



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 05:46 AM UTC

PDB ID : 4JTH / pdb_00004jth
Title : Crystal structure of F114R/R117Q mutant of 3-deoxy-D-manno-octulosonate 8-phosphate synthase (KDO8PS) from *Neisseria meningitidis*
Authors : Allison, T.M.; Cochrane, F.C.; Jameson, G.B.; Parker, E.J.
Deposited on : 2013-03-23
Resolution : 2.00 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

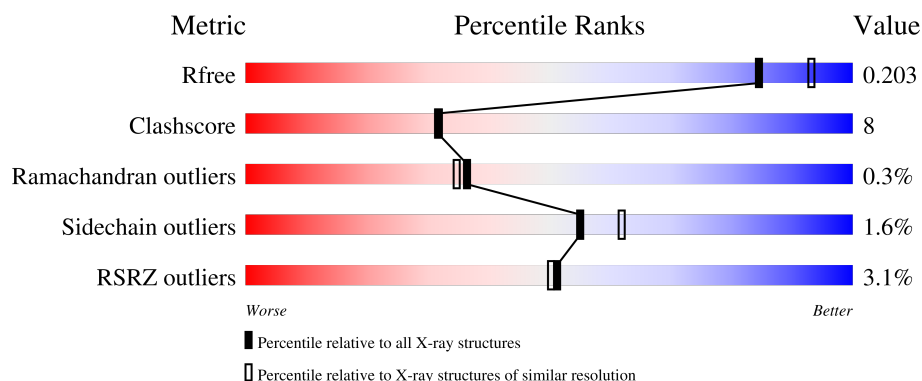
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	3% 82% 7% 11%
1	B	280	2% 75% 12% 10%
1	C	280	2% 82% 9% 8%
1	D	280	4% 80% 10% 9%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8774 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	250	Total	C	N	O	S	0	5	0
			1953	1254	326	362	11			
1	B	251	Total	C	N	O	S	0	10	0
			1994	1280	336	366	12			
1	C	259	Total	C	N	O	S	0	6	0
			2020	1296	341	372	11			
1	D	256	Total	C	N	O	S	0	6	0
			2000	1281	341	367	11			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	114	ARG	PHE	engineered mutation	UNP Q9JZ55
A	117	GLN	ARG	engineered mutation	UNP Q9JZ55
B	114	ARG	PHE	engineered mutation	UNP Q9JZ55
B	117	GLN	ARG	engineered mutation	UNP Q9JZ55
C	114	ARG	PHE	engineered mutation	UNP Q9JZ55
C	117	GLN	ARG	engineered mutation	UNP Q9JZ55
D	114	ARG	PHE	engineered mutation	UNP Q9JZ55
D	117	GLN	ARG	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	2	Total	Cl	0	0
			2	2		
2	D	2	Total	Cl	0	0
			2	2		

- 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 GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

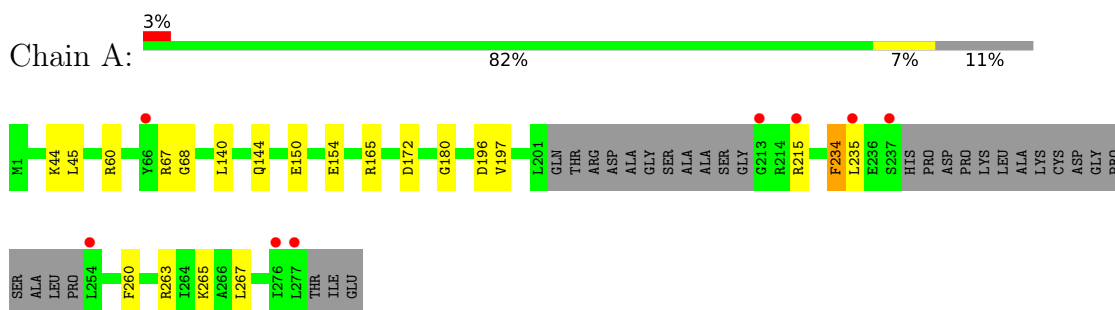
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	162	Total	O	0	0
			162	162		
5	B	224	Total	O	0	0
			224	224		
5	C	215	Total	O	0	0
			215	215		
5	D	194	Total	O	0	0
			194	194		

