



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 10:24 pm BST

PDB ID : 1O60
Title : Crystal structure of KDO-8-phosphate synthase
Authors : Structural GenomiX
Deposited on : 2003-10-23
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
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

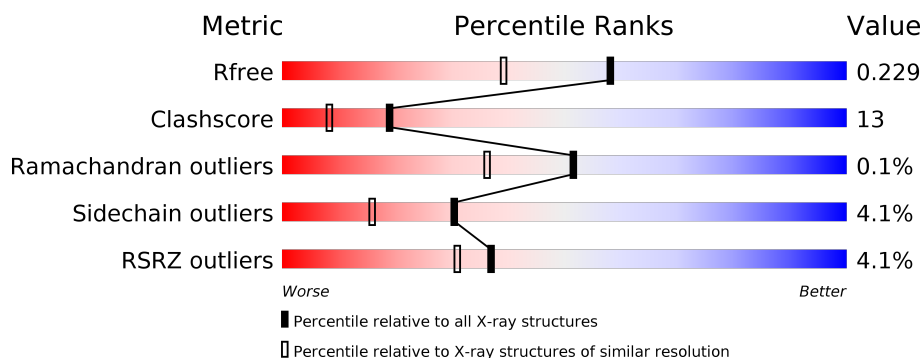
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (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 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	292	<div> <div>0%</div> <div> <div>75%</div> <div>16%</div> <div>8%</div> </div> </div>
1	B	292	<div> <div>4%</div> <div> <div>71%</div> <div>20%</div> <div>8%</div> </div> </div>
1	C	292	<div> <div>5%</div> <div> <div>67%</div> <div>21%</div> <div>8%</div> </div> </div>
1	D	292	<div> <div>5%</div> <div> <div>69%</div> <div>22%</div> <div>8%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8811 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	269	Total	C	N	O	S	Se	0	3	0
			2067	1317	345	391	5	9			
1	B	269	Total	C	N	O	S	Se	0	0	0
			2059	1311	344	390	5	9			
1	C	269	Total	C	N	O	S	Se	0	2	0
			2066	1316	345	391	5	9			
1	D	269	Total	C	N	O	S	Se	0	1	0
			2051	1308	343	386	5	9			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP P45251
A	25	MSE	MET	MODIFIED RESIDUE	UNP P45251
A	33	MSE	MET	MODIFIED RESIDUE	UNP P45251
A	35	MSE	MET	MODIFIED RESIDUE	UNP P45251
A	74	MSE	MET	MODIFIED RESIDUE	UNP P45251
A	128	MSE	MET	MODIFIED RESIDUE	UNP P45251
A	148	MSE	MET	MODIFIED RESIDUE	UNP P45251
A	181	MSE	MET	MODIFIED RESIDUE	UNP P45251
A	187	MSE	MET	MODIFIED RESIDUE	UNP P45251
A	267	MSE	MET	MODIFIED RESIDUE	UNP P45251
A	285	GLY	-	cloning artifact	UNP P45251
A	286	SER	-	cloning artifact	UNP P45251
A	287	HIS	-	cloning artifact	UNP P45251
A	288	HIS	-	cloning artifact	UNP P45251
A	289	HIS	-	cloning artifact	UNP P45251
A	290	HIS	-	cloning artifact	UNP P45251
A	291	HIS	-	cloning artifact	UNP P45251
A	292	HIS	-	cloning artifact	UNP P45251
B	1	MSE	MET	MODIFIED RESIDUE	UNP P45251
B	25	MSE	MET	MODIFIED RESIDUE	UNP P45251
B	33	MSE	MET	MODIFIED RESIDUE	UNP P45251

