



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:46 AM UTC

PDB ID : 1X2G / pdb_00001x2g
Title : Crystal Structure of Lipate-Protein Ligase A from Escherichia coli
Authors : Fujiwara, K.; Toma, S.; Okamura-Ikeda, K.; Motokawa, Y.; Nakagawa, A.;
Taniguchi, H.
Deposited on : 2005-04-23
Resolution : 2.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

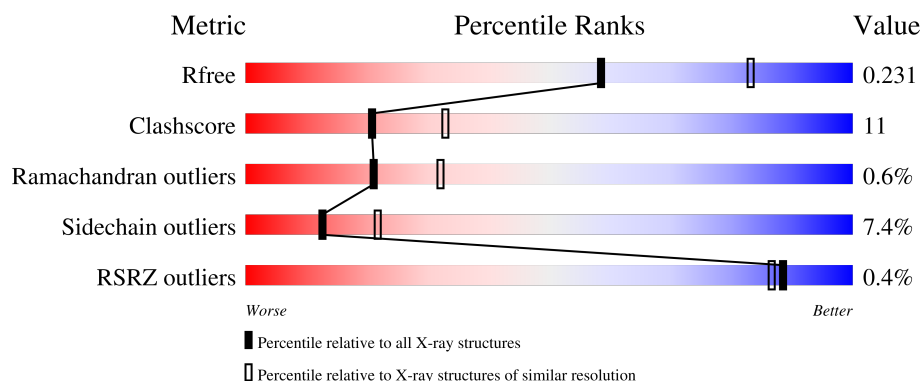
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	337	% 77% 19% ..
1	B	337	% 78% 16% ...
1	C	337	 74% 20% 6% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8425 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 Lipoate-protein ligase A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	Se	0	0	0
			2618	1646	467	494	6	5			
1	B	330	Total	C	N	O	S	Se	0	0	0
			2614	1644	466	493	6	5			
1	C	337	Total	C	N	O	S	Se	0	0	0
			2663	1673	476	503	6	5			

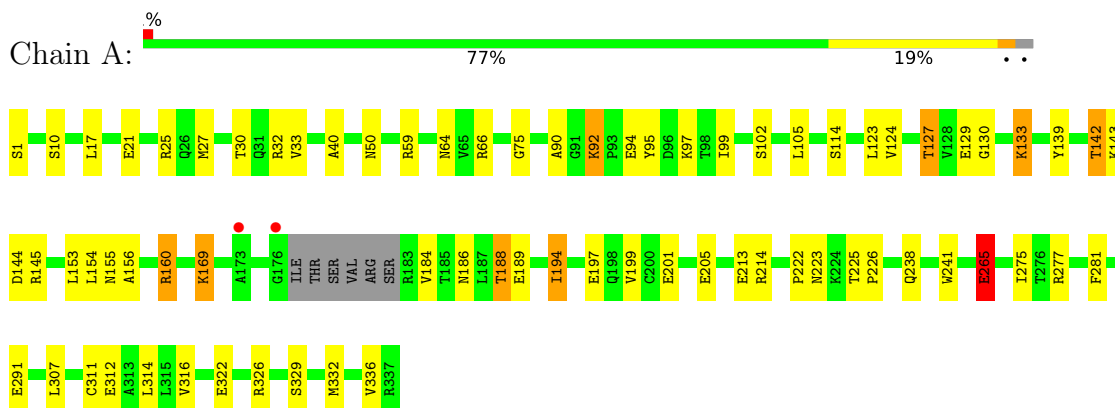
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	177	Total	O	0	0
			177	177		
2	B	175	Total	O	0	0
			175	175		
2	C	178	Total	O	0	0
			178	178		

