



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 06:05 am BST

PDB ID : 3QQ1
Title : Crystal structure of a double mutant [A58P, DEL(N59)] 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 : 2011-02-14
Resolution : 2.70 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

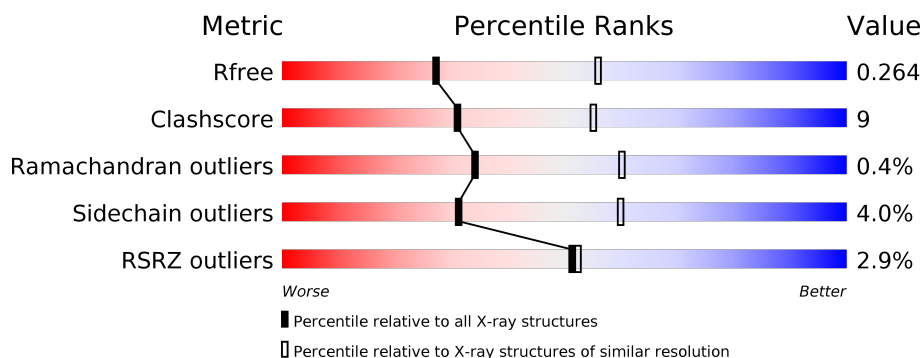
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 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	279	<div> <div>3%</div> <div> <div></div> <div>75%</div> <div>12%</div> <div>•</div> <div>11%</div> </div> </div>
1	B	279	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>13%</div> <div>•</div> <div>10%</div> </div> </div>
1	C	279	<div> <div>%</div> <div> <div></div> <div>76%</div> <div>13%</div> <div></div> <div>11%</div> </div> </div>
1	D	279	<div> <div>4%</div> <div> <div></div> <div>69%</div> <div>18%</div> <div>•</div> <div>10%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7850 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	249	Total	C	N	O	S	0	0	0
			1889	1219	311	348	11			
1	B	251	Total	C	N	O	S	0	5	0
			1940	1253	322	354	11			
1	C	249	Total	C	N	O	S	0	2	0
			1939	1250	325	353	11			
1	D	251	Total	C	N	O	S	0	2	0
			1943	1250	325	357	11			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	ALA	DELETION	UNP Q9JZ55
A	58	PRO	ASN	ENGINEERED MUTATION	UNP Q9JZ55
B	?	-	ALA	DELETION	UNP Q9JZ55
B	58	PRO	ASN	ENGINEERED MUTATION	UNP Q9JZ55
C	?	-	ALA	DELETION	UNP Q9JZ55
C	58	PRO	ASN	ENGINEERED MUTATION	UNP Q9JZ55
D	?	-	ALA	DELETION	UNP Q9JZ55
D	58	PRO	ASN	ENGINEERED MUTATION	UNP Q9JZ55

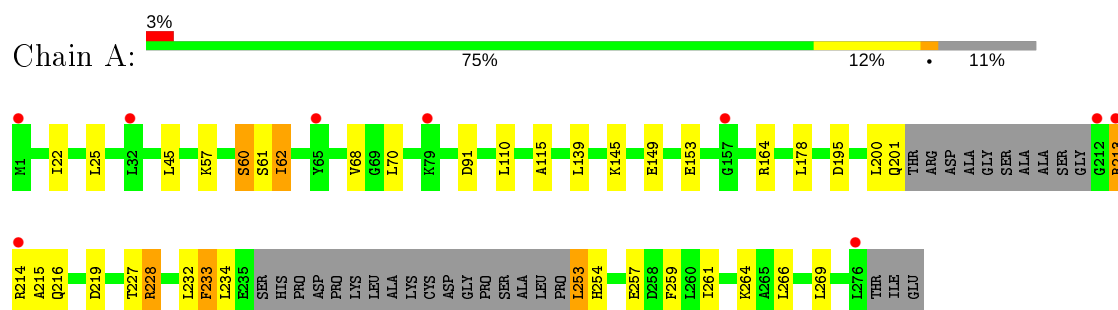
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	30	Total	O	0	0
			30	30		
2	B	38	Total	O	0	0
			38	38		
2	C	37	Total	O	0	0
			37	37		
2	D	34	Total	O	0	0
			34	34		

