



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:28 PM UTC

PDB ID : 3K8E / pdb_00003k8e
Title : Crystal structure of E. coli lipopolysaccharide specific CMP-KDO synthetase
Authors : Heyes, D.J.; Levy, C.W.; Lafite, P.; Scrutton, N.S.; Leys, D.
Deposited on : 2009-10-14
Resolution : 2.51 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
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

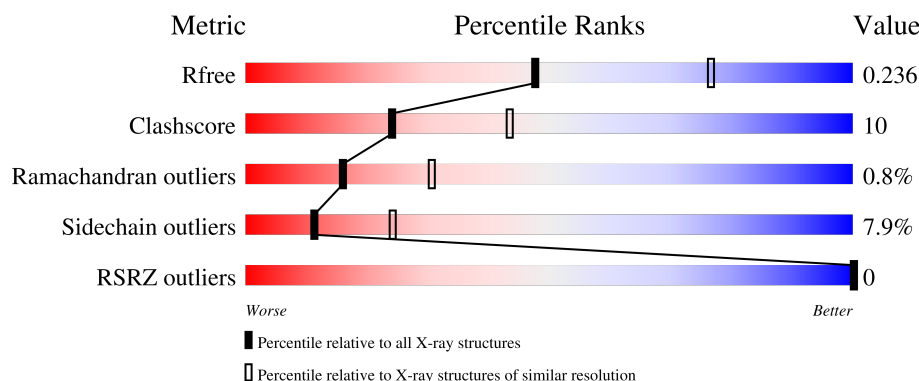
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	264	
1	B	264	
1	C	264	
1	D	264	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7657 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 3-deoxy-manno-octulosonate cytidyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	244	Total	C	N	O	S	0	0	0
			1895	1192	342	355	6			
1	D	238	Total	C	N	O	S	0	0	0
			1832	1161	324	339	8			
1	A	242	Total	C	N	O	S	0	0	0
			1882	1185	340	350	7			
1	B	246	Total	C	N	O	S	0	0	0
			1906	1200	344	355	7			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-15	HIS	-	expression tag	UNP P04951
C	-14	HIS	-	expression tag	UNP P04951
C	-13	HIS	-	expression tag	UNP P04951
C	-12	HIS	-	expression tag	UNP P04951
C	-11	HIS	-	expression tag	UNP P04951
C	-10	HIS	-	expression tag	UNP P04951
C	-9	SER	-	expression tag	UNP P04951
C	-8	SER	-	expression tag	UNP P04951
C	-7	GLY	-	expression tag	UNP P04951
C	-6	LEU	-	expression tag	UNP P04951
C	-5	VAL	-	expression tag	UNP P04951
C	-4	PRO	-	expression tag	UNP P04951
C	-3	ARG	-	expression tag	UNP P04951
C	-2	GLY	-	expression tag	UNP P04951
C	-1	SER	-	expression tag	UNP P04951
C	0	HIS	-	expression tag	UNP P04951
D	-15	HIS	-	expression tag	UNP P04951
D	-14	HIS	-	expression tag	UNP P04951
D	-13	HIS	-	expression tag	UNP P04951
D	-12	HIS	-	expression tag	UNP P04951
D	-11	HIS	-	expression tag	UNP P04951

