



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:14 PM GMT

PDB ID : 4R9M
Title : Crystal structure of spermidine N-acetyltransferase from Escherichia coli
Authors : Filippova, E.V.; Minasov, G.; Kiryukhina, O.; Shuvalova, L.; Grimshaw, S.; Wolfe, A.J.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2014-09-05
Resolution : 2.90 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

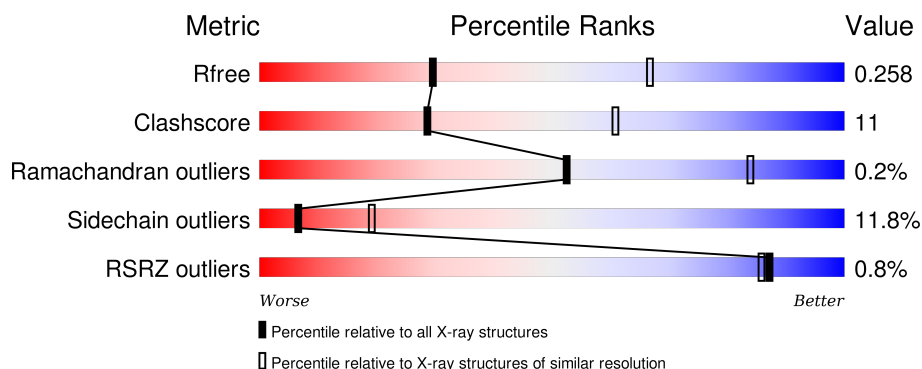
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION




The reported resolution of this entry is 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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	210	 60% 18% • 20%
1	B	210	 53% 20% • 22%
1	C	210	 54% 19% 5% 22%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4201 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 Spermidine N(1)-acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	169	Total	C	N	O	S	0	0	0
			1417	910	245	255	7			
1	B	163	Total	C	N	O	S	0	0	0
			1377	887	239	245	6			
1	C	164	Total	C	N	O	S	0	0	0
			1385	891	240	248	6			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	EXPRESSION TAG	UNP P0A951
A	-22	HIS	-	EXPRESSION TAG	UNP P0A951
A	-21	HIS	-	EXPRESSION TAG	UNP P0A951
A	-20	HIS	-	EXPRESSION TAG	UNP P0A951
A	-19	HIS	-	EXPRESSION TAG	UNP P0A951
A	-18	HIS	-	EXPRESSION TAG	UNP P0A951
A	-17	HIS	-	EXPRESSION TAG	UNP P0A951
A	-16	SER	-	EXPRESSION TAG	UNP P0A951
A	-15	SER	-	EXPRESSION TAG	UNP P0A951
A	-14	GLY	-	EXPRESSION TAG	UNP P0A951
A	-13	VAL	-	EXPRESSION TAG	UNP P0A951
A	-12	ASP	-	EXPRESSION TAG	UNP P0A951
A	-11	LEU	-	EXPRESSION TAG	UNP P0A951
A	-10	GLY	-	EXPRESSION TAG	UNP P0A951
A	-9	THR	-	EXPRESSION TAG	UNP P0A951
A	-8	GLU	-	EXPRESSION TAG	UNP P0A951
A	-7	ASN	-	EXPRESSION TAG	UNP P0A951
A	-6	LEU	-	EXPRESSION TAG	UNP P0A951
A	-5	TYR	-	EXPRESSION TAG	UNP P0A951
A	-4	PHE	-	EXPRESSION TAG	UNP P0A951
A	-3	GLN	-	EXPRESSION TAG	UNP P0A951
A	-2	SER	-	EXPRESSION TAG	UNP P0A951
A	-1	ASN	-	EXPRESSION TAG	UNP P0A951