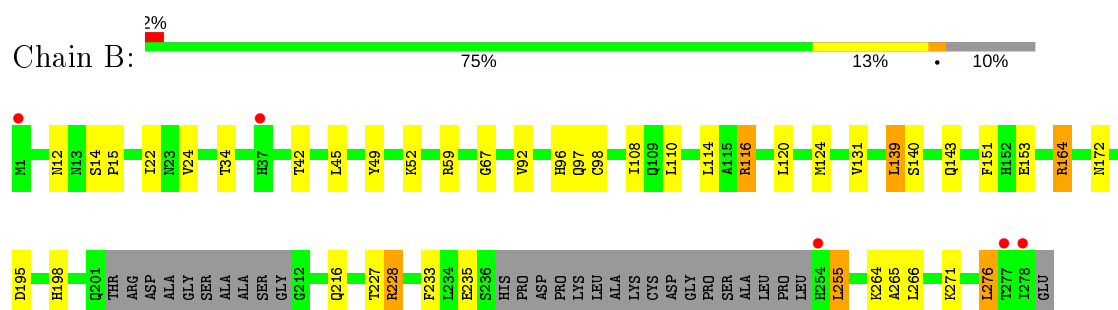
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

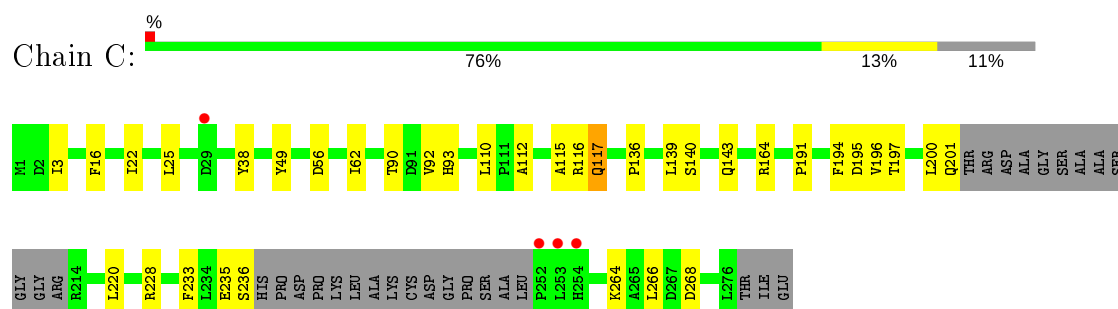
- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase



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

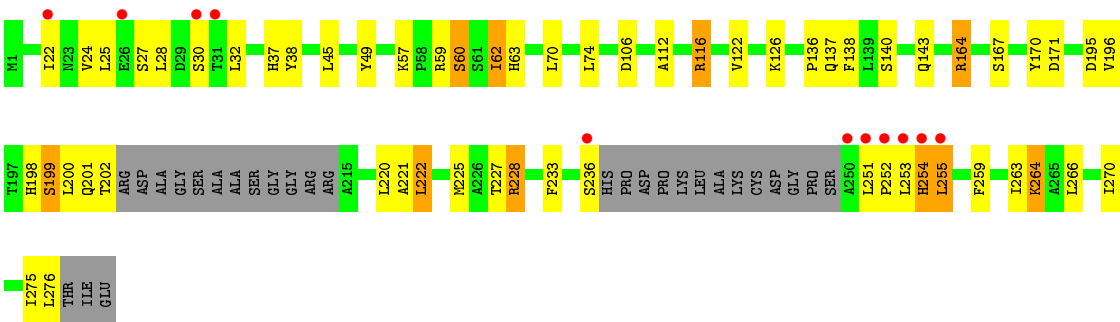


- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase



- 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	81.82Å 85.15Å 163.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.89 – 2.70 47.87 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.89-2.70) 99.6 (47.87-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.97 (at 2.69Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.211 , 0.275 0.207 , 0.264	Depositor DCC
R_{free} test set	1562 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	36.8	Xtriage
Anisotropy	0.295	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 31.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.025 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7850	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% 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	0.78	0/1922	0.67	2/2602 (0.1%)
1	B	0.66	0/1985	0.63	0/2688
1	C	0.60	1/1976 (0.1%)	0.57	0/2670
1	D	0.79	1/1980 (0.1%)	0.67	0/2678
All	All	0.71	2/7863 (0.0%)	0.64	2/10638 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	92	VAL	CB-CG2	-5.17	1.42	1.52
1	D	138	PHE	CD2-CE2	-5.01	1.29	1.39