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-10	HIS	-	expression tag	UNP P04951
D	-9	SER	-	expression tag	UNP P04951
D	-8	SER	-	expression tag	UNP P04951
D	-7	GLY	-	expression tag	UNP P04951
D	-6	LEU	-	expression tag	UNP P04951
D	-5	VAL	-	expression tag	UNP P04951
D	-4	PRO	-	expression tag	UNP P04951
D	-3	ARG	-	expression tag	UNP P04951
D	-2	GLY	-	expression tag	UNP P04951
D	-1	SER	-	expression tag	UNP P04951
D	0	HIS	-	expression tag	UNP P04951
A	-15	HIS	-	expression tag	UNP P04951
A	-14	HIS	-	expression tag	UNP P04951
A	-13	HIS	-	expression tag	UNP P04951
A	-12	HIS	-	expression tag	UNP P04951
A	-11	HIS	-	expression tag	UNP P04951
A	-10	HIS	-	expression tag	UNP P04951
A	-9	SER	-	expression tag	UNP P04951
A	-8	SER	-	expression tag	UNP P04951
A	-7	GLY	-	expression tag	UNP P04951
A	-6	LEU	-	expression tag	UNP P04951
A	-5	VAL	-	expression tag	UNP P04951
A	-4	PRO	-	expression tag	UNP P04951
A	-3	ARG	-	expression tag	UNP P04951
A	-2	GLY	-	expression tag	UNP P04951
A	-1	SER	-	expression tag	UNP P04951
A	0	HIS	-	expression tag	UNP P04951
B	-15	HIS	-	expression tag	UNP P04951
B	-14	HIS	-	expression tag	UNP P04951
B	-13	HIS	-	expression tag	UNP P04951
B	-12	HIS	-	expression tag	UNP P04951
B	-11	HIS	-	expression tag	UNP P04951
B	-10	HIS	-	expression tag	UNP P04951
B	-9	SER	-	expression tag	UNP P04951
B	-8	SER	-	expression tag	UNP P04951
B	-7	GLY	-	expression tag	UNP P04951
B	-6	LEU	-	expression tag	UNP P04951
B	-5	VAL	-	expression tag	UNP P04951
B	-4	PRO	-	expression tag	UNP P04951
B	-3	ARG	-	expression tag	UNP P04951
B	-2	GLY	-	expression tag	UNP P04951
B	-1	SER	-	expression tag	UNP P04951

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	HIS	-	expression tag	UNP P04951

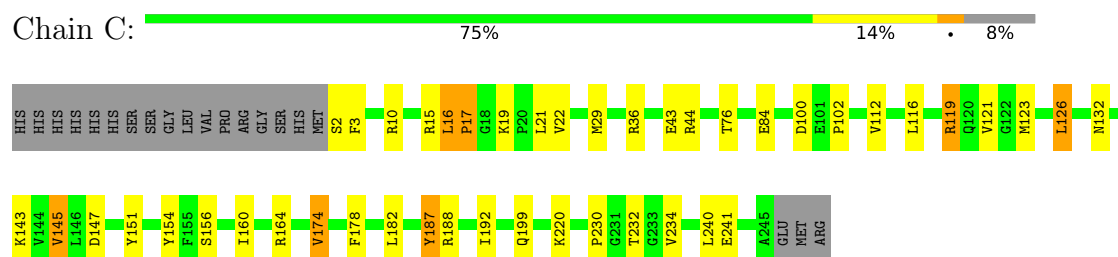
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	30	Total 30	O 30	0	0
2	D	39	Total 39	O 39	0	0
2	A	36	Total 36	O 36	0	0
2	B	37	Total 37	O 37	0	0

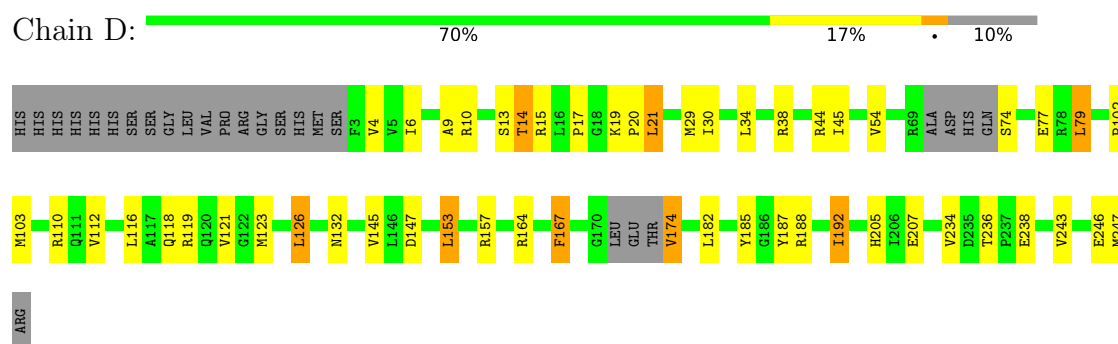
3 Residue-property plots

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.