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Chain	Residue	Modelled	Actual	Comment	Reference
B	35	MSE	MET	MODIFIED RESIDUE	UNP P45251
B	74	MSE	MET	MODIFIED RESIDUE	UNP P45251
B	128	MSE	MET	MODIFIED RESIDUE	UNP P45251
B	148	MSE	MET	MODIFIED RESIDUE	UNP P45251
B	181	MSE	MET	MODIFIED RESIDUE	UNP P45251
B	187	MSE	MET	MODIFIED RESIDUE	UNP P45251
B	267	MSE	MET	MODIFIED RESIDUE	UNP P45251
B	285	GLY	-	cloning artifact	UNP P45251
B	286	SER	-	cloning artifact	UNP P45251
B	287	HIS	-	cloning artifact	UNP P45251
B	288	HIS	-	cloning artifact	UNP P45251
B	289	HIS	-	cloning artifact	UNP P45251
B	290	HIS	-	cloning artifact	UNP P45251
B	291	HIS	-	cloning artifact	UNP P45251
B	292	HIS	-	cloning artifact	UNP P45251
C	1	MSE	MET	MODIFIED RESIDUE	UNP P45251
C	25	MSE	MET	MODIFIED RESIDUE	UNP P45251
C	33	MSE	MET	MODIFIED RESIDUE	UNP P45251
C	35	MSE	MET	MODIFIED RESIDUE	UNP P45251
C	74	MSE	MET	MODIFIED RESIDUE	UNP P45251
C	128	MSE	MET	MODIFIED RESIDUE	UNP P45251
C	148	MSE	MET	MODIFIED RESIDUE	UNP P45251
C	181	MSE	MET	MODIFIED RESIDUE	UNP P45251
C	187	MSE	MET	MODIFIED RESIDUE	UNP P45251
C	267	MSE	MET	MODIFIED RESIDUE	UNP P45251
C	285	GLY	-	cloning artifact	UNP P45251
C	286	SER	-	cloning artifact	UNP P45251
C	287	HIS	-	cloning artifact	UNP P45251
C	288	HIS	-	cloning artifact	UNP P45251
C	289	HIS	-	cloning artifact	UNP P45251
C	290	HIS	-	cloning artifact	UNP P45251
C	291	HIS	-	cloning artifact	UNP P45251
C	292	HIS	-	cloning artifact	UNP P45251
D	1	MSE	MET	MODIFIED RESIDUE	UNP P45251
D	25	MSE	MET	MODIFIED RESIDUE	UNP P45251
D	33	MSE	MET	MODIFIED RESIDUE	UNP P45251
D	35	MSE	MET	MODIFIED RESIDUE	UNP P45251
D	74	MSE	MET	MODIFIED RESIDUE	UNP P45251
D	128	MSE	MET	MODIFIED RESIDUE	UNP P45251
D	148	MSE	MET	MODIFIED RESIDUE	UNP P45251
D	181	MSE	MET	MODIFIED RESIDUE	UNP P45251
D	187	MSE	MET	MODIFIED RESIDUE	UNP P45251

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Chain	Residue	Modelled	Actual	Comment	Reference
D	267	MSE	MET	MODIFIED RESIDUE	UNP P45251
D	285	GLY	-	cloning artifact	UNP P45251
D	286	SER	-	cloning artifact	UNP P45251
D	287	HIS	-	cloning artifact	UNP P45251
D	288	HIS	-	cloning artifact	UNP P45251
D	289	HIS	-	cloning artifact	UNP P45251
D	290	HIS	-	cloning artifact	UNP P45251
D	291	HIS	-	cloning artifact	UNP P45251
D	292	HIS	-	cloning artifact	UNP P45251

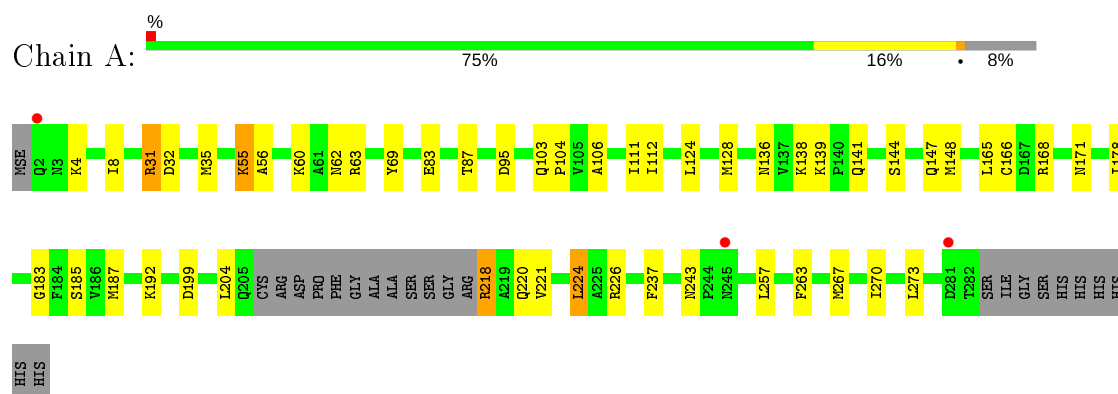
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	168	Total	O	0	0
			168	168		
2	B	137	Total	O	0	0
			137	137		
2	C	143	Total	O	0	0
			143	143		
2	D	120	Total	O	0	0
			120	120		

