



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:00 AM UTC

PDB ID : 3FYO / pdb_00003fyo
Title : Crystal structure of the triple mutant (N23C/D247E/P249A) of 3-deoxy-D-manno-octulosonate 8-phosphate synthase (KDO8PS) from *Neisseria meningitidis*
Authors : Jameson, G.B.; Parker, E.J.; Cochrane, F.P.; Patchett, M.L.
Deposited on : 2009-01-22
Resolution : 1.90 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

<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
Mogul : 2022.3.0, CSD as543be (2022)
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

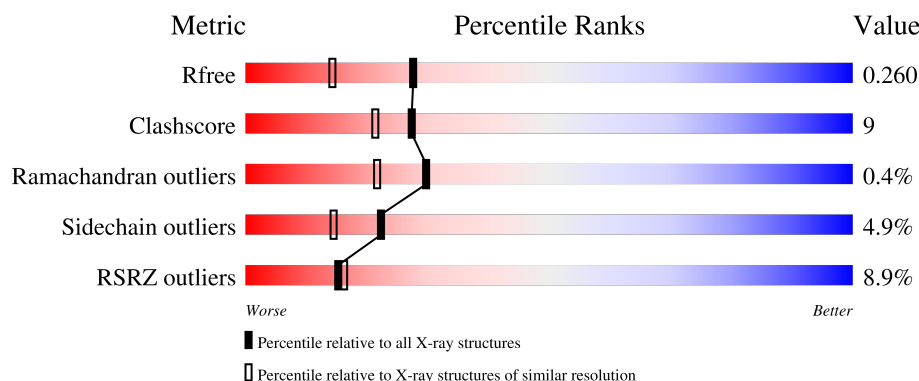
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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 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	280	12% 73% 16% • 9%
1	B	280	9% 79% 13% • 6%
1	C	280	4% 79% 12% • 9%
1	D	280	8% 72% 16% • 7%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8610 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-D-manno-octulosonic acid 8-phosphate synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	0	10	0
			1995	1287	335	361	12			
1	B	262	Total	C	N	O	S	0	11	0
			2082	1337	352	381	12			
1	C	256	Total	C	N	O	S	0	5	0
			2002	1290	336	364	12			
1	D	261	Total	C	N	O	S	0	12	0
			2064	1327	348	374	15			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	CYS	ASN	engineered mutation	UNP Q9JZ55
A	247	GLU	ASP	engineered mutation	UNP Q9JZ55
A	249	ALA	PRO	engineered mutation	UNP Q9JZ55
B	23	CYS	ASN	engineered mutation	UNP Q9JZ55
B	247	GLU	ASP	engineered mutation	UNP Q9JZ55
B	249	ALA	PRO	engineered mutation	UNP Q9JZ55
C	23	CYS	ASN	engineered mutation	UNP Q9JZ55
C	247	GLU	ASP	engineered mutation	UNP Q9JZ55
C	249	ALA	PRO	engineered mutation	UNP Q9JZ55
D	23	CYS	ASN	engineered mutation	UNP Q9JZ55
D	247	GLU	ASP	engineered mutation	UNP Q9JZ55
D	249	ALA	PRO	engineered mutation	UNP Q9JZ55

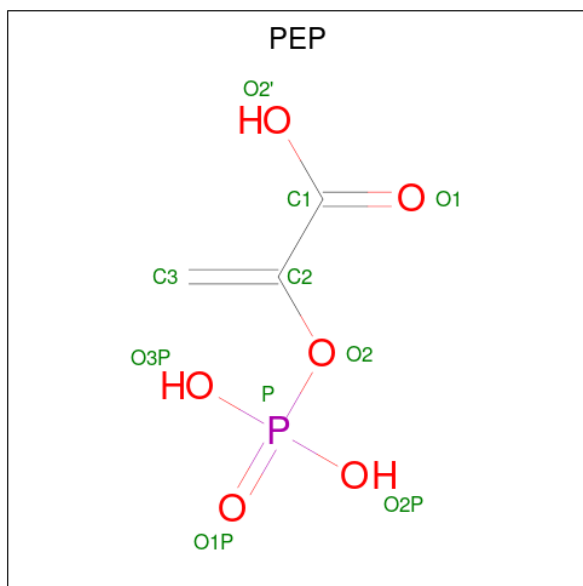
- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		
2	C	1	Total	Cl	0	0
			1	1		