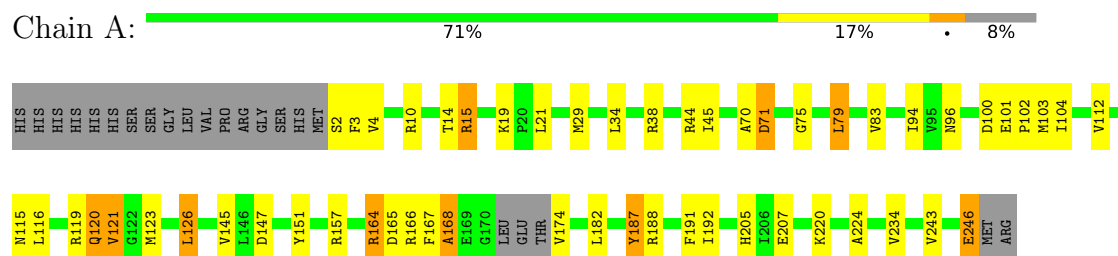
- Molecule 1: 3-deoxy-manno-octulosonate cytidyltransferase



- Molecule 1: 3-deoxy-manno-octulosonate cytidyltransferase

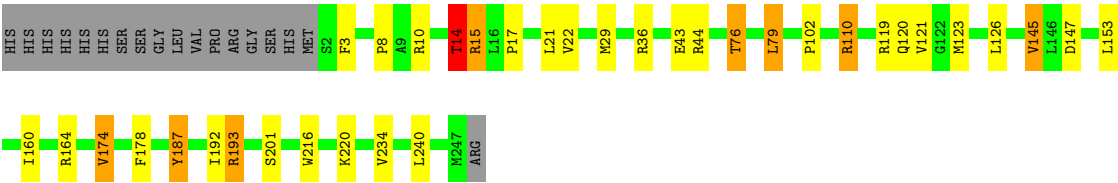


- Molecule 1: 3-deoxy-manno-octulosonate cytidyltransferase



- Molecule 1: 3-deoxy-manno-octulosonate cytidyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	124.47Å 77.18Å 143.63Å 90.00° 90.03° 90.00°	Depositor
Resolution (Å)	47.00 – 2.51 47.00 – 2.51	Depositor EDS
% Data completeness (in resolution range)	98.1 (47.00-2.51) 98.1 (47.00-2.51)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.82 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.199 , 0.234 0.206 , 0.236	Depositor DCC
R_{free} test set	2360 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	54.5	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.480 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7657	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.08% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.72	0/1919	0.85	1/2609 (0.0%)
1	B	0.75	0/1945	0.87	2/2648 (0.1%)
1	C	0.75	1/1933 (0.1%)	0.87	2/2631 (0.1%)
1	D	0.76	0/1868	0.87	0/2542
All	All	0.74	1/7665 (0.0%)	0.86	5/10430 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	2	SER	C-N	-5.84	1.25	1.33

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	3	PHE	N-CA-C	7.00	119.81	109.24
1	C	2	SER	O-C-N	-6.70	112.28	123.00
1	B	3	PHE	N-CA-C	5.77	117.85	109.07
1	B	14	THR	N-CA-C	-5.74	102.55	110.35
1	A	3	PHE	N-CA-C	5.05	116.75	109.07

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1882	0	1851	47	0
1	B	1906	0	1871	28	0
1	C	1895	0	1858	32	0
1	D	1832	0	1798	42	0
2	A	36	0	0	1	0
2	B	37	0	0	1	0
2	C	30	0	0	0	0
2	D	39	0	0	1	0
All	All	7657	0	7378	149	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 10.

