



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 24, 2024 – 08:10 AM EST

PDB ID : 1XV2
Title : Crystal Structure of a Protein of Unknown Function Similar to Alpha-acetolactate Decarboxylase from Staphylococcus aureus
Authors : Duke, N.E.C.; Zhang, R.; Quartey, P.; Collart, F.; Holzle, D.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2004-10-26
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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

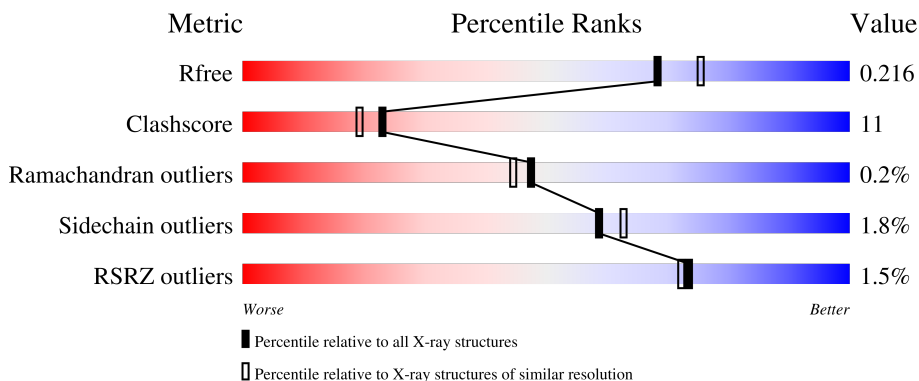
1 Overall quality at a glance

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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (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	237	 76% 20% ..
1	B	237	 72% 25% ..
1	C	237	 81% 17% ..
1	D	237	 82% 15% ..

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 8068 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 hypothetical protein, similar to alpha-acetolactate decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	232	1850	1177	313	355	5	0	0	0
1	B	232	1850	1177	313	355	5	0	0	0
1	C	233	1857	1181	314	357	5	0	0	0
1	D	232	1850	1177	313	355	5	0	0	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	cloning artifact	UNP Q7A3A7
A	-1	ASN	-	cloning artifact	UNP Q7A3A7
A	0	ALA	-	cloning artifact	UNP Q7A3A7
A	1	MSE	MET	modified residue	UNP Q7A3A7
A	15	MSE	MET	modified residue	UNP Q7A3A7
A	103	MSE	MET	modified residue	UNP Q7A3A7
A	121	MSE	MET	modified residue	UNP Q7A3A7
A	125	MSE	MET	modified residue	UNP Q7A3A7
A	126	MSE	MET	modified residue	UNP Q7A3A7
B	-2	SER	-	cloning artifact	UNP Q7A3A7
B	-1	ASN	-	cloning artifact	UNP Q7A3A7
B	0	ALA	-	cloning artifact	UNP Q7A3A7
B	1	MSE	MET	modified residue	UNP Q7A3A7
B	15	MSE	MET	modified residue	UNP Q7A3A7
B	103	MSE	MET	modified residue	UNP Q7A3A7
B	121	MSE	MET	modified residue	UNP Q7A3A7
B	125	MSE	MET	modified residue	UNP Q7A3A7
B	126	MSE	MET	modified residue	UNP Q7A3A7
C	-2	SER	-	cloning artifact	UNP Q7A3A7
C	-1	ASN	-	cloning artifact	UNP Q7A3A7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	0	ALA	-	cloning artifact	UNP Q7A3A7
C	1	MSE	MET	modified residue	UNP Q7A3A7
C	15	MSE	MET	modified residue	UNP Q7A3A7
C	103	MSE	MET	modified residue	UNP Q7A3A7
C	121	MSE	MET	modified residue	UNP Q7A3A7
C	125	MSE	MET	modified residue	UNP Q7A3A7
C	126	MSE	MET	modified residue	UNP Q7A3A7
D	-2	SER	-	cloning artifact	UNP Q7A3A7
D	-1	ASN	-	cloning artifact	UNP Q7A3A7
D	0	ALA	-	cloning artifact	UNP Q7A3A7
D	1	MSE	MET	modified residue	UNP Q7A3A7
D	15	MSE	MET	modified residue	UNP Q7A3A7
D	103	MSE	MET	modified residue	UNP Q7A3A7
D	121	MSE	MET	modified residue	UNP Q7A3A7
D	125	MSE	MET	modified residue	UNP Q7A3A7
D	126	MSE	MET	modified residue	UNP Q7A3A7

