



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:07 PM UTC

PDB ID : 1N3B / pdb_00001n3b
Title : Crystal Structure of Dephosphocoenzyme A kinase from Escherichia coli
Authors : O'Toole, N.; Barbosa, J.A.R.G.; Li, Y.; Hung, L.-W.; Matte, A.; Cygler, M.;
Montreal-Kingston Bacterial Structural Genomics Initiative (BSGI)
Deposited on : 2002-10-25
Resolution : 1.80 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

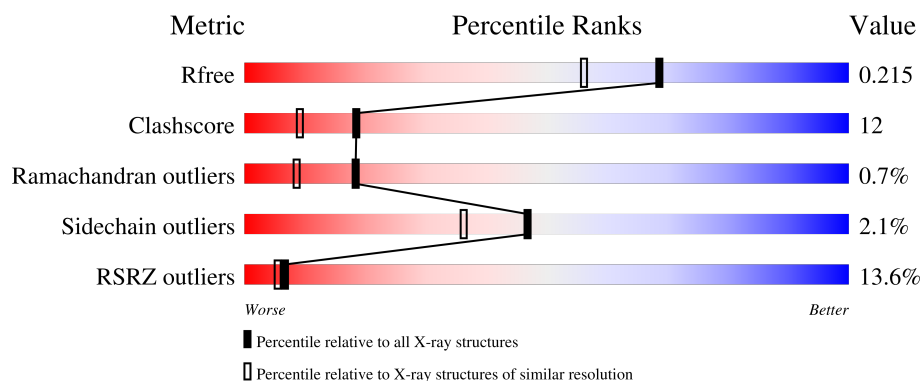
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	216	18% 60% 20% 19%
1	B	216	2% 79% 16% 5%
1	C	216	17% 72% 19% 5%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5046 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 Dephospho-CoA kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	176	Total	C	N	O	Se	0	0	0
			1359	853	244	260	2			
1	B	206	Total	C	N	O	Se	0	0	0
			1594	999	292	300	3			
1	C	206	Total	C	N	O	Se	0	0	0
			1594	999	292	300	3			

There are 39 discrepancies between the modelled and reference sequences:

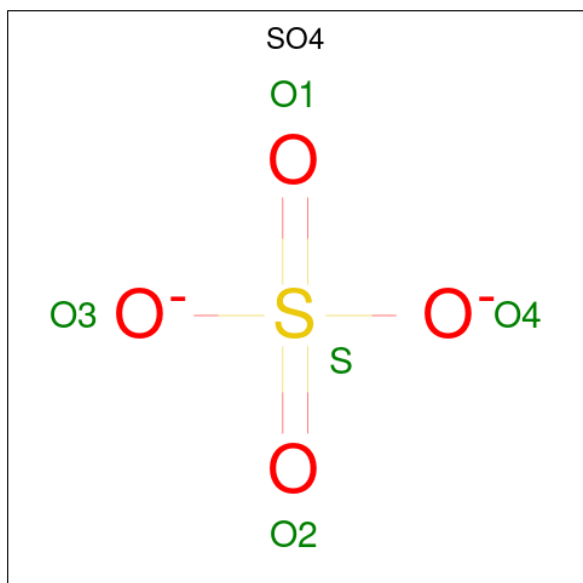
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP P0A6I9
A	58	MSE	MET	modified residue	UNP P0A6I9
A	142	MSE	MET	modified residue	UNP P0A6I9
A	207	GLY	-	expression tag	UNP P0A6I9
A	208	SER	-	expression tag	UNP P0A6I9
A	209	HIS	-	expression tag	UNP P0A6I9
A	210	HIS	-	expression tag	UNP P0A6I9
A	211	HIS	-	expression tag	UNP P0A6I9
A	212	HIS	-	expression tag	UNP P0A6I9
A	213	HIS	-	expression tag	UNP P0A6I9
A	214	HIS	-	expression tag	UNP P0A6I9
A	215	HIS	-	expression tag	UNP P0A6I9
A	216	HIS	-	expression tag	UNP P0A6I9
B	1	MSE	MET	modified residue	UNP P0A6I9
B	58	MSE	MET	modified residue	UNP P0A6I9
B	142	MSE	MET	modified residue	UNP P0A6I9
B	207	GLY	-	expression tag	UNP P0A6I9
B	208	SER	-	expression tag	UNP P0A6I9
B	209	HIS	-	expression tag	UNP P0A6I9
B	210	HIS	-	expression tag	UNP P0A6I9
B	211	HIS	-	expression tag	UNP P0A6I9
B	212	HIS	-	expression tag	UNP P0A6I9
B	213	HIS	-	expression tag	UNP P0A6I9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	214	HIS	-	expression tag	UNP P0A6I9
B	215	HIS	-	expression tag	UNP P0A6I9
B	216	HIS	-	expression tag	UNP P0A6I9
C	1	MSE	MET	modified residue	UNP P0A6I9
C	58	MSE	MET	modified residue	UNP P0A6I9
C	142	MSE	MET	modified residue	UNP P0A6I9
C	207	GLY	-	expression tag	UNP P0A6I9
C	208	SER	-	expression tag	UNP P0A6I9
C	209	HIS	-	expression tag	UNP P0A6I9
C	210	HIS	-	expression tag	UNP P0A6I9
C	211	HIS	-	expression tag	UNP P0A6I9
C	212	HIS	-	expression tag	UNP P0A6I9
C	213	HIS	-	expression tag	UNP P0A6I9
C	214	HIS	-	expression tag	UNP P0A6I9
C	215	HIS	-	expression tag	UNP P0A6I9
C	216	HIS	-	expression tag	UNP P0A6I9