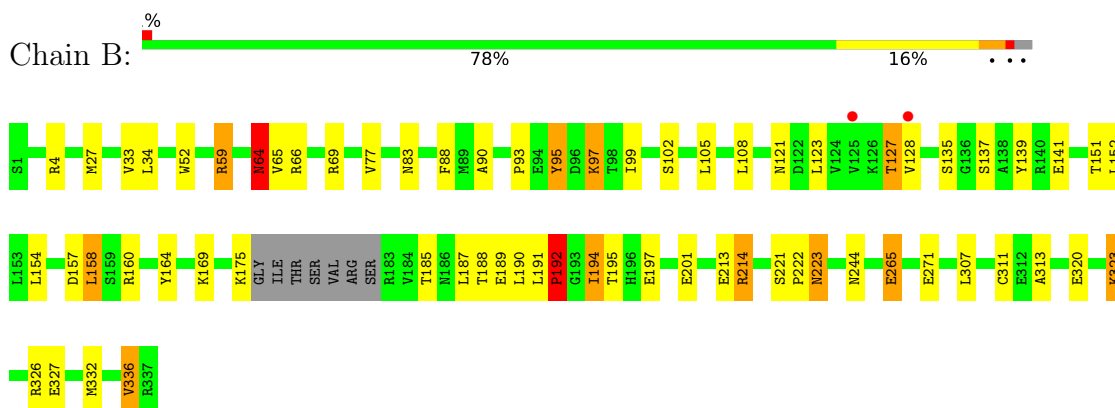
3 Residue-property plots [i](#)

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.

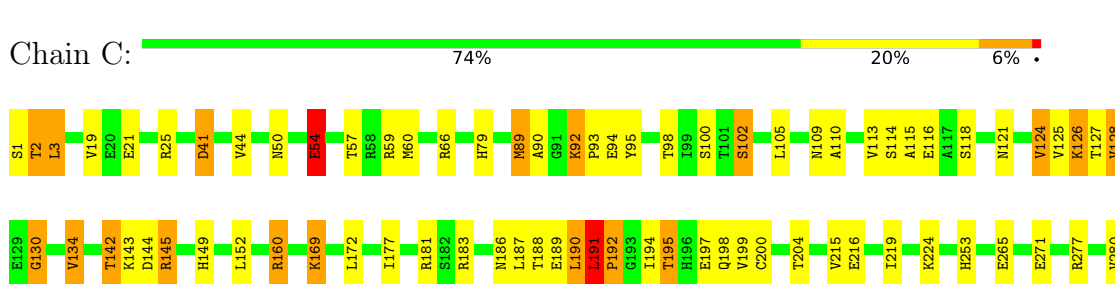
• Molecule 1: Lipoate-protein ligase A



• Molecule 1: Lipoate-protein ligase A



• Molecule 1: Lipoate-protein ligase A



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	82.00Å 112.80Å 289.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.60 – 2.40 43.60 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.6 (43.60-2.40) 99.5 (43.60-2.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.71 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.171 , 0.233 0.170 , 0.231	Depositor DCC
R_{free} test set	2698 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	43.0	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8425	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.38% 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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	1.17	3/2669 (0.1%)	1.14	4/3606 (0.1%)
1	B	1.16	5/2665 (0.2%)	1.14	1/3601 (0.0%)
1	C	1.15	6/2715 (0.2%)	1.15	10/3670 (0.3%)
All	All	1.16	14/8049 (0.2%)	1.14	15/10877 (0.1%)

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	89	MSE	SE-CE	-11.77	1.60	1.95
1	C	54	GLU	CB-CG	7.17	1.74	1.52
1	B	33	VAL	CA-CB	6.24	1.62	1.54
1	B	64	ASN	CA-C	6.14	1.60	1.53
1	B	27	MSE	SE-CE	5.96	2.13	1.95
1	A	75	GLY	N-CA	5.84	1.50	1.44
1	C	54	GLU	CD-OE2	5.46	1.35	1.25
1	A	275	ILE	CA-CB	5.42	1.60	1.53
1	C	280	VAL	CA-CB	5.27	1.60	1.54
1	B	265	GLU	CD-OE1	5.27	1.35	1.25
1	C	265	GLU	CD-OE1	5.25	1.35	1.25
1	A	265	GLU	CD-OE1	5.19	1.35	1.25
1	B	214	ARG	N-CA	-5.07	1.39	1.45
1	C	19	VAL	CA-CB	-5.01	1.48	1.54