All (149) 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:29:MET:HE2	1:B:102:PRO:HD3	1.35	1.09
1:B:193:ARG:HH11	1:B:193:ARG:HG3	0.96	1.07
1:D:29:MET:HE1	1:D:234:VAL:HB	1.42	1.01
1:A:29:MET:HE1	1:A:234:VAL:HB	1.42	1.01
1:A:123:MET:HE3	1:A:191:PHE:HE2	1.30	0.97
1:B:193:ARG:HH11	1:B:193:ARG:CG	1.78	0.94
1:D:103:MET:HE2	1:D:247:MET:HG3	1.47	0.94
1:B:29:MET:HE1	1:B:234:VAL:HB	1.46	0.94
1:C:29:MET:HE1	1:C:234:VAL:HB	1.47	0.93
1:B:193:ARG:HG3	1:B:193:ARG:NH1	1.75	0.93
1:A:79:LEU:HD11	1:A:96:ASN:ND2	1.92	0.85
1:A:116:LEU:HD11	1:A:188:ARG:HD2	1.59	0.84
1:A:123:MET:HE3	1:A:191:PHE:CE2	2.17	0.79
1:A:29:MET:HE2	1:A:102:PRO:HD3	1.68	0.76
1:A:29:MET:CE	1:A:102:PRO:HD3	2.16	0.75
1:A:167:PHE:O	1:A:168:ALA:CB	2.34	0.75
1:D:103:MET:HE3	1:D:243:VAL:HG13	1.69	0.74
1:A:123:MET:HE2	1:A:187:TYR:CE2	2.23	0.74
1:B:76:THR:CG2	1:B:192:ILE:HD11	2.17	0.73
1:C:29:MET:CE	1:C:102:PRO:HD3	2.20	0.72
1:D:174:VAL:N	2:D:256:HOH:O	2.21	0.72
1:A:112:VAL:HG22	1:A:126:LEU:HD13	1.71	0.71
1:D:29:MET:CE	1:D:102:PRO:HD3	2.20	0.71
1:C:29:MET:HE2	1:C:102:PRO:HD3	1.72	0.70
1:C:123:MET:HE2	1:C:187:TYR:CE2	2.26	0.70
1:A:116:LEU:CD1	1:A:188:ARG:HD2	2.21	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:10:ARG:O	1:C:10:ARG:HG2	1.92	0.70
1:D:29:MET:HE2	1:D:102:PRO:HD3	1.76	0.68
1:C:16:LEU:O	1:C:17:PRO:CB	2.42	0.67
1:D:103:MET:HE3	1:D:243:VAL:HG22	1.76	0.67
1:D:103:MET:CE	1:D:243:VAL:HG13	2.26	0.66
1:D:14:THR:O	1:D:15:ARG:CG	2.44	0.65
1:B:121:VAL:HG22	1:B:220:LYS:O	1.95	0.65
1:A:123:MET:HE2	1:A:187:TYR:CD2	2.32	0.64
1:C:29:MET:HE1	1:C:234:VAL:CB	2.26	0.63
1:B:29:MET:CE	1:B:102:PRO:HD3	2.21	0.63
1:D:103:MET:HE2	1:D:247:MET:CG	2.26	0.63
1:A:103:MET:HE1	1:A:246:GLU:C	2.23	0.63
1:A:29:MET:CE	1:A:234:VAL:HB	2.25	0.62
1:D:112:VAL:HG22	1:D:126:LEU:HD13	1.83	0.61
1:D:116:LEU:HD11	1:D:188:ARG:HG2	1.83	0.60
1:C:22:VAL:HG11	1:C:240:LEU:HD21	1.83	0.59
1:D:164:ARG:C	1:D:167:PHE:O	2.45	0.59
1:B:14:THR:O	1:B:15:ARG:NE	2.36	0.59
1:A:167:PHE:O	1:A:168:ALA:HB2	2.02	0.59
1:B:123:MET:HE2	1:B:187:TYR:CD2	2.38	0.59
1:C:145:VAL:HG12	1:C:160:ILE:HD12	1.84	0.58
1:A:101:GLU:OE2	1:A:104:ILE:HD12	2.02	0.58
1:C:76:THR:CG2	1:C:192:ILE:HD11	2.33	0.58
1:B:29:MET:HE1	1:B:234:VAL:CB	2.27	0.58
1:B:29:MET:HE2	1:B:102:PRO:CD	2.22	0.58