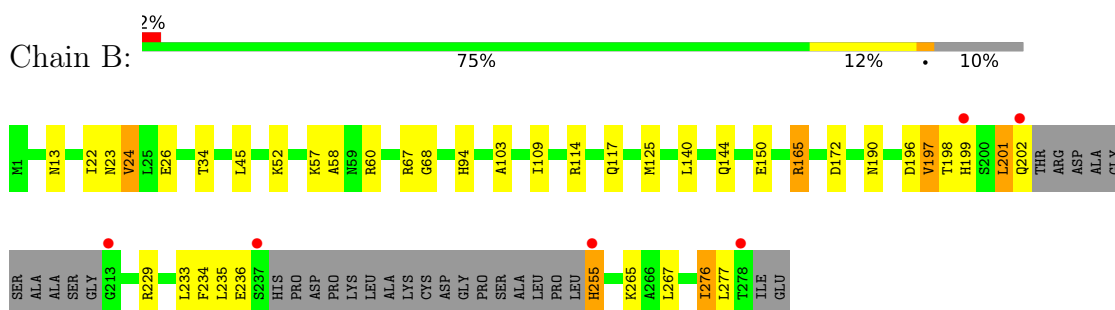
3 Residue-property plots [i](#)

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.

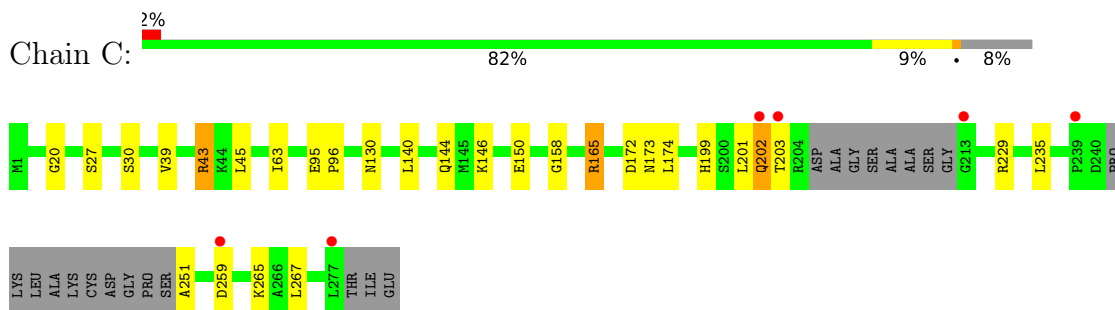
- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase



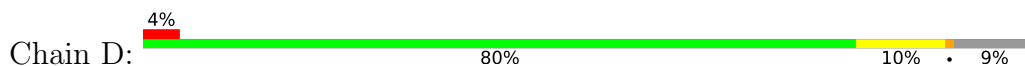
- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase

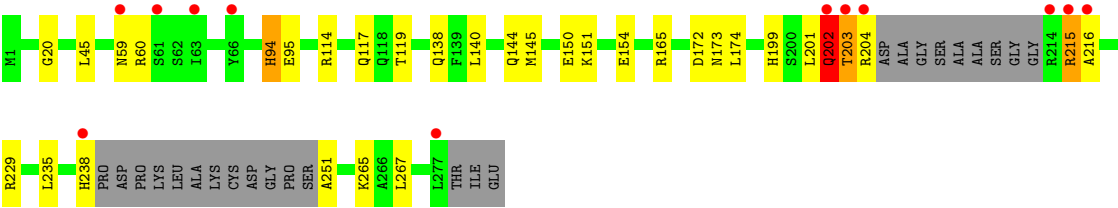


- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase



- 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	82.15Å 85.99Å 163.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.95 – 2.00 45.95 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (45.95-2.00) 100.0 (45.95-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.81 (at 2.00Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.159 , 0.194 0.167 , 0.203	Depositor DCC
R_{free} test set	3931 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	22.8	Xtriage
Anisotropy	0.410	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.023 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8774	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% 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: GOL, CL, NA