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	233	PHE	CB-CA-C	-5.67	99.06	110.40
1	A	234	LEU	CA-CB-CG	5.05	126.91	115.30

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	1889	0	1912	34	0
1	B	1940	0	1988	36	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1939	0	2001	22	0
1	D	1943	0	1989	57	0
2	A	30	0	0	0	0
2	B	38	0	0	1	0
2	C	37	0	0	0	0
2	D	34	0	0	0	0
All	All	7850	0	7890	142	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 9.

All (142) 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:A:70:LEU:O	1:A:70:LEU:HD23	1.34	1.24
1:A:70:LEU:C	1:A:70:LEU:HD23	1.54	1.23
1:D:70:LEU:HD12	1:D:70:LEU:O	1.50	1.12
1:D:251:LEU:HD12	1:D:252:PRO:HD2	1.25	1.11
1:D:62:ILE:HG23	1:D:63:HIS:CD2	1.86	1.10
1:D:253:LEU:O	1:D:254:HIS:HB3	1.33	1.07
1:D:62:ILE:CG2	1:D:63:HIS:CD2	2.43	1.02
1:C:200:LEU:O	1:C:201:GLN:HB2	1.60	1.00
1:D:195:ASP:OD2	1:D:198:HIS:ND1	1.94	1.00
1:A:70:LEU:C	1:A:70:LEU:CD2	2.30	0.99
1:A:22:ILE:HD11	1:A:25:LEU:HD23	1.56	0.87
1:D:253:LEU:O	1:D:254:HIS:CB	2.18	0.87
1:D:252:PRO:HB2	1:D:255:LEU:CD1	2.03	0.87
1:D:251:LEU:HD12	1:D:252:PRO:CD	2.06	0.86
1:D:252:PRO:HB2	1:D:255:LEU:HD13	1.55	0.86
1:B:195:ASP:HA	1:B:233:PHE:HB3	1.58	0.84
1:D:62:ILE:CG2	1:D:63:HIS:HD2	1.90	0.82
1:D:70:LEU:HD12	1:D:70:LEU:C	1.94	0.81
1:D:195:ASP:O	1:D:199:SER:OG	1.99	0.79
1:A:253:LEU:HB3	1:A:254:HIS:HD2	1.47	0.78
1:D:57:LYS:O	1:D:60:SER:OG	2.00	0.78
1:C:136:PRO:HG2	1:C:139:LEU:HB2	1.64	0.77
1:D:70:LEU:CD1	1:D:70:LEU:O	2.30	0.75
1:A:195:ASP:HA	1:A:233:PHE:HB2	1.71	0.73
1:A:253:LEU:HB3	1:A:254:HIS:CD2	2.23	0.73
1:C:200:LEU:O	1:C:201:GLN:CB	2.37	0.72
1:A:45:LEU:O	1:A:264:LYS:HE3	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:LEU:HD13	1:B:124:MET:HE3	1.73	0.70
1:D:195:ASP:HA	1:D:233:PHE:HB3	1.74	0.69
1:D:236:SER:HB2	1:D:251:LEU:O	1.95	0.67
1:A:227:THR:O	1:A:228:ARG:HB3	1.94	0.67
1:D:222:LEU:HD13	1:D:263:ILE:HD13	1.79	0.65
1:C:195:ASP:HA	1:C:233:PHE:HB3	1.80	0.64
1:B:110:LEU:HD13	1:B:124:MET:CE	2.28	0.64
1:C:264:LYS:NZ	1:C:268:ASP:OD2	2.29	0.63
1:A:213:ARG:CG	1:A:214:ARG:N	2.61	0.63
1:A:227:THR:O	1:A:228:ARG:CB	2.48	0.62
1:D:252:PRO:HB2	1:D:255:LEU:HD12	1.80	0.62
1:A:214:ARG:O	1:A:259:PHE:HZ	1.81	0.62
1:D:254:HIS:O	1:D:254:HIS:CD2	2.52	0.61
1:D:255:LEU:O	1:D:259:PHE:HB2	2.02	0.60
1:D:254:HIS:HD2	1:D:254:HIS:O	1.83	0.60