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0

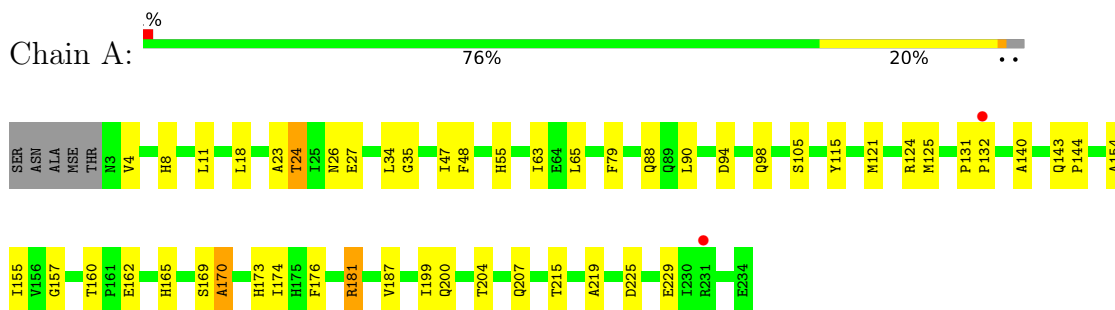
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	107	Total O 107 107	0	0
3	B	82	Total O 82 82	0	0
3	C	246	Total O 246 246	0	0
3	D	222	Total O 222 222	0	0

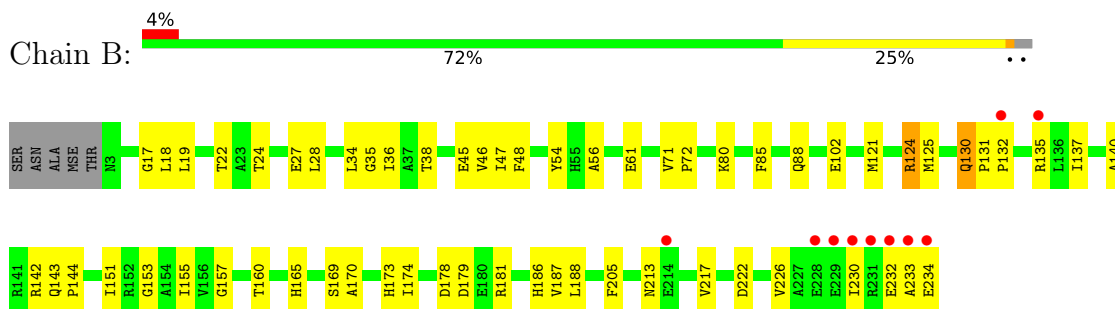
3 Residue-property plots [i](#)

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.

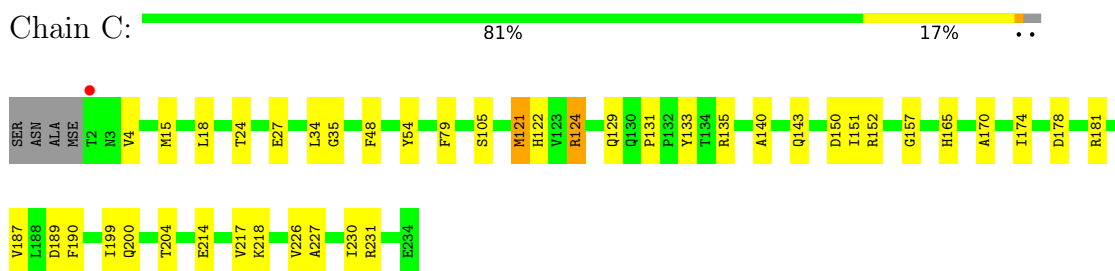
- Molecule 1: hypothetical protein, similar to alpha-acetolactate decarboxylase