1:A:79:LEU:CD1	1:A:96:ASN:ND2	2.66	0.57
1:A:120:GLN:O	1:A:120:GLN:CG	2.52	0.57
1:A:121:VAL:HG21	1:A:151:TYR:HE1	1.69	0.57
1:D:123:MET:HE3	1:D:185:TYR:CD2	2.40	0.56
1:A:116:LEU:HD11	1:A:188:ARG:CD	2.33	0.56
1:A:123:MET:HE2	1:A:187:TYR:HE2	1.71	0.56
1:A:70:ALA:O	1:A:71:ASP:C	2.49	0.55
1:A:75:GLY:O	1:A:79:LEU:HD12	2.06	0.55
1:C:29:MET:HE2	1:C:100:ASP:C	2.31	0.55
1:D:14:THR:O	1:D:15:ARG:HG3	2.05	0.55
1:C:43:GLU:O	1:C:44:ARG:HG3	2.07	0.54
1:C:121:VAL:HG22	1:C:220:LYS:O	2.07	0.54
1:B:43:GLU:O	1:B:44:ARG:HG3	2.08	0.53
1:D:103:MET:HE1	1:D:243:VAL:HA	1.92	0.52
1:C:174:VAL:HG22	1:C:178:PHE:CE2	2.45	0.51
1:D:38:ARG:HG3	1:D:45:ILE:HD13	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:ARG:HB3	1:A:121:VAL:HG23	1.92	0.51
1:B:123:MET:HE2	1:B:187:TYR:CE2	2.45	0.51
1:B:110:ARG:HH11	1:B:110:ARG:HG2	1.75	0.51
1:B:145:VAL:CG1	1:B:160:ILE:HD12	2.40	0.51
1:C:29:MET:HE3	1:C:102:PRO:HD3	1.92	0.51
1:C:119:ARG:HG2	1:C:151:TYR:OH	2.10	0.51
1:C:22:VAL:HG11	1:C:240:LEU:CD2	2.41	0.51
1:D:21:LEU:HD13	1:D:54:VAL:HG22	1.93	0.50
1:D:29:MET:HE1	1:D:234:VAL:CB	2.30	0.50
1:B:123:MET:HG3	1:B:187:TYR:HD2	1.75	0.50
1:C:123:MET:HE2	1:C:187:TYR:HE2	1.76	0.50
1:D:29:MET:HE3	1:D:102:PRO:HD3	1.94	0.50
1:D:74:SER:CB	1:D:77:GLU:OE1	2.60	0.50
1:D:123:MET:HE3	1:D:185:TYR:HD2	1.76	0.50
1:D:14:THR:O	1:D:15:ARG:HG2	2.12	0.49
1:C:143:LYS:N	1:C:143:LYS:HD2	2.28	0.49
1:C:121:VAL:HG23	1:C:220:LYS:HB2	1.94	0.49
1:D:4:VAL:HG22	1:D:44:ARG:HB2	1.94	0.49
1:A:29:MET:HE2	1:A:100:ASP:HA	1.95	0.49
1:A:29:MET:HE3	1:A:102:PRO:HD3	1.90	0.49
1:C:147:ASP:C	1:C:147:ASP:OD1	2.56	0.48
1:A:165:ASP:O	1:A:166:ARG:C	2.53	0.48
1:B:147:ASP:OD1	1:B:147:ASP:C	2.57	0.48
1:D:19:LYS:HB3	1:D:20:PRO:HD3	1.95	0.48
1:C:29:MET:CE	1:C:234:VAL:HB	2.33	0.47
1:C:123:MET:HE2	1:C:187:TYR:CD2	2.49	0.47
1:D:164:ARG:O	1:D:167:PHE:O	2.32	0.47
1:A:121:VAL:HG13	1:A:220:LYS:HB2	1.96	0.47
1:D:29:MET:CE	1:D:234:VAL:HB	2.28	0.47
1:B:145:VAL:HG12	1:B:160:ILE:HD12	1.97	0.47
1:D:9:ALA:HB2	1:D:30:ILE:HD13	1.97	0.47
1:D:34:LEU:HD12	1:D:45:ILE:HG21	1.97	0.47
1:B:201:SER:HB3	1:B:216:TRP:CG	2.51	0.46
1:A:120:GLN:O	1:A:120:GLN:HG3	2.15	0.46
1:C:29:MET:HE2	1:C:100:ASP:HA	1.97	0.46
1:A:157:ARG:HD2	1:A:207:GLU:O	2.16	0.46
1:A:38:ARG:HG2	2:A:278:HOH:O	2.16	0.45