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	1.15	1/1987 (0.1%)	1.00	5/2686 (0.2%)
1	B	1.27	5/2038 (0.2%)	1.06	3/2754 (0.1%)
1	C	1.25	3/2066 (0.1%)	1.08	0/2793
1	D	1.15	2/2049 (0.1%)	1.03	1/2769 (0.0%)
All	All	1.21	11/8140 (0.1%)	1.05	9/11002 (0.1%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	173	ASN	CA-CB	-6.41	1.42	1.53
1	C	43	ARG	C-O	-5.85	1.17	1.24
1	A	263	ARG	C-O	-5.71	1.17	1.24
1	C	158	GLY	CA-C	5.47	1.57	1.51
1	D	145	MET	C-O	-5.33	1.17	1.24
1	B	197[A]	VAL	C-O	-5.21	1.18	1.24
1	B	197[B]	VAL	C-O	-5.21	1.18	1.24
1	B	276	ILE	C-O	-5.17	1.18	1.24
1	B	229	ARG	CZ-NH2	-5.12	1.26	1.33
1	B	23	ASN	N-CA	-5.08	1.40	1.46
1	D	95	GLU	CA-C	-5.07	1.46	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	234	PHE	CB-CA-C	-6.44	99.12	109.75
1	B	201	LEU	CB-CA-C	-6.36	99.33	111.48
1	D	172	ASP	CB-CA-C	-6.32	101.53	111.51
1	A	172	ASP	CB-CA-C	-5.87	102.24	111.51
1	B	172	ASP	CB-CA-C	-5.66	102.57	111.51
1	A	67	ARG	N-CA-CB	-5.37	102.63	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	60	ARG	N-CA-C	-5.13	101.79	109.95
1	A	68	GLY	N-CA-C	-5.05	105.11	112.17
1	B	165	ARG	CB-CA-C	-5.02	101.42	110.35

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	1953	0	1993	12	0
1	B	1994	0	2048	60	0
1	C	2020	0	2078	23	0
1	D	2000	0	2055	40	0
2	B	1	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	1	0
3	C	1	0	0	0	0
4	D	6	0	8	0	0
5	A	162	0	0	2	0
5	B	224	0	0	4	0
5	C	215	0	0	5	0
5	D	194	0	0	15	0
All	All	8774	0	8182	131	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 8.