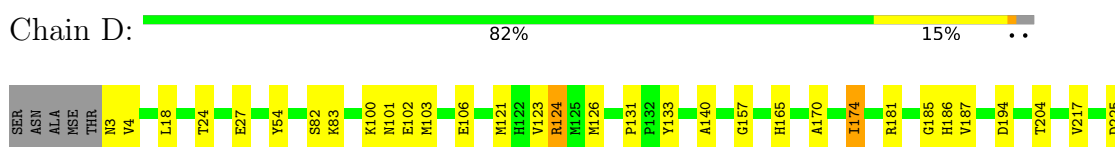
- Molecule 1: hypothetical protein, similar to alpha-acetolactate decarboxylase



- Molecule 1: hypothetical protein, similar to alpha-acetolactate decarboxylase



- Molecule 1: hypothetical protein, similar to alpha-acetolactate decarboxylase



V226	A227	E228	E229	I230	R231	E232	A233	E234
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	48.43Å 160.00Å 68.06Å 90.00° 107.55° 90.00°	Depositor
Resolution (Å)	30.24 – 2.00 30.24 – 1.66	Depositor EDS
% Data completeness (in resolution range)	94.3 (30.24-2.00) 55.2 (30.24-1.66)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 1.66Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.191 , 0.236 0.193 , 0.216	Depositor DCC
R_{free} test set	1588 reflections (1.25%)	wwPDB-VP
Wilson B-factor (Å ²)	12.5	Xtrriage
Anisotropy	0.714	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8068	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% 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:
ZN

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	0.31	0/1887	0.58	0/2540
1	B	0.31	0/1887	0.56	0/2540
1	C	0.35	0/1894	0.62	0/2550
1	D	0.37	0/1887	0.62	0/2540
All	All	0.34	0/7555	0.60	0/10170

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1850	0	1781	38	0
1	B	1850	0	1781	46	0
1	C	1857	0	1788	42	0
1	D	1850	0	1781	42	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	107	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	82	0	0	1	0
3	C	246	0	0	3	0
3	D	222	0	0	2	0
All	All	8068	0	7131	165	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 11.

