



Full wwPDB X-ray Structure Validation Report

Mar 31, 2014 – 01:52 PM BST

PDB ID : 4MRS
Title : Structure of a bacterial Atm1-family ABC transporter
Authors : Lee, J.Y.; Yang, J.G.; Zhitnitsky, D.; Lewinson, O.; Rees, D.C.
Deposited on : 2013-09-17
Resolution : 2.35 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

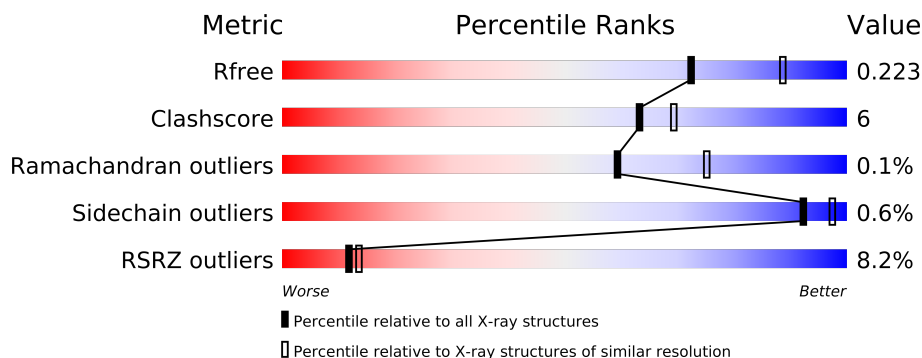
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable23004
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable23004

1 Overall quality at a glance

The reported resolution of this entry is 2.35 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}	66092	3327 (2.40-2.32)
Clashscore	79885	1064 (2.38-2.34)
Ramachandran outliers	78287	1048 (2.38-2.34)
Sidechain outliers	78261	1049 (2.38-2.34)
RSRZ outliers	66119	3330 (2.40-2.32)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	614	
1	B	614	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	LDA	A	701	-	X
2	LDA	A	706	-	X
2	LDA	A	707	-	X
2	LDA	A	708	-	X
2	LDA	A	709	-	X
2	LDA	B	701	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
2	LDA	B	702	-	X
2	LDA	B	706	-	X
2	LDA	B	707	-	X
2	LDA	B	708	-	X
2	LDA	B	709	-	X
3	GDS	A	703	-	X
3	GDS	B	703	-	X
4	PO4	B	704	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9931 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

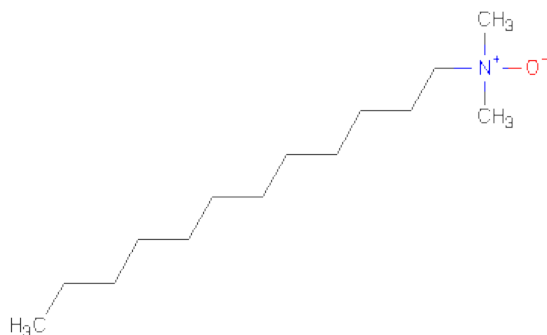
- Molecule 1 is a protein called ABC transporter related protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	600	Total	C	N	O	Se	0	0	0
			4658	2970	828	848	12			
1	B	598	Total	C	N	O	Se	0	0	0
			4645	2963	825	845	12			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	609	HIS	-	EXPRESSION TAG	UNP Q2G506
A	610	HIS	-	EXPRESSION TAG	UNP Q2G506
A	611	HIS	-	EXPRESSION TAG	UNP Q2G506
A	612	HIS	-	EXPRESSION TAG	UNP Q2G506
A	613	HIS	-	EXPRESSION TAG	UNP Q2G506
A	614	HIS	-	EXPRESSION TAG	UNP Q2G506
B	609	HIS	-	EXPRESSION TAG	UNP Q2G506
B	610	HIS	-	EXPRESSION TAG	UNP Q2G506
B	611	HIS	-	EXPRESSION TAG	UNP Q2G506
B	612	HIS	-	EXPRESSION TAG	UNP Q2G506
B	613	HIS	-	EXPRESSION TAG	UNP Q2G506
B	614	HIS	-	EXPRESSION TAG	UNP Q2G506