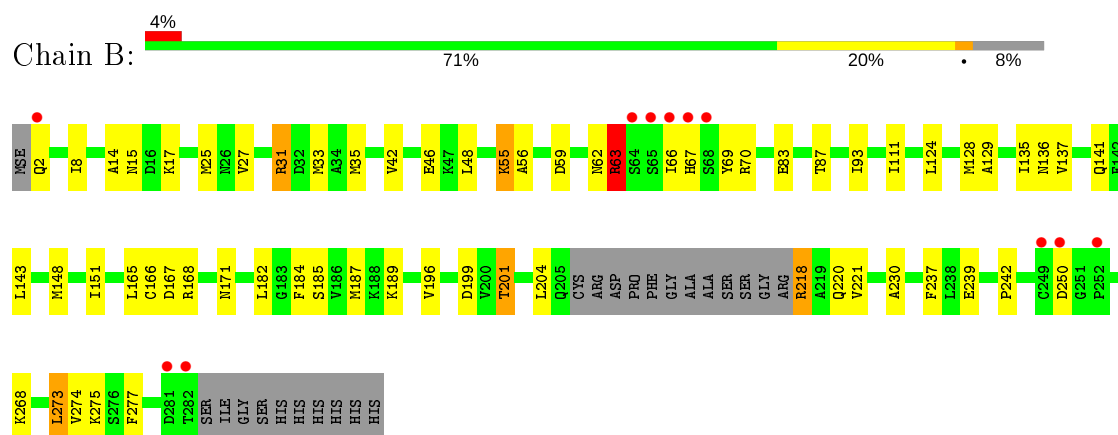
3 Residue-property plots

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.

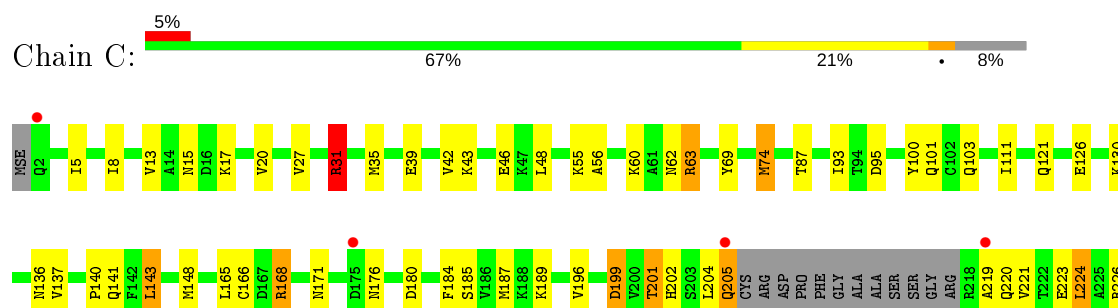
- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase

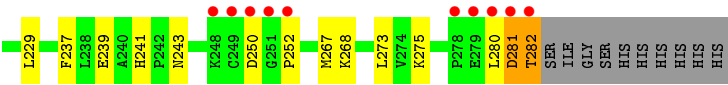


- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase

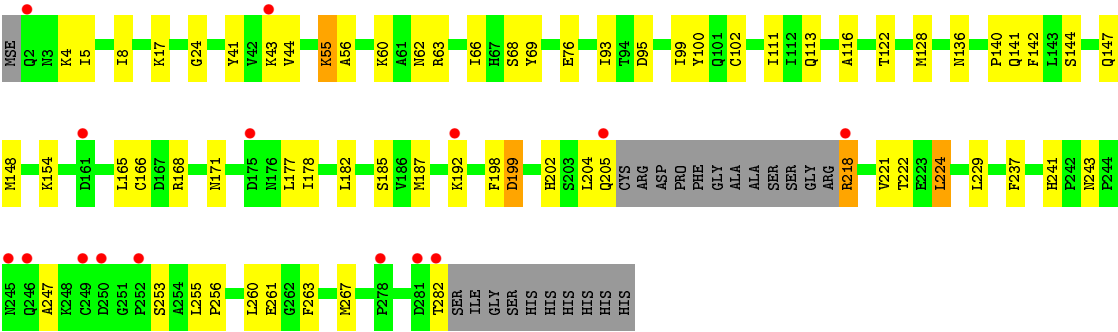


- Molecule 1: 2-dehydro-3-deoxyphosphooctonate aldolase





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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.97Å 92.89Å 82.30Å 90.00° 116.82° 90.00°	Depositor
Resolution (Å)	39.22 – 1.80 38.78 – 1.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (39.22-1.80) 97.5 (38.78-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.23 (at 1.81Å)	Xtriage
Refinement program	REFMAC 4.0	Depositor
R, R_{free}	0.207 , 0.242 0.197 , 0.229	Depositor DCC
R_{free} test set	9540 reflections (9.97%)	wwPDB-VP
Wilson B-factor (Å ²)	20.9	Xtriage
Anisotropy	0.334	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 56.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.017 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8811	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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	0.58	0/2103	1.08	4/2822 (0.1%)
1	B	0.56	0/2084	0.98	2/2798 (0.1%)
1	C	0.57	0/2099	1.07	8/2818 (0.3%)
1	D	0.54	0/2080	0.95	4/2793 (0.1%)
All	All	0.56	0/8366	1.02	18/11231 (0.2%)