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	1	Total	Na	0	0
			1	1		

- Molecule 4 is PHOSPHOENOLPYRUVATE (CCD ID: PEP) (formula: C₃H₅O₆P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	D	1	Total	C	O	P	0	0
			10	3	6	1		

- Molecule 5 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	1	Total	Mn	0	0
			1	1		

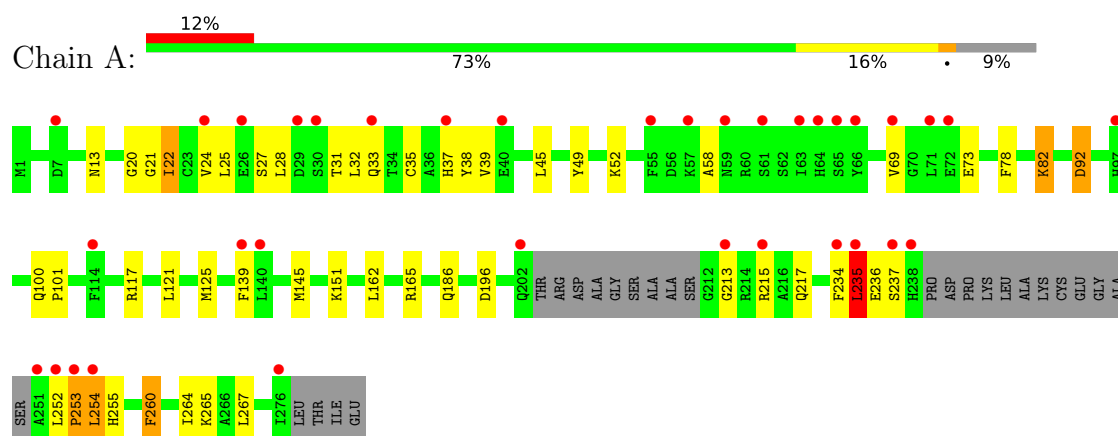
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	86	Total	O	0	0
			86	86		
6	B	122	Total	O	0	0
			122	122		
6	C	123	Total	O	0	0
			123	123		
6	D	122	Total	O	0	0
			122	122		

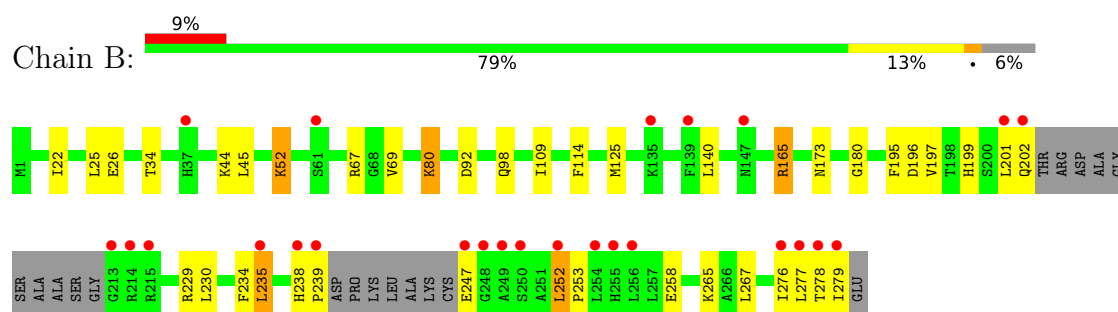
3 Residue-property plots [i](#)

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.