1:A:34:LEU:HD12	1:A:45:ILE:HG21	1.98	0.45
1:D:6:ILE:HD12	1:D:79:LEU:HD23	1.98	0.45
1:B:201:SER:HB3	1:B:216:TRP:CD1	2.50	0.45
1:A:4:VAL:HB	1:A:94:ILE:HD13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:118:GLN:HE21	1:D:119:ARG:NH2	2.15	0.45
1:A:167:PHE:O	1:A:168:ALA:HB3	2.14	0.45
1:D:192:ILE:HD13	1:D:192:ILE:HA	1.75	0.44
1:D:145:VAL:HG23	1:D:153:LEU:HB2	1.99	0.44
1:A:19:LYS:HZ1	1:A:100:ASP:HB3	1.81	0.44
1:B:8:PRO:HG3	1:B:79:LEU:CD1	2.46	0.44
1:B:121:VAL:HG23	1:B:220:LYS:HB2	1.99	0.44
1:D:103:MET:HE3	1:D:243:VAL:CG2	2.47	0.44
1:D:123:MET:CE	1:D:185:TYR:CD2	3.00	0.44
1:C:22:VAL:CG1	1:C:240:LEU:HD21	2.48	0.44
1:D:103:MET:HE3	1:D:243:VAL:CG1	2.43	0.44
1:C:116:LEU:HD11	1:C:188:ARG:HG3	2.00	0.44
1:C:154:TYR:CE2	1:C:156:SER:HB2	2.53	0.43
1:A:29:MET:HE3	1:A:102:PRO:CD	2.48	0.43
1:A:14:THR:O	1:A:15:ARG:CB	2.67	0.43
1:B:201:SER:CB	1:B:216:TRP:CG	3.02	0.43
1:A:14:THR:O	1:A:15:ARG:HB2	2.17	0.43
1:C:123:MET:HG2	1:C:187:TYR:HD2	1.84	0.42
1:D:103:MET:CE	1:D:243:VAL:HA	2.49	0.42
1:C:76:THR:HG23	1:C:192:ILE:HD11	2.00	0.42
1:A:192:ILE:HD13	1:A:192:ILE:HA	1.87	0.42
1:D:15:ARG:HD3	1:D:236:THR:HG22	2.01	0.42
1:C:192:ILE:HA	1:C:192:ILE:HD12	1.87	0.42
1:D:147:ASP:HB3	1:D:153:LEU:HD13	2.02	0.42
1:A:4:VAL:HG22	1:A:44:ARG:HB2	2.01	0.42
1:A:19:LYS:NZ	1:A:100:ASP:HB3	2.34	0.41
1:A:147:ASP:OD1	1:A:147:ASP:C	2.63	0.41
1:A:29:MET:HE2	1:A:100:ASP:C	2.46	0.41
1:A:103:MET:HE3	1:A:243:VAL:HG13	2.03	0.41
1:A:164:ARG:O	1:A:167:PHE:O	2.39	0.41
1:B:192:ILE:HD12	1:B:192:ILE:HA	1.91	0.41
1:D:157:ARG:HD2	1:D:207:GLU:O	2.21	0.41
1:B:22:VAL:HG11	1:B:240:LEU:CD2	2.51	0.41
1:D:103:MET:HE1	1:D:246:GLU:HB2	2.02	0.41
1:D:21:LEU:HD11	1:D:54:VAL:HA	2.03	0.40
1:A:10:ARG:HH12	1:A:19:LYS:HZ1	1.69	0.40
1:C:112:VAL:HG22	1:C:126:LEU:HD13	2.03	0.40
1:C:182:LEU:HD21	1:C:230:PRO:CG	2.52	0.40
1:B:10:ARG:NH2	2:B:252:HOH:O	2.53	0.40
1:B:174:VAL:HG22	1:B:178:PHE:CE2	2.55	0.40
1:A:115:ASN:HD21	1:A:224:ALA:CB	2.34	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	238/264 (90%)	227 (95%)	8 (3%)	3 (1%)	9	18
1	B	244/264 (92%)	235 (96%)	8 (3%)	1 (0%)	30	49
1	C	242/264 (92%)	232 (96%)	8 (3%)	2 (1%)	16	31
1	D	232/264 (88%)	220 (95%)	10 (4%)	2 (1%)	14	27
All	All	956/1056 (90%)	914 (96%)	34 (4%)	8 (1%)	16	31