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	160	ARG	N-CA-C	7.11	119.93	111.33
1	C	2	THR	N-CA-C	7.08	121.86	111.87
1	C	188	THR	N-CA-C	-6.58	105.11	113.01
1	C	54	GLU	CB-CG-CD	6.50	123.64	112.60
1	C	3	LEU	N-CA-C	-6.29	100.39	110.14
1	C	2	THR	CB-CA-C	5.88	120.56	111.28
1	C	124	VAL	N-CA-C	5.85	116.91	108.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	145	ARG	N-CA-C	5.83	117.64	108.96
1	C	160	ARG	N-CA-C	5.79	118.34	111.33
1	C	128	VAL	N-CA-C	-5.37	108.27	113.53
1	A	291	GLU	N-CA-C	-5.29	105.59	111.36
1	C	92	LYS	CB-CA-C	5.23	116.99	109.09
1	A	314	LEU	N-CA-C	5.10	117.74	111.82
1	B	313	ALA	N-CA-C	-5.10	106.32	112.54
1	C	219	ILE	CB-CA-C	5.08	117.10	110.96

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	2618	0	2554	56	0
1	B	2614	0	2551	43	0
1	C	2663	0	2605	80	0
2	A	177	0	0	7	0
2	B	175	0	0	8	0
2	C	178	0	0	16	0
All	All	8425	0	7710	173	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 (173) 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:89:MSE:HE2	1:C:145:ARG:HB2	1.31	1.11
1:C:169:LYS:H	1:C:169:LYS:HD3	1.01	1.08
1:C:337:ARG:HG2	2:C:515:HOH:O	1.56	1.05
1:C:126:LYS:HG3	1:C:130:GLY:O	1.58	1.04
1:C:142:THR:HG22	1:C:144:ASP:H	1.26	1.00
1:C:1:SER:HB3	2:C:493:HOH:O	1.63	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:ARG:H	1:B:244:ASN:HD22	1.12	0.96
1:C:50:ASN:O	1:C:54:GLU:HG3	1.67	0.94
1:A:197:GLU:O	1:A:201:GLU:HG3	1.69	0.92
1:C:169:LYS:H	1:C:169:LYS:CD	1.84	0.89
1:C:169:LYS:HD3	1:C:169:LYS:N	1.87	0.88
1:C:215:VAL:HG11	2:C:376:HOH:O	1.74	0.87
1:A:169:LYS:HD2	1:A:169:LYS:H	1.41	0.86
1:C:127:THR:H	1:C:130:GLY:HA3	1.39	0.85
1:C:89:MSE:HE2	1:C:145:ARG:CB	2.09	0.83
1:A:1:SER:HB3	1:A:213:GLU:OE2	1.77	0.83
1:A:155:ASN:ND2	1:A:188:THR:HG21	1.94	0.82
1:C:191:LEU:H	1:C:192:PRO:HD3	1.46	0.81
1:C:54:GLU:HB3	2:C:505:HOH:O	1.81	0.80
1:C:134:VAL:HG21	1:C:187:LEU:HD13	1.67	0.75
1:A:265:GLU:HG3	2:A:493:HOH:O	1.87	0.75
1:A:155:ASN:HD22	1:A:188:THR:HG21	1.50	0.74
1:C:323:LYS:HA	1:C:326:ARG:HG2	1.70	0.73
1:C:89:MSE:CE	1:C:145:ARG:HB2	2.16	0.72
1:B:271:GLU:HG2	2:B:497:HOH:O	1.90	0.71
1:A:17:LEU:HG	2:A:471:HOH:O	1.89	0.71
1:A:194:ILE:HD11	1:A:199:VAL:CG2	2.21	0.71
1:B:307:LEU:CD1	1:B:336:VAL:HG21	2.21	0.69