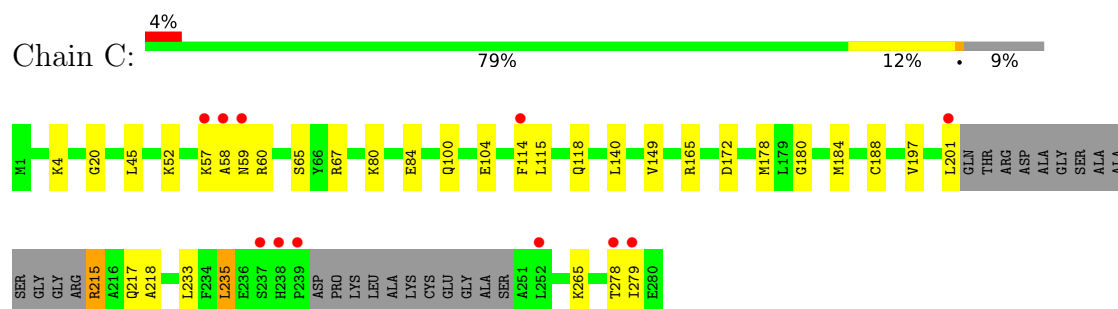
- Molecule 1: 3-deoxy-D-manno-octulosonic acid 8-phosphate synthetase



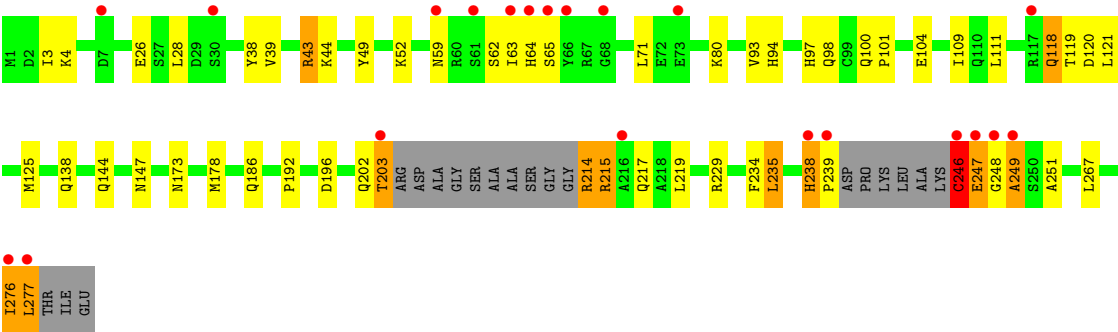
- Molecule 1: 3-deoxy-D-manno-octulosonic acid 8-phosphate synthetase



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- Molecule 1: 3-deoxy-D-manno-octulosonic acid 8-phosphate synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.74Å 86.20Å 163.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.94 – 1.90 33.94 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.4 (33.94-1.90) 99.4 (33.94-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 1.89Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.216 , 0.257 0.217 , 0.260	Depositor DCC
R_{free} test set	4644 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	32.8	Xtriage
Anisotropy	0.066	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	0.001 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8610	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.67% 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

5.1 Standard geometry

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

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	0.60	0/2042	0.80	2/2763 (0.1%)
1	B	0.63	0/2134	0.72	0/2884
1	C	0.62	0/2043	0.76	1/2760 (0.0%)
1	D	0.60	0/2137	0.74	1/2889 (0.0%)
All	All	0.61	0/8356	0.75	4/11296 (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	C	0	1
1	D	0	1
All	All	0	2

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	255	HIS	N-CA-C	-6.19	104.43	112.23
1	A	235	LEU	N-CA-C	5.32	115.56	108.38
1	C	235	LEU	N-CA-C	5.29	117.11	109.07
1	D	235	LEU	N-CA-C	5.14	116.65	108.79

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	215	ARG	Peptide
1	D	246	CYS	Peptide