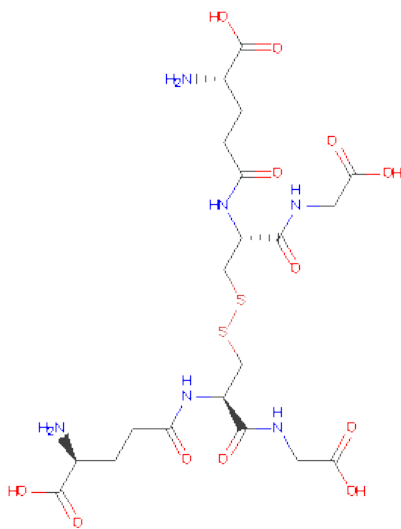
- Molecule 2 is LAURYL DIMETHYLAMINE-N-OXIDE (three-letter code: LDA) (formula: C₁₄H₃₁NO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			16	14	1	1		
2	A	1	Total	C	N	O	0	0
			16	14	1	1		
2	A	1	Total	C	N	O	0	0
			16	14	1	1		
2	A	1	Total	C			0	0
			6	6				
2	A	1	Total	C			0	0
			5	5				
2	A	1	Total	C			0	0
			6	6				
2	A	1	Total	C			0	0
			4	4				
2	B	1	Total	C	N	O	0	0
			16	14	1	1		
2	B	1	Total	C	N	O	0	0
			16	14	1	1		
2	B	1	Total	C			0	0
			7	7				
2	B	1	Total	C			0	0
			6	6				
2	B	1	Total	C	N	O	0	0
			11	9	1	1		
2	B	1	Total	C			0	0
			5	5				

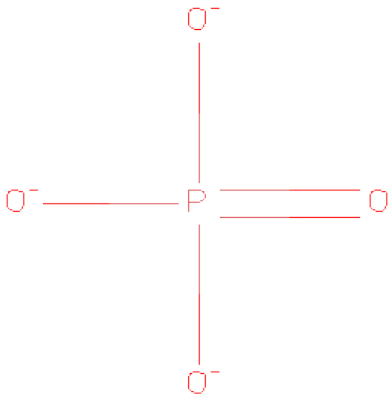
- Molecule 3 is OXIDIZED GLUTATHIONE DISULFIDE (three-letter code: GDS) (formula:

C₂₀H₃₂N₆O₁₂S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			40	20	6	12	2		
3	B	1	Total	C	N	O	S	0	0
			40	20	6	12	2		

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



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

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

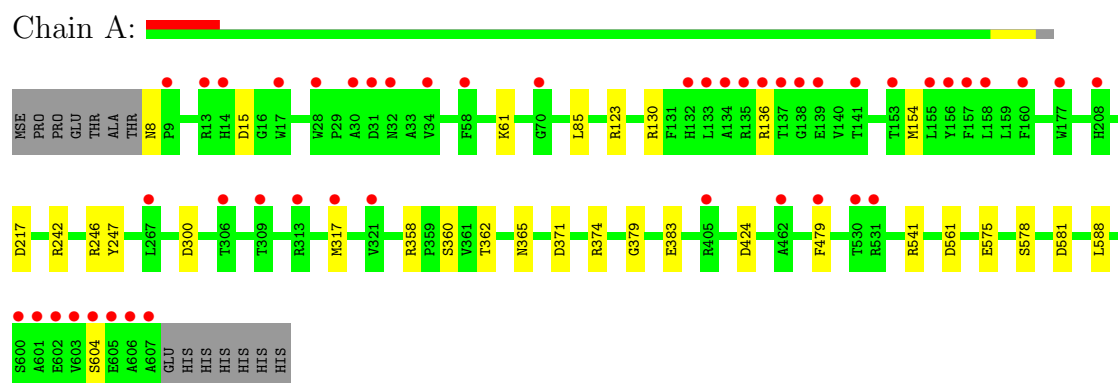
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	197	Total	O	0	0
			197	197		
5	B	201	Total	O	0	0
			201	201		