1:B:83:ASN:HD22	1:B:151:THR:HG21	1.57	0.69
1:B:69:ARG:H	1:B:244:ASN:ND2	1.88	0.68
1:B:59:ARG:HH11	1:B:59:ARG:HG3	1.59	0.68
1:A:142:THR:HG22	1:A:144:ASP:H	1.59	0.68
1:A:66:ARG:HH21	1:B:64:ASN:HD22	1.42	0.67
1:C:304:ALA:O	1:C:308:GLN:HG3	1.94	0.67
1:A:27:MSE:HE1	1:A:33:VAL:CG2	2.24	0.67
1:A:27:MSE:HE1	1:A:33:VAL:HG21	1.77	0.67
1:C:59:ARG:HH11	1:C:59:ARG:HG3	1.59	0.66
1:A:1:SER:CB	1:A:213:GLU:OE2	2.43	0.66
1:C:90:ALA:HB3	1:C:95:TYR:CD1	2.31	0.66
1:C:89:MSE:CE	1:C:145:ARG:CB	2.73	0.65
1:B:265:GLU:HG3	2:B:373:HOH:O	1.96	0.65
1:A:127:THR:HB	1:A:130:GLY:O	1.97	0.64
1:C:127:THR:H	1:C:130:GLY:CA	2.09	0.64
1:C:89:MSE:HE3	1:C:145:ARG:NE	2.13	0.64
1:B:64:ASN:HB3	2:B:446:HOH:O	1.97	0.64
1:C:57:THR:HA	1:C:60:MSE:HE2	1.80	0.63
1:A:238:GLN:HG3	2:A:452:HOH:O	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:SER:HA	2:A:437:HOH:O	1.99	0.62
1:B:323:LYS:O	1:B:327:GLU:HG3	2.00	0.62
1:C:125:VAL:HG21	1:C:190:LEU:HD12	1.82	0.62
1:A:307:LEU:CD1	1:A:336:VAL:HG21	2.31	0.61
1:C:224:LYS:HD2	2:C:472:HOH:O	2.00	0.60
1:C:191:LEU:N	1:C:192:PRO:HD3	2.13	0.60
1:A:133:LYS:HB3	1:A:184:VAL:HG22	1.83	0.60
1:B:59:ARG:HH11	1:B:59:ARG:CG	2.13	0.60
1:C:41:ASP:OD2	1:C:160:ARG:NH2	2.35	0.59
1:B:83:ASN:HD22	1:B:151:THR:CG2	2.15	0.59
1:C:1:SER:H2	1:C:215:VAL:HG22	1.68	0.58
1:A:316:VAL:HB	2:C:486:HOH:O	2.03	0.58
1:C:66:ARG:HG3	2:C:503:HOH:O	2.03	0.58
1:C:190:LEU:O	1:C:191:LEU:HB2	2.03	0.58
1:C:253:HIS:CE1	2:C:498:HOH:O	2.56	0.58
1:A:154:LEU:O	1:A:188:THR:HB	2.04	0.57
1:C:307:LEU:HD13	1:C:336:VAL:HG21	1.85	0.57
1:C:152:LEU:HD13	1:C:199:VAL:HG11	1.85	0.57
1:C:271:GLU:HG2	2:C:496:HOH:O	2.05	0.57
1:C:44:VAL:HG23	1:C:79:HIS:HD2	1.70	0.56
1:C:253:HIS:HE1	2:C:498:HOH:O	1.88	0.56
1:C:297:LEU:HD22	1:C:307:LEU:HD23	1.87	0.56
1:C:50:ASN:O	1:C:54:GLU:CG	2.48	0.56
1:B:139:TYR:CE2	1:B:141:GLU:HG3	2.42	0.55
1:C:118:SER:OG	1:C:124:VAL:HG23	2.07	0.55
1:C:200:CYS:O	1:C:204:THR:HG23	2.06	0.55
1:B:195:THR:HG22	1:B:197:GLU:OE1	2.07	0.55
1:C:191:LEU:H	1:C:192:PRO:CD	2.16	0.55
1:B:69:ARG:N	1:B:244:ASN:HD22	1.94	0.55
1:C:115:ALA:O	1:C:116:GLU:HG3	2.07	0.55
1:C:169:LYS:HD2	2:C:480:HOH:O	2.07	0.54
1:A:307:LEU:HD13	1:A:336:VAL:HG21	1.89	0.54