All (165) 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:121:MSE:HE3	1:C:190:PHE:HB3	1.20	1.17
1:C:121:MSE:HE2	1:C:122:HIS:H	1.10	1.08
1:D:174:ILE:HG12	1:D:187:VAL:CG1	1.87	1.04
1:D:186:HIS:HE2	1:D:234:GLU:HG3	1.27	0.99
1:C:121:MSE:HE1	1:C:187:VAL:HG13	1.46	0.97
1:C:121:MSE:HE2	1:C:122:HIS:N	1.82	0.94
1:A:24:THR:HG22	1:A:27:GLU:H	1.36	0.87
1:C:121:MSE:HE3	1:C:190:PHE:CB	2.03	0.85
1:C:24:THR:HG22	1:C:27:GLU:HG3	1.57	0.84
1:D:174:ILE:HG12	1:D:187:VAL:HG11	1.59	0.82
1:D:174:ILE:O	1:D:187:VAL:HG12	1.82	0.80
1:C:24:THR:HG23	1:C:27:GLU:H	1.46	0.78
1:B:24:THR:HG22	1:B:27:GLU:HG3	1.67	0.77
1:D:24:THR:HG23	1:D:27:GLU:H	1.50	0.77
1:C:24:THR:HG22	1:C:27:GLU:CG	2.17	0.74
1:C:121:MSE:CE	1:C:190:PHE:HB3	2.10	0.73
1:D:100:LYS:HA	1:D:103:MSE:HE2	1.72	0.71
1:D:24:THR:HG22	1:D:27:GLU:HG3	1.73	0.71
1:B:165:HIS:CD2	1:B:170:ALA:HA	2.26	0.71
1:D:186:HIS:HE2	1:D:234:GLU:CG	2.03	0.71
1:C:121:MSE:CE	1:C:189:ASP:O	2.40	0.70
1:D:186:HIS:NE2	1:D:234:GLU:HG3	2.04	0.69
1:D:174:ILE:HD13	1:D:174:ILE:N	2.07	0.69
1:C:152:ARG:HD2	3:C:927:HOH:O	1.92	0.69
1:C:24:THR:CG2	1:C:27:GLU:HG3	2.23	0.68
1:A:165:HIS:CD2	1:A:170:ALA:HA	2.30	0.67
1:C:24:THR:CG2	1:C:27:GLU:H	2.08	0.66
1:D:170:ALA:HB3	1:D:231:ARG:NH1	2.12	0.65
1:D:126:MSE:HE1	1:D:140:ALA:HB2	1.77	0.65
1:B:24:THR:HG23	1:B:27:GLU:H	1.61	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:105:SER:HB2	1:D:217:VAL:HG13	1.79	0.64
1:B:28:LEU:HD21	1:B:71:VAL:HG21	1.80	0.63
1:A:160:THR:HB	1:A:169:SER:O	1.99	0.63
1:D:123:VAL:HG23	1:D:186:HIS:O	2.00	0.62
1:C:121:MSE:HE2	1:C:121:MSE:HA	1.82	0.62
1:D:157:GLY:HA3	1:D:174:ILE:HG22	1.81	0.62
1:B:24:THR:HG22	1:B:27:GLU:CG	2.29	0.61
1:B:135:ARG:HH22	1:B:222:ASP:HB3	1.66	0.61
1:D:24:THR:HG22	1:D:27:GLU:CG	2.30	0.61
1:A:121:MSE:HE1	1:A:155:ILE:HD11	1.82	0.61
1:C:121:MSE:CE	1:C:121:MSE:HA	2.31	0.61
1:D:174:ILE:HD13	1:D:174:ILE:H	1.63	0.61
1:B:130:GLN:HG3	1:B:131:PRO:HD2	1.84	0.60
1:C:121:MSE:CE	1:C:122:HIS:H	2.01	0.60
1:B:174:ILE:HG13	1:B:187:VAL:HB	1.83	0.59
1:A:79:PHE:HA	1:A:200:GLN:HE22	1.67	0.59
1:D:123:VAL:HG21	1:D:185:GLY:CA	2.32	0.59
1:B:24:THR:CG2	1:B:27:GLU:HG3	2.31	0.59
1:D:24:THR:CG2	1:D:27:GLU:HG3	2.33	0.57
1:D:123:VAL:HG22	1:D:124:ARG:N	2.20	0.57
1:A:225:ASP:O	1:A:229:GLU:HG3	2.04	0.56
1:B:151:ILE:HD11	1:B:178:ASP:HB3	1.87	0.56
1:A:4:VAL:HG22	1:A:204:THR:HB	1.88	0.56
1:A:35:GLY:N	1:A:48:PHE:HB2	2.21	0.56
1:B:135:ARG:HB3	1:B:137:ILE:HG22	1.88	0.55
1:D:100:LYS:CA	1:D:103:MSE:HE2	2.35	0.55
1:B:47:ILE:HD12	1:B:47:ILE:N	2.21	0.55