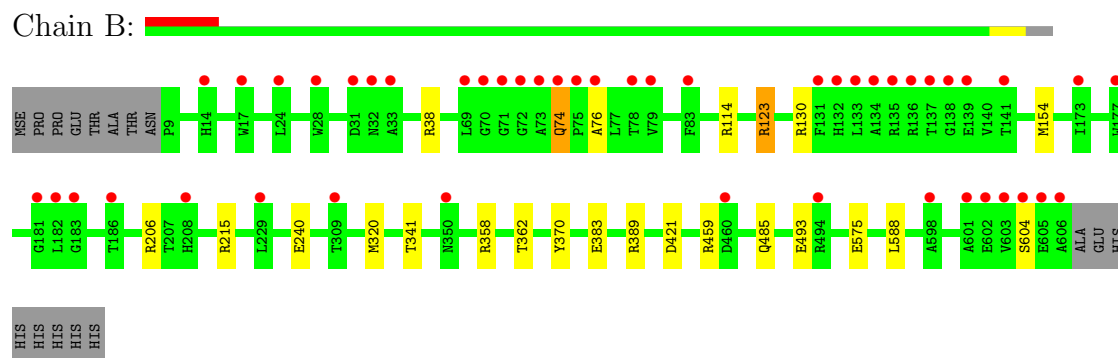
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: ABC transporter related protein



- Molecule 1: ABC transporter related protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	322.17Å 95.40Å 79.40Å 90.00° 101.81° 90.00°	Depositor
Resolution (Å)	39.66 – 2.35 39.66 – 2.34	Depositor EDS
% Data completeness (in resolution range)	96.1 (39.66-2.35) 83.7 (39.66-2.34)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.22 (at 2.34Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.193 , 0.223 0.194 , 0.223	Depositor DCC
R_{free} test set	4133 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	48.9	Xtriage
Anisotropy	0.357	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 42.1	EDS
Estimated twinning fraction	0.024 for -h-2*1,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 95159 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9931	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.26% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, GDS, LDA

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	0.45	0/4732	0.60	0/6408
1	B	0.45	0/4719	0.63	0/6389
All	All	0.45	0/9451	0.62	0/12797

