



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:33 PM GMT

PDB ID : 1KUT
Title : Structural Genomics, Protein TM1243, (SAICAR synthetase)
Authors : Zhang, R.; Skarina, T.; Beasley, S.; Edwards, A.; Joachimiak, A.; Savchenko, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2002-01-22
Resolution : 2.20 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

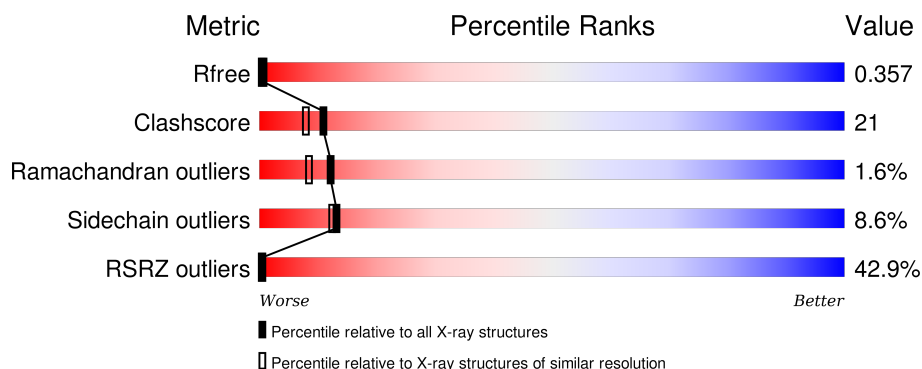
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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. 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	230	
1	B	230	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3583 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 Phosphoribosylaminoimidazole-succinocarboxamide synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	213	Total	C	N	O	S	Se	0	0	0
			1718	1109	286	316	3	4			
1	B	222	Total	C	N	O	S	Se	0	0	0
			1784	1149	298	330	3	4			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q9X0X0
A	51	MSE	MET	MODIFIED RESIDUE	UNP Q9X0X0
A	79	MSE	MET	MODIFIED RESIDUE	UNP Q9X0X0
A	124	MSE	MET	MODIFIED RESIDUE	UNP Q9X0X0
A	144	MSE	MET	MODIFIED RESIDUE	UNP Q9X0X0
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q9X0X0
B	51	MSE	MET	MODIFIED RESIDUE	UNP Q9X0X0
B	79	MSE	MET	MODIFIED RESIDUE	UNP Q9X0X0
B	124	MSE	MET	MODIFIED RESIDUE	UNP Q9X0X0
B	144	MSE	MET	MODIFIED RESIDUE	UNP Q9X0X0

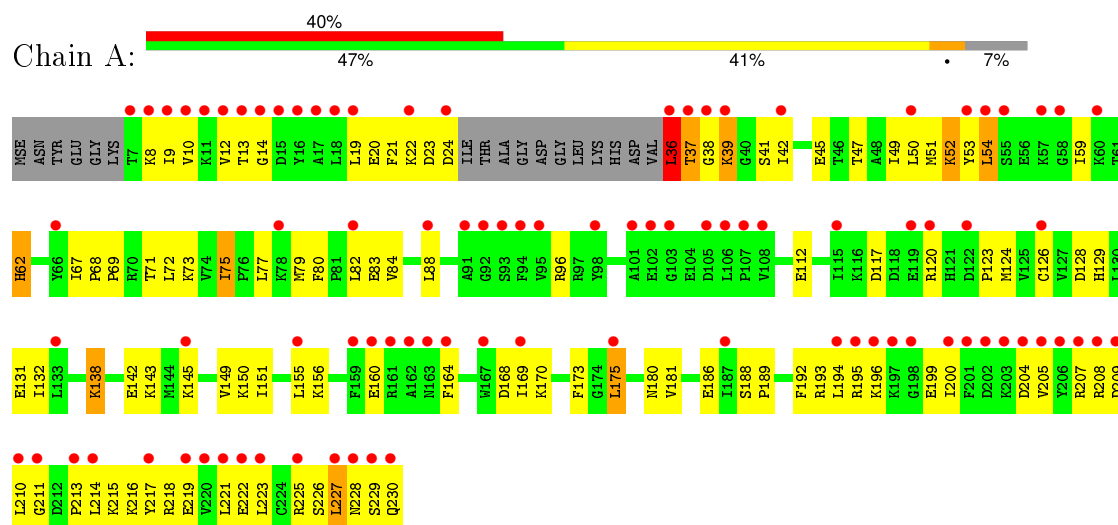
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	42	Total	O	0	0
			42	42		
2	B	39	Total	O	0	0
			39	39		