All (131) 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:C:259:ASP:HB2	5:C:601:HOH:O	1.45	1.17
1:B:199[B]:HIS:CE1	1:B:236[B]:GLU:HG3	1.91	1.05
1:D:154:GLU:HG2	5:D:585:HOH:O	1.65	0.97
1:C:146:LYS:HD2	5:C:604:HOH:O	1.65	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109[B]:ILE:HD12	1:B:125[B]:MET:HG3	1.51	0.92
1:D:265:LYS:CE	5:D:534:HOH:O	2.22	0.86
1:B:199[B]:HIS:NE2	1:B:236[B]:GLU:CG	2.40	0.85
1:B:196:ASP:OD2	1:B:199[B]:HIS:HD2	1.61	0.84
1:C:43:ARG:HD3	5:C:482:HOH:O	1.78	0.84
1:B:117:GLN:OE1	1:D:114:ARG:NH1	2.10	0.83
1:D:202:GLN:HE21	1:D:202:GLN:HA	1.45	0.82
1:D:202:GLN:HA	1:D:202:GLN:NE2	1.92	0.82
1:B:190:ASN:HB2	5:B:600:HOH:O	1.81	0.81
1:D:154:GLU:CG	5:D:585:HOH:O	2.26	0.78
1:B:109[A]:ILE:HG22	1:B:125[A]:MET:HE3	1.65	0.78
1:D:150:GLU:HG2	5:D:587:HOH:O	1.84	0.78
1:B:103:ALA:HB2	1:B:109[A]:ILE:HD11	1.68	0.76
1:B:13:ASN:HD21	1:B:276:ILE:HD11	1.51	0.75
1:B:199[B]:HIS:NE2	1:B:236[B]:GLU:HG3	1.98	0.75
1:A:196:ASP:HA	1:A:234:PHE:HB2	1.66	0.75
1:B:199[A]:HIS:HB2	1:B:236[A]:GLU:OE2	1.87	0.75
1:D:150:GLU:HB3	5:D:587:HOH:O	1.87	0.73
1:B:52:LYS:HE2	1:B:234:PHE:CZ	2.25	0.72
1:C:165:ARG:HH22	1:C:202:GLN:HE22	1.37	0.72
1:B:199[A]:HIS:O	1:B:199[A]:HIS:CD2	2.44	0.71
1:B:199[B]:HIS:CE1	1:B:236[B]:GLU:CG	2.71	0.71
1:A:44:LYS:HB2	5:A:365:HOH:O	1.90	0.70
1:B:165:ARG:NH1	1:B:199[A]:HIS:CE1	2.61	0.69
1:B:190:ASN:CB	5:B:600:HOH:O	2.40	0.68
1:D:265:LYS:HE3	5:D:534:HOH:O	1.87	0.68
1:B:109[B]:ILE:CD1	1:B:125[B]:MET:HG3	2.23	0.67
1:B:199[B]:HIS:NE2	1:B:236[B]:GLU:HG2	2.08	0.67
1:D:150:GLU:CB	5:D:587:HOH:O	2.42	0.66
1:D:138:GLN:HG2	2:D:302:CL:CL	2.32	0.66
1:A:197[A]:VAL:HG12	1:A:234:PHE:O	1.95	0.66
1:B:199[A]:HIS:O	1:B:199[A]:HIS:CG	2.49	0.65
1:B:198:THR:HG21	1:B:236[A]:GLU:HB2	1.78	0.64
1:B:255:HIS:N	1:B:255:HIS:HD1	1.95	0.64
1:B:114:ARG:NH2	1:D:117:GLN:OE1	2.30	0.64
1:B:13:ASN:ND2	1:B:276:ILE:HD11	2.13	0.64
1:B:52:LYS:HE2	1:B:234:PHE:CE1	2.31	0.64
1:C:201:LEU:O	1:C:203:THR:N	2.27	0.63
1:A:154[B]:GLU:HG3	1:C:63:ILE:HD11	1.78	0.63
1:B:190:ASN:CG	5:B:600:HOH:O	2.41	0.62
1:B:109[A]:ILE:CG2	1:B:125[A]:MET:HE3	2.30	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:154:GLU:CB	5:D:585:HOH:O	2.48	0.61
1:D:45:LEU:O	1:D:265:LYS:HE3	2.00	0.61
1:B:26:GLU:O	1:B:26:GLU:HG3	2.00	0.60
1:D:150:GLU:CG	5:D:587:HOH:O	2.49	0.59
1:C:146:LYS:CD	5:C:604:HOH:O	2.37	0.59
1:D:173:ASN:HA	1:D:203:THR:HG21	1.85	0.58
1:D:174:LEU:HD12	1:D:202:GLN:CB	2.33	0.58
1:D:238:HIS:C	5:D:525:HOH:O	2.46	0.58
1:B:198:THR:HG21	1:B:236[B]:GLU:HB3	1.85	0.58
1:B:196:ASP:OD2	1:B:199[B]:HIS:CD2	2.48	0.57
1:D:60:ARG:HD2	5:D:495:HOH:O	2.03	0.57
1:D:215:ARG:NH2	1:D:215:ARG:HG3	2.18	0.57
1:B:45:LEU:O	1:B:265:LYS:HE3	2.06	0.55
1:D:174:LEU:HD12	1:D:202:GLN:HB2	1.86	0.55
1:C:165:ARG:HH22	1:C:202:GLN:NE2	2.04	0.55
1:B:198:THR:HG21	1:B:236[B]:GLU:CB	2.38	0.54
1:D:94[B]:HIS:HE1	5:D:540:HOH:O	1.91	0.54
1:B:52:LYS:CE	1:B:234:PHE:CZ	2.91	0.53
1:D:251:ALA:HB2	5:D:502:HOH:O	2.07	0.53
1:D:265:LYS:NZ	5:D:534:HOH:O	2.27	0.53
1:A:215:ARG:O	1:A:260:PHE:HZ	1.90	0.53
1:D:215:ARG:HG2	1:D:216:ALA:N	2.23	0.53
1:B:255:HIS:N	1:B:255:HIS:ND1	2.55	0.52