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		

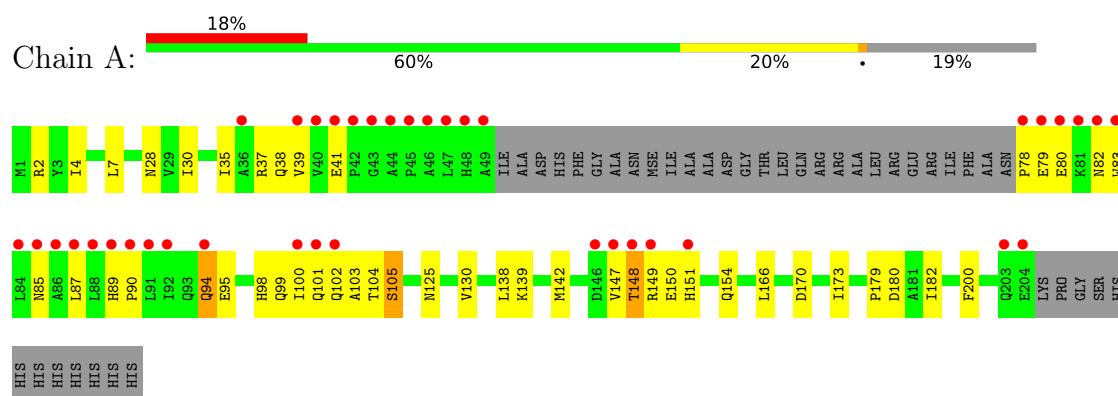
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	107	Total	O	0	0
			107	107		
3	B	189	Total	O	0	0
			189	189		
3	C	153	Total	O	0	0
			153	153		

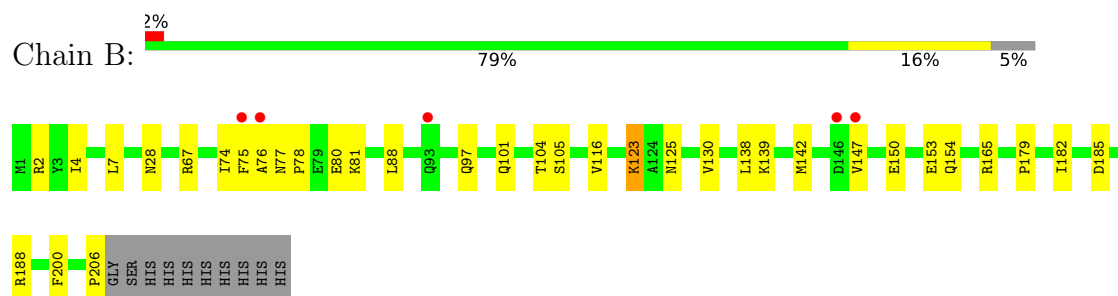
3 Residue-property plots [i](#)

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.