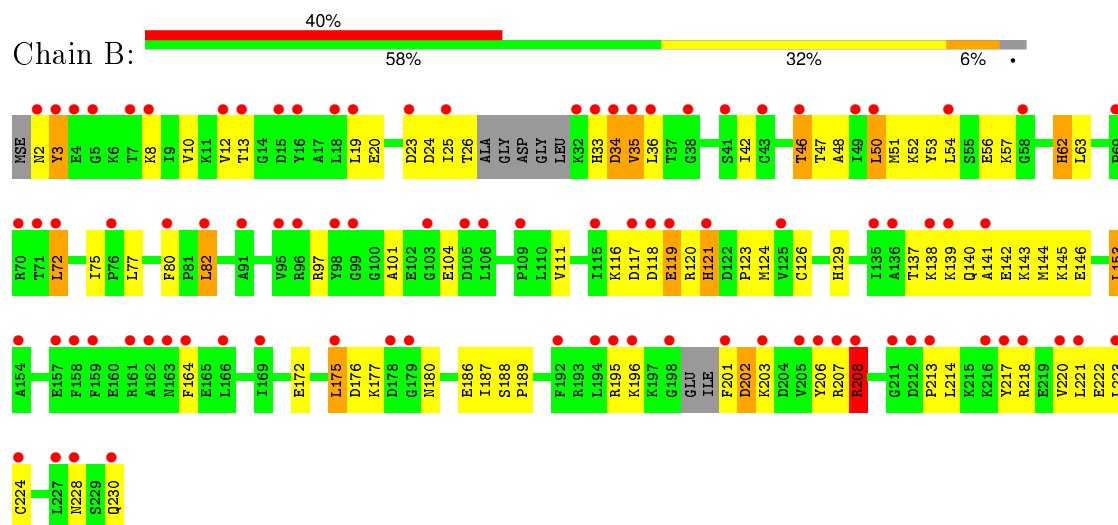
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ($\text{RSRZ} > 2$). 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: Phosphoribosylaminoimidazole-succinocarboxamide synthase



- Molecule 1: Phosphoribosylaminoimidazole-succinocarboxamide synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.51Å 43.07Å 80.22Å 90.00° 92.30° 90.00°	Depositor
Resolution (Å)	10.00 – 2.20 40.08 – 2.01	Depositor EDS
% Data completeness (in resolution range)	82.0 (10.00-2.20) 92.2 (40.08-2.01)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.01Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.245 , 0.281 0.363 , 0.357	Depositor DCC
R_{free} test set	1094 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	30.1	Xtriage
Anisotropy	0.369	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 54.2	EDS
Estimated twinning fraction	0.048 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 27003 reflections	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	3583	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 9.21% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.44	0/1744	0.72	3/2334 (0.1%)
1	B	0.46	0/1811	1.37	10/2425 (0.4%)
All	All	0.45	0/3555	1.10	13/4759 (0.3%)

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	B	0	1

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	208	ARG	NE-CZ-NH1	-34.03	103.29	120.30
1	B	208	ARG	NE-CZ-NH2	28.74	134.67	120.30
1	B	207	ARG	NE-CZ-NH2	-25.71	107.44	120.30
1	B	207	ARG	NE-CZ-NH1	24.04	132.32	120.30
1	A	36	LEU	CB-CG-CD2	9.85	127.75	111.00
1	A	36	LEU	CB-CG-CD1	-9.23	95.31	111.00
1	B	207	ARG	CD-NE-CZ	7.87	134.62	123.60
1	B	208	ARG	CA-CB-CG	6.53	127.76	113.40
1	B	207	ARG	CG-CD-NE	6.42	125.27	111.80
1	B	208	ARG	CG-CD-NE	6.31	125.05	111.80
1	B	207	ARG	CA-CB-CG	6.19	127.01	113.40
1	B	208	ARG	CB-CG-CD	-6.05	95.86	111.60
1	A	36	LEU	CA-CB-CG	5.78	128.59	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	208	ARG	Sidechain

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	1718	0	1775	81	0
1	B	1784	0	1823	72	0
2	A	42	0	0	1	0
2	B	39	0	0	3	0
All	All	3583	0	3598	149	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 21.