1:C:35:GLY:N	1:C:48:PHE:HB2	2.23	0.54
1:D:233:ALA:O	1:D:234:GLU:HB3	2.08	0.54
1:A:169:SER:OG	1:A:173:HIS:HB2	2.08	0.53
1:C:231:ARG:HG2	1:C:231:ARG:HH11	1.73	0.53
1:B:124:ARG:HD2	1:B:124:ARG:C	2.29	0.53
1:B:137:ILE:HD11	1:B:232:GLU:OE1	2.08	0.52
1:D:174:ILE:CG1	1:D:187:VAL:HG11	2.35	0.52
1:A:140:ALA:O	1:A:143:GLN:HG2	2.09	0.52
1:B:56:ALA:HA	1:B:61:GLU:O	2.09	0.52
1:C:165:HIS:HB2	1:C:170:ALA:HB2	1.91	0.52
1:C:131:PRO:HG3	1:C:133:TYR:CZ	2.45	0.52
1:C:121:MSE:HE2	1:C:121:MSE:CA	2.40	0.52
1:B:124:ARG:HB3	1:B:188:LEU:HD11	1.93	0.51
1:B:45:GLU:HG3	1:B:186:HIS:H	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:ILE:C	1:A:63:ILE:HD12	2.31	0.50
1:A:157:GLY:HA3	1:A:174:ILE:HG22	1.92	0.50
1:C:150:ASP:HB3	3:C:996:HOH:O	2.11	0.50
1:B:233:ALA:O	1:B:234:GLU:HG2	2.11	0.49
1:D:54:TYR:CZ	1:D:181:ARG:HD2	2.47	0.49
1:A:174:ILE:HG13	1:A:187:VAL:HB	1.95	0.49
1:B:131:PRO:HA	1:B:132:PRO:C	2.33	0.49
1:D:24:THR:CG2	1:D:27:GLU:H	2.22	0.49
1:D:123:VAL:HG23	1:D:186:HIS:C	2.33	0.48
1:A:24:THR:CG2	1:A:27:GLU:HG3	2.43	0.48
1:B:121:MSE:HE1	1:B:155:ILE:HD11	1.95	0.48
1:B:132:PRO:HG2	3:B:867:HOH:O	2.12	0.48
1:C:4:VAL:HG22	1:C:204:THR:HB	1.96	0.48
1:B:22:THR:HG22	1:B:213:ASN:ND2	2.28	0.48
1:D:174:ILE:N	1:D:174:ILE:CD1	2.77	0.48
1:D:131:PRO:HG3	1:D:133:TYR:CZ	2.49	0.47
1:D:18:LEU:N	1:D:18:LEU:HD12	2.29	0.47
1:D:100:LYS:CB	1:D:103:MSE:HE2	2.44	0.47
1:B:46:VAL:C	1:B:47:ILE:HD12	2.35	0.47
1:D:3:ASN:N	3:D:907:HOH:O	2.47	0.47
1:A:18:LEU:HD12	1:A:18:LEU:N	2.30	0.46
1:B:226:VAL:O	1:B:230:ILE:HG13	2.16	0.46
1:A:24:THR:HG22	1:A:27:GLU:N	2.18	0.46
1:A:63:ILE:HD12	1:A:63:ILE:O	2.16	0.46
1:B:160:THR:HB	1:B:169:SER:O	2.16	0.46
1:A:8:HIS:CE1	1:A:23:ALA:HB2	2.50	0.46
1:B:54:TYR:OH	1:B:181:ARG:NH1	2.49	0.46
1:B:153:GLY:HA2	1:B:179:ASP:OD1	2.16	0.46
1:A:94:ASP:O	1:A:98:GLN:HG2	2.16	0.46
1:B:233:ALA:O	1:B:234:GLU:CG	2.64	0.46
1:C:15:MSE:HE2	1:C:230:ILE:HA	1.98	0.46
1:C:199:ILE:HG22	1:C:200:GLN:N	2.31	0.46
1:A:131:PRO:HA	1:A:132:PRO:C	2.36	0.46
1:B:140:ALA:O	1:B:143:GLN:HG2	2.15	0.45
1:B:54:TYR:CZ	1:B:181:ARG:HD2	2.51	0.45
1:D:123:VAL:HG21	1:D:185:GLY:HA3	1.96	0.45
1:D:225:ASP:O	1:D:229:GLU:HG3	2.16	0.45
1:D:174:ILE:HG12	1:D:187:VAL:HG13	1.88	0.45
1:A:121:MSE:HE3	1:A:176:PHE:CD1	2.52	0.45
1:D:226:VAL:HG13	1:D:227:ALA:N	2.32	0.45
1:A:11:LEU:HD22	1:A:169:SER:HB3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:GLN:O	1:A:90:LEU:HG	2.17	0.44
1:C:135:ARG:HD3	3:C:1009:HOH:O	2.16	0.44
1:A:162:GLU:HA	1:A:165:HIS:NE2	2.33	0.44