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

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	226	ARG	NE-CZ-NH2	17.49	129.04	120.30
1	A	226	ARG	NE-CZ-NH1	-15.15	112.73	120.30
1	C	226	ARG	NE-CZ-NH2	14.00	127.30	120.30
1	C	226	ARG	NE-CZ-NH1	-12.74	113.93	120.30
1	C	226	ARG	CD-NE-CZ	12.17	140.64	123.60
1	A	95	ASP	CB-CG-OD1	6.72	124.34	118.30
1	C	31	ARG	CA-CB-CG	6.54	127.79	113.40
1	A	226	ARG	CD-NE-CZ	6.42	132.59	123.60
1	D	199	ASP	CB-CG-OD1	6.41	124.07	118.30
1	C	95	ASP	CB-CG-OD1	6.14	123.83	118.30
1	C	199	ASP	CB-CG-OD1	5.80	123.52	118.30
1	D	55	LYS	CB-CG-CD	5.36	125.52	111.60
1	C	31	ARG	NE-CZ-NH1	-5.34	117.63	120.30
1	B	63	ARG	CD-NE-CZ	5.28	130.99	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	199	ASP	CB-CG-OD2	-5.27	113.56	118.30
1	D	95	ASP	CB-CG-OD1	5.12	122.90	118.30
1	B	63	ARG	NE-CZ-NH1	-5.09	117.75	120.30
1	D	199	ASP	CB-CG-OD2	-5.06	113.75	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	252	PRO	Mainchain

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	2067	0	2091	43	0
1	B	2059	0	2089	49	0
1	C	2066	0	2093	71	0
1	D	2051	0	2080	66	0
2	A	168	0	0	3	0
2	B	137	0	0	1	0
2	C	143	0	0	2	0
2	D	120	0	0	1	0
All	All	8811	0	8353	215	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 13.