All (149) 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:51:MSE:HE3	1:B:63:LEU:HB2	1.53	0.91
1:A:37:THR:HG22	1:A:214:LEU:HD13	1.58	0.85
1:B:46:THR:HG22	1:B:221:LEU:HD13	1.58	0.84
1:B:121:HIS:O	1:B:123:PRO:HD3	1.77	0.83
1:B:126:CYS:H	1:B:129:HIS:HD2	1.24	0.82
1:B:36:LEU:HD11	1:B:213:PRO:HB2	1.61	0.82
1:B:42:ILE:O	1:B:46:THR:HG23	1.80	0.82
1:A:37:THR:CG2	1:A:214:LEU:HD13	2.10	0.81
1:B:47:THR:HG22	1:B:51:MSE:HE2	1.62	0.81
1:A:156:LYS:O	1:A:160:GLU:HG3	1.79	0.81
1:B:42:ILE:HD12	1:B:214:LEU:HD11	1.63	0.80
1:A:168:ASP:OD1	1:A:193:ARG:HG3	1.82	0.80
1:A:189:PRO:HB3	1:A:216:LYS:HE2	1.61	0.79
1:A:54:LEU:HD21	1:A:151:ILE:HG23	1.68	0.76
1:B:42:ILE:HG21	1:B:217:TYR:HB3	1.67	0.76
1:A:218:ARG:O	1:A:222:GLU:HG3	1.86	0.76
1:B:52:LYS:O	1:B:56:GLU:HG2	1.87	0.74
1:B:51:MSE:CE	1:B:63:LEU:HB2	2.17	0.74
1:A:227:LEU:C	1:A:229:SER:H	1.90	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:CYS:H	1:B:129:HIS:CD2	2.05	0.73
1:A:84:VAL:HG11	1:A:173:PHE:HE1	1.54	0.72
1:A:225:ARG:O	1:A:230:GLN:HB2	1.91	0.71
1:A:209:ASP:OD1	1:A:215:LYS:HD3	1.91	0.71
1:A:10:VAL:HG22	1:A:19:LEU:HD13	1.73	0.70
1:B:142:GLU:O	1:B:146:GLU:HG3	1.92	0.68
1:A:225:ARG:C	1:A:230:GLN:HB2	2.15	0.68
1:A:42:ILE:HG23	1:A:221:LEU:HD22	1.77	0.67
1:B:218:ARG:O	1:B:222:GLU:HG3	1.95	0.66
1:A:83:GLU:HG2	1:A:170:LYS:HE2	1.77	0.65
1:A:128:ASP:O	1:A:132:ILE:HD13	1.97	0.65
1:B:189:PRO:HG3	1:B:220:VAL:HG21	1.79	0.64
1:A:138:LYS:O	1:A:142:GLU:HB2	1.99	0.63
1:A:145:LYS:O	1:A:149:VAL:HG23	2.00	0.62
1:A:39:LYS:HB3	1:A:214:LEU:HD11	1.82	0.62
1:A:169:ILE:HG13	1:A:192:PHE:HB3	1.82	0.62
1:A:96:ARG:O	1:B:120:ARG:HD3	1.99	0.62
1:B:3:TYR:HB3	1:B:10:VAL:HG12	1.81	0.61
1:A:208:ARG:HG3	1:A:209:ASP:N	2.15	0.61
1:A:84:VAL:HG11	1:A:173:PHE:CE1	2.36	0.61
1:B:34:ASP:CG	1:B:35:VAL:H	2.03	0.61
1:A:131:GLU:CD	1:A:138:LYS:HD3	2.20	0.61
1:B:23:ASP:HB3	1:B:36:LEU:O	2.00	0.59
1:B:25:ILE:HG22	1:B:26:THR:N	2.17	0.59
1:B:137:THR:OG1	1:B:140:GLN:HG3	2.02	0.59
1:A:8:LYS:HB2	1:A:20:GLU:O	2.04	0.58
1:B:62:HIS:HB3	1:B:77:LEU:HD11	1.86	0.58
1:A:200:ILE:O	1:A:200:ILE:HG13	2.04	0.57
1:A:52:LYS:O	1:A:52:LYS:HD3	2.04	0.57
1:A:205:VAL:O	1:A:209:ASP:HB3	2.05	0.56
1:B:2:ASN:O	1:B:3:TYR:HB2	2.05	0.56