1:B:22:THR:HG22	1:B:213:ASN:HD22	1.82	0.44
1:C:34:LEU:C	1:C:34:LEU:HD23	2.37	0.44
1:C:157:GLY:HA3	1:C:174:ILE:HG22	1.99	0.44
1:C:226:VAL:HG13	1:C:227:ALA:N	2.32	0.44
1:A:34:LEU:C	1:A:34:LEU:HD23	2.38	0.44
1:B:85:PHE:CE2	1:B:102:GLU:HG3	2.53	0.44
1:D:4:VAL:HG22	1:D:204:THR:HB	1.99	0.44
1:C:54:TYR:OH	1:C:181:ARG:NH1	2.51	0.43
1:D:103:MSE:HE3	1:D:106:GLU:HB3	2.00	0.43
1:A:24:THR:HB	1:A:27:GLU:HG3	1.99	0.43
1:A:181:ARG:NH2	3:A:874:HOH:O	2.51	0.43
1:C:214:GLU:HA	1:C:217:VAL:HG22	2.00	0.43
1:C:217:VAL:HG23	1:C:218:LYS:HG3	2.00	0.43
1:D:165:HIS:HB2	1:D:170:ALA:HB2	2.00	0.43
1:D:170:ALA:O	1:D:231:ARG:NH1	2.50	0.43
1:A:47:ILE:N	1:A:47:ILE:HD12	2.33	0.43
1:B:35:GLY:N	1:B:48:PHE:HB2	2.33	0.43
1:B:45:GLU:OE1	1:B:186:HIS:HB2	2.19	0.43
1:C:231:ARG:HG2	1:C:231:ARG:NH1	2.32	0.43
1:D:82:SER:HB3	3:D:865:HOH:O	2.18	0.43
1:A:79:PHE:HA	1:A:200:GLN:NE2	2.33	0.42
1:B:143:GLN:HA	1:B:144:PRO:HD3	1.89	0.42
1:A:125:MSE:O	1:A:144:PRO:HD2	2.18	0.42
1:B:157:GLY:HA3	1:B:174:ILE:HG22	2.01	0.42
1:C:121:MSE:HE2	1:C:189:ASP:O	2.19	0.42
1:C:151:ILE:HD11	1:C:178:ASP:HB3	2.02	0.42
1:B:17:GLY:O	1:B:19:LEU:HG	2.19	0.42
1:B:135:ARG:HH11	1:B:135:ARG:HG2	1.83	0.42
1:C:54:TYR:CZ	1:C:181:ARG:HD2	2.54	0.42
1:B:38:THR:O	1:B:72:PRO:HD2	2.19	0.42
1:A:55:HIS:HB2	1:A:65:LEU:HD21	2.02	0.41
1:B:18:LEU:HD12	1:B:18:LEU:N	2.35	0.41
1:B:80:LYS:HD3	1:B:80:LYS:HA	1.95	0.41
1:A:24:THR:HG22	1:A:27:GLU:HG3	2.03	0.41
1:C:135:ARG:H	1:C:135:ARG:HG2	1.68	0.41
1:A:215:THR:O	1:A:219:ALA:HB2	2.21	0.41
1:B:125:MSE:O	1:B:144:PRO:HD2	2.21	0.41
1:D:123:VAL:CG2	1:D:124:ARG:N	2.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:SER:HB2	1:B:217:VAL:HG13	2.03	0.41
1:A:199:ILE:HG22	1:A:200:GLN:N	2.36	0.41
1:C:140:ALA:O	1:C:143:GLN:HG2	2.20	0.41
1:B:34:LEU:C	1:B:34:LEU:HD23	2.42	0.41
1:B:36:ILE:HD12	1:B:173:HIS:CE1	2.56	0.41
1:C:79:PHE:HA	1:C:200:GLN:HE22	1.85	0.41
1:A:24:THR:HG23	1:A:26:ASN:H	1.85	0.40
1:A:115:TYR:HD1	1:A:154:ALA:HB2	1.86	0.40
1:A:207:GLN:HB3	1:B:205:PHE:HB3	2.03	0.40
1:D:83:LYS:HE3	1:D:102:GLU:O	2.20	0.40
1:C:18:LEU:HD12	1:C:18:LEU:N	2.36	0.40
1:C:124:ARG:HD2	1:C:124:ARG:C	2.41	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	230/237 (97%)	221 (96%)	8 (4%)	1 (0%)	34	30
1	B	230/237 (97%)	221 (96%)	8 (4%)	1 (0%)	34	30
1	C	231/237 (98%)	226 (98%)	5 (2%)	0	100	100
1	D	230/237 (97%)	223 (97%)	7 (3%)	0	100	100
All	All	921/948 (97%)	891 (97%)	28 (3%)	2 (0%)	47	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	142	ARG
1	A	170	ALA

