



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 10:07 PM UTC

PDB ID : 3STG / pdb_00003stg
Title : Crystal structure of A58P, DEL(N59), and loop 7 truncated mutant of 3-deoxy-D-manno-octulosonate 8-phosphate synthase (KDO8PS) from *Neisseria meningitidis*
Authors : Allison, T.M.; Jameson, G.B.; Parker, E.J.
Deposited on : 2011-07-09
Resolution : 2.20 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

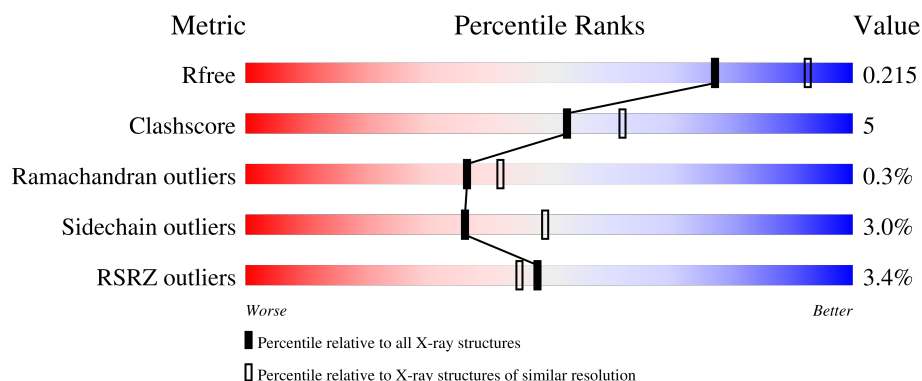
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	268	
1	B	268	
1	C	268	
1	D	268	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8729 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 2-dehydro-3-deoxyphosphooctonate aldolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	260	Total	C	N	O	S	0	2	0
			2020	1298	340	371	11			
1	B	268	Total	C	N	O	S	0	4	0
			2080	1337	349	382	12			
1	C	260	Total	C	N	O	S	0	4	0
			2030	1307	342	369	12			
1	D	258	Total	C	N	O	S	0	3	0
			2013	1296	339	366	12			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	58	PRO	ALA	engineered mutation	UNP Q9JZ55
A	?	-	ASN	deletion	UNP Q9JZ55
A	?	-	GLN	deletion	UNP Q9JZ55
A	?	-	THR	deletion	UNP Q9JZ55
A	?	-	ARG	deletion	UNP Q9JZ55
A	?	-	ASP	deletion	UNP Q9JZ55
A	?	-	ALA	deletion	UNP Q9JZ55
A	?	-	GLY	deletion	UNP Q9JZ55
A	?	-	SER	deletion	UNP Q9JZ55
A	?	-	ALA	deletion	UNP Q9JZ55
A	?	-	ALA	deletion	UNP Q9JZ55
A	?	-	SER	deletion	UNP Q9JZ55
A	?	-	GLY	deletion	UNP Q9JZ55
B	58	PRO	ALA	engineered mutation	UNP Q9JZ55
B	?	-	ASN	deletion	UNP Q9JZ55
B	?	-	GLN	deletion	UNP Q9JZ55
B	?	-	THR	deletion	UNP Q9JZ55
B	?	-	ARG	deletion	UNP Q9JZ55
B	?	-	ASP	deletion	UNP Q9JZ55
B	?	-	ALA	deletion	UNP Q9JZ55
B	?	-	GLY	deletion	UNP Q9JZ55