1:A:227:LEU:C	1:A:229:SER:N	2.56	0.56
1:A:9:ILE:HG23	1:A:22:LYS:NZ	2.21	0.55
1:A:194:LEU:C	1:A:195:ARG:HD3	2.28	0.55
1:B:177:LYS:HG3	2:B:315:HOH:O	2.08	0.54
1:B:119:GLU:C	1:B:121:HIS:H	2.11	0.54
1:B:33:HIS:ND1	1:B:34:ASP:N	2.55	0.54
1:B:139:LYS:HE3	1:B:143:LYS:NZ	2.23	0.54
1:A:8:LYS:HB3	1:A:21:PHE:HA	1.89	0.54
1:A:41:SER:HA	1:A:69:PRO:O	2.07	0.54
1:A:188:SER:HB2	1:A:189:PRO:HD2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:GLU:OE1	1:A:138:LYS:HD3	2.07	0.54
1:B:82:LEU:O	1:B:172:GLU:HG2	2.08	0.54
1:B:3:TYR:HB3	1:B:10:VAL:CG1	2.37	0.54
1:A:143:LYS:HB3	1:A:181:VAL:HG21	1.89	0.54
1:A:67:ILE:HD11	1:A:73:LYS:HB2	1.89	0.54
1:A:188:SER:HB2	1:A:189:PRO:CD	2.37	0.54
1:B:195:ARG:HH11	1:B:195:ARG:HG3	1.70	0.54
1:A:45:GLU:O	1:A:49:ILE:HG13	2.08	0.54
1:B:8:LYS:HD2	1:B:19:LEU:HD11	1.90	0.53
1:A:123:PRO:HG3	1:B:97:ARG:O	2.09	0.53
1:A:126:CYS:H	1:A:129:HIS:CD2	2.26	0.53
1:A:112:GLU:OE2	1:B:129:HIS:HE1	1.91	0.53
1:A:156:LYS:HE2	1:A:160:GLU:OE2	2.09	0.53
1:B:75:ILE:O	1:B:75:ILE:HG13	2.09	0.53
1:A:42:ILE:HG23	1:A:221:LEU:CD2	2.38	0.52
1:A:62:HIS:HA	1:A:75:ILE:HG23	1.91	0.52
1:A:12:VAL:HG12	1:A:14:GLY:H	1.75	0.52
1:A:47:THR:O	1:A:51:MSE:HB2	2.10	0.51
1:B:63:LEU:HD13	1:B:72:LEU:HG	1.93	0.51
1:A:189:PRO:HB2	1:A:217:TYR:CE2	2.46	0.50
1:B:2:ASN:HB2	1:B:10:VAL:O	2.12	0.50
1:A:208:ARG:HH11	1:A:208:ARG:HG3	1.76	0.50
1:A:62:HIS:HB3	1:A:77:LEU:HD11	1.94	0.50
1:A:214:LEU:HA	1:A:217:TYR:HD1	1.76	0.50
1:B:53:TYR:O	1:B:57:LYS:HG3	2.12	0.49
1:A:169:ILE:CG1	1:A:192:PHE:HB3	2.42	0.49
1:B:175:LEU:HA	1:B:180:ASN:O	2.12	0.49
1:B:117:ASP:O	1:B:121:HIS:O	2.31	0.49
1:A:208:ARG:NH1	1:A:208:ARG:HG3	2.27	0.49
1:B:118:ASP:N	1:B:118:ASP:OD2	2.46	0.48
1:B:201:PHE:O	1:B:202:ASP:C	2.50	0.48
1:A:207:ARG:O	1:A:210:LEU:HB3	2.14	0.48
1:B:223:LEU:HD23	1:B:223:LEU:C	2.34	0.48
1:B:24:ASP:O	1:B:25:ILE:HD13	2.13	0.48
1:B:46:THR:HG22	1:B:221:LEU:CD1	2.40	0.47
1:B:46:THR:HG21	2:B:310:HOH:O	2.13	0.47
1:A:226:SER:HA	1:A:230:GLN:HB3	1.96	0.47
1:B:195:ARG:CZ	1:B:195:ARG:HB2	2.44	0.47
1:A:117:ASP:OD2	1:A:120:ARG:HD2	2.14	0.47
1:B:111:VAL:HG11	1:B:145:LYS:HG2	1.96	0.46
1:B:13:THR:O	1:B:13:THR:HG23	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:ARG:O	1:A:210:LEU:CB	2.63	0.46