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	1995	0	2023	49	0
1	B	2082	0	2129	41	0
1	C	2002	0	2056	27	0
1	D	2064	0	2126	47	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	C	1	0	0	0	0
4	D	10	0	2	0	0
5	D	1	0	0	0	0
6	A	86	0	0	5	0
6	B	122	0	0	4	0
6	C	123	0	0	4	0
6	D	122	0	0	1	0
All	All	8610	0	8336	151	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 (151) 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:D:118:GLN:NE2	1:D:120:ASP:OD2	1.70	1.25
1:A:235:LEU:HD12	1:A:235:LEU:C	1.66	1.20
1:C:165[B]:ARG:HG3	1:C:165[B]:ARG:HH11	0.96	1.12
1:B:109:ILE:HG22	1:B:125:MET:HE3	1.22	1.10
1:A:139:PHE:O	6:A:306:HOH:O	1.70	1.08
1:B:52:LYS:HD3	1:B:234:PHE:HZ	1.20	1.04
1:C:165[B]:ARG:HG3	1:C:165[B]:ARG:NH1	1.65	0.99
1:C:165[B]:ARG:HH11	1:C:165[B]:ARG:CG	1.76	0.97
1:A:235:LEU:C	1:A:235:LEU:CD1	2.35	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:235:LEU:CD1	1:A:236:GLU:N	2.30	0.95
1:D:246:CYS:O	1:D:247:GLU:CG	2.15	0.95
1:A:260[B]:PHE:CE2	1:A:264[B]:ILE:HD11	2.02	0.94
1:B:52:LYS:HD3	1:B:234:PHE:CZ	2.00	0.94
1:B:109:ILE:CG2	1:B:125:MET:HE3	1.98	0.92
1:B:235:LEU:C	1:B:235:LEU:HD13	1.97	0.89
1:D:246:CYS:O	1:D:247:GLU:HG3	1.76	0.86
1:A:260[B]:PHE:CE2	1:A:264[B]:ILE:CD1	2.59	0.85
1:A:235:LEU:HD13	1:A:236:GLU:N	1.91	0.84
1:D:93:VAL:HG22	1:D:125[B]:MET:HE2	1.58	0.84
1:B:98:GLN:OE1	6:B:409:HOH:O	1.96	0.83
1:A:58:ALA:HB1	1:C:118:GLN:HE22	1.45	0.82
1:B:235:LEU:HD13	1:B:235:LEU:O	1.80	0.80
1:D:93:VAL:CG2	1:D:125[B]:MET:HE2	2.12	0.80
1:B:109:ILE:HG22	1:B:125:MET:CE	2.10	0.77
1:B:52:LYS:CD	1:B:234:PHE:CZ	2.59	0.76
1:C:172:ASP:OD1	6:C:671:HOH:O	2.03	0.76
1:A:237:SER:HB3	1:A:252:LEU:O	1.88	0.74
1:A:267:LEU:HD13	1:B:267[A]:LEU:CD1	2.22	0.70
1:D:238[B]:HIS:ND1	1:D:239:PRO:HD2	2.07	0.69
1:A:235:LEU:HD12	1:A:236:GLU:N	1.97	0.69
1:B:199[A]:HIS:O	1:B:202:GLN:N	2.23	0.69
1:B:235:LEU:O	1:B:235:LEU:CD1	2.39	0.69
1:A:37[A]:HIS:CD2	6:A:282:HOH:O	2.46	0.69
1:B:165[A]:ARG:NH2	1:B:199[A]:HIS:CE1	2.61	0.69
1:A:22:ILE:HD11	1:A:25:LEU:HD23	1.74	0.68
1:B:252:LEU:HD22	1:B:253:PRO:HD2	1.75	0.68
1:B:247:GLU:HA	6:B:577:HOH:O	1.95	0.67
1:A:27:SER:O	1:A:31:THR:HG23	1.95	0.66
1:D:235:LEU:C	1:D:235:LEU:HD12	2.19	0.66
1:D:247:GLU:HB3	1:D:249:ALA:H	1.61	0.65
1:A:260[B]:PHE:CD2	1:A:264[B]:ILE:HD12	2.31	0.65
1:B:92[A]:ASP:OD1	6:B:549:HOH:O	2.15	0.65
1:C:58:ALA:O	1:C:59:ASN:CB	2.43	0.64
1:D:246:CYS:O	1:D:247:GLU:CB	2.45	0.64