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	196/194 (101%)	193 (98%)	3 (2%)	65	69
1	B	196/194 (101%)	193 (98%)	3 (2%)	65	69
1	C	197/194 (102%)	194 (98%)	3 (2%)	65	69
1	D	196/194 (101%)	191 (97%)	5 (3%)	46	48
All	All	785/776 (101%)	771 (98%)	14 (2%)	59	63

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	THR
1	A	124	ARG
1	A	181	ARG
1	B	88	GLN
1	B	124	ARG
1	B	130	GLN
1	C	121	MSE
1	C	124	ARG
1	C	129	GLN
1	D	101	ASN
1	D	121	MSE
1	D	124	ARG
1	D	174	ILE
1	D	194	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	ASN
1	A	26	ASN
1	A	92	GLN
1	A	101	ASN
1	A	149	GLN

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Mol	Chain	Res	Type
1	A	200	GLN
1	B	3	ASN
1	B	26	ASN
1	B	88	GLN
1	B	101	ASN
1	B	130	GLN
1	B	200	GLN
1	C	3	ASN
1	C	26	ASN
1	C	88	GLN
1	C	98	GLN
1	C	101	ASN
1	C	120	HIS
1	C	200	GLN
1	C	206	GLN
1	D	26	ASN
1	D	31	HIS
1	D	98	GLN
1	D	149	GLN
1	D	200	GLN
1	D	206	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	227/237 (95%)	-0.40	2 (0%) 84 83	12, 21, 33, 39	0
1	B	227/237 (95%)	-0.07	10 (4%) 34 33	10, 24, 48, 59	0
1	C	228/237 (96%)	-0.83	1 (0%) 92 92	3, 10, 20, 36	0
1	D	227/237 (95%)	-0.73	1 (0%) 92 92	3, 10, 20, 40	0
All	All	909/948 (95%)	-0.51	14 (1%) 73 72	3, 16, 36, 59	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	233	ALA	5.0
1	A	132	PRO	3.6
1	B	230	ILE	3.4
1	C	2	THR	3.1
1	B	214	GLU	3.1
1	B	234	GLU	2.9
1	B	231	ARG	2.9
1	D	234	GLU	2.8
1	A	231	ARG	2.6
1	B	228	GLU	2.6
1	B	132	PRO	2.5
1	B	135	ARG	2.4
1	B	232	GLU	2.3
1	B	229	GLU	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 monosaccharides 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	ZN	B	802	1/1	0.98	0.06	29,29,29,29	0
2	ZN	A	801	1/1	1.00	0.04	17,17,17,17	0
2	ZN	C	803	1/1	1.00	0.04	10,10,10,10	0
2	ZN	D	804	1/1	1.00	0.05	12,12,12,12	0

6.5 Other polymers [i](#)

There are no such residues in this entry.