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Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ALA	-	EXPRESSION TAG	UNP P0A951
B	-23	MET	-	EXPRESSION TAG	UNP P0A951
B	-22	HIS	-	EXPRESSION TAG	UNP P0A951
B	-21	HIS	-	EXPRESSION TAG	UNP P0A951
B	-20	HIS	-	EXPRESSION TAG	UNP P0A951
B	-19	HIS	-	EXPRESSION TAG	UNP P0A951
B	-18	HIS	-	EXPRESSION TAG	UNP P0A951
B	-17	HIS	-	EXPRESSION TAG	UNP P0A951
B	-16	SER	-	EXPRESSION TAG	UNP P0A951
B	-15	SER	-	EXPRESSION TAG	UNP P0A951
B	-14	GLY	-	EXPRESSION TAG	UNP P0A951
B	-13	VAL	-	EXPRESSION TAG	UNP P0A951
B	-12	ASP	-	EXPRESSION TAG	UNP P0A951
B	-11	LEU	-	EXPRESSION TAG	UNP P0A951
B	-10	GLY	-	EXPRESSION TAG	UNP P0A951
B	-9	THR	-	EXPRESSION TAG	UNP P0A951
B	-8	GLU	-	EXPRESSION TAG	UNP P0A951
B	-7	ASN	-	EXPRESSION TAG	UNP P0A951
B	-6	LEU	-	EXPRESSION TAG	UNP P0A951
B	-5	TYR	-	EXPRESSION TAG	UNP P0A951
B	-4	PHE	-	EXPRESSION TAG	UNP P0A951
B	-3	GLN	-	EXPRESSION TAG	UNP P0A951
B	-2	SER	-	EXPRESSION TAG	UNP P0A951
B	-1	ASN	-	EXPRESSION TAG	UNP P0A951
B	0	ALA	-	EXPRESSION TAG	UNP P0A951
C	-23	MET	-	EXPRESSION TAG	UNP P0A951
C	-22	HIS	-	EXPRESSION TAG	UNP P0A951
C	-21	HIS	-	EXPRESSION TAG	UNP P0A951
C	-20	HIS	-	EXPRESSION TAG	UNP P0A951
C	-19	HIS	-	EXPRESSION TAG	UNP P0A951
C	-18	HIS	-	EXPRESSION TAG	UNP P0A951
C	-17	HIS	-	EXPRESSION TAG	UNP P0A951
C	-16	SER	-	EXPRESSION TAG	UNP P0A951
C	-15	SER	-	EXPRESSION TAG	UNP P0A951
C	-14	GLY	-	EXPRESSION TAG	UNP P0A951
C	-13	VAL	-	EXPRESSION TAG	UNP P0A951
C	-12	ASP	-	EXPRESSION TAG	UNP P0A951
C	-11	LEU	-	EXPRESSION TAG	UNP P0A951
C	-10	GLY	-	EXPRESSION TAG	UNP P0A951
C	-9	THR	-	EXPRESSION TAG	UNP P0A951
C	-8	GLU	-	EXPRESSION TAG	UNP P0A951
C	-7	ASN	-	EXPRESSION TAG	UNP P0A951

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-6	LEU	-	EXPRESSION TAG	UNP P0A951
C	-5	TYR	-	EXPRESSION TAG	UNP P0A951
C	-4	PHE	-	EXPRESSION TAG	UNP P0A951
C	-3	GLN	-	EXPRESSION TAG	UNP P0A951
C	-2	SER	-	EXPRESSION TAG	UNP P0A951
C	-1	ASN	-	EXPRESSION TAG	UNP P0A951
C	0	ALA	-	EXPRESSION TAG	UNP P0A951

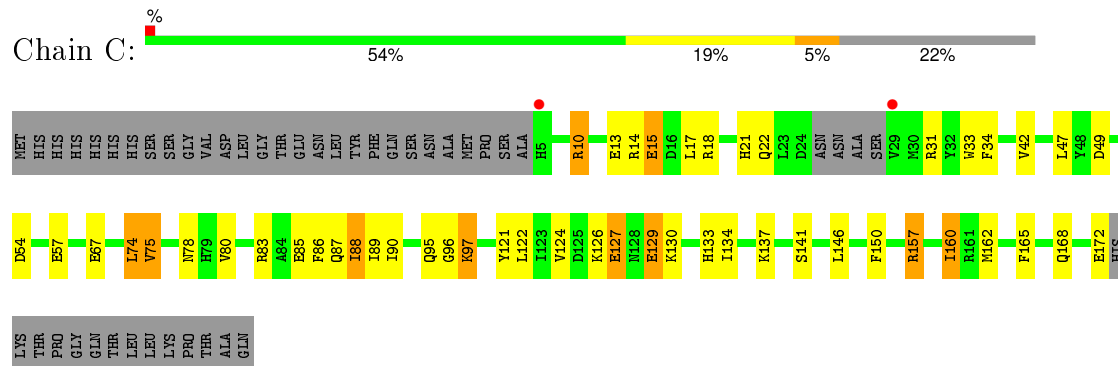
- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	15	Total O 15 15	0	0
3	B	3	Total O 3 3	0	0
3	C	3	Total O 3 3	0	0