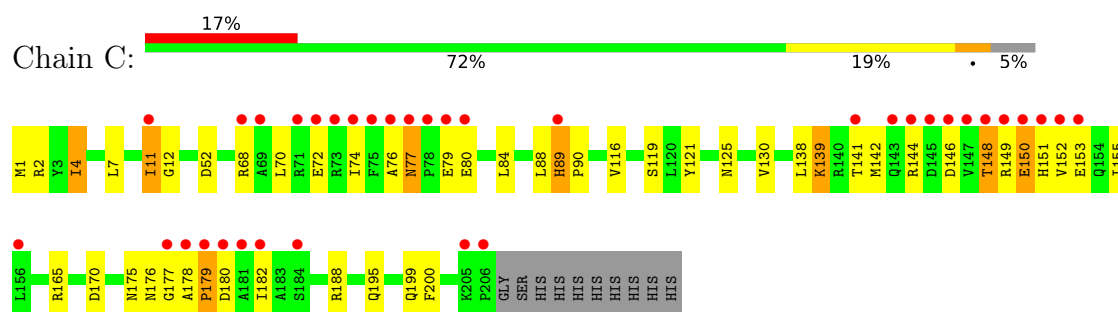
• Molecule 1: Dephospho-CoA kinase



• Molecule 1: Dephospho-CoA kinase



• Molecule 1: Dephospho-CoA kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.50Å 82.41Å 76.00Å 90.00° 94.77° 90.00°	Depositor
Resolution (Å)	40.00 – 1.80 40.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	86.8 (40.00-1.80) 87.5 (40.00-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 1.75Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.217 , 0.246 0.221 , 0.215	Depositor DCC
R_{free} test set	5806 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	20.1	Xtriage
Anisotropy	0.340	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.7	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	5046	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% 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: SO4

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	0.49	0/1380	0.90	4/1880 (0.2%)
1	B	0.51	0/1619	0.91	4/2202 (0.2%)
1	C	0.49	0/1619	0.93	8/2202 (0.4%)
All	All	0.50	0/4618	0.91	16/6284 (0.3%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	200	PHE	N-CA-C	9.25	121.36	111.28
1	A	200	PHE	N-CA-C	7.42	120.31	111.33
1	C	170	ASP	N-CA-C	-6.55	103.35	112.45
1	C	200	PHE	N-CA-C	6.43	119.11	111.71
1	C	176	ASN	N-CA-C	-6.41	104.51	112.90
1	A	170	ASP	N-CA-C	-6.21	103.82	112.45
1	B	67	ARG	N-CA-C	6.04	118.36	111.11
1	C	4	ILE	N-CA-C	5.62	115.94	107.80
1	C	146	ASP	N-CA-C	5.59	115.95	108.23
1	C	89	HIS	CA-C-N	5.50	125.15	119.05
1	C	89	HIS	C-N-CA	5.50	125.15	119.05
1	B	147	VAL	N-CA-C	5.33	117.30	108.99
1	C	88	LEU	N-CA-C	5.28	117.95	111.82
1	A	105	SER	CA-C-N	5.28	125.34	119.32
1	A	105	SER	C-N-CA	5.28	125.34	119.32
1	B	88	LEU	N-CA-C	5.25	118.14	111.69

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	1359	0	1365	48	0
1	B	1594	0	1603	24	0
1	C	1594	0	1603	41	0
2	A	15	0	0	0	0
2	B	15	0	0	0	0
2	C	20	0	0	0	0
3	A	107	0	0	0	0
3	B	189	0	0	0	0
3	C	153	0	0	0	0
All	All	5046	0	4571	110	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 12.