1:B:188:SER:HB2	1:B:189:PRO:HD2	1.97	0.46
1:B:224:CYS:O	1:B:228:ASN:HB2	2.16	0.46
1:A:227:LEU:O	1:A:229:SER:N	2.48	0.45
1:A:80:PHE:CE2	1:A:175:LEU:HD13	2.51	0.45
1:A:164:PHE:HB3	1:A:194:LEU:HD11	1.98	0.45
1:A:8:LYS:NZ	1:A:186:GLU:OE1	2.47	0.45
1:B:141:ALA:HA	1:B:144:MSE:HE3	1.97	0.45
1:A:59:ILE:HD11	1:A:150:LYS:HE2	1.98	0.45
1:A:37:THR:HG21	1:A:214:LEU:HD13	1.96	0.45
1:A:42:ILE:HG23	1:A:221:LEU:HB2	1.98	0.45
1:B:80:PHE:HE1	1:B:175:LEU:HD13	1.82	0.45
1:B:164:PHE:CD1	1:B:223:LEU:HD11	2.52	0.45
1:B:188:SER:HB2	1:B:189:PRO:CD	2.46	0.44
1:B:164:PHE:CE2	1:B:196:LYS:HD2	2.53	0.44
1:A:204:ASP:O	1:A:208:ARG:HG2	2.18	0.44
1:B:50:LEU:HD23	1:B:187:ILE:HG22	2.00	0.44
1:A:164:PHE:CE2	1:A:196:LYS:HB3	2.53	0.44
1:B:101:ALA:HB3	1:B:104:GLU:HB2	1.99	0.43
1:A:124:MSE:O	1:B:124:MSE:HE1	2.18	0.43
1:B:177:LYS:HB2	1:B:177:LYS:HE3	1.76	0.43
1:B:25:ILE:HG22	1:B:26:THR:H	1.83	0.43
1:A:71:THR:HG22	1:A:72:LEU:N	2.34	0.43
1:A:22:LYS:HE2	1:A:24:ASP:OD2	2.19	0.43
1:A:216:LYS:HG3	1:A:217:TYR:N	2.34	0.42
1:B:203:LYS:HG2	1:B:203:LYS:H	1.61	0.42
1:B:47:THR:CG2	1:B:51:MSE:HE2	2.43	0.42
1:A:68:PRO:HA	1:A:69:PRO:HA	1.74	0.42
1:B:20:GLU:HB2	2:B:372:HOH:O	2.20	0.42
1:B:164:PHE:CE1	1:B:223:LEU:HD12	2.56	0.41
1:A:196:LYS:NZ	1:A:199:GLU:HG3	2.35	0.41
1:B:34:ASP:CG	1:B:35:VAL:N	2.72	0.41
1:A:47:THR:OG1	1:A:186:GLU:HB2	2.19	0.41
1:B:153:LEU:HD12	1:B:153:LEU:HA	1.91	0.41
1:B:220:VAL:O	1:B:223:LEU:HB3	2.19	0.41
1:A:219:GLU:O	1:A:223:LEU:HB2	2.20	0.41
1:A:126:CYS:HB3	2:A:305:HOH:O	2.20	0.41
1:A:39:LYS:HB3	1:A:214:LEU:CD1	2.50	0.41
1:A:188:SER:CB	1:A:189:PRO:CD	2.97	0.41
1:B:33:HIS:ND1	1:B:33:HIS:C	2.74	0.41
1:B:176:ASP:OD2	1:B:180:ASN:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:48:ALA:HA	1:B:63:LEU:HD22	2.03	0.40
1:A:211:GLY:C	1:A:213:PRO:HD3	2.42	0.40
1:A:227:LEU:HD12	1:A:227:LEU:HA	1.87	0.40
1:A:145:LYS:HE3	1:A:145:LYS:HB3	1.87	0.40
1:B:25:ILE:CG2	1:B:26:THR:N	2.84	0.40
1:A:23:ASP:O	1:A:36:LEU:HD23	2.22	0.40
1:B:8:LYS:NZ	1:B:186:GLU:OE1	2.53	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	209/230 (91%)	202 (97%)	4 (2%)	3 (1%)	14	10
1	B	216/230 (94%)	201 (93%)	11 (5%)	4 (2%)	10	6
All	All	425/460 (92%)	403 (95%)	15 (4%)	7 (2%)	12	8