- Molecule 1: Spermidine N(1)-acetyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	110.92Å 110.92Å 108.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.61 – 2.90 29.60 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.9 (29.61-2.90) 99.1 (29.60-2.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.10 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.176 , 0.259 0.174 , 0.258	Depositor DCC
R_{free} test set	774 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	77.3	Xtriage
Anisotropy	0.131	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 56.0	EDS
Estimated twinning fraction	0.010 for -h,-l,-k 0.000 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 15415 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4201	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.87% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 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	0.76	0/1450	0.91	0/1951
1	B	0.73	0/1409	0.94	1/1895 (0.1%)
1	C	0.75	0/1417	0.93	1/1906 (0.1%)
All	All	0.75	0/4276	0.93	2/5752 (0.0%)

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	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	71	LEU	CB-CG-CD1	-5.79	101.15	111.00
1	C	160	ILE	CG1-CB-CG2	-5.09	100.19	111.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	23	LEU	Peptide

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	1417	0	1386	25	0
1	B	1377	0	1348	35	0
1	C	1385	0	1352	38	0
2	A	1	0	0	0	0
3	A	15	0	0	0	0
3	B	3	0	0	0	0
3	C	3	0	0	0	0
All	All	4201	0	4086	95	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 11.

All (95) 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:B:31:ARG:HG3	1:B:31:ARG:HH11	1.17	1.05
1:C:10:ARG:HG2	1:C:10:ARG:HH11	1.35	0.89
1:A:30:MET:CE	1:A:30:MET:HA	2.09	0.83
1:B:31:ARG:CG	1:B:31:ARG:HH11	1.95	0.80
1:A:88:ILE:HD11	1:A:100:ALA:HA	1.63	0.79
1:C:130:LYS:HE3	1:C:134:ILE:HD11	1.68	0.75
1:B:169:TYR:O	1:B:172:GLU:HG2	1.87	0.75
1:C:21:HIS:O	1:C:31:ARG:NH2	2.21	0.74
1:C:13:GLU:HB3	1:C:15:GLU:OE1	1.88	0.74
1:C:33:TRP:HB3	1:C:47:LEU:HD13	1.71	0.72
1:C:10:ARG:HG2	1:C:10:ARG:NH1	2.02	0.70
1:B:49:ASP:OD1	1:B:49:ASP:N	2.25	0.69
1:B:31:ARG:NH1	1:B:31:ARG:HG3	1.97	0.68
1:A:30:MET:HE3	1:A:30:MET:HA	1.75	0.67
1:C:34:PHE:CE2	1:C:54:ASP:HB3	2.30	0.66
1:B:126:LYS:HA	1:B:160:ILE:HD13	1.78	0.66
1:B:52:ILE:HD11	1:C:42:VAL:HG12	1.80	0.63
1:A:30:MET:HE2	1:A:30:MET:HA	1.82	0.61
1:C:165:PHE:H	1:C:168:GLN:NE2	1.99	0.61
1:B:23:LEU:N	1:B:23:LEU:HD12	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:ILE:HD11	1:B:90:ILE:HD11	1.84	0.60
1:B:150:PHE:HB3	1:B:152:ILE:CD1	2.33	0.58
1:A:123:ILE:HG23	1:A:159:ALA:HB1	1.86	0.58
1:A:49:ASP:N	1:A:49:ASP:OD1	2.37	0.57
1:C:89:ILE:C	1:C:90:ILE:HD12	2.25	0.56
1:A:125:ASP:OD2	1:A:157:ARG:NH2	2.38	0.56
1:A:117:LEU:HD12	1:A:120:LEU:HD21	1.87	0.56
1:C:88:ILE:HG12	1:C:90:ILE:CD1	2.36	0.55
1:B:134:ILE:O	1:B:137:LYS:HB2	2.08	0.54
1:A:55:GLN:OE1	1:A:55:GLN:HA	2.07	0.54