All (110) 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:4:ILE:H	1:B:125:ASN:HD22	1.18	0.89
1:A:138:LEU:HD13	1:A:149:ARG:NH1	1.87	0.89
1:C:89:HIS:HB3	1:C:90:PRO:HD3	1.54	0.89
1:C:4:ILE:H	1:C:125:ASN:HD22	1.17	0.87
1:A:4:ILE:H	1:A:125:ASN:HD22	1.22	0.87
1:A:28:ASN:HD21	1:A:104:THR:H	1.30	0.78
1:A:148:THR:HG23	1:A:151:HIS:HB2	1.65	0.78
1:B:150:GLU:O	1:B:154:GLN:HG2	1.84	0.77
1:A:138:LEU:O	1:A:142:MSE:HG3	1.84	0.76
1:C:149:ARG:O	1:C:153:GLU:HG3	1.86	0.76
1:B:139:LYS:HA	1:B:142:MSE:HE2	1.71	0.73
1:C:11:ILE:HD13	1:C:12:GLY:N	2.03	0.73
1:B:76:ALA:O	1:B:78:PRO:HD3	1.89	0.73
1:C:7:LEU:HD11	1:C:130:VAL:CG2	2.20	0.72
1:A:149:ARG:HH11	1:A:149:ARG:HG3	1.54	0.71
1:C:179:PRO:HB2	1:C:182:ILE:HB	1.72	0.70
1:A:99:GLN:HA	1:A:102:GLN:HE21	1.57	0.69
1:C:138:LEU:O	1:C:142:MSE:HG3	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:7:LEU:HD11	1:C:130:VAL:HG23	1.73	0.68
1:A:7:LEU:HD11	1:A:130:VAL:HG23	1.76	0.67
1:A:82:ASN:HA	1:A:85:ASN:HD22	1.62	0.65
1:A:30:ILE:HD12	1:A:100:ILE:HG12	1.79	0.64
1:B:77:ASN:O	1:B:81:LYS:HG3	1.98	0.64
1:A:28:ASN:ND2	1:A:104:THR:H	1.95	0.64
1:A:89:HIS:HB2	1:A:90:PRO:HD3	1.80	0.63
1:A:7:LEU:HD11	1:A:130:VAL:CG2	2.30	0.62
1:A:139:LYS:HA	1:A:142:MSE:HE3	1.82	0.62
1:A:148:THR:HG23	1:A:151:HIS:CB	2.30	0.62
1:C:139:LYS:HA	1:C:142:MSE:HE2	1.80	0.62
1:A:139:LYS:HA	1:A:142:MSE:CE	2.31	0.61
1:A:85:ASN:HB3	1:A:89:HIS:NE2	2.16	0.60
1:C:4:ILE:H	1:C:125:ASN:ND2	1.96	0.60
1:A:150:GLU:O	1:A:154:GLN:HG2	2.03	0.58
1:C:175:ASN:ND2	1:C:182:ILE:HD11	2.18	0.58
1:A:28:ASN:HD22	1:A:103:ALA:HB1	1.69	0.58
1:C:11:ILE:HD12	1:C:141:THR:OG1	2.03	0.57
1:A:138:LEU:HD13	1:A:149:ARG:HH12	1.67	0.57
1:A:28:ASN:ND2	1:A:105:SER:H	2.03	0.57
1:A:2:ARG:NH2	1:A:103:ALA:O	2.38	0.56
1:A:147:VAL:HG12	1:A:148:THR:H	1.70	0.56
1:C:70:LEU:O	1:C:74:ILE:HG13	2.06	0.56
1:A:28:ASN:HD21	1:A:104:THR:N	2.01	0.56
1:C:195:GLN:O	1:C:199:GLN:HG2	2.05	0.56
1:C:11:ILE:HD11	1:C:144:ARG:NH2	2.21	0.56
1:C:116:VAL:HG21	1:C:165:ARG:HG2	1.88	0.55
1:B:138:LEU:O	1:B:142:MSE:HG3	2.07	0.55
1:C:179:PRO:HB3	1:C:182:ILE:HD12	1.88	0.55
1:C:188:ARG:HG3	1:C:188:ARG:HH11	1.70	0.55
1:C:139:LYS:HA	1:C:142:MSE:CE	2.38	0.54
1:B:74:ILE:HB	1:B:81:LYS:HG2	1.88	0.54
1:B:97:GLN:HE21	1:B:101:GLN:HE21	1.55	0.54
1:A:147:VAL:HG11	1:A:151:HIS:CD2	2.44	0.53
1:C:178:ALA:HB3	1:C:180:ASP:N	2.22	0.53
1:A:94:GLN:HG2	1:A:95:GLU:N	2.21	0.53
1:B:153:GLU:OE2	1:C:1:MSE:HG3	2.08	0.53
1:A:79:GLU:O	1:A:83:TRP:HD1	1.92	0.52
1:C:150:GLU:O	1:C:151:HIS:HB3	2.10	0.52