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4658	0	0	24	0
1	B	4645	0	0	21	0
2	A	69	0	119	3	0
2	B	61	0	104	2	0
3	A	40	0	28	9	0
3	B	40	0	28	2	0
4	A	10	0	0	0	0
4	B	10	0	0	2	0
5	A	197	0	0	13	0
5	B	201	0	0	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	9931	0	279	54	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (54) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:374:ARG:NH1	5:A:957:HOH:O	1.89	1.05
1:A:365:ASN:ND2	5:A:903:HOH:O	2.00	0.93
1:B:130:ARG:NH2	1:B:370:TYR:O	2.07	0.87
3:A:703:GDS:HA31	3:A:703:GDS:HB21	1.63	0.79
1:B:358:ARG:NH2	5:B:920:HOH:O	2.16	0.78
1:B:485:GLN:NE2	5:B:890:HOH:O	2.17	0.78
1:A:15:ASP:OD2	5:A:949:HOH:O	2.02	0.77
1:A:358:ARG:NH1	5:A:991:HOH:O	2.17	0.77
1:A:379:GLY:O	5:A:884:HOH:O	2.03	0.77
1:A:581:ASP:OD2	5:A:980:HOH:O	2.05	0.73
4:B:705:PO4:O1	5:B:925:HOH:O	2.07	0.71
3:A:703:GDS:HA42	3:A:703:GDS:HB51	1.73	0.70
1:A:362:THR:OG1	1:A:383:GLU:OE2	2.12	0.68
1:A:604:SER:OG	1:B:604:SER:OG	2.12	0.67
1:B:362:THR:OG1	1:B:383:GLU:OE2	2.12	0.67
1:A:136:ARG:NH2	1:B:493:GLU:OE1	2.29	0.65
1:A:578:SER:N	5:A:980:HOH:O	2.30	0.64
1:B:114:ARG:NH1	4:B:705:PO4:O4	2.34	0.61
1:B:215:ARG:NE	5:B:854:HOH:O	2.32	0.61
3:A:703:GDS:CB2	3:A:703:GDS:HA31	2.31	0.59
1:B:74:GLN:O	1:B:76:ALA:N	2.36	0.58
1:A:317:MSE:SE	3:B:703:GDS:HA41	2.54	0.58
1:B:421:ASP:O	5:B:814:HOH:O	2.17	0.58
1:A:8:ASN:N	5:A:935:HOH:O	2.35	0.58
1:A:479:PHE:N	5:A:929:HOH:O	2.37	0.57
1:B:421:ASP:OD2	5:B:980:HOH:O	2.17	0.56
1:A:61:LYS:NZ	1:A:300:ASP:OD1	2.40	0.54
1:A:360:SER:OG	5:A:898:HOH:O	2.18	0.54
3:A:703:GDS:OE1	1:B:206:ARG:NH2	2.41	0.54
1:B:38:ARG:NH1	5:B:902:HOH:O	2.41	0.53
1:B:459:ARG:NH1	5:B:977:HOH:O	2.41	0.53
1:A:247:TYR:OH	2:A:702:LDA:H22	2.09	0.53
1:B:114:ARG:NE	5:B:880:HOH:O	2.42	0.52
1:A:130:ARG:NH2	1:A:371:ASP:OD1	2.43	0.51
1:A:424:ASP:OD2	5:A:808:HOH:O	2.20	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:389:ARG:NH1	5:B:860:HOH:O	2.45	0.48
3:A:703:GDS:HB22	1:B:320:MSE:CE	2.44	0.48
1:A:246:ARG:NH1	5:A:810:HOH:O	2.46	0.48
3:A:703:GDS:HB12	3:A:703:GDS:OE3	2.14	0.47
3:A:703:GDS:CA1	3:A:703:GDS:HN2	2.27	0.47
2:B:706:LDA:H92	2:B:707:LDA:H41	1.97	0.47
3:A:703:GDS:HN2	3:A:703:GDS:HA1	1.79	0.46
1:B:240:GLU:OE1	1:B:240:GLU:N	2.48	0.46
1:A:246:ARG:NH1	5:A:908:HOH:O	2.48	0.46
1:B:459:ARG:NH2	5:B:977:HOH:O	2.48	0.45
1:A:300:ASP:OD2	2:A:706:LDA:HM11	2.17	0.44
1:B:123:ARG:NH1	1:B:341:THR:O	2.51	0.43
2:B:708:LDA:H42	2:B:709:LDA:H42	2.01	0.42
1:B:575:GLU:OE1	1:B:588:LEU:N	2.53	0.41
1:A:575:GLU:OE2	1:A:588:LEU:N	2.54	0.41
1:A:541:ARG:NE	1:A:561:ASP:OD2	2.54	0.41
3:A:703:GDS:HA41	3:A:703:GDS:HB62	2.02	0.41
1:A:217:ASP:OD1	2:A:702:LDA:O1	2.39	0.41
3:B:703:GDS:HA5	3:B:703:GDS:HA42	1.79	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	598/614 (97%)	587 (98%)	11 (2%)	0	100	100
1	B	596/614 (97%)	588 (99%)	7 (1%)	1 (0%)	56	70
All	All	1194/1228 (97%)	1175 (98%)	18 (2%)	1 (0%)	59	75

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	74	GLN

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	478/478 (100%)	474 (99%)	4 (1%)	89	95
1	B	477/478 (100%)	475 (100%)	2 (0%)	95	98
All	All	955/956 (100%)	949 (99%)	6 (1%)	92	97

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	85	LEU
1	A	123	ARG
1	A	154	MSE
1	A	242	ARG
1	B	123	ARG
1	B	154	MSE

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 chains 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 ⓘ