1:C:215:ARG:N	6:C:611:HOH:O	2.30	0.64
1:A:58:ALA:HB1	1:C:118:GLN:NE2	2.12	0.64
1:A:260[B]:PHE:CD2	1:A:264[B]:ILE:CD1	2.82	0.63
1:A:78:PHE:O	1:A:82:LYS:HG3	1.98	0.63
1:A:235:LEU:HD12	1:A:235:LEU:O	1.98	0.63
1:D:111:LEU:HA	1:D:125[B]:MET:HE1	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:26:GLU:O	1:B:69:VAL:HG11	1.98	0.63
1:A:13[B]:ASN:HB2	6:A:441:HOH:O	1.98	0.62
1:B:26:GLU:O	1:B:69:VAL:CG1	2.47	0.62
1:B:52:LYS:CD	1:B:234:PHE:HZ	1.96	0.62
1:D:121:LEU:O	1:D:125[A]:MET:HG3	1.99	0.62
1:A:196:ASP:HA	1:A:234:PHE:HB3	1.81	0.61
1:B:67:ARG:NH2	1:D:119:THR:HG22	2.15	0.60
1:A:100:GLN:N	1:A:101:PRO:HD2	2.16	0.60
1:B:80:LYS:N	1:B:80:LYS:HD3	2.16	0.60
1:A:52[B]:LYS:NZ	1:A:236:GLU:OE2	2.30	0.60
1:B:196:ASP:HA	1:B:234:PHE:HB3	1.83	0.59
1:C:52:LYS:C	1:C:52:LYS:HD3	2.28	0.59
1:C:100:GLN:O	1:C:104:GLU:HG3	2.03	0.58
1:A:267:LEU:CD1	1:B:267[A]:LEU:HD13	2.35	0.56
1:D:173[A]:ASN:OD1	1:D:203:THR:CG2	2.53	0.56
1:C:279:ILE:HD11	1:D:219:LEU:HB2	1.87	0.55
1:B:235:LEU:C	1:B:235:LEU:CD1	2.69	0.55
1:B:196:ASP:OD2	1:B:199[B]:HIS:HD2	1.90	0.55
1:D:111:LEU:CA	1:D:125[B]:MET:HE1	2.37	0.55
1:A:254[A]:LEU:HD23	1:A:254[A]:LEU:O	2.06	0.55
1:D:109:ILE:HD12	1:D:125[B]:MET:HG2	1.89	0.55
1:D:100:GLN:HB3	1:D:101:PRO:HD3	1.90	0.54
1:A:237:SER:CB	1:A:252:LEU:O	2.56	0.54
1:B:22:ILE:CG2	1:B:34:THR:HG21	2.37	0.54
1:A:25:LEU:HD22	1:A:31:THR:HG21	1.90	0.53
1:B:278:THR:O	1:B:279:ILE:C	2.51	0.53
1:A:28:LEU:O	1:A:32:LEU:HG	2.10	0.52
1:D:100:GLN:O	1:D:104:GLU:HG2	2.09	0.52
1:C:178:MET:HG3	1:D:178:MET:HE3	1.91	0.52
1:A:254[B]:LEU:O	1:A:254[B]:LEU:HD23	2.09	0.52
1:D:97:HIS:HD2	6:D:572:HOH:O	1.93	0.52
1:A:235:LEU:HD13	1:A:236:GLU:H	1.73	0.52
1:A:267:LEU:CD1	1:B:267[A]:LEU:CD1	2.88	0.51
1:C:197:VAL:HG21	1:C:233:LEU:HD11	1.92	0.51
1:D:202:GLN:O	1:D:203:THR:HG23	2.11	0.51
1:C:197:VAL:CG2	1:C:233:LEU:HD11	2.40	0.51
1:A:69:VAL:HG22	1:A:73:GLU:CB	2.41	0.50
1:D:39:VAL:O	1:D:43:ARG:HG2	2.12	0.50
1:B:238:HIS:HB2	1:B:239:PRO:HD2	1.94	0.49
1:D:248:GLY:O	1:D:249:ALA:HB3	2.13	0.49
1:B:67:ARG:NH2	1:D:119:THR:CG2	2.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:63:ILE:HG13	1:D:64:HIS:CD2	2.48	0.48
1:A:117:ARG:O	1:C:59:ASN:HA	2.14	0.48
1:A:215:ARG:O	1:A:260[A]:PHE:HZ	1.96	0.48
1:D:52:LYS:HD3	1:D:52:LYS:C	2.38	0.48
1:A:213:GLY:O	1:A:217:GLN:N	2.39	0.48
1:A:151[A]:LYS:NZ	6:A:486:HOH:O	2.28	0.48
1:B:195:PHE:CE2	1:B:197[B]:VAL:HG22	2.48	0.47