1:C:45:LEU:O	1:C:265:LYS:HE3	2.10	0.52
1:B:198:THR:HB	1:B:236[A]:GLU:HG3	1.91	0.52
1:D:215:ARG:HG3	1:D:215:ARG:HH21	1.73	0.51
1:B:199[A]:HIS:CB	1:B:236[A]:GLU:OE2	2.57	0.51
1:D:119:THR:HG22	1:D:151:LYS:NZ	2.26	0.51
1:B:199[A]:HIS:CD2	1:B:199[A]:HIS:C	2.89	0.50
1:C:165:ARG:NH2	1:C:202:GLN:HE22	2.05	0.50
1:C:172:ASP:O	1:C:203:THR:OG1	2.30	0.49
1:C:199:HIS:O	1:C:202:GLN:HB2	2.11	0.49
1:C:267[B]:LEU:C	1:C:267[B]:LEU:HD23	2.36	0.49
1:D:203:THR:O	1:D:203:THR:OG1	2.30	0.49
1:B:199[B]:HIS:HE1	1:B:236[B]:GLU:HG3	1.64	0.49
1:C:27:SER:OG	1:C:30:SER:HB3	2.13	0.49
1:B:24[B]:VAL:CG1	1:B:68:GLY:HA2	2.42	0.48
1:C:165:ARG:HH12	1:C:202:GLN:NE2	2.11	0.48
1:A:45:LEU:O	1:A:265:LYS:HE3	2.15	0.47
1:D:173:ASN:HA	1:D:203:THR:CG2	2.45	0.47
1:D:199:HIS:O	1:D:202:GLN:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:165:ARG:O	1:B:199[A]:HIS:HD2	1.98	0.46
1:C:39:VAL:O	1:C:43:ARG:HG3	2.16	0.46
1:B:198:THR:HG21	1:B:236[A]:GLU:O	2.16	0.45
1:A:197[A]:VAL:CG1	1:A:235:LEU:HD23	2.46	0.45
1:C:201:LEU:O	1:C:203:THR:HG22	2.16	0.45
1:C:140:LEU:HG	1:C:144:GLN:HB2	2.00	0.44
1:A:150[A]:GLU:CD	1:A:150[A]:GLU:C	2.86	0.44
1:B:197[B]:VAL:HG21	1:B:233:LEU:HD11	1.98	0.44
1:B:198:THR:CG2	1:B:236[A]:GLU:HB2	2.46	0.44
1:B:201:LEU:C	1:B:202:GLN:HG3	2.42	0.44
1:C:251:ALA:HB2	5:C:590:HOH:O	2.18	0.44
1:B:267:LEU:HD23	1:B:267:LEU:C	2.42	0.44
1:D:267:LEU:C	1:D:267:LEU:HD23	2.43	0.44
1:D:140:LEU:HG	1:D:144:GLN:HB2	1.99	0.43
1:A:267[B]:LEU:C	1:A:267[B]:LEU:HD23	2.44	0.43
1:B:57:LYS:HD3	1:B:60:ARG:HG2	2.01	0.43
1:D:20:GLY:HA2	1:D:235:LEU:O	2.18	0.43
1:B:52:LYS:CD	1:B:234:PHE:CZ	3.01	0.43
1:B:235:LEU:O	5:B:573:HOH:O	2.20	0.43
1:D:235:LEU:C	1:D:235:LEU:HD12	2.43	0.43
1:D:265:LYS:CD	5:D:534:HOH:O	2.60	0.43
1:B:103:ALA:CB	1:B:109[A]:ILE:HD11	2.46	0.43
1:B:117:GLN:O	1:D:59:ASN:HA	2.18	0.43
1:B:201:LEU:O	1:B:202:GLN:CB	2.67	0.43
1:B:22:ILE:HG23	1:B:34:THR:HG21	2.00	0.43
1:B:165:ARG:HB3	1:B:199[A]:HIS:CD2	2.53	0.43
1:D:201:LEU:O	1:D:203:THR:N	2.47	0.42
1:A:140:LEU:HG	1:A:144:GLN:HB2	2.00	0.42
1:B:196:ASP:HA	1:B:234:PHE:HB3	2.01	0.42
1:A:180:GLY:HA3	5:A:315:HOH:O	2.19	0.42
1:B:140:LEU:HG	1:B:144:GLN:HB2	2.01	0.41
1:D:204:ARG:HH11	1:D:204:ARG:HD2	1.65	0.41
1:B:24[B]:VAL:CG1	1:B:68:GLY:CA	2.98	0.41
1:D:202:GLN:HE21	1:D:202:GLN:CA	2.25	0.41
1:B:198:THR:HG21	1:B:236[B]:GLU:O	2.19	0.41
1:D:215:ARG:HH21	1:D:215:ARG:CG	2.33	0.41
1:B:58:ALA:HB2	1:B:94:HIS:CE1	2.56	0.41
1:B:201:LEU:O	1:B:202:GLN:CG	2.69	0.41
1:A:215:ARG:O	1:A:260:PHE:CZ	2.73	0.41
1:C:172:ASP:C	1:C:203:THR:HG1	2.29	0.41
1:C:174:LEU:N	1:C:203:THR:HG21	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24[B]:VAL:HG13	1:B:68:GLY:CA	2.51	0.40
1:B:57:LYS:O	1:B:67:ARG:HG2	2.22	0.40
1:C:95:GLU:HA	1:C:96:PRO:HD3	1.97	0.40
1:C:20:GLY:HA2	1:C:235:LEU:O	2.22	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	249/280 (89%)	244 (98%)	5 (2%)	0	100	100
1	B	255/280 (91%)	250 (98%)	5 (2%)	0	100	100
1	C	259/280 (92%)	252 (97%)	6 (2%)	1 (0%)	30	27
1	D	256/280 (91%)	252 (98%)	2 (1%)	2 (1%)	16	11
All	All	1019/1120 (91%)	998 (98%)	18 (2%)	3 (0%)	36	35