All (215) 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:220[B]:GLN:NE2	1:C:223[B]:GLU:CD	2.03	1.12
1:C:220[B]:GLN:NE2	1:C:223[B]:GLU:OE1	1.82	1.09
1:B:137:VAL:HG11	1:B:151:ILE:HD12	1.45	0.98
1:A:165:LEU:HB3	1:A:187:MSE:HE1	1.51	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:148:MSE:HE2	1:C:165:LEU:HD22	1.52	0.91
1:D:44:VAL:HG22	1:D:261:GLU:HG3	1.52	0.91
1:C:148:MSE:HG3	1:C:187:MSE:HE3	1.54	0.90
1:B:141:GLN:HG3	1:B:171:ASN:HD21	1.38	0.88
1:A:148:MSE:HG3	1:A:187:MSE:HE3	1.54	0.88
1:D:148:MSE:HG3	1:D:187:MSE:HE3	1.56	0.87
1:B:201:THR:HG21	1:B:239:GLU:H	1.42	0.84
1:D:141:GLN:HG3	1:D:171:ASN:HD21	1.42	0.84
1:C:137:VAL:HG11	1:C:148:MSE:HE1	1.59	0.83
1:C:220[B]:GLN:NE2	1:C:223[B]:GLU:OE2	2.08	0.83
1:C:201:THR:HG21	1:C:239:GLU:H	1.44	0.81
1:C:220[B]:GLN:OE1	1:D:185[B]:SER:OG	1.99	0.81
1:A:165:LEU:HB3	1:A:187:MSE:CE	2.11	0.80
1:C:165:LEU:HB3	1:C:187:MSE:HE1	1.63	0.79
1:B:35:MSE:HE1	1:B:87:THR:HG21	1.64	0.78
1:C:241:HIS:HD2	1:C:243:ASN:H	1.27	0.78
1:A:55:LYS:HD2	1:A:56:ALA:N	1.99	0.78
1:D:55:LYS:HD2	1:D:56:ALA:N	1.99	0.77
1:D:165:LEU:HB3	1:D:187:MSE:HE1	1.68	0.76
1:B:141:GLN:HG3	1:B:171:ASN:ND2	2.02	0.74
1:C:148:MSE:CG	1:C:187:MSE:HE3	2.19	0.73
1:D:165:LEU:HD13	1:D:187:MSE:HE2	1.70	0.73
1:D:141:GLN:HG3	1:D:171:ASN:ND2	2.04	0.72
1:D:241:HIS:HD2	1:D:243:ASN:H	1.37	0.72
1:C:140:PRO:HG2	1:C:143:LEU:HD23	1.71	0.72
1:C:74:MSE:CE	1:C:101:GLN:HG2	2.19	0.72
1:B:204:LEU:CD1	1:B:221:VAL:HG22	2.20	0.71
1:D:165:LEU:HB3	1:D:187:MSE:CE	2.21	0.71
1:A:141:GLN:OE1	1:A:171:ASN:ND2	2.22	0.71
1:A:185:SER:HB2	2:A:363:HOH:O	1.91	0.70
1:D:148:MSE:HE3	1:D:165:LEU:HD22	1.73	0.68
1:B:204:LEU:HD13	1:B:221:VAL:HG22	1.76	0.68
1:C:74:MSE:HE1	1:C:100:TYR:CE1	2.28	0.68
1:D:148:MSE:CG	1:D:187:MSE:HE3	2.23	0.68
1:B:148:MSE:O	1:B:151:ILE:HG12	1.94	0.67
1:B:25:MSE:HE1	1:B:33:MSE:HG2	1.76	0.67
1:A:148:MSE:CG	1:A:187:MSE:HE3	2.24	0.67
1:B:148:MSE:HE3	1:B:165:LEU:HD22	1.76	0.66
1:A:165:LEU:HD13	1:A:187:MSE:HE2	1.78	0.66
1:C:165:LEU:HB3	1:C:187:MSE:CE	2.24	0.66
1:A:139:LYS:HB2	1:A:148:MSE:HE2	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:144:SER:OG	1:D:147:GLN:HG2	1.95	0.65
1:A:204:LEU:CD1	1:A:221:VAL:HG22	2.27	0.65
1:D:204:LEU:HD12	1:D:253:SER:HB2	1.77	0.65
1:B:8:ILE:HD13	1:B:111:ILE:HD11	1.78	0.65
1:B:148:MSE:SE	1:B:151:ILE:HD11	2.46	0.65
1:C:224:LEU:HG	1:D:182:LEU:HD21	1.79	0.65
1:D:218:ARG:O	1:D:221:VAL:HG23	1.96	0.64
1:B:184:PHE:HD1	1:B:196:VAL:HG11	1.63	0.64
1:C:220[B]:GLN:HE22	1:C:223[B]:GLU:CD	1.84	0.64
1:C:184:PHE:HD1	1:C:196:VAL:HG11	1.61	0.64
1:A:8:ILE:HD13	1:A:111:ILE:HD11	1.80	0.63
1:C:141:GLN:OE1	1:C:171:ASN:ND2	2.32	0.62
1:D:199:ASP:OD2	1:D:202:HIS:HD2	1.82	0.62
1:A:220[B]:GLN:NE2	1:B:185:SER:OG	2.33	0.62
1:C:165:LEU:HD13	1:C:187:MSE:HE2	1.80	0.62
1:C:280:LEU:HD23	1:D:222:THR:HG21	1.82	0.62
1:C:74:MSE:HE2	1:C:101:GLN:HG2	1.82	0.61
1:C:5:ILE:HD13	1:C:17:LYS:NZ	2.15	0.61
1:A:35:MSE:HE1	1:A:87:THR:HG21	1.83	0.61
1:A:63:ARG:HD3	1:A:69:TYR:O	2.01	0.61
1:C:241:HIS:CD2	1:C:243:ASN:H	2.15	0.60
1:D:148:MSE:HE2	1:D:187:MSE:HE1	1.83	0.60
1:C:55:LYS:HD2	1:C:56:ALA:N	2.15	0.60
1:C:63:ARG:HD3	1:C:69:TYR:O	2.02	0.60
1:D:8:ILE:HD13	1:D:111:ILE:HD11	1.84	0.60
1:B:137:VAL:HG11	1:B:151:ILE:CD1	2.27	0.59
1:B:148:MSE:SE	1:B:151:ILE:CD1	3.00	0.59