1:C:67:ARG:HG3	6:C:662:HOH:O	2.13	0.47
1:B:25:LEU:O	1:B:69:VAL:HG13	2.14	0.47
1:C:178:MET:HE3	1:D:178:MET:HG3	1.96	0.47
1:D:267:LEU:C	1:D:267:LEU:HD23	2.40	0.47
1:D:246:CYS:O	1:D:247:GLU:HG2	2.10	0.47
1:D:173[B]:ASN:HA	1:D:203:THR:HG22	1.97	0.46
1:D:173[A]:ASN:HA	1:D:203:THR:HG22	1.97	0.45
1:A:20:GLY:HA2	1:A:235:LEU:O	2.16	0.45
1:D:215:ARG:NH2	1:D:251:ALA:O	2.42	0.45
1:C:235:LEU:C	1:C:235:LEU:HD12	2.41	0.45
1:C:201:LEU:HD11	1:C:218:ALA:HA	1.98	0.45
1:A:35:CYS:O	1:A:39:VAL:HG23	2.16	0.44
1:D:276:ILE:O	1:D:277:LEU:CB	2.65	0.44
1:A:236:GLU:C	1:A:237:SER:HG	2.20	0.44
1:A:267:LEU:HD13	1:B:267[A]:LEU:HD11	1.98	0.43
1:A:37[A]:HIS:CG	6:A:282:HOH:O	2.70	0.43
1:B:229:ARG:CD	1:B:276:ILE:HD11	2.49	0.43
1:C:114:PHE:CE2	1:C:115:LEU:HD21	2.54	0.43
1:B:44:LYS:HE2	1:B:258:GLU:OE2	2.19	0.43
1:A:38:TYR:HB3	1:A:49:TYR:CZ	2.53	0.42
1:A:92:ASP:OD1	1:A:92:ASP:N	2.36	0.42
1:B:199[A]:HIS:C	1:B:201:LEU:N	2.77	0.42
1:D:39:VAL:O	1:D:43:ARG:CG	2.66	0.42
1:A:121:LEU:O	1:A:125:MET:HG3	2.19	0.42
1:D:196:ASP:HA	1:D:234:PHE:HB3	2.01	0.42
1:D:235:LEU:C	1:D:235:LEU:CD1	2.90	0.42
1:D:38:TYR:HB3	1:D:49:TYR:CZ	2.55	0.42
1:A:252:LEU:O	1:A:253:PRO:O	2.38	0.42
1:D:71:LEU:CD1	1:D:98:GLN:HG2	2.49	0.42
1:C:180:GLY:HA3	6:C:417:HOH:O	2.19	0.42
1:A:145:MET:HE3	1:A:162:LEU:HD22	2.02	0.42
1:D:3:ILE:HD12	1:D:192:PRO:HG2	2.01	0.42
1:B:180:GLY:HA3	6:B:406:HOH:O	2.19	0.41
1:C:80:LYS:NZ	1:C:84:GLU:OE1	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:LEU:O	1:B:265:LYS:HE3	2.20	0.41
1:C:184:MET:O	1:C:188:CYS:HB2	2.20	0.41
1:C:215:ARG:C	1:C:217:GLN:N	2.76	0.41
1:D:28:LEU:HD13	1:D:80:LYS:HD2	2.02	0.41
1:C:165[B]:ARG:NH1	1:C:165[B]:ARG:CG	2.46	0.41
1:D:173[B]:ASN:OD1	1:D:203:THR:HG22	2.20	0.41
1:B:114:PHE:HE1	1:D:118:GLN:HG3	1.85	0.41
1:B:199[A]:HIS:O	1:B:201:LEU:N	2.54	0.41
1:C:45:LEU:O	1:C:265:LYS:HE3	2.21	0.41
1:D:214:ARG:O	1:D:217:GLN:N	2.39	0.41
1:D:144:GLN:OE1	1:D:147:ASN:ND2	2.54	0.40
1:A:45:LEU:O	1:A:265:LYS:HE3	2.21	0.40
1:B:252:LEU:HD22	1:B:253:PRO:CD	2.48	0.40
1:C:20:GLY:HA2	1:C:235:LEU:O	2.22	0.40
1:D:62:SER:HB3	1:D:65:SER:OG	2.22	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	259/280 (92%)	249 (96%)	9 (4%)	1 (0%)	30	22
1	B	267/280 (95%)	258 (97%)	9 (3%)	0	100	100
1	C	255/280 (91%)	250 (98%)	5 (2%)	0	100	100
1	D	266/280 (95%)	257 (97%)	6 (2%)	3 (1%)	11	4
All	All	1047/1120 (94%)	1014 (97%)	29 (3%)	4 (0%)	30	22