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	SER	deletion	UNP Q9JZ55
B	?	-	ALA	deletion	UNP Q9JZ55
B	?	-	ALA	deletion	UNP Q9JZ55
B	?	-	SER	deletion	UNP Q9JZ55
B	?	-	GLY	deletion	UNP Q9JZ55
C	58	PRO	ALA	engineered mutation	UNP Q9JZ55
C	?	-	ASN	deletion	UNP Q9JZ55
C	?	-	GLN	deletion	UNP Q9JZ55
C	?	-	THR	deletion	UNP Q9JZ55
C	?	-	ARG	deletion	UNP Q9JZ55
C	?	-	ASP	deletion	UNP Q9JZ55
C	?	-	ALA	deletion	UNP Q9JZ55
C	?	-	GLY	deletion	UNP Q9JZ55
C	?	-	SER	deletion	UNP Q9JZ55
C	?	-	ALA	deletion	UNP Q9JZ55
C	?	-	ALA	deletion	UNP Q9JZ55
C	?	-	SER	deletion	UNP Q9JZ55
C	?	-	GLY	deletion	UNP Q9JZ55
D	58	PRO	ALA	engineered mutation	UNP Q9JZ55
D	?	-	ASN	deletion	UNP Q9JZ55
D	?	-	GLN	deletion	UNP Q9JZ55
D	?	-	THR	deletion	UNP Q9JZ55
D	?	-	ARG	deletion	UNP Q9JZ55
D	?	-	ASP	deletion	UNP Q9JZ55
D	?	-	ALA	deletion	UNP Q9JZ55
D	?	-	GLY	deletion	UNP Q9JZ55
D	?	-	SER	deletion	UNP Q9JZ55
D	?	-	ALA	deletion	UNP Q9JZ55
D	?	-	ALA	deletion	UNP Q9JZ55
D	?	-	SER	deletion	UNP Q9JZ55
D	?	-	GLY	deletion	UNP Q9JZ55

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total Cl 3 3	0	0
2	B	4	Total Cl 4 4	0	0
2	C	2	Total Cl 2 2	0	0
2	D	3	Total Cl 3 3	0	0

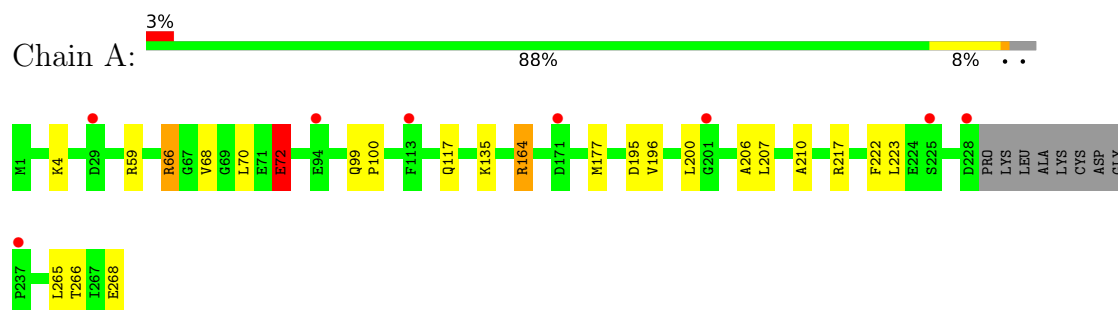
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	150	Total 150	O 150	0	0
3	B	138	Total 138	O 138	0	0
3	C	135	Total 135	O 135	0	0
3	D	151	Total 151	O 151	0	0

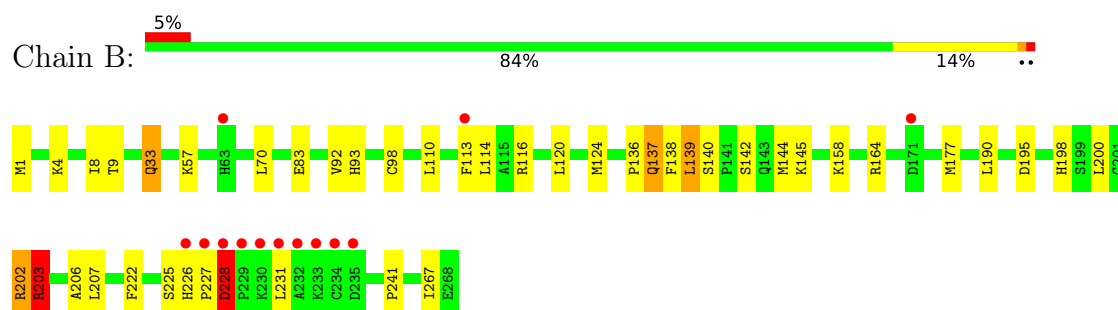
3 Residue-property plots [i](#)

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.