1:D:222:LEU:CD1	1:D:263:ILE:HD13	2.32	0.59
1:D:37:HIS:NE2	1:D:253:LEU:CB	2.66	0.59
1:D:62:ILE:HG23	1:D:63:HIS:CG	2.35	0.59
1:D:140:SER:OG	1:D:143[A]:GLN:HG2	2.03	0.59
1:B:45:LEU:O	1:B:264:LYS:HE3	2.03	0.58
1:D:62:ILE:HG21	1:D:63:HIS:CD2	2.37	0.58
1:B:164:ARG:HG3	1:B:164:ARG:HH11	1.67	0.58
1:B:139:LEU:HG	1:B:143:GLN:HG3	1.86	0.58
1:D:137:GLN:HG3	1:D:167:SER:HB2	1.85	0.58
1:B:108:ILE:HD12	1:B:124:MET:HG2	1.87	0.56
1:D:252:PRO:CB	1:D:255:LEU:HD13	2.33	0.56
1:D:62:ILE:HG21	1:D:63:HIS:HD2	1.70	0.56
1:D:45:LEU:O	1:D:264:LYS:HE2	2.06	0.56
1:C:266[A]:LEU:HD13	1:D:266:LEU:HD13	1.89	0.55
1:D:62:ILE:CG2	1:D:63:HIS:N	2.70	0.55
1:B:120:LEU:HG	1:B:124:MET:CE	2.36	0.55
1:B:227:THR:O	1:B:228:ARG:HB3	2.07	0.54
1:A:57:LYS:O	1:A:60:SER:OG	2.26	0.53
1:C:110:LEU:HD21	1:C:115:ALA:HA	1.90	0.53
1:C:3:ILE:HD12	1:C:191:PRO:HG2	1.92	0.52
1:D:254:HIS:C	1:D:254:HIS:CD2	2.83	0.52
1:A:25:LEU:O	1:A:68:VAL:HG13	2.09	0.52
1:B:52:LYS:HE2	1:B:233:PHE:CZ	2.44	0.52
1:C:112:ALA:HB1	1:C:136:PRO:HA	1.91	0.52
1:A:22:ILE:HD11	1:A:25:LEU:CD2	2.34	0.51
1:B:96:HIS:HB3	2:B:292:HOH:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:114:LEU:HD13	1:B:120:LEU:CD2	2.40	0.51
1:B:164:ARG:HG3	1:B:164:ARG:NH1	2.23	0.51
1:A:214:ARG:O	1:A:259:PHE:CZ	2.63	0.50
1:A:213:ARG:C	1:A:215:ALA:N	2.62	0.50
1:D:227:THR:O	1:D:228:ARG:CB	2.59	0.50
1:D:22:ILE:HD12	1:D:24:VAL:O	2.11	0.50
1:C:56:ASP:OD1	1:C:56:ASP:O	2.30	0.50
1:D:266:LEU:HD12	1:D:270:ILE:HG12	1.94	0.50
1:B:92[B]:VAL:HG11	1:B:98:CYS:SG	2.52	0.49
1:B:120:LEU:HG	1:B:124:MET:HE1	1.94	0.49
1:A:110:LEU:HD21	1:A:115:ALA:HA	1.94	0.49
1:D:164:ARG:O	1:D:164:ARG:HD3	2.13	0.49
1:D:201:GLN:O	1:D:202:THR:OG1	2.29	0.48
1:B:114:LEU:HD13	1:B:120:LEU:HD22	1.95	0.48
1:C:56:ASP:C	1:C:56:ASP:OD1	2.51	0.48
1:D:227:THR:O	1:D:228:ARG:HB3	2.13	0.48
1:B:227:THR:O	1:B:228:ARG:CB	2.62	0.48
1:A:153:GLU:HG3	1:C:62:ILE:HD11	1.96	0.48
1:D:170:TYR:O	1:D:171:ASP:OD2	2.32	0.47
1:A:213:ARG:O	1:A:216:GLN:HB2	2.13	0.47
1:C:194:PHE:CE2	1:C:196[B]:VAL:HG12	2.50	0.47
1:D:275:ILE:O	1:D:276:LEU:CB	2.62	0.47
1:B:12:ASN:O	1:B:271:LYS:HE2	2.15	0.46
1:D:38:TYR:HB3	1:D:49:TYR:CZ	2.50	0.46
1:A:200:LEU:O	1:A:201:GLN:CB	2.62	0.46
1:B:92[B]:VAL:HG13	1:B:97:GLN:HB2	1.97	0.46
1:D:116:ARG:HG3	1:D:116:ARG:O	2.16	0.46
1:A:22:ILE:CD1	1:A:25:LEU:HD23	2.36	0.45