1:A:194:ILE:HD11	1:A:199:VAL:HG22	1.90	0.54
1:A:225:THR:HG22	1:A:226:PRO:O	2.08	0.53
1:B:4:ARG:NH2	2:B:410:HOH:O	2.41	0.53
1:B:307:LEU:HD12	1:B:336:VAL:HG21	1.88	0.53
1:C:144:ASP:OD1	1:C:145:ARG:HD2	2.08	0.53
1:C:1:SER:N	1:C:215:VAL:HG22	2.23	0.53
1:A:10:SER:HB2	1:A:222:PRO:HD3	1.90	0.52
1:A:142:THR:CG2	1:A:143:LYS:N	2.72	0.52
1:B:97:LYS:HG3	1:B:139:TYR:CD1	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:187:LEU:C	1:B:194:ILE:HD11	2.35	0.52
1:C:125:VAL:HG21	1:C:190:LEU:CD1	2.40	0.52
1:A:59:ARG:HD3	2:A:348:HOH:O	2.09	0.52
1:B:121:ASN:HB3	1:B:137:SER:O	2.10	0.51
1:C:311:CYS:HB2	1:C:332:MSE:HE1	1.94	0.50
1:C:98:THR:O	1:C:102:SER:HB3	2.11	0.50
1:A:205:GLU:OE1	1:A:214:ARG:NH2	2.44	0.50
1:A:27:MSE:CE	1:A:33:VAL:HG21	2.41	0.49
1:C:215:VAL:HG12	1:C:216:GLU:N	2.27	0.49
1:C:326:ARG:HD3	2:C:514:HOH:O	2.13	0.49
1:A:127:THR:HG22	1:A:129:GLU:N	2.28	0.49
1:A:194:ILE:HD11	1:A:199:VAL:HG23	1.95	0.48
1:C:59:ARG:HD3	2:C:388:HOH:O	2.13	0.48
1:C:194:ILE:HG23	1:C:198:GLN:HE21	1.78	0.48
1:A:156:ALA:H	1:A:186:ASN:ND2	2.12	0.48
1:B:97:LYS:HG3	1:B:139:TYR:CG	2.49	0.47
1:B:195:THR:HG21	2:B:392:HOH:O	2.14	0.47
1:B:93:PRO:HD2	2:B:466:HOH:O	2.14	0.47
1:A:156:ALA:H	1:A:186:ASN:HD21	1.62	0.47
1:A:142:THR:HG22	1:A:144:ASP:N	2.26	0.47
1:C:41:ASP:CG	1:C:160:ARG:NH2	2.72	0.47
1:B:123:LEU:HB2	1:B:135:SER:HB3	1.97	0.47
1:A:97:LYS:HD2	1:A:139:TYR:HB2	1.97	0.46
1:B:188:THR:N	1:B:194:ILE:HD11	2.30	0.46
1:C:127:THR:N	1:C:130:GLY:HA3	2.20	0.46
1:A:194:ILE:CD1	1:A:199:VAL:CG2	2.92	0.46
1:C:142:THR:CG2	1:C:143:LYS:N	2.79	0.46
1:A:127:THR:HG23	1:A:129:GLU:OE1	2.16	0.46
1:C:118:SER:OG	1:C:124:VAL:CG2	2.63	0.46
1:A:92:LYS:NZ	1:A:143:LYS:O	2.49	0.45
1:B:83:ASN:HB3	1:B:151:THR:HG23	1.99	0.45
1:B:105:LEU:HD23	1:B:105:LEU:HA	1.80	0.45
1:C:25:ARG:NH1	2:C:500:HOH:O	2.49	0.45
1:A:90:ALA:HB3	1:A:95:TYR:CG	2.52	0.45
1:B:34:LEU:HD13	1:B:88:PHE:CE1	2.51	0.45
1:C:326:ARG:CD	2:C:514:HOH:O	2.64	0.45
1:B:90:ALA:HB3	1:B:95:TYR:CG	2.52	0.45
1:C:50:ASN:HB2	1:C:281:PHE:CG	2.53	0.44
1:C:66:ARG:HA	1:C:66:ARG:HD3	1.73	0.44
1:B:213:GLU:HG2	1:B:214:ARG:N	2.32	0.44
1:A:160:ARG:HD2	2:A:353:HOH:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:GLU:O	1:A:25:ARG:HG2	2.18	0.43