1:B:199:ASP:HA	1:B:237:PHE:HB3	1.83	0.59
1:C:280:LEU:HD23	1:D:222:THR:CG2	2.32	0.59
1:C:241:HIS:HD2	1:C:243:ASN:N	1.99	0.59
1:D:136:ASN:HD21	1:D:166:CYS:HB2	1.68	0.58
1:C:137:VAL:CG1	1:C:148:MSE:HE1	2.32	0.58
1:A:141:GLN:HG3	1:A:171:ASN:OD1	2.03	0.57
1:D:60:LYS:HE2	1:D:60:LYS:HA	1.87	0.57
1:D:204:LEU:CD1	1:D:221:VAL:HG22	2.34	0.57
1:A:218:ARG:O	1:A:221:VAL:HG23	2.05	0.57
1:D:218:ARG:N	1:D:253:SER:HA	2.19	0.57
1:C:280:LEU:O	1:C:281:ASP:C	2.43	0.57
1:C:204:LEU:CD1	1:C:221:VAL:HG22	2.35	0.57
1:C:126:GLU:HG2	1:C:130:LYS:HE3	1.87	0.57
1:A:183:GLY:HA3	2:A:316:HOH:O	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:270:ILE:HD13	1:B:273:LEU:HD13	1.86	0.56
1:D:205:GLN:HB2	2:D:350:HOH:O	2.04	0.56
1:B:33:MSE:HE2	1:B:242:PRO:HG2	1.87	0.55
1:C:199:ASP:OD2	1:C:202:HIS:HD2	1.89	0.55
1:C:31:ARG:C	1:C:31:ARG:HD3	2.26	0.55
1:A:139:LYS:HE2	2:A:319:HOH:O	2.07	0.54
1:C:220[B]:GLN:HE21	1:C:223[B]:GLU:CD	2.00	0.54
1:B:31:ARG:HD2	1:B:83:GLU:OE1	2.07	0.54
1:A:204:LEU:HD13	1:A:221:VAL:HG22	1.89	0.54
1:B:136:ASN:HD21	1:B:166:CYS:HB2	1.72	0.54
1:B:184:PHE:HD1	1:B:196:VAL:CG1	2.21	0.54
1:B:59:ASP:OD1	1:B:70:ARG:HD3	2.08	0.53
1:D:148:MSE:HG3	1:D:187:MSE:CE	2.34	0.53
1:C:5:ILE:HD13	1:C:17:LYS:HZ2	1.72	0.53
1:C:180:ASP:CG	1:D:178:ILE:HD12	2.29	0.53
1:C:168:ARG:HH11	1:C:168:ARG:HG2	1.74	0.53
1:C:199:ASP:HA	1:C:237:PHE:HB3	1.91	0.53
1:A:199:ASP:HA	1:A:237:PHE:HB3	1.91	0.52
1:C:8:ILE:HD13	1:C:111:ILE:HD11	1.91	0.52
1:C:136:ASN:HD21	1:C:166:CYS:HB2	1.74	0.52
1:D:204:LEU:HD13	1:D:221:VAL:HG22	1.92	0.52
1:D:5:ILE:HD13	1:D:17:LYS:NZ	2.25	0.52
1:A:148:MSE:HG3	1:A:187:MSE:CE	2.32	0.52
1:B:201:THR:HG21	1:B:239:GLU:N	2.20	0.52
1:B:63:ARG:HE	1:B:63:ARG:HA	1.74	0.52
1:B:141:GLN:NE2	1:B:171:ASN:OD1	2.28	0.51
1:B:218:ARG:O	1:B:221:VAL:HG23	2.10	0.51
1:B:55:LYS:HB2	1:B:93:ILE:CG2	2.40	0.51
1:C:148:MSE:HG3	1:C:187:MSE:CE	2.36	0.51
1:C:74:MSE:HE1	1:C:101:GLN:HG2	1.90	0.51
1:B:66:ILE:HG22	1:D:122:THR:HG22	1.91	0.51
1:A:165:LEU:HD13	1:A:187:MSE:CE	2.39	0.51
1:A:112:ILE:HG22	1:A:128:MSE:HE3	1.91	0.51
1:C:35:MSE:HE1	1:C:87:THR:HG21	1.92	0.50
1:C:184:PHE:HD1	1:C:196:VAL:CG1	2.25	0.50
1:C:27:VAL:HG22	1:C:60:LYS:HE3	1.94	0.50
1:D:177:LEU:O	1:D:178:ILE:HD13	2.11	0.50
1:D:241:HIS:CD2	1:D:243:ASN:H	2.24	0.50
1:A:31:ARG:NH2	1:A:83:GLU:OE2	2.45	0.50
1:C:201:THR:HG21	1:C:239:GLU:N	2.20	0.50
1:C:185:SER:O	1:C:189:LYS:HG3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:LYS:NZ	1:A:168:ARG:HH11	2.11	0.49
1:A:55:LYS:C	1:A:55:LYS:HD2	2.32	0.49
1:A:106:ALA:HB2	1:A:112:ILE:HD11	1.93	0.49
1:B:14:ALA:HB3	1:B:17:LYS:HG3	1.95	0.49
1:D:63:ARG:HD2	1:D:69:TYR:O	2.13	0.49
1:D:199:ASP:HA	1:D:237:PHE:HB3	1.93	0.49
1:D:241:HIS:CE1	1:D:247:ALA:HA	2.47	0.49
1:D:43:LYS:NZ	1:D:43:LYS:HB2	2.28	0.48
1:B:42:VAL:O	1:B:46:GLU:HG2	2.14	0.48
1:B:124:LEU:O	1:B:128:MSE:HG3	2.12	0.48
1:B:8:ILE:CD1	1:B:111:ILE:HD11	2.43	0.48
1:C:148:MSE:CB	1:C:187:MSE:HE3	2.44	0.48
1:B:63:ARG:HD3	1:B:69:TYR:O	2.14	0.47
1:D:99:ILE:HG13	1:D:100:TYR:N	2.28	0.47
1:D:241:HIS:HD2	1:D:243:ASN:N	2.10	0.47
1:C:141:GLN:HG3	1:C:171:ASN:OD1	2.15	0.47
1:C:55:LYS:HB2	1:C:93:ILE:HG23	1.95	0.47
1:D:204:LEU:O	1:D:205:GLN:C	2.54	0.47
1:B:55:LYS:HB2	1:B:93:ILE:HG23	1.97	0.46
1:A:31:ARG:HG3	1:A:32:ASP:N	2.29	0.46