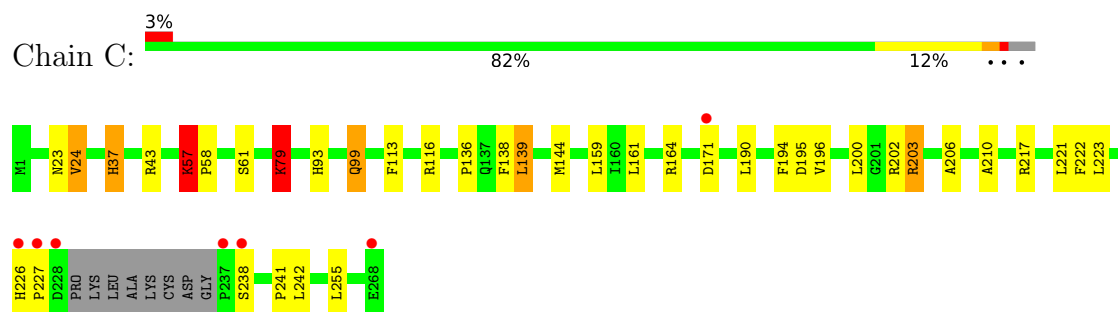
- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase



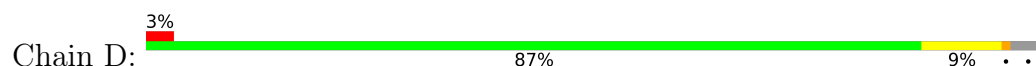
- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase

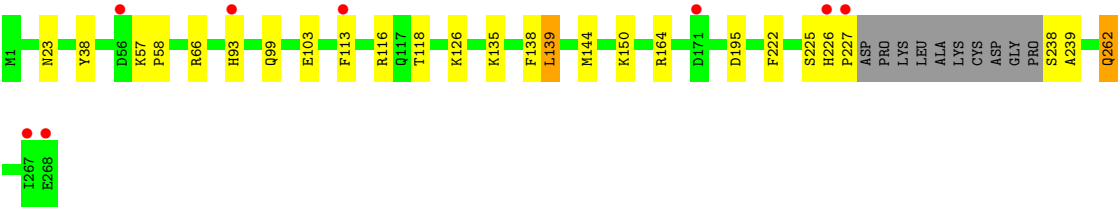


- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase



- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.69Å 104.05Å 149.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.99 – 2.20 35.99 – 2.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (35.99-2.20) 100.0 (35.99-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.77 (at 2.20Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.197 , 0.238 (Not available) , 0.215	Depositor DCC
R_{free} test set	3250 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 36.2	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	0.94	EDS
Total number of atoms	8729	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.13% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CL

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	1/2061 (0.0%)	0.79	1/2789 (0.0%)
1	B	0.87	7/2127 (0.3%)	0.94	9/2884 (0.3%)
1	C	0.77	7/2071 (0.3%)	0.78	0/2802
1	D	0.68	2/2053 (0.1%)	0.77	1/2777 (0.0%)
All	All	0.76	17/8312 (0.2%)	0.82	11/11252 (0.1%)

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	99	GLN	C-O	-8.70	1.16	1.24
1	D	23	ASN	C-O	-7.96	1.14	1.24
1	B	140	SER	C-O	-6.97	1.15	1.23
1	B	137	GLN	C-O	-6.68	1.15	1.24
1	B	225	SER	C-O	-6.27	1.15	1.23
1	C	203	ARG	C-O	-5.97	1.16	1.24
1	B	139	LEU	C-O	-5.75	1.17	1.24
1	D	262	GLN	C-O	-5.60	1.16	1.24
1	B	138	PHE	C-O	-5.55	1.16	1.23
1	B	93	HIS	CA-CB	-5.54	1.46	1.54
1	B	198	HIS	C-O	-5.36	1.16	1.24
1	A	72	GLU	C-O	-5.33	1.17	1.24
1	C	79	LYS	C-O	-5.27	1.18	1.24
1	C	57	LYS	C-O	-5.25	1.18	1.24
1	C	37[A]	HIS	C-O	-5.13	1.18	1.24
1	C	37[B]	HIS	C-O	-5.13	1.18	1.24
1	C	61	SER	C-O	-5.04	1.16	1.23