1:C:136:PRO:HG2	1:C:139:LEU:CB	2.40	0.45
1:C:140:SER:OG	1:C:143:GLN:HG3	2.16	0.45
1:D:253:LEU:C	1:D:255:LEU:H	2.19	0.45
1:D:32:LEU:HA	1:D:32:LEU:HD23	1.65	0.45
1:C:266[B]:LEU:C	1:C:266[B]:LEU:HD23	2.36	0.45
1:C:117:GLN:HB3	1:C:117:GLN:HE21	1.51	0.45
1:A:213:ARG:HG3	1:A:214:ARG:N	2.27	0.45
1:A:178:LEU:HD22	1:B:216:GLN:HG2	1.99	0.45
1:D:70:LEU:HD11	1:D:74:LEU:CD1	2.47	0.45
1:D:28:LEU:O	1:D:28:LEU:HG	2.17	0.44
1:B:164:ARG:HH11	1:B:164:ARG:CG	2.28	0.44
1:B:276:LEU:HD23	1:B:276:LEU:HA	1.72	0.44
1:C:197:THR:HG21	1:C:235:GLU:HB2	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:116:ARG:CG	1:D:116:ARG:O	2.65	0.44
1:D:200:LEU:O	1:D:201:GLN:C	2.53	0.44
1:A:213:ARG:O	1:A:216:GLN:N	2.42	0.44
1:A:145:LYS:O	1:A:149:GLU:HB2	2.17	0.44
1:D:112:ALA:HB1	1:D:136:PRO:HA	2.00	0.44
1:C:22:ILE:HD11	1:C:25:LEU:HD23	2.00	0.44
1:D:221:ALA:O	1:D:225:MET:HG2	2.18	0.44
1:A:91:ASP:N	1:A:91:ASP:OD1	2.50	0.43
1:B:255:LEU:H	1:B:255:LEU:HD23	1.83	0.43
1:B:195:ASP:OD2	1:B:198:HIS:ND1	2.36	0.43
1:A:269:LEU:HD11	1:B:265:ALA:HB1	1.99	0.43
1:A:213:ARG:C	1:A:215:ALA:H	2.20	0.43
1:A:257:GLU:O	1:A:261:ILE:HG13	2.19	0.43
1:D:222:LEU:HD13	1:D:263:ILE:CD1	2.49	0.43
1:B:24[B]:VAL:HG13	1:B:67:GLY:HA2	2.01	0.42
1:B:42:THR:HG21	1:B:49:TYR:HB2	1.99	0.42
1:C:93:HIS:CD2	1:C:93:HIS:N	2.87	0.42
1:D:22:ILE:HD11	1:D:25:LEU:HD23	2.01	0.42
1:C:38:TYR:HB3	1:C:49:TYR:CZ	2.54	0.42
1:B:22:ILE:HG23	1:B:34:THR:HG21	2.01	0.42
1:B:140:SER:O	1:B:143:GLN:HG2	2.19	0.41
1:B:24[A]:VAL:HG21	1:B:59:ARG:HH21	1.85	0.41
1:A:266:LEU:HD13	1:B:266[B]:LEU:HD21	2.02	0.41
1:D:196:VAL:HG23	1:D:220:LEU:HD23	2.03	0.41
1:C:196[A]:VAL:HG23	1:C:220:LEU:HD23	2.02	0.41
1:B:116:ARG:HB3	1:D:59:ARG:HD2	2.02	0.41
1:B:235:GLU:HG3	1:B:235:GLU:O	2.20	0.41
1:A:266:LEU:CD1	1:B:266[B]:LEU:HD21	2.50	0.41
1:D:122:VAL:O	1:D:126:LYS:HG2	2.21	0.41
1:A:213:ARG:HG2	1:A:214:ARG:N	2.34	0.40
1:A:62:ILE:HD13	1:A:62:ILE:HG21	1.63	0.40
1:B:131:VAL:HG11	1:B:151:PHE:CZ	2.56	0.40
1:B:14:SER:O	1:B:15:PRO:C	2.58	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	243/279 (87%)	237 (98%)	5 (2%)	1 (0%)	34	60
1	B	250/279 (90%)	245 (98%)	4 (2%)	1 (0%)	34	60
1	C	245/279 (88%)	239 (98%)	5 (2%)	1 (0%)	34	60
1	D	247/279 (88%)	241 (98%)	5 (2%)	1 (0%)	34	60
All	All	985/1116 (88%)	962 (98%)	19 (2%)	4 (0%)	34	60