19 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	LDA	A	701	-	15,15,15	3.85	2 (13%)	17,17,17	0.91	1 (5%)
2	LDA	A	702	-	15,15,15	3.86	1 (6%)	17,17,17	0.78	1 (5%)
3	GDS	A	703	-	39,39,39	2.21	6 (15%)	50,50,50	1.47	10 (20%)
4	PO4	A	704	-	4,4,4	0.25	0	6,6,6	0.31	0
4	PO4	A	705	-	4,4,4	0.31	0	6,6,6	0.31	0
2	LDA	A	706	-	15,15,15	3.83	2 (13%)	17,17,17	0.62	1 (5%)
2	LDA	A	707	-	5,5,15	3.03	2 (40%)	4,4,17	0.64	0
2	LDA	A	708	-	4,4,15	2.94	2 (50%)	3,3,17	0.74	0
2	LDA	A	709	-	5,5,15	2.82	2 (40%)	4,4,17	0.81	0
2	LDA	A	710	-	3,3,15	3.83	2 (66%)	2,2,17	1.22	0
2	LDA	B	701	-	15,15,15	3.83	2 (13%)	17,17,17	0.64	0
2	LDA	B	702	-	15,15,15	3.82	2 (13%)	17,17,17	0.74	0
3	GDS	B	703	-	39,39,39	2.04	6 (15%)	50,50,50	1.44	9 (18%)
4	PO4	B	704	-	4,4,4	0.27	0	6,6,6	0.31	0
4	PO4	B	705	-	4,4,4	0.27	0	6,6,6	0.31	0
2	LDA	B	706	-	6,6,15	2.71	2 (33%)	5,5,17	0.48	0
2	LDA	B	707	-	5,5,15	3.35	2 (40%)	4,4,17	0.83	0
2	LDA	B	708	-	10,10,15	4.74	3 (30%)	12,12,17	1.21	1 (8%)
2	LDA	B	709	-	4,4,15	3.70	2 (50%)	3,3,17	1.16	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	LDA	A	701	-	-	0/13/13/13	0/0/0/0
2	LDA	A	702	-	-	0/13/13/13	0/0/0/0
3	GDS	A	703	-	-	0/51/51/51	0/0/0/0
4	PO4	A	704	-	-	0/0/0/0	0/0/0/0
4	PO4	A	705	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	LDA	A	706	-	-	0/13/13/13	0/0/0/0
2	LDA	A	707	-	-	0/3/3/13	0/0/0/0
2	LDA	A	708	-	-	0/2/2/13	0/0/0/0
2	LDA	A	709	-	-	0/3/3/13	0/0/0/0
2	LDA	A	710	-	-	0/1/1/13	0/0/0/0
2	LDA	B	701	-	-	0/13/13/13	0/0/0/0
2	LDA	B	702	-	-	0/13/13/13	0/0/0/0
3	GDS	B	703	-	-	0/51/51/51	0/0/0/0
4	PO4	B	704	-	-	0/0/0/0	0/0/0/0
4	PO4	B	705	-	-	0/0/0/0	0/0/0/0
2	LDA	B	706	-	-	0/4/4/13	0/0/0/0
2	LDA	B	707	-	-	0/3/3/13	0/0/0/0
2	LDA	B	708	-	-	0/8/8/13	0/0/0/0
2	LDA	B	709	-	-	0/2/2/13	0/0/0/0

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	702	LDA	O1-N1	-14.66	1.25	1.39
2	A	701	LDA	O1-N1	-14.56	1.25	1.39
2	A	706	LDA	O1-N1	-14.51	1.25	1.39
2	B	702	LDA	O1-N1	-14.50	1.25	1.39
2	B	701	LDA	O1-N1	-14.50	1.25	1.39
2	B	708	LDA	O1-N1	-14.09	1.26	1.39
3	A	703	GDS	CD1-N2	6.59	1.47	1.34
3	A	703	GDS	C2-N3	6.11	1.46	1.33
3	A	703	GDS	C5-N4	6.05	1.46	1.33
3	B	703	GDS	CD1-N2	5.95	1.46	1.34
3	A	703	GDS	CD6-N5	5.69	1.45	1.34
3	B	703	GDS	CD6-N5	5.67	1.45	1.34
3	B	703	GDS	C2-N3	5.58	1.45	1.33
2	B	707	LDA	C8-C7	-5.47	1.53	1.55
2	B	709	LDA	C7-C6	-5.32	1.53	1.55
3	B	703	GDS	C5-N4	5.29	1.45	1.33
2	A	707	LDA	C3-C4	-5.15	1.53	1.55
2	B	709	LDA	C3-C4	-5.13	1.53	1.55
2	A	710	LDA	C7-C8	-5.12	1.53	1.55
2	B	707	LDA	C3-C4	-5.08	1.53	1.55
2	B	706	LDA	C4-C5	-4.82	1.53	1.55
2	A	708	LDA	C3-C4	-4.57	1.53	1.55
2	A	709	LDA	C3-C4	-4.51	1.53	1.55
2	B	706	LDA	C10-C9	-4.50	1.53	1.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	708	LDA	C7-C6	-4.40	1.53	1.55
2	A	707	LDA	C8-C7	-4.37	1.53	1.55
2	A	709	LDA	C8-C7	-4.36	1.53	1.55
2	A	710	LDA	C10-C9	-4.22	1.53	1.55
2	A	708	LDA	C7-C6	-3.68	1.53	1.55
2	A	701	LDA	C1-N1	-2.73	1.46	1.51
3	A	703	GDS	CG1-CD1	2.62	1.56	1.51
2	A	706	LDA	C1-N1	-2.48	1.46	1.51
2	B	701	LDA	C1-N1	-2.45	1.46	1.51
2	B	702	LDA	C1-N1	-2.39	1.46	1.51
3	B	703	GDS	CG1-CD1	2.37	1.56	1.51
3	A	703	GDS	CG6-CD6	2.20	1.55	1.51
3	B	703	GDS	CG6-CD6	2.09	1.55	1.51
2	B	708	LDA	C1-N1	-2.04	1.47	1.51