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	231	LEU	N-CA-C	11.61	123.67	111.14
1	B	203	ARG	N-CA-C	6.58	121.31	113.28
1	B	228[A]	ASP	CA-C-N	6.42	126.76	119.83
1	B	228[A]	ASP	C-N-CA	6.42	126.76	119.83
1	B	228[B]	ASP	CA-C-N	6.42	126.76	119.83
1	B	228[B]	ASP	C-N-CA	6.42	126.76	119.83
1	B	228[A]	ASP	N-CA-C	-6.36	96.82	109.60
1	B	228[B]	ASP	N-CA-C	-6.36	96.82	109.60
1	B	8	ILE	N-CA-C	5.73	116.97	109.58
1	A	223	LEU	N-CA-C	5.40	117.28	109.07
1	D	239	ALA	N-CA-C	5.09	117.67	110.50

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	2020	0	2073	17	0
1	B	2080	0	2120	30	0
1	C	2030	0	2087	26	0
1	D	2013	0	2076	17	0
2	A	3	0	0	1	0
2	B	4	0	0	2	0
2	C	2	0	0	0	0
2	D	3	0	0	1	0
3	A	150	0	0	1	0
3	B	138	0	0	3	0
3	C	135	0	0	2	0
3	D	151	0	0	3	0
All	All	8729	0	8356	82	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 5.