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	202	GLN
1	C	229	ARG
1	D	229	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	216/236 (92%)	215 (100%)	1 (0%)	81	87
1	B	223/236 (94%)	218 (98%)	5 (2%)	45	50
1	C	226/236 (96%)	222 (98%)	4 (2%)	51	58
1	D	224/236 (95%)	218 (97%)	6 (3%)	39	42
All	All	889/944 (94%)	873 (98%)	16 (2%)	55	58

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	165	ARG
1	B	24[A]	VAL
1	B	24[B]	VAL
1	B	150	GLU
1	B	255	HIS
1	B	277	LEU
1	C	130	ASN
1	C	150	GLU
1	C	165	ARG
1	C	202	GLN
1	D	94[A]	HIS
1	D	94[B]	HIS
1	D	165	ARG
1	D	202	GLN
1	D	203	THR
1	D	215	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	37	HIS
1	A	59	ASN
1	A	64	HIS
1	A	94	HIS
1	B	37	HIS
1	B	118	GLN
1	B	147	ASN
1	B	186	GLN
1	B	274	GLN

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Mol	Chain	Res	Type
1	C	100	GLN
1	C	110	GLN
1	C	133	ASN
1	C	202	GLN
1	C	274	GLN
1	D	59	ASN
1	D	64	HIS
1	D	202	GLN
1	D	217	GLN

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 7 ligands modelled in this entry, 6 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	GOL	D	303	-	5,5,5	0.22	0	5,5,5	0.72	0

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	GOL	D	303	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	303	GOL	C1-C2-C3-O3
4	D	303	GOL	O2-C2-C3-O3

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	250/280 (89%)	-0.06	8 (3%)	50	49	7, 27, 57, 80	6 (2%)
1	B	251/280 (89%)	-0.27	6 (2%)	59	59	8, 21, 41, 84	10 (3%)
1	C	259/280 (92%)	-0.35	6 (2%)	61	60	8, 20, 41, 83	6 (2%)
1	D	256/280 (91%)	-0.12	12 (4%)	36	35	9, 26, 54, 74	6 (2%)
All	All	1016/1120 (90%)	-0.20	32 (3%)	51	50	7, 23, 53, 84	28 (2%)

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	254	LEU	4.7
1	B	278	THR	4.5
1	A	66	TYR	4.3
1	D	277	LEU	4.2
1	A	235	LEU	4.2
1	D	203	THR	3.7
1	D	66	TYR	3.3
1	A	213	GLY	3.2
1	B	255	HIS	3.1
1	D	61	SER	3.0
1	B	199[A]	HIS	3.0
1	C	203	THR	2.9
1	D	59	ASN	2.8
1	C	239	PRO	2.7
1	D	214	ARG	2.7
1	D	216	ALA	2.7
1	D	238	HIS	2.5
1	C	202	GLN	2.5
1	A	277	LEU	2.5
1	D	215	ARG	2.4
1	B	213	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	276	ILE	2.4
1	D	202	GLN	2.4
1	C	213	GLY	2.3
1	C	259	ASP	2.3
1	B	237	SER	2.2
1	D	204	ARG	2.2
1	A	237	SER	2.2
1	C	277	LEU	2.2
1	D	63	ILE	2.1
1	A	215	ARG	2.1
1	B	202	GLN	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
4	GOL	D	303	6/6	0.79	0.22	41,49,51,54	6
2	CL	D	302	1/1	0.95	0.12	47,47,47,47	0
3	NA	C	303	1/1	0.96	0.10	28,28,28,28	0
2	CL	B	301	1/1	0.97	0.06	25,25,25,25	0
2	CL	C	301	1/1	0.98	0.06	24,24,24,24	0
2	CL	C	302	1/1	0.98	0.08	32,32,32,32	0
2	CL	D	301	1/1	0.98	0.05	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.