1:D:165:LEU:CD1	1:D:187:MSE:HE2	2.43	0.46
1:D:148:MSE:CB	1:D:187:MSE:HE3	2.45	0.46
1:D:55:LYS:HB2	1:D:93:ILE:CG2	2.45	0.46
1:C:42:VAL:O	1:C:46:GLU:HB2	2.16	0.46
1:A:136:ASN:HD21	1:A:166:CYS:HB2	1.80	0.46
1:C:224:LEU:HG	1:D:182:LEU:CD2	2.45	0.46
1:B:129:ALA:HA	1:B:135:ILE:HD11	1.97	0.46
1:C:8:ILE:CD1	1:C:111:ILE:HD11	2.45	0.46
1:A:144:SER:H	1:A:147[B]:GLN:HE21	1.63	0.46
1:A:144:SER:OG	1:A:147[B]:GLN:HG3	2.17	0.45
1:B:167:ASP:HB2	1:B:187:MSE:HE1	1.98	0.45
1:B:63:ARG:NE	1:B:63:ARG:HA	2.31	0.45
1:D:199:ASP:OD2	1:D:202:HIS:CD2	2.66	0.45
1:C:219:ALA:HA	1:D:282:THR:HG23	1.98	0.45
1:C:48:LEU:O	1:C:268:LYS:HE3	2.16	0.45
1:B:66:ILE:HG13	1:B:67:HIS:CD2	2.52	0.45
1:D:41:TYR:CD1	1:D:260:LEU:HD21	2.52	0.45
1:A:224:LEU:HG	1:B:182:LEU:HD21	1.99	0.45
1:B:55:LYS:HD2	1:B:56:ALA:N	2.32	0.45
1:C:199:ASP:OD2	1:C:202:HIS:CD2	2.70	0.45
1:B:218:ARG:HG3	1:B:220:GLN:H	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:165:LEU:HB3	1:D:187:MSE:HE2	1.99	0.45
1:D:4:LYS:HE3	1:D:192:LYS:O	2.16	0.45
1:B:273:LEU:HD22	1:B:277:PHE:CZ	2.53	0.44
1:C:202:HIS:HE1	2:C:435:HOH:O	2.00	0.44
1:A:62:ASN:OD1	1:C:121:GLN:NE2	2.50	0.44
1:C:165:LEU:CD1	1:C:187:MSE:HE2	2.48	0.44
1:D:63:ARG:HD3	1:D:68:SER:HB3	1.98	0.44
1:D:165:LEU:HD13	1:D:187:MSE:CE	2.43	0.44
1:A:165:LEU:CD1	1:A:187:MSE:HE2	2.46	0.44
1:C:165:LEU:HD13	1:C:187:MSE:CE	2.46	0.44
1:D:122:THR:HG23	1:D:154:LYS:CE	2.48	0.44
1:C:204:LEU:O	1:C:205:GLN:C	2.56	0.43
1:B:171:ASN:ND2	1:D:142:PHE:HB2	2.33	0.43
1:A:204:LEU:HD11	1:A:221:VAL:HG22	2.01	0.43
1:D:5:ILE:HD13	1:D:17:LYS:HZ1	1.83	0.43
1:A:8:ILE:CD1	1:A:111:ILE:HD11	2.48	0.43
1:A:263:PHE:O	1:A:267:MSE:HG2	2.18	0.43
1:C:60:LYS:HE2	2:C:340:HOH:O	2.19	0.43
1:B:63:ARG:CA	1:B:63:ARG:HE	2.32	0.42
1:C:15:ASN:O	1:C:275:LYS:HE2	2.20	0.42
1:D:198:PHE:HZ	1:D:224:LEU:HD13	1.84	0.42
1:A:124:LEU:O	1:A:128:MSE:HG3	2.19	0.42
1:A:4:LYS:HE3	1:A:192:LYS:O	2.19	0.42
2:B:373:HOH:O	1:D:66:ILE:HG23	2.18	0.42
1:B:230:ALA:HA	1:B:274:VAL:HG21	2.02	0.42
1:D:141:GLN:NE2	1:D:171:ASN:OD1	2.53	0.42
1:D:116:ALA:HB1	1:D:140:PRO:HA	2.02	0.42
1:B:15:ASN:O	1:B:275:LYS:HE2	2.20	0.42
1:D:204:LEU:CD1	1:D:253:SER:HB2	2.46	0.42
1:C:39:GLU:HG2	1:C:43:LYS:NZ	2.35	0.41
1:A:103:GLN:N	1:A:104:PRO:HD2	2.35	0.41
1:D:113:GLN:HE22	1:D:237:PHE:HE2	1.66	0.41
1:D:255:LEU:HA	1:D:256:PRO:HD3	1.94	0.41
1:D:24:GLY:HA3	1:D:55:LYS:O	2.20	0.41
1:A:178:ILE:HD13	1:A:178:ILE:HG21	1.91	0.41
1:C:176:ASN:HA	1:C:205:GLN:HG3	2.03	0.41
1:D:55:LYS:HB2	1:D:93:ILE:HG23	2.02	0.41
1:C:229:LEU:CD1	1:C:267:MSE:HB3	2.51	0.40
1:C:280:LEU:O	1:C:282:THR:N	2.54	0.40
1:D:229:LEU:CD1	1:D:267:MSE:HB3	2.51	0.40
1:B:48:LEU:O	1:B:268:LYS:HE3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:25:MSE:HE2	1:B:27:VAL:O	2.21	0.40
1:C:13:VAL:HG22	1:C:20:VAL:HG21	2.02	0.40
1:A:60:LYS:HE2	1:A:60:LYS:HA	2.02	0.40
1:D:102:CYS:SG	1:D:128:MSE:HG2	2.61	0.40
1:C:280:LEU:CD2	1:D:263:PHE:HA	2.50	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	268/292 (92%)	262 (98%)	6 (2%)	0	100	100
1	B	265/292 (91%)	257 (97%)	8 (3%)	0	100	100
1	C	267/292 (91%)	260 (97%)	6 (2%)	1 (0%)	34	21
1	D	266/292 (91%)	257 (97%)	9 (3%)	0	100	100
All	All	1066/1168 (91%)	1036 (97%)	29 (3%)	1 (0%)	51	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	281	ASP