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	228	ARG
1	D	228	ARG
1	B	228	ARG
1	C	228	ARG

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	206/236 (87%)	197 (96%)	9 (4%)	28	56
1	B	215/236 (91%)	208 (97%)	7 (3%)	38	67
1	C	217/236 (92%)	211 (97%)	6 (3%)	43	73
1	D	216/236 (92%)	204 (94%)	12 (6%)	21	45
All	All	854/944 (90%)	820 (96%)	34 (4%)	31	60

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	SER
1	A	61	SER
1	A	62	ILE
1	A	139	LEU
1	A	164	ARG
1	A	213	ARG
1	A	219	ASP
1	A	232	LEU
1	A	253	LEU
1	B	116	ARG
1	B	139	LEU
1	B	153	GLU
1	B	164	ARG
1	B	172	ASN
1	B	255	LEU
1	B	276	LEU
1	C	16	PHE
1	C	90	THR
1	C	116	ARG
1	C	117	GLN
1	C	164	ARG
1	C	236	SER
1	D	27	SER
1	D	30	SER
1	D	60	SER
1	D	62	ILE
1	D	106	ASP
1	D	116	ARG
1	D	164	ARG
1	D	199	SER
1	D	222	LEU
1	D	254	HIS
1	D	255	LEU
1	D	264	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	254	HIS
1	B	93	HIS

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Mol	Chain	Res	Type
1	B	117	GLN
1	B	216	GLN
1	B	273	GLN
1	C	37	HIS
1	C	63	HIS
1	C	93	HIS
1	C	99	GLN
1	C	117	GLN
1	D	63	HIS
1	D	254	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 carbohydrates 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 ⓘ

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(Å²)	Q<0.9	
1	A	249/279 (89%)	0.07	9 (3%)	42	42	15, 32, 60, 82	1 (0%)
1	B	251/279 (89%)	0.01	5 (1%)	65	67	16, 31, 52, 63	0
1	C	249/279 (89%)	-0.07	4 (1%)	72	74	17, 32, 54, 69	0
1	D	251/279 (89%)	0.13	11 (4%)	34	33	17, 33, 64, 83	0
All	All	1000/1116 (89%)	0.04	29 (2%)	51	52	15, 32, 58, 83	1 (0%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	252	PRO	7.8
1	D	251	LEU	7.0
1	B	277	THR	5.3
1	A	276	LEU	4.7
1	D	254	HIS	4.2
1	A	214	ARG	4.1
1	D	253	LEU	4.1
1	C	253	LEU	3.8
1	D	22	ILE	3.7
1	A	213	ARG	3.6
1	B	1	MET	3.4
1	B	254	HIS	3.3
1	A	65	TYR	3.1
1	D	250	ALA	2.9
1	D	236	SER	2.9
1	A	212	GLY	2.8
1	A	32	LEU	2.7
1	D	255	LEU	2.5
1	D	26	GLU	2.5
1	B	278	ILE	2.5
1	C	252	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	31	THR	2.3
1	A	79	LYS	2.2
1	A	157	GLY	2.2
1	B	37	HIS	2.1
1	C	254	HIS	2.1
1	C	29	ASP	2.1
1	D	30	SER	2.1
1	A	1	MET	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 carbohydrates 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.