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	16	LEU
1	C	17	PRO
1	A	168	ALA
1	D	17	PRO
1	A	71	ASP
1	D	13	SER
1	A	15	ARG
1	B	17	PRO

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	194/217 (89%)	180 (93%)	14 (7%)	13	28
1	B	196/217 (90%)	180 (92%)	16 (8%)	10	23
1	C	195/217 (90%)	180 (92%)	15 (8%)	12	25
1	D	187/217 (86%)	171 (91%)	16 (9%)	10	21
All	All	772/868 (89%)	711 (92%)	61 (8%)	11	24

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	15	ARG
1	C	19	LYS
1	C	21	LEU
1	C	36	ARG
1	C	84	GLU
1	C	119	ARG
1	C	126	LEU
1	C	132	ASN
1	C	145	VAL
1	C	164	ARG
1	C	174	VAL
1	C	187	TYR
1	C	199	GLN
1	C	232	THR
1	C	241	GLU
1	D	10	ARG
1	D	14	THR
1	D	21	LEU
1	D	79	LEU
1	D	110	ARG
1	D	121	VAL
1	D	126	LEU
1	D	132	ASN
1	D	153	LEU
1	D	167	PHE
1	D	174	VAL
1	D	182	LEU
1	D	187	TYR
1	D	192	ILE
1	D	205	HIS
1	D	238	GLU
1	A	2	SER
1	A	21	LEU

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Mol	Chain	Res	Type
1	A	79	LEU
1	A	83	VAL
1	A	120	GLN
1	A	121	VAL
1	A	126	LEU
1	A	145	VAL
1	A	164	ARG
1	A	174	VAL
1	A	182	LEU
1	A	187	TYR
1	A	205	HIS
1	A	246	GLU
1	B	14	THR
1	B	15	ARG
1	B	21	LEU
1	B	36	ARG
1	B	76	THR
1	B	79	LEU
1	B	110	ARG
1	B	119	ARG
1	B	120	GLN
1	B	126	LEU
1	B	145	VAL
1	B	153	LEU
1	B	164	ARG
1	B	174	VAL
1	B	187	TYR
1	B	193	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	51	HIS
1	D	51	HIS
1	D	115	ASN
1	D	118	GLN
1	D	140	ASN
1	D	222	HIS
1	A	115	ASN
1	A	197	ASN
1	B	51	HIS
1	B	73	GLN

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Mol	Chain	Res	Type
1	B	205	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](#)

There are no ligands in this entry.

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	242/264 (91%)	-1.16	0 100 100	47, 54, 65, 70	0
1	B	246/264 (93%)	-1.16	0 100 100	48, 54, 61, 68	0
1	C	244/264 (92%)	-1.15	0 100 100	20, 54, 60, 66	0
1	D	238/264 (90%)	-1.19	0 100 100	48, 54, 63, 66	0
All	All	970/1056 (91%)	-1.16	0 100 100	20, 54, 63, 70	0

There are no RSRZ outliers to report.

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.