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	708	LDA	C2-C1-N1	-3.70	107.43	113.80
3	A	703	GDS	CB5-SG5-SG2	3.40	110.65	103.87
2	A	701	LDA	C2-C1-N1	-3.39	107.96	113.80
3	B	703	GDS	CG1-CB1-CA1	-3.35	108.18	114.43
3	A	703	GDS	CA5-C5-N4	3.13	122.83	116.72
3	A	703	GDS	CB2-SG2-SG5	3.12	110.10	103.87
3	B	703	GDS	CB2-SG2-SG5	3.10	110.06	103.87
3	B	703	GDS	CB6-CA6-C6	-3.01	103.85	110.98
3	B	703	GDS	CB5-SG5-SG2	2.80	109.45	103.87
3	A	703	GDS	CA2-C2-N3	2.61	121.82	116.72
3	B	703	GDS	C4-CA4-N4	-2.57	106.67	112.83
3	B	703	GDS	C3-CA3-N3	-2.47	106.91	112.83
3	A	703	GDS	O5-C5-N4	-2.45	118.07	123.05
3	A	703	GDS	CG1-CD1-N2	2.39	120.36	115.83
2	A	702	LDA	CM2-N1-CM1	-2.33	106.19	108.85
3	A	703	GDS	O1-CD1-CG1	-2.20	117.56	121.92
3	A	703	GDS	CG6-CB6-CA6	-2.18	110.37	114.43
3	B	703	GDS	OE8-C6-CA6	2.14	121.69	116.88
3	A	703	GDS	OE4-C3-CA3	2.07	120.49	112.98
3	A	703	GDS	OE6-C4-CA4	2.05	120.44	112.98
2	A	706	LDA	C2-C1-N1	-2.03	110.30	113.80
3	B	703	GDS	OE2-C1-CA1	2.01	121.40	116.88
3	B	703	GDS	OE2-C1-OE1	-2.00	119.54	124.07

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2		OWAB(Å ²)	Q < 0.9
1	A	600/614 (97%)	0.46	47 (7%)	13 15	33, 60, 104, 146	0
1	B	598/614 (97%)	0.42	47 (7%)	13 15	31, 56, 107, 165	0
All	All	1198/1228 (97%)	0.44	94 (7%)	12 15	31, 58, 105, 165	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	71	GLY	19.8
1	B	70	GLY	13.9
1	B	72	GLY	11.2
1	A	138	GLY	9.9
1	A	137	THR	8.6
1	B	604	SER	7.8
1	A	133	LEU	6.8
1	B	69	LEU	6.7
1	B	133	LEU	6.5
1	A	601	ALA	6.4
1	A	603	VAL	6.4
1	B	135	ARG	6.3
1	B	73	ALA	6.3
1	B	603	VAL	6.1
1	A	607	ALA	6.0
1	A	139	GLU	5.3
1	B	74	GLN	5.2
1	A	604	SER	5.2
1	A	605	GLU	5.1
1	B	138	GLY	5.0
1	B	606	ALA	4.7
1	B	134	ALA	4.7
1	B	605	GLU	4.6
1	B	137	THR	4.6