1:B:7:LEU:HD11	1:B:130:VAL:CG2	2.41	0.51
1:A:138:LEU:C	1:A:142:MSE:HE2	2.36	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:LEU:HD11	1:B:130:VAL:HG23	1.93	0.51
1:B:80:GLU:HA	1:B:80:GLU:OE1	2.09	0.51
1:A:28:ASN:ND2	1:A:104:THR:N	2.57	0.51
1:A:28:ASN:ND2	1:A:103:ALA:HB1	2.26	0.51
1:A:37:ARG:HH11	1:A:37:ARG:HG3	1.76	0.50
1:B:75:PHE:O	1:B:76:ALA:HB3	2.12	0.50
1:A:138:LEU:CD1	1:A:149:ARG:NH1	2.70	0.49
1:B:123:LYS:HZ2	1:B:123:LYS:HB3	1.77	0.49
1:B:116:VAL:HG21	1:B:165:ARG:HG2	1.95	0.48
1:B:179:PRO:O	1:B:182:ILE:HG12	2.13	0.48
1:A:138:LEU:CD1	1:A:149:ARG:HH12	2.25	0.48
1:B:153:GLU:CD	1:C:1:MSE:HG3	2.38	0.48
1:C:119:SER:HA	1:C:121:TYR:CE2	2.48	0.48
1:C:142:MSE:SE	1:C:148:THR:C	3.07	0.48
1:C:68:ARG:O	1:C:72:GLU:HG3	2.14	0.47
1:A:147:VAL:HG12	1:A:148:THR:N	2.29	0.47
1:B:97:GLN:HE21	1:B:101:GLN:NE2	2.13	0.47
1:C:11:ILE:HD11	1:C:144:ARG:HH22	1.80	0.47
1:C:188:ARG:HG3	1:C:188:ARG:NH1	2.29	0.46
1:A:78:PRO:O	1:A:79:GLU:HB2	2.14	0.46
1:C:74:ILE:HD11	1:C:84:LEU:CD2	2.46	0.46
1:A:35:ILE:O	1:A:39:VAL:HG23	2.15	0.46
1:B:97:GLN:OE1	1:B:97:GLN:HA	2.16	0.46
1:A:149:ARG:HH11	1:A:149:ARG:CG	2.25	0.45
1:C:179:PRO:HB2	1:C:182:ILE:CB	2.43	0.45
1:C:80:GLU:HA	1:C:80:GLU:OE1	2.17	0.44
1:A:80:GLU:HA	1:A:83:TRP:CD1	2.52	0.44
1:A:173:ILE:HD11	1:A:182:ILE:HB	1.98	0.44
1:B:28:ASN:HD22	1:B:105:SER:H	1.65	0.43
1:C:138:LEU:HD21	1:C:153:GLU:HG2	2.00	0.43
1:A:83:TRP:CD1	1:A:83:TRP:H	2.36	0.43
1:B:185:ASP:OD1	1:B:188:ARG:NH2	2.49	0.43
1:C:139:LYS:CB	1:C:139:LYS:NZ	2.82	0.43
1:A:37:ARG:HG3	1:A:37:ARG:NH1	2.34	0.42
1:C:178:ALA:N	1:C:179:PRO:HA	2.35	0.42
1:A:85:ASN:C	1:A:87:LEU:N	2.77	0.42
1:A:98:HIS:O	1:A:101:GLN:HG2	2.20	0.42
1:C:11:ILE:HD13	1:C:11:ILE:C	2.44	0.42
1:A:149:ARG:NH1	1:A:149:ARG:HG3	2.27	0.42
1:B:28:ASN:ND2	1:B:104:THR:H	2.17	0.42
1:C:2:ARG:HA	1:C:2:ARG:HE	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:152:VAL:O	1:C:155:ILE:HB	2.21	0.41
1:C:179:PRO:CB	1:C:182:ILE:HD12	2.49	0.41
1:C:178:ALA:H	1:C:179:PRO:HA	1.86	0.41
1:A:166:LEU:HD13	1:B:206:PRO:HB3	2.03	0.41
1:C:11:ILE:HD13	1:C:12:GLY:H	1.83	0.41
1:C:76:ALA:O	1:C:77:ASN:HB2	2.20	0.41
1:B:4:ILE:H	1:B:125:ASN:ND2	2.00	0.40
1:A:38:GLN:HA	1:A:41:GLU:HG3	2.03	0.40
1:A:179:PRO:O	1:A:182:ILE:HG12	2.21	0.40
1:A:149:ARG:NH1	1:A:149:ARG:CG	2.84	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	172/216 (80%)	166 (96%)	6 (4%)	0	100	100
1	B	204/216 (94%)	202 (99%)	2 (1%)	0	100	100
1	C	204/216 (94%)	191 (94%)	9 (4%)	4 (2%)	6	1
All	All	580/648 (90%)	559 (96%)	17 (3%)	4 (1%)	18	8