All (82) 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:92[A]:VAL:HG23	1:B:124:MET:HE2	1.42	0.99
1:B:92[A]:VAL:CG2	1:B:124:MET:HE2	2.02	0.88
1:A:68:VAL:HB	1:A:72:GLU:HG2	1.62	0.81
1:B:120:LEU:HG	1:B:124:MET:HE3	1.69	0.73
1:D:150:LYS:NZ	3:D:613:HOH:O	2.22	0.68
1:C:202:ARG:NH2	3:C:547:HOH:O	2.27	0.66
1:B:70:LEU:C	1:B:70:LEU:HD13	2.20	0.66
1:D:38:TYR:OH	1:D:225:SER:HB2	2.00	0.61
1:B:33:GLN:NE2	1:B:228[B]:ASP:OD2	2.33	0.61
1:C:194:PHE:CE2	1:C:196[B]:VAL:HG12	2.34	0.61
1:B:4:LYS:HE2	1:B:9:THR:OG1	2.01	0.61
1:B:139:LEU:HD23	1:B:144:MET:HE2	1.83	0.61
1:C:226:HIS:CB	1:C:227:PRO:HD2	2.31	0.60
1:C:255[B]:LEU:C	1:C:255[B]:LEU:HD23	2.25	0.60
1:B:110:LEU:HA	1:B:124:MET:HE1	1.84	0.60
1:A:177:MET:HG3	1:B:177:MET:HE3	1.84	0.59
1:A:177:MET:HE3	1:B:177:MET:HG3	1.87	0.57
1:A:196[B]:VAL:HG11	1:A:210:ALA:HB2	1.87	0.56
1:B:92[B]:VAL:HG11	1:B:98:CYS:SG	2.47	0.55
1:A:195:ASP:HA	1:A:222:PHE:HB3	1.90	0.54
1:C:136:PRO:HG2	1:C:139:LEU:HD22	1.88	0.54
1:B:57:LYS:NZ	3:B:348:HOH:O	2.40	0.53
1:A:70:LEU:C	1:A:70:LEU:HD23	2.36	0.51
1:C:138:PHE:CD2	1:C:139:LEU:HD13	2.46	0.51
1:A:117:GLN:NE2	1:C:113:PHE:CD1	2.79	0.51
1:A:59:ARG:HB2	2:A:271:CL:CL	2.49	0.50
1:C:138:PHE:CE2	1:C:139:LEU:HD13	2.47	0.49
1:A:207:LEU:HB2	1:B:267:ILE:HD11	1.94	0.49
1:C:23:ASN:OD1	1:C:57:LYS:NZ	2.37	0.49
1:B:116:ARG:HG3	1:D:113:PHE:CE2	2.48	0.48
1:B:113[B]:PHE:CD1	1:B:114:LEU:HG	2.48	0.48
1:B:113[A]:PHE:HZ	1:D:116:ARG:HH11	1.60	0.48
1:C:79:LYS:HA	1:C:79:LYS:HD2	1.64	0.47
1:D:139:LEU:HD23	1:D:144[B]:MET:HE2	1.96	0.47
1:C:139:LEU:HD23	1:C:144[B]:MET:CE	2.45	0.46
1:B:57:LYS:HG3	2:B:271:CL:CL	2.53	0.46
1:D:118:THR:OG1	1:D:150:LYS:HE2	2.16	0.46
1:D:116:ARG:HA	3:D:377:HOH:O	2.15	0.46
1:C:203:ARG:HD3	1:C:241:PRO:HD3	1.97	0.45
1:C:58:PRO:HG3	1:C:93:HIS:CD2	2.51	0.45
1:B:1:MET:HE2	1:B:190:LEU:HD23	1.99	0.45
1:C:37[B]:HIS:HD2	1:C:242:LEU:HG	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:ASP:HA	1:B:222:PHE:HB3	1.97	0.45
1:D:226:HIS:CD2	3:D:274:HOH:O	2.70	0.45
1:C:200:LEU:HD11	1:C:206:ALA:HA	1.99	0.45
3:A:329:HOH:O	1:C:171:ASP:CB	2.64	0.44
1:D:195:ASP:HA	1:D:222:PHE:HB3	1.98	0.44
1:B:136:PRO:HD2	1:B:144:MET:HE1	2.00	0.44
1:A:68:VAL:HB	1:A:72:GLU:CG	2.42	0.44
1:A:265:LEU:HD23	1:B:207:LEU:HD22	2.00	0.44
1:C:195:ASP:HA	1:C:222:PHE:HB3	2.00	0.44
1:D:139:LEU:HD23	1:D:144[B]:MET:CE	2.48	0.43
1:A:200:LEU:HD11	1:A:206:ALA:HA	1.99	0.43
1:A:99:GLN:N	1:A:100:PRO:HD2	2.34	0.43
1:D:138:PHE:CE2	1:D:139:LEU:HD13	2.54	0.43
1:B:137:GLN:HG2	2:B:272:CL:CL	2.55	0.43
1:B:226:HIS:HD2	1:B:241:PRO:HA	1.82	0.43
1:C:226:HIS:CB	1:C:227:PRO:CD	2.97	0.43
1:B:70:LEU:HD13	1:B:70:LEU:O	2.18	0.42
1:D:57:LYS:HE3	2:D:271:CL:CL	2.56	0.42
1:D:58:PRO:HA	1:D:66:ARG:HH21	1.85	0.42
1:B:203:ARG:H	1:B:203:ARG:HG3	1.59	0.42
1:B:70:LEU:C	1:B:70:LEU:CD1	2.90	0.42
1:B:142:SER:O	1:B:145:LYS:HG2	2.20	0.42
1:C:159:LEU:HB2	1:C:190:LEU:HD13	2.00	0.42
1:A:66:ARG:HH21	1:A:66:ARG:HD3	1.59	0.42
1:A:135:LYS:O	1:A:164:ARG:HD2	2.20	0.41
1:D:99:GLN:O	1:D:103:GLU:HG2	2.20	0.41
1:A:66:ARG:HE	1:A:66:ARG:HB3	1.74	0.41
1:A:59:ARG:HG2	1:C:116:ARG:HB2	2.02	0.41
1:D:226:HIS:HA	1:D:227:PRO:HD3	1.93	0.41
1:C:202:ARG:NH2	3:C:548:HOH:O	2.54	0.41
1:B:200:LEU:HD11	1:B:206:ALA:HA	2.03	0.41
1:C:24:VAL:HG21	1:C:57:LYS:HG3	2.03	0.41
1:C:144[A]:MET:HE3	1:C:161:LEU:HD22	2.03	0.41
1:B:57:LYS:HE2	3:B:600:HOH:O	2.20	0.40
1:B:202:ARG:O	3:B:601:HOH:O	2.22	0.40
1:C:37[B]:HIS:CD2	1:C:242:LEU:HG	2.57	0.40
1:C:196[B]:VAL:HG11	1:C:210:ALA:HA	2.03	0.40
1:D:58:PRO:HG2	1:D:93:HIS:CD2	2.57	0.40
1:D:135:LYS:HG3	1:D:144[B]:MET:HE3	2.03	0.40
1:C:196[B]:VAL:CG1	1:C:221:LEU:HD11	2.52	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	258/268 (96%)	253 (98%)	4 (2%)	1 (0%)	30	34
1	B	270/268 (101%)	265 (98%)	4 (2%)	1 (0%)	30	34
1	C	260/268 (97%)	254 (98%)	5 (2%)	1 (0%)	30	34
1	D	257/268 (96%)	253 (98%)	4 (2%)	0	100	100
All	All	1045/1072 (98%)	1025 (98%)	17 (2%)	3 (0%)	36	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	227	PRO
1	A	217	ARG
1	C	217	ARG