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	37	THR
1	B	3	TYR
1	A	228	ASN
1	B	34	ASP
1	B	119	GLU
1	A	38	GLY
1	B	35	VAL

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	188/196 (96%)	171 (91%)	17 (9%)	12	11
1	B	194/196 (99%)	178 (92%)	16 (8%)	14	13
All	All	382/392 (97%)	349 (91%)	33 (9%)	13	12

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	THR
1	A	36	LEU
1	A	39	LYS
1	A	50	LEU
1	A	52	LYS
1	A	53	TYR
1	A	54	LEU
1	A	62	HIS
1	A	75	ILE
1	A	79	MSE
1	A	82	LEU
1	A	88	LEU
1	A	138	LYS
1	A	155	LEU
1	A	175	LEU
1	A	180	ASN
1	A	227	LEU
1	B	12	VAL
1	B	46	THR
1	B	50	LEU
1	B	54	LEU
1	B	62	HIS
1	B	72	LEU
1	B	82	LEU
1	B	116	LYS
1	B	121	HIS
1	B	138	LYS

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Mol	Chain	Res	Type
1	B	153	LEU
1	B	175	LEU
1	B	202	ASP
1	B	206	TYR
1	B	208	ARG
1	B	230	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	129	HIS
1	A	228	ASN
1	B	2	ASN
1	B	129	HIS
1	B	230	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 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 ⓘ

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	209/230 (90%)	2.26	92 (44%)  	15, 27, 48, 55	0
1	B	218/230 (94%)	2.00	91 (41%)  	15, 29, 47, 54	0
All	All	427/460 (92%)	2.13	183 (42%)  	15, 28, 48, 55	0

All (183) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	32	LYS	8.5
1	A	205	VAL	7.9
1	A	204	ASP	7.9
1	A	206	TYR	7.7
1	A	200	ILE	7.6
1	A	210	LEU	7.5
1	B	220	VAL	6.3
1	A	7	THR	6.0
1	A	208	ARG	6.0
1	B	211	GLY	5.9
1	B	25	ILE	5.8
1	B	19	LEU	5.7
1	A	9	ILE	5.7
1	A	13	THR	5.7
1	A	36	LEU	5.6
1	B	16	TYR	5.5
1	A	197	LYS	5.5
1	A	211	GLY	5.4
1	B	227	LEU	5.0
1	B	195	ARG	5.0
1	A	12	VAL	5.0
1	A	230	GLN	5.0
1	B	121	HIS	5.0
1	B	3	TYR	5.0

