
1:A:476:GLU:H	1:A:476:GLU:CD	2.20	0.48
1:A:164:MET:HE2	1:A:214:MET:HE1	1.95	0.48
1:A:741:LEU:C	1:A:741:LEU:HD12	2.38	0.48
1:B:178:PRO:O	1:B:179:TRP:HB2	2.13	0.48
1:A:159:THR:HG22	1:A:161:VAL:HG23	1.96	0.48
1:B:741:LEU:HD12	1:B:741:LEU:C	2.39	0.48
1:A:325:TRP:CD2	1:A:441:GLU:HB2	2.49	0.47
1:B:510:HIS:NE2	1:B:580:ARG:NH1	2.63	0.47
1:B:785:ARG:NH2	1:B:829:ASP:OD1	2.40	0.47
1:A:159:THR:HG22	1:A:161:VAL:CG2	2.45	0.46
1:A:826:LEU:O	1:A:843:HIS:HA	2.16	0.46
1:A:655:GLU:CD	1:A:659:ARG:HH22	2.23	0.46
1:B:270:ARG:HA	1:A:494:PHE:CE2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:321:ARG:HH11	4:B:912:MPD:H4	1.79	0.46
1:A:703:PRO:HG2	1:A:706:PRO:HB3	1.98	0.46
1:B:706:PRO:O	1:B:707:PHE:C	2.59	0.46
1:A:745:TYR:CE2	1:A:755:GLU:HA	2.51	0.45
1:B:849:THR:HG22	7:B:1005:HOH:O	2.16	0.45
1:A:655:GLU:CG	1:A:659:ARG:HH22	2.29	0.45
1:A:820:ARG:HD3	7:A:1582:HOH:O	2.17	0.44
1:B:367:GLN:NE2	1:B:370:ARG:HH22	2.16	0.44
1:A:476:GLU:CD	1:A:476:GLU:N	2.76	0.44
1:B:159:THR:HG22	1:B:161[A]:VAL:HG23	2.00	0.43
1:A:164:MET:CA	1:A:164:MET:HE3	2.45	0.43
1:A:339:ALA:HB1	1:A:486:ARG:CZ	2.49	0.43
1:A:81[A]:ARG:NH1	7:A:1008:HOH:O	2.40	0.43
1:A:476:GLU:O	1:A:531:ARG:HD3	2.19	0.43
1:B:8:ILE:HG21	1:B:15:VAL:HG21	2.01	0.42
1:A:558:ILE:C	1:A:558:ILE:HD12	2.44	0.42
1:A:587:ASP:OD2	1:A:590:ARG:HD2	2.19	0.42
1:B:322:ARG:HD3	7:B:1346:HOH:O	2.19	0.42
1:B:798:ARG:CG	1:B:820:ARG:HH22	2.33	0.42
1:A:430:TRP:CE2	1:A:438:GLY:HA3	2.55	0.42
1:B:234:TRP:HB3	1:B:235:PRO:HD2	2.01	0.42
1:B:342:LEU:O	1:B:346:LEU:HG	2.20	0.42
1:A:52:ARG:NH1	1:A:88:ASP:CG	2.77	0.42
1:A:178:PRO:O	1:A:179:TRP:HB2	2.21	0.41
1:A:325:TRP:HA	4:A:910:MPD:H53	2.01	0.41
1:A:565:PRO:HD2	1:A:629:TYR:OH	2.21	0.41
1:A:808:GLY:HA3	1:A:820:ARG:O	2.20	0.41
1:B:719:HIS:O	1:B:723:VAL:HG23	2.21	0.41
1:A:606:GLU:OE1	1:A:657:HIS:ND1	2.46	0.41
1:A:147:ARG:HH11	1:A:147:ARG:HD2	1.71	0.41
1:A:356:VAL:CG2	1:A:415:GLU:HG2	2.51	0.41
1:B:526:TRP:HB3	1:B:527:PRO:HD3	2.03	0.41
1:B:573:ALA:HB2	1:B:652:LEU:HD11	2.03	0.40
1:B:638:LEU:HD11	1:B:661:ALA:HB3	2.04	0.40
1:A:164:MET:HE2	1:A:164:MET:HB3	1.87	0.40
1:A:564:GLU:HB3	1:A:617:GLN:HG3	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	854/867 (98%)	835 (98%)	19 (2%)	0	100	100
1	B	856/867 (99%)	832 (97%)	24 (3%)	0	100	100
All	All	1710/1734 (99%)	1667 (98%)	43 (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	662/673 (98%)	656 (99%)	6 (1%)	70	60
1	B	664/673 (99%)	658 (99%)	6 (1%)	70	60
All	All	1326/1346 (98%)	1314 (99%)	12 (1%)	70	60

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	2	SER
1	B	151	PRO
1	B	196	ARG
1	B	512	GLU
1	B	616	ARG
1	B	741	LEU
1	A	196	ARG
1	A	275	ASP