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	148	THR
1	C	177	GLY
1	C	77	ASN
1	C	179	PRO

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	145/173 (84%)	142 (98%)	3 (2%)	47	36
1	B	167/173 (96%)	165 (99%)	2 (1%)	63	57
1	C	167/173 (96%)	162 (97%)	5 (3%)	36	24
All	All	479/519 (92%)	469 (98%)	10 (2%)	47	36

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94	GLN
1	A	148	THR
1	A	180	ASP
1	B	2	ARG
1	B	123	LYS
1	C	11	ILE
1	C	52	ASP
1	C	79	GLU
1	C	139	LYS
1	C	150	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	38	GLN
1	A	85	ASN
1	A	97	GLN
1	A	102	GLN
1	A	125	ASN
1	A	137	GLN
1	A	151	HIS
1	A	175	ASN
1	B	28	ASN
1	B	82	ASN

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Mol	Chain	Res	Type
1	B	101	GLN
1	B	125	ASN
1	B	137	GLN
1	B	151	HIS
1	B	175	ASN
1	B	192	HIS
1	B	195	GLN
1	B	199	GLN
1	C	57	ASN
1	C	77	ASN
1	C	85	ASN
1	C	94	GLN
1	C	125	ASN
1	C	137	GLN
1	C	175	ASN
1	C	199	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](#)

10 ligands are modelled in this entry.

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	$\# Z > 2$
2	SO4	B	509	-	4,4,4	0.36	0	6,6,6	0.14	0
2	SO4	A	505	-	4,4,4	0.38	0	6,6,6	0.14	0
2	SO4	A	501	-	4,4,4	0.29	0	6,6,6	0.08	0
2	SO4	C	508	-	4,4,4	0.38	0	6,6,6	0.09	0
2	SO4	C	507	-	4,4,4	0.36	0	6,6,6	0.07	0
2	SO4	B	506	-	4,4,4	0.38	0	6,6,6	0.09	0
2	SO4	A	510	-	4,4,4	0.37	0	6,6,6	0.11	0
2	SO4	C	503	-	4,4,4	0.34	0	6,6,6	0.26	0
2	SO4	C	504	-	4,4,4	0.40	0	6,6,6	0.07	0
2	SO4	B	502	-	4,4,4	0.36	0	6,6,6	0.22	0