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	253	PRO
1	D	249	ALA
1	D	247	GLU
1	D	229	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	217/235 (92%)	205 (94%)	12 (6%)	19	11
1	B	232/235 (99%)	222 (96%)	10 (4%)	26	18
1	C	221/235 (94%)	214 (97%)	7 (3%)	34	27
1	D	232/235 (99%)	213 (92%)	19 (8%)	10	5
All	All	902/940 (96%)	854 (95%)	48 (5%)	22	12

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ILE
1	A	24	VAL
1	A	33	GLN
1	A	82	LYS
1	A	92	ASP
1	A	165	ARG
1	A	186	GLN
1	A	235	LEU
1	A	254[A]	LEU
1	A	254[B]	LEU
1	A	260[A]	PHE
1	A	260[B]	PHE
1	B	52	LYS
1	B	80	LYS
1	B	140	LEU
1	B	165[A]	ARG
1	B	165[B]	ARG
1	B	173	ASN

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Mol	Chain	Res	Type
1	B	230	LEU
1	B	235	LEU
1	B	252	LEU
1	B	277	LEU
1	C	4	LYS
1	C	57	LYS
1	C	60	ARG
1	C	65	SER
1	C	140	LEU
1	C	149	VAL
1	C	278	THR
1	D	4	LYS
1	D	26	GLU
1	D	43	ARG
1	D	44	LYS
1	D	59	ASN
1	D	94[A]	HIS
1	D	94[B]	HIS
1	D	118	GLN
1	D	138[A]	GLN
1	D	138[B]	GLN
1	D	186	GLN
1	D	203	THR
1	D	214	ARG
1	D	215	ARG
1	D	238[A]	HIS
1	D	238[B]	HIS
1	D	246	CYS
1	D	276	ILE
1	D	277	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	100	GLN
1	A	118	GLN
1	B	94	HIS
1	B	98	GLN
1	B	147	ASN
1	B	173	ASN
1	B	274	GLN

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Mol	Chain	Res	Type
1	C	100	GLN
1	C	110	GLN
1	C	118	GLN
1	C	133	ASN
1	D	64	HIS
1	D	217	GLN
1	D	255	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
4	PEP	D	281	-	9,9,9	2.65	1 (11%)	11,13,13	2.21	5 (45%)

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
4	PEP	D	281	-	-	1/9/9/9	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	281	PEP	C3-C2	7.37	1.51	1.31

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	281	PEP	C3-C2-C1	-4.36	114.66	122.73
4	D	281	PEP	O2-C2-C3	-3.67	117.83	124.85
4	D	281	PEP	O2'-C1-C2	2.67	118.47	113.91
4	D	281	PEP	O1-C1-C2	-2.30	118.32	121.79
4	D	281	PEP	O2-P-O1P	-2.09	102.79	109.47