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Mol	Chain	Res	Type
1	A	512	GLU
1	A	616	ARG
1	A	741	LEU
1	A	844	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	625	HIS
1	A	26	ASN
1	A	625	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 23 ligands modelled in this entry, 16 are monoatomic - leaving 7 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	MPD	A	911	-	7,7,7	0.30	0	9,10,10	0.59	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	MLI	B	911	-	6,6,6	3.24	4 (66%)	7,7,7	1.08	0
4	MPD	A	910	-	7,7,7	0.76	0	9,10,10	1.52	2 (22%)
3	TRS	A	909	-	7,7,7	0.29	0	9,9,9	0.54	0
4	MPD	B	912	-	7,7,7	0.79	0	9,10,10	1.18	1 (11%)
4	MPD	B	910	-	7,7,7	0.24	0	9,10,10	0.41	0
3	TRS	B	909	-	7,7,7	0.43	0	9,9,9	0.73	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	MPD	A	911	-	-	0/5/5/5	-
5	MLI	B	911	-	-	0/4/4/4	-
4	MPD	A	910	-	-	4/5/5/5	-
3	TRS	A	909	-	-	5/9/9/9	-
4	MPD	B	912	-	-	4/5/5/5	-
4	MPD	B	910	-	-	1/5/5/5	-
3	TRS	B	909	-	-	3/9/9/9	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	911	MLI	O6-C2	5.41	1.39	1.22
5	B	911	MLI	O8-C3	4.33	1.36	1.22
5	B	911	MLI	C1-C3	3.05	1.55	1.51
5	B	911	MLI	C1-C2	2.31	1.54	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	910	MPD	O2-C2-CM	-3.14	98.21	107.99
4	A	910	MPD	O4-C4-C3	2.43	121.01	111.35
4	B	912	MPD	O4-C4-C3	2.24	120.26	111.35

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	909	TRS	N-C-C1-O1
3	A	909	TRS	N-C-C3-O3
4	B	912	MPD	C2-C3-C4-O4
3	A	909	TRS	N-C-C2-O2
4	B	910	MPD	O2-C2-C3-C4
4	A	910	MPD	C2-C3-C4-O4
3	B	909	TRS	C3-C-C1-O1
3	A	909	TRS	C1-C-C2-O2
3	A	909	TRS	C3-C-C2-O2
3	A	909	TRS	C2-C-C3-O3
4	B	912	MPD	O2-C2-C3-C4
4	A	910	MPD	O2-C2-C3-C4
4	B	912	MPD	C1-C2-C3-C4
4	B	912	MPD	CM-C2-C3-C4
4	A	910	MPD	C1-C2-C3-C4
4	A	910	MPD	CM-C2-C3-C4
3	B	909	TRS	C2-C-C1-O1

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	910	MPD	2	0
4	B	912	MPD	2	0
4	B	910	MPD	2	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	853/867 (98%)	-0.45	4 (0%) 87 90	10, 19, 35, 70	3 (0%)
1	B	855/867 (98%)	-0.43	7 (0%) 82 86	11, 20, 37, 79	3 (0%)
All	All	1708/1734 (98%)	-0.44	11 (0%) 85 89	10, 19, 36, 79	6 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	855	THR	3.9
1	B	339	ALA	3.3
1	B	856	LEU	3.2
1	B	274	PRO	2.6
1	A	844	LEU	2.4
1	A	176	ASP	2.3
1	B	271	ASP	2.3
1	B	340	PRO	2.2
1	A	843	HIS	2.2
1	B	2	SER	2.2
1	B	277	PRO	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MPD	A	910	8/8	0.88	0.12	29,31,36,39	0
4	MPD	B	912	8/8	0.89	0.12	26,29,33,35	0
5	MLI	B	911	7/7	0.89	0.10	30,33,36,37	0
4	MPD	B	910	8/8	0.93	0.10	27,34,40,42	0
2	MG	A	907	1/1	0.96	0.15	35,35,35,35	0
4	MPD	A	911	8/8	0.96	0.08	25,31,33,34	0
3	TRS	A	909	8/8	0.96	0.06	16,19,23,24	0
2	MG	B	907	1/1	0.97	0.09	38,38,38,38	0
2	MG	B	908	1/1	0.97	0.11	40,40,40,40	0
2	MG	A	903	1/1	0.97	0.17	32,32,32,32	0
2	MG	A	905	1/1	0.97	0.07	31,31,31,31	0
2	MG	B	906	1/1	0.97	0.06	33,33,33,33	0
3	TRS	B	909	8/8	0.97	0.05	16,20,22,22	0
2	MG	B	902	1/1	0.98	0.03	17,17,17,17	0
2	MG	B	905	1/1	0.98	0.06	33,33,33,33	0
2	MG	A	904	1/1	0.98	0.04	22,22,22,22	0
2	MG	B	903	1/1	0.99	0.05	22,22,22,22	0
2	MG	B	904	1/1	0.99	0.04	18,18,18,18	0
2	MG	B	901	1/1	0.99	0.03	23,23,23,23	0
2	MG	A	906	1/1	0.99	0.03	20,20,20,20	0
2	MG	A	901	1/1	0.99	0.02	19,19,19,19	0
2	MG	A	902	1/1	0.99	0.05	20,20,20,20	0
6	CL	A	908	1/1	0.99	0.04	22,22,22,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.