1:A:57:GLU:HG2	1:A:75:VAL:HA	1.90	0.54
1:A:122:LEU:CD1	1:A:124:VAL:HG13	2.36	0.54
1:B:88:ILE:HD13	1:B:103:ALA:CB	2.37	0.54
1:C:10:ARG:CG	1:C:10:ARG:HH11	2.14	0.54
1:B:19:TYR:O	1:B:23:LEU:HD12	2.09	0.53
1:B:170:LEU:C	1:B:172:GLU:H	2.11	0.53
1:B:73:GLU:O	1:B:86:PHE:HA	2.09	0.53
1:B:150:PHE:HB3	1:B:152:ILE:HD12	1.91	0.52
1:C:96:GLY:O	1:C:97:LYS:HG3	2.10	0.52
1:A:13:GLU:HB3	1:A:15:GLU:OE1	2.10	0.52
1:C:165:PHE:H	1:C:168:GLN:HE21	1.55	0.52
1:B:152:ILE:HB	1:B:157:ARG:HG3	1.92	0.52
1:C:141:SER:O	1:C:162:MET:HA	2.10	0.52
1:C:126:LYS:HA	1:C:160:ILE:HD13	1.91	0.51
1:C:150:PHE:O	1:C:157:ARG:HG2	2.10	0.51
1:B:40:ALA:HB3	1:B:43:GLU:HG3	1.92	0.51
1:B:170:LEU:C	1:B:172:GLU:N	2.63	0.50
1:C:85:GLU:HB2	1:C:121:TYR:CE2	2.47	0.49
1:B:52:ILE:CD1	1:C:42:VAL:HG12	2.42	0.49
1:C:75:VAL:HG12	1:C:85:GLU:HB3	1.93	0.49
1:B:23:LEU:HD12	1:B:23:LEU:H	1.76	0.49
1:C:88:ILE:HD13	1:C:88:ILE:O	2.13	0.49
1:C:122:LEU:HD12	1:C:122:LEU:C	2.34	0.48
1:C:34:PHE:CE2	1:C:57:GLU:HB2	2.49	0.47
1:A:57:GLU:HG2	1:A:75:VAL:HG22	1.96	0.47
1:B:88:ILE:HD13	1:B:103:ALA:HB1	1.94	0.47
1:C:34:PHE:CZ	1:C:57:GLU:HB2	2.49	0.47
1:A:78:ASN:C	1:A:78:ASN:OD1	2.53	0.46
1:A:134:ILE:O	1:A:137:LYS:HB2	2.15	0.46
1:B:7:VAL:HA	1:B:63:GLU:O	2.16	0.46
1:A:15:GLU:H	1:A:15:GLU:CD	2.18	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:14:ARG:HH11	1:C:49:ASP:CG	2.19	0.46
1:B:34:PHE:HA	1:B:57:GLU:OE1	2.15	0.46
1:A:60:PHE:CD1	1:A:60:PHE:N	2.84	0.46
1:B:170:LEU:O	1:B:172:GLU:N	2.49	0.45
1:B:34:PHE:CE2	1:B:54:ASP:HB3	2.52	0.45
1:C:127:GLU:O	1:C:129:GLU:HG2	2.16	0.45
1:C:133:HIS:CE1	1:C:137:LYS:HE3	2.52	0.45
1:B:14:ARG:HB2	1:C:42:VAL:HG21	1.99	0.45
1:A:85:GLU:HB2	1:A:121:TYR:CE2	2.52	0.44
1:C:17:LEU:HA	1:C:17:LEU:HD12	1.87	0.44
1:A:117:LEU:CD1	1:A:120:LEU:HD21	2.47	0.43
1:C:78:ASN:OD1	1:C:80:VAL:N	2.51	0.43
1:B:59:ARG:HG2	1:B:73:GLU:HG2	2.01	0.43
1:A:99:LEU:HD23	1:A:99:LEU:HA	1.81	0.42
1:B:45:SER:O	1:B:49:ASP:OD1	2.38	0.42
1:C:90:ILE:O	1:C:95:GLN:NE2	2.51	0.42
1:C:33:TRP:HB3	1:C:47:LEU:CD1	2.46	0.42
1:C:96:GLY:C	1:C:97:LYS:HG3	2.40	0.42
1:C:14:ARG:O	1:C:17:LEU:HB2	2.20	0.42
1:A:23:LEU:O	1:A:24:ASP:HB3	2.20	0.42
1:C:157:ARG:HG2	1:C:157:ARG:H	1.72	0.41
1:B:150:PHE:CB	1:B:152:ILE:CD1	2.98	0.41
1:C:90:ILE:N	1:C:90:ILE:HD12	2.34	0.41
1:B:122:LEU:O	1:B:161:ARG:HA	2.21	0.41
1:A:106:LEU:HD23	1:A:106:LEU:HA	1.83	0.41
1:C:57:GLU:HA	1:C:74:LEU:O	2.21	0.41
1:C:86:PHE:C	1:C:87:GLN:HG3	2.41	0.41
1:B:99:LEU:O	1:B:100:ALA:C	2.59	0.41
1:A:164:ILE:HG13	1:A:165:PHE:H	1.86	0.40
1:B:55:GLN:HA	1:B:55:GLN:NE2	2.36	0.40
1:B:164:ILE:HG12	1:B:165:PHE:N	2.36	0.40
1:A:17:LEU:HD12	1:A:17:LEU:HA	1.83	0.40
1:B:13:GLU:O	1:B:16:ASP:HB2	2.22	0.40
1:A:143:GLU:HG3	1:A:161:ARG:HG2	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	165/210 (79%)	155 (94%)	10 (6%)	0	100	100
1	B	159/210 (76%)	148 (93%)	10 (6%)	1 (1%)	30	67
1	C	160/210 (76%)	144 (90%)	16 (10%)	0	100	100
All	All	484/630 (77%)	447 (92%)	36 (7%)	1 (0%)	52	84