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	229/236 (97%)	222 (97%)	7 (3%)	40	25
1	B	228/236 (97%)	216 (95%)	12 (5%)	22	9
1	C	229/236 (97%)	216 (94%)	13 (6%)	20	8
1	D	226/236 (96%)	221 (98%)	5 (2%)	52	39
All	All	912/944 (97%)	875 (96%)	37 (4%)	30	16

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	ARG
1	A	55	LYS
1	A	218	ARG
1	A	224	LEU
1	A	243	ASN
1	A	257	LEU
1	A	273	LEU
1	B	2	GLN
1	B	31	ARG
1	B	55	LYS
1	B	62	ASN
1	B	63	ARG
1	B	143	LEU
1	B	168	ARG
1	B	189	LYS
1	B	201	THR
1	B	218	ARG
1	B	250	ASP
1	B	273	LEU
1	C	31	ARG
1	C	62	ASN
1	C	63	ARG
1	C	74	MSE
1	C	103	GLN
1	C	143	LEU
1	C	168	ARG
1	C	201	THR
1	C	205	GLN
1	C	224	LEU
1	C	250	ASP
1	C	273	LEU
1	C	282	THR
1	D	62	ASN

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Mol	Chain	Res	Type
1	D	76	GLU
1	D	168	ARG
1	D	218	ARG
1	D	224	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	ASN
1	A	113	GLN
1	A	136	ASN
1	A	150	ASN
1	B	62	ASN
1	B	67	HIS
1	B	97	HIS
1	B	113	GLN
1	B	136	ASN
1	B	141	GLN
1	B	171	ASN
1	C	113	GLN
1	C	121	GLN
1	C	136	ASN
1	C	202	HIS
1	C	205	GLN
1	C	241	HIS
1	D	36	GLN
1	D	113	GLN
1	D	136	ASN
1	D	141	GLN
1	D	171	ASN
1	D	202	HIS
1	D	241	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 [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(Å ²)	Q<0.9
1	A	260/292 (89%)	-0.08	3 (1%) 79 76	11, 19, 34, 44	0
1	B	260/292 (89%)	0.18	11 (4%) 36 30	11, 20, 46, 63	0
1	C	260/292 (89%)	0.23	14 (5%) 25 20	14, 22, 43, 66	0
1	D	260/292 (89%)	0.28	15 (5%) 23 18	14, 25, 45, 64	0
All	All	1040/1168 (89%)	0.15	43 (4%) 37 31	11, 21, 43, 66	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	66	ILE	7.2
1	C	282	THR	7.1
1	C	250	ASP	6.7
1	D	282	THR	6.1
1	B	249	CYS	5.6
1	D	249	CYS	5.5
1	C	280	LEU	5.3
1	C	252	PRO	5.3
1	B	250	ASP	5.1
1	B	282	THR	5.1
1	D	250	ASP	4.8
1	B	67	HIS	4.5
1	C	205	GLN	4.4
1	B	252	PRO	4.3
1	D	245	ASN	4.1
1	B	2	GLN	4.0
1	D	252	PRO	3.9
1	D	2	GLN	3.8
1	C	2	GLN	3.7
1	C	249	CYS	3.6
1	B	65	SER	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	281	ASP	3.6
1	C	281	ASP	3.6
1	A	245	ASN	3.5
1	B	68	SER	3.4
1	D	281	ASP	3.4
1	C	279	GLU	3.0
1	A	2	GLN	2.9
1	B	64	SER	2.9
1	C	251	GLY	2.8
1	D	205	GLN	2.8
1	D	175	ASP	2.8
1	D	218	ARG	2.7
1	C	278	PRO	2.7
1	D	43	LYS	2.6
1	D	246	GLN	2.5
1	C	175	ASP	2.4
1	D	192	LYS	2.3
1	D	161	ASP	2.3
1	D	278	PRO	2.2
1	A	281	ASP	2.2
1	C	219	ALA	2.1
1	C	248	LYS	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 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.