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Mol	Chain	Res	Type	RSRZ
1	A	93	SER	4.9
1	A	207	ARG	4.9
1	A	225	ARG	4.9
1	B	33	HIS	4.8
1	B	91	ALA	4.6
1	A	16	TYR	4.6
1	A	228	ASN	4.5
1	A	14	GLY	4.5
1	A	10	VAL	4.5
1	A	223	LEU	4.5
1	A	24	ASP	4.5
1	A	229	SER	4.4
1	B	198	GLY	4.4
1	B	2	ASN	4.4
1	A	203	LYS	4.4
1	A	209	ASP	4.3
1	A	11	LYS	4.3
1	A	15	ASP	4.3
1	A	222	GLU	4.3
1	A	37	THR	4.1
1	A	155	LEU	4.1
1	B	36	LEU	4.1
1	B	206	TYR	4.1
1	A	101	ALA	4.0
1	A	227	LEU	4.0
1	A	160	GLU	4.0
1	A	214	LEU	4.0
1	B	224	CYS	3.9
1	A	22	LYS	3.9
1	B	35	VAL	3.9
1	A	201	PHE	3.9
1	B	70	ARG	3.9
1	B	58	GLY	3.8
1	B	223	LEU	3.8
1	B	15	ASP	3.8
1	B	161	ARG	3.7
1	B	203	LYS	3.7
1	A	163	ASN	3.6
1	B	208	ARG	3.6
1	B	118	ASP	3.6
1	B	221	LEU	3.5
1	A	161	ARG	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	5	GLY	3.5
1	B	201	PHE	3.4
1	B	7	THR	3.4
1	B	164	PHE	3.4
1	A	219	GLU	3.4
1	A	213	PRO	3.3
1	B	162	ALA	3.3
1	A	103	GLY	3.2
1	A	60	LYS	3.2
1	B	8	LYS	3.2
1	B	95	VAL	3.2
1	A	8	LYS	3.2
1	A	39	LYS	3.1
1	B	205	VAL	3.1
1	B	213	PRO	3.1
1	B	98	TYR	3.1
1	B	217	TYR	3.1
1	B	163	ASN	3.1
1	A	220	VAL	3.1
1	A	202	ASP	3.0
1	A	94	PHE	3.0
1	A	164	PHE	3.0
1	A	18	LEU	3.0
1	B	158	PHE	2.9
1	A	196	LYS	2.9
1	A	195	ARG	2.9
1	B	46	THR	2.9
1	A	102	GLU	2.9
1	B	139	LYS	2.9
1	A	95	VAL	2.9
1	B	34	ASP	2.9
1	B	136	ALA	2.8
1	B	72	LEU	2.8
1	B	125	VAL	2.8
1	A	50	LEU	2.8
1	A	38	GLY	2.8
1	B	54	LEU	2.8
1	A	159	PHE	2.8
1	A	167	TRP	2.8
1	A	108	VAL	2.7
1	B	207	ARG	2.7
1	B	106	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	57	LYS	2.7
1	B	159	PHE	2.7
1	B	179	GLY	2.7
1	A	162	ALA	2.7
1	A	105	ASP	2.7
1	B	154	ALA	2.7
1	B	230	GLN	2.7
1	A	78	LYS	2.6
1	A	19	LEU	2.6
1	B	218	ARG	2.6
1	A	217	TYR	2.5
1	B	194	LEU	2.5
1	A	106	LEU	2.5
1	B	76	PRO	2.5
1	B	23	ASP	2.5
1	B	196	LYS	2.4
1	A	198	GLY	2.4
1	B	119	GLU	2.4
1	B	43	CYS	2.4
1	A	145	LYS	2.4
1	B	212	ASP	2.4
1	A	98	TYR	2.3
1	A	175	LEU	2.3
1	A	120	ARG	2.3
1	A	107	PRO	2.3
1	A	88	LEU	2.3
1	B	96	ARG	2.3
1	B	18	LEU	2.3
1	B	105	ASP	2.3
1	B	228	ASN	2.3
1	A	187	ILE	2.3
1	B	12	VAL	2.3
1	B	115	ILE	2.3
1	B	135	ILE	2.3
1	B	99	GLY	2.3
1	B	13	THR	2.3
1	A	54	LEU	2.3
1	B	50	LEU	2.3
1	A	92	GLY	2.2
1	B	49	ILE	2.2
1	A	122	ASP	2.2
1	A	53	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	82	LEU	2.2
1	B	109	PRO	2.2
1	A	42	ILE	2.2
1	B	4	GLU	2.2
1	B	138	LYS	2.2
1	B	41	SER	2.2
1	A	91	ALA	2.1
1	B	169	ILE	2.1
1	A	82	LEU	2.1
1	B	178	ASP	2.1
1	B	141	ALA	2.1
1	A	58	GLY	2.1
1	A	119	GLU	2.1
1	B	192	PHE	2.1
1	A	133	LEU	2.1
1	B	175	LEU	2.1
1	A	17	ALA	2.1
1	B	103	GLY	2.1
1	A	169	ILE	2.1
1	B	69	PRO	2.1
1	A	55	SER	2.1
1	A	221	LEU	2.1
1	B	157	GLU	2.1
1	B	166	LEU	2.1
1	B	38	GLY	2.0
1	A	126	CYS	2.0
1	B	71	THR	2.0
1	A	115	ILE	2.0
1	B	117	ASP	2.0
1	B	216	LYS	2.0
1	A	194	LEU	2.0
1	B	80	PHE	2.0
1	A	66	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no 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.