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 ⓘ

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	174/216 (80%)	0.83	38 (21%) 2 2	12, 23, 69, 89	0
1	B	203/216 (93%)	0.19	5 (2%) 58 58	10, 21, 40, 50	0
1	C	203/216 (93%)	0.78	36 (17%) 4 3	12, 24, 67, 76	0
All	All	580/648 (89%)	0.59	79 (13%) 7 5	10, 22, 65, 89	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	83	TRP	6.7
1	C	75	PHE	6.7
1	C	178	ALA	5.8
1	C	147	VAL	5.8
1	C	76	ALA	5.6
1	C	148	THR	5.6
1	A	47	LEU	5.3
1	A	43	GLY	4.9
1	A	84	LEU	4.8
1	A	87	LEU	4.8
1	C	74	ILE	4.7
1	A	78	PRO	4.7
1	A	49	ALA	4.5
1	B	146	ASP	4.4
1	C	180	ASP	4.1
1	A	88	LEU	4.0
1	C	151	HIS	3.9
1	A	48	HIS	3.9
1	A	147	VAL	3.8
1	A	86	ALA	3.8
1	A	40	VAL	3.8
1	C	179	PRO	3.7
1	C	77	ASN	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	76	ALA	3.5
1	A	82	ASN	3.5
1	A	204	GLU	3.5
1	C	145	ASP	3.5
1	A	42	PRO	3.4
1	C	152	VAL	3.4
1	A	46	ALA	3.3
1	C	141	THR	3.3
1	A	149	ARG	3.3
1	C	146	ASP	3.2
1	A	79	GLU	3.2
1	A	80	GLU	3.1
1	A	39	VAL	3.1
1	A	45	PRO	3.1
1	C	71	ARG	3.1
1	C	206	PRO	3.1
1	C	150	GLU	3.0
1	C	72	GLU	2.9
1	A	85	ASN	2.9
1	C	181	ALA	2.9
1	C	182	ILE	2.9
1	C	153	GLU	2.9
1	C	78	PRO	2.8
1	C	149	ARG	2.8
1	A	81	LYS	2.7
1	A	89	HIS	2.7
1	A	203	GLN	2.7
1	A	91	LEU	2.7
1	A	36	ALA	2.6
1	C	80	GLU	2.6
1	B	93	GLN	2.6
1	A	90	PRO	2.6
1	A	146	ASP	2.6
1	C	79	GLU	2.6
1	A	44	ALA	2.6
1	C	69	ALA	2.6
1	A	100	ILE	2.6
1	C	73	ARG	2.5
1	C	143	GLN	2.5
1	C	156	LEU	2.5
1	A	151	HIS	2.5
1	A	41	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	75	PHE	2.4
1	A	102	GLN	2.4
1	A	92	ILE	2.3
1	C	177	GLY	2.3
1	C	205	LYS	2.3
1	C	144	ARG	2.3
1	A	101	GLN	2.3
1	B	147	VAL	2.3
1	A	94	GLN	2.2
1	C	184	SER	2.1
1	C	68	ARG	2.1
1	C	11	ILE	2.1
1	A	148	THR	2.1
1	C	89	HIS	2.1

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	SO4	B	509	5/5	0.73	0.19	84,84,84,85	0
2	SO4	B	506	5/5	0.74	0.17	82,82,82,82	0
2	SO4	A	510	5/5	0.75	0.20	82,82,82,83	0
2	SO4	A	505	5/5	0.76	0.14	77,77,78,78	0
2	SO4	C	508	5/5	0.80	0.21	87,87,88,88	0
2	SO4	C	507	5/5	0.87	0.11	66,67,67,67	0
2	SO4	C	504	5/5	0.87	0.14	45,45,48,49	0
2	SO4	C	503	5/5	0.95	0.08	34,35,36,37	0
2	SO4	A	501	5/5	0.97	0.06	23,24,25,26	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	B	502	5/5	0.98	0.06	18,19,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.