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	171	ALA

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	150/185 (81%)	132 (88%)	18 (12%)	6	18
1	B	145/185 (78%)	127 (88%)	18 (12%)	6	17
1	C	146/185 (79%)	130 (89%)	16 (11%)	8	23
All	All	441/555 (80%)	389 (88%)	52 (12%)	6	19

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	15	GLU

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Mol	Chain	Res	Type
1	A	18	ARG
1	A	22	GLN
1	A	30	MET
1	A	31	ARG
1	A	35	GLU
1	A	49	ASP
1	A	67	GLU
1	A	74	LEU
1	A	75	VAL
1	A	88	ILE
1	A	91	SER
1	A	127	GLU
1	A	130	LYS
1	A	146	LEU
1	A	157	ARG
1	A	172	GLU
1	B	6	SER
1	B	17	LEU
1	B	18	ARG
1	B	22	GLN
1	B	31	ARG
1	B	44	LEU
1	B	47	LEU
1	B	49	ASP
1	B	55	GLN
1	B	71	LEU
1	B	75	VAL
1	B	88	ILE
1	B	122	LEU
1	B	124	VAL
1	B	127	GLU
1	B	145	GLU
1	B	152	ILE
1	B	160	ILE
1	C	10	ARG
1	C	15	GLU
1	C	18	ARG
1	C	22	GLN
1	C	67	GLU
1	C	74	LEU
1	C	75	VAL
1	C	83	ARG

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Mol	Chain	Res	Type
1	C	88	ILE
1	C	97	LYS
1	C	124	VAL
1	C	127	GLU
1	C	129	GLU
1	C	146	LEU
1	C	157	ARG
1	C	172	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	53	HIS
1	A	148	HIS
1	A	153	ASN
1	B	55	GLN
1	C	133	HIS
1	C	155	GLN
1	C	168	GLN

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 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [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	169/210 (80%)	-0.28	0 100 100	64, 78, 110, 126	0
1	B	163/210 (77%)	-0.27	2 (1%) 81 78	61, 79, 106, 129	0
1	C	164/210 (78%)	-0.24	2 (1%) 81 78	62, 81, 110, 148	0
All	All	496/630 (78%)	-0.26	4 (0%) 87 86	61, 80, 110, 148	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	29	VAL	4.1
1	C	5	HIS	2.7
1	B	127	GLU	2.4
1	B	129	GLU	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 carbohydrates 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. 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(\AA^2)	Q<0.9
2	MG	A	201	1/1	0.65	0.20	-	80,80,80,80	1

6.5 Other polymers [i](#)

There are no such residues in this entry.