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	226/230 (98%)	220 (97%)	6 (3%)	39	53
1	B	231/230 (100%)	223 (96%)	8 (4%)	32	43
1	C	226/230 (98%)	217 (96%)	9 (4%)	28	38
1	D	225/230 (98%)	220 (98%)	5 (2%)	45	61
All	All	908/920 (99%)	880 (97%)	28 (3%)	36	48

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LYS
1	A	66	ARG
1	A	72	GLU
1	A	164	ARG
1	A	266	THR
1	A	268	GLU
1	B	33	GLN
1	B	83	GLU
1	B	158	LYS
1	B	164	ARG
1	B	202	ARG
1	B	203	ARG
1	B	228[A]	ASP
1	B	228[B]	ASP
1	C	24	VAL
1	C	43	ARG
1	C	57	LYS
1	C	79	LYS
1	C	99	GLN
1	C	139	LEU
1	C	164	ARG
1	C	223	LEU
1	C	238	SER
1	D	126	LYS
1	D	139	LEU
1	D	164	ARG
1	D	238	SER
1	D	262	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	63	HIS
1	A	262	GLN
1	B	226	HIS
1	C	93	HIS
1	C	96	HIS
1	C	243	HIS
1	D	146	ASN
1	D	226	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 12 ligands modelled in this entry, 12 are 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 [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	260/268 (97%)	-0.14	8 (3%)	51	48	9, 21, 38, 60	2 (0%)
1	B	268/268 (100%)	0.14	13 (4%)	35	31	9, 24, 45, 57	4 (1%)
1	C	260/268 (97%)	0.01	7 (2%)	56	53	9, 24, 45, 84	4 (1%)
1	D	258/268 (96%)	-0.18	8 (3%)	51	48	9, 20, 37, 76	3 (1%)
All	All	1046/1072 (97%)	-0.04	36 (3%)	48	45	9, 22, 42, 84	13 (1%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	232	ALA	12.4
1	B	229	PRO	7.0
1	B	233	LYS	6.4
1	C	237	PRO	6.3
1	C	227	PRO	6.0
1	B	234	CYS	4.9
1	B	113[A]	PHE	4.8
1	C	226	HIS	4.5
1	C	171	ASP	4.4
1	B	228[A]	ASP	4.3
1	B	171	ASP	4.3
1	A	171	ASP	4.1
1	B	227	PRO	4.0
1	C	228	ASP	3.9
1	A	237	PRO	3.4
1	D	226	HIS	3.4
1	B	231	LEU	3.3
1	D	268	GLU	3.0
1	C	238	SER	2.9
1	C	268	GLU	2.9
1	D	93	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	113	PHE	2.8
1	D	171	ASP	2.7
1	D	113	PHE	2.7
1	D	56	ASP	2.6
1	A	201	GLY	2.5
1	D	227	PRO	2.5
1	B	230	LYS	2.3
1	A	94	GLU	2.2
1	A	228	ASP	2.2
1	A	29[A]	ASP	2.1
1	B	226	HIS	2.1
1	D	267	ILE	2.1
1	A	225	SER	2.1
1	B	235	ASP	2.1
1	B	63	HIS	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 [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
2	CL	C	270	1/1	0.93	0.15	35,35,35,35	0
2	CL	B	271	1/1	0.96	0.10	31,31,31,31	0
2	CL	A	271	1/1	0.96	0.12	33,33,33,33	0
2	CL	D	271	1/1	0.96	0.09	35,35,35,35	0
2	CL	A	270	1/1	0.97	0.05	27,27,27,27	0
2	CL	A	269	1/1	0.97	0.09	22,22,22,22	0
2	CL	B	270	1/1	0.97	0.05	22,22,22,22	0
2	CL	B	269	1/1	0.98	0.14	28,28,28,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CL	B	272	1/1	0.98	0.05	30,30,30,30	0
2	CL	D	269	1/1	0.99	0.06	24,24,24,24	0
2	CL	D	270	1/1	0.99	0.05	26,26,26,26	0
2	CL	C	269	1/1	0.99	0.06	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.