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	281	PEP	O2'-C1-C2-C3

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

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	255/280 (91%)	0.81	35 (13%) 6 7	12, 27, 58, 63	11 (4%)
1	B	262/280 (93%)	0.36	25 (9%) 14 14	8, 23, 47, 66	17 (6%)
1	C	256/280 (91%)	0.10	11 (4%) 40 42	10, 24, 42, 60	7 (2%)
1	D	261/280 (93%)	0.46	21 (8%) 18 19	11, 25, 46, 58	15 (5%)
All	All	1034/1120 (92%)	0.43	92 (8%) 15 16	8, 25, 47, 66	50 (4%)

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	248	GLY	9.6
1	D	246	CYS	8.3
1	B	247	GLU	7.7
1	D	248	GLY	6.5
1	B	249	ALA	5.9
1	B	279	ILE	5.7
1	A	251	ALA	5.3
1	A	235	LEU	4.8
1	D	277	LEU	4.8
1	A	66	TYR	4.7
1	B	277	LEU	4.7
1	D	238[A]	HIS	4.5
1	C	58	ALA	4.4
1	D	30	SER	4.3
1	B	239	PRO	4.3
1	B	37[A]	HIS	4.1
1	B	238	HIS	4.1
1	A	238	HIS	3.9
1	B	254	LEU	3.9
1	C	239	PRO	3.8
1	A	114[A]	PHE	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	214	ARG	3.7
1	D	66	TYR	3.7
1	C	238	HIS	3.6
1	C	279	ILE	3.6
1	D	249	ALA	3.5
1	A	26	GLU	3.5
1	B	202	GLN	3.5
1	D	64	HIS	3.4
1	A	237	SER	3.4
1	D	59	ASN	3.3
1	B	252	LEU	3.3
1	B	255	HIS	3.2
1	C	237	SER	3.2
1	D	216	ALA	3.1
1	A	63	ILE	3.0
1	A	24	VAL	3.0
1	A	61	SER	3.0
1	B	213	GLY	3.0
1	A	234	PHE	2.9
1	A	69	VAL	2.9
1	A	252	LEU	2.9
1	B	215	ARG	2.8
1	A	254[A]	LEU	2.8
1	B	278	THR	2.8
1	D	203	THR	2.8
1	A	7	ASP	2.8
1	D	247	GLU	2.7
1	B	147	ASN	2.7
1	B	201	LEU	2.7
1	C	57	LYS	2.7
1	A	213	GLY	2.6
1	A	139	PHE	2.6
1	B	256	LEU	2.6
1	A	215	ARG	2.5
1	A	253	PRO	2.5
1	B	139	PHE	2.5
1	D	7	ASP	2.5
1	A	276	ILE	2.5
1	C	278	THR	2.5
1	A	97	HIS	2.5
1	D	239	PRO	2.4
1	C	252	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	63	ILE	2.4
1	D	276	ILE	2.4
1	B	235	LEU	2.3
1	A	202	GLN	2.3
1	A	64	HIS	2.3
1	D	73	GLU	2.2
1	B	250	SER	2.2
1	A	55	PHE	2.2
1	B	135	LYS	2.2
1	A	71	LEU	2.2
1	D	61	SER	2.1
1	D	65	SER	2.1
1	A	140	LEU	2.1
1	A	37[A]	HIS	2.1
1	A	29	ASP	2.1
1	B	61	SER	2.1
1	D	68	GLY	2.1
1	D	117[A]	ARG	2.1
1	C	114	PHE	2.1
1	A	40	GLU	2.1
1	A	33	GLN	2.1
1	A	30	SER	2.0
1	A	65	SER	2.0
1	B	276	ILE	2.0
1	C	59	ASN	2.0
1	C	201	LEU	2.0
1	A	72	GLU	2.0
1	A	57	LYS	2.0
1	A	59	ASN	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 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
5	MN	D	282	1/1	0.88	0.18	27,27,27,27	1
4	PEP	D	281	10/10	0.92	0.11	19,21,26,27	10
2	CL	C	281	1/1	0.94	0.12	24,24,24,24	1
2	CL	B	281	1/1	0.97	0.04	15,15,15,15	1
3	NA	C	282	1/1	0.98	0.11	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.