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Mol	Chain	Res	Type	RSRZ
1	A	177	TRP	4.4
1	A	17	TRP	4.3
1	A	600	SER	4.2
1	A	135	ARG	4.1
1	B	177	TRP	4.1
1	B	132	HIS	4.0
1	B	14	HIS	3.9
1	A	602	GLU	3.9
1	B	76	ALA	3.8
1	B	75	PRO	3.8
1	A	134	ALA	3.8
1	A	136	ARG	3.7
1	B	229	LEU	3.6
1	A	14	HIS	3.6
1	A	132	HIS	3.6
1	A	606	ALA	3.5
1	A	31	ASP	3.5
1	B	131	PHE	3.5
1	A	28	TRP	3.4
1	A	70	GLY	3.4
1	B	24	LEU	3.4
1	A	479	PHE	3.4
1	B	17	TRP	3.3
1	A	9	PRO	3.1
1	A	141	THR	3.0
1	A	267	LEU	3.0
1	A	531	ARG	3.0
1	B	79	VAL	2.9
1	A	30	ALA	2.9
1	B	494	ARG	2.9
1	B	460	ASP	2.8
1	A	58	PHE	2.8
1	A	309	THR	2.7
1	A	32	ASN	2.7
1	B	78	THR	2.7
1	A	157	PHE	2.6
1	B	601	ALA	2.6
1	B	28	TRP	2.6
1	B	31	ASP	2.5
1	A	208	HIS	2.5
1	A	156	TYR	2.5
1	B	139	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	462	ALA	2.5
1	B	183	GLY	2.4
1	B	181	GLY	2.4
1	A	160	PHE	2.4
1	A	158	LEU	2.4
1	B	136	ARG	2.4
1	B	182	LEU	2.3
1	A	317	MSE	2.3
1	B	350	ASN	2.3
1	B	602	GLU	2.3
1	A	306	THR	2.3
1	A	153	THR	2.3
1	A	155	LEU	2.2
1	B	598	ALA	2.2
1	B	83	PHE	2.2
1	A	13	ARG	2.2
1	A	313	ARG	2.2
1	B	309	THR	2.1
1	A	34	VAL	2.1
1	B	33	ALA	2.1
1	B	141	THR	2.1
1	B	173	ILE	2.1
1	B	32	ASN	2.0
1	A	321	VAL	2.0
1	A	530	THR	2.0
1	B	208	HIS	2.0
1	A	405	ARG	2.0
1	B	186	THR	2.0

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

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	LDA	A	707	6/16	0.56	20.10	63,72,74,74	0
2	LDA	B	709	5/16	0.35	15.45	53,58,62,64	0
2	LDA	A	708	5/16	0.61	14.09	68,72,77,78	0
2	LDA	B	706	7/16	0.57	12.30	60,69,74,76	0
2	LDA	B	702	16/16	0.39	7.67	60,78,101,108	0
2	LDA	B	701	16/16	0.41	7.67	58,72,82,86	0
2	LDA	A	701	16/16	0.54	7.46	55,66,101,111	0
2	LDA	A	706	16/16	0.45	5.79	57,80,111,118	0
3	GDS	A	703	40/40	0.37	4.32	79,117,136,137	0
3	GDS	B	703	40/40	0.35	4.26	42,84,111,133	0
2	LDA	A	709	6/16	0.25	3.96	56,59,64,65	0
2	LDA	B	707	6/16	0.36	3.69	84,89,91,91	0
2	LDA	B	708	11/16	0.39	3.67	61,71,89,89	0
4	PO4	B	704	5/5	0.16	2.28	134,136,137,140	0
2	LDA	A	702	16/16	0.40	1.94	55,67,89,92	0
4	PO4	B	705	5/5	0.13	1.56	133,136,138,140	0
4	PO4	A	704	5/5	0.12	0.63	150,151,151,151	0
2	LDA	A	710	4/16	0.23	0.51	49,55,58,61	0
4	PO4	A	705	5/5	0.13	0.04	79,80,85,88	5

6.5 Other polymers ⓘ

There are no such residues in this entry.