1:C:191:LEU:N	1:C:192:PRO:CD	2.79	0.43
1:A:142:THR:HG22	1:A:143:LYS:N	2.34	0.43
1:A:153:LEU:HD13	1:A:184:VAL:HG13	2.00	0.43
1:B:158:LEU:H	1:B:158:LEU:HD12	1.83	0.43
1:C:60:MSE:HE2	1:C:60:MSE:HB2	1.92	0.43
1:C:194:ILE:HG22	1:C:195:THR:N	2.34	0.43
1:B:189:GLU:O	1:B:192:PRO:HD3	2.19	0.43
1:C:109:ASN:O	1:C:110:ALA:C	2.60	0.43
1:B:157:ASP:HB3	1:B:160:ARG:HG3	2.01	0.43
1:C:224:LYS:CD	2:C:472:HOH:O	2.63	0.43
1:C:92:LYS:HD3	1:C:143:LYS:O	2.19	0.42
1:C:194:ILE:HG23	1:C:198:GLN:NE2	2.34	0.42
1:C:195:THR:HG22	1:C:197:GLU:OE1	2.19	0.42
1:A:40:ALA:HB1	2:B:512:HOH:O	2.19	0.42
1:A:105:LEU:HD21	1:A:123:LEU:HD22	2.02	0.42
1:B:65:VAL:HG21	1:B:164:TYR:CZ	2.54	0.42
1:C:190:LEU:O	1:C:191:LEU:CB	2.68	0.42
1:A:153:LEU:CD1	1:A:184:VAL:HG13	2.50	0.42
1:C:92:LYS:HB2	1:C:93:PRO:HA	2.02	0.42
1:A:50:ASN:HB2	1:A:281:PHE:CD1	2.55	0.41
1:A:64:ASN:ND2	1:B:66:ARG:HH12	2.18	0.41
1:A:223:ASN:HB3	1:B:52:TRP:CZ3	2.55	0.41
1:A:64:ASN:HB3	1:B:64:ASN:HB2	2.03	0.41
1:B:326:ARG:HD2	2:B:477:HOH:O	2.20	0.41
1:C:59:ARG:HG3	1:C:59:ARG:NH1	2.32	0.41
1:C:109:ASN:HA	1:C:113:VAL:O	2.20	0.41
1:B:108:LEU:HD21	1:B:152:LEU:HD22	2.03	0.41
1:B:127:THR:HB	1:B:128:VAL:H	1.69	0.41
1:A:27:MSE:HE1	1:A:33:VAL:HG22	2.03	0.41
1:A:241:TRP:CZ2	1:B:222:PRO:HB2	2.56	0.41
1:B:221:SER:OG	1:B:223:ASN:ND2	2.54	0.41
1:C:315:LEU:HD11	1:C:322:GLU:HG2	2.02	0.41
1:A:311:CYS:HB2	1:A:332:MSE:HE1	2.02	0.41
1:C:50:ASN:HB2	1:C:281:PHE:CD1	2.56	0.41
1:A:277:ARG:HB3	1:C:277:ARG:HB3	2.03	0.40
1:A:312:GLU:OE2	1:A:329:SER:OG	2.35	0.40
1:B:311:CYS:HB2	1:B:332:MSE:HE1	2.03	0.40
1:A:322:GLU:HG3	2:A:514:HOH:O	2.22	0.40
1:C:21:GLU:CD	1:C:25:ARG:NH1	2.80	0.40
1:C:44:VAL:CG2	1:C:79:HIS:CD2	3.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:THR:HG21	1:A:144:ASP:OD1	2.20	0.40
1:C:149:HIS:CD2	1:C:149:HIS:N	2.89	0.40
1:C:149:HIS:CD2	1:C:149:HIS:H	2.39	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	327/337 (97%)	318 (97%)	9 (3%)	0	100	100
1	B	326/337 (97%)	313 (96%)	12 (4%)	1 (0%)	36	50
1	C	335/337 (99%)	314 (94%)	16 (5%)	5 (2%)	8	12
All	All	988/1011 (98%)	945 (96%)	37 (4%)	6 (1%)	21	32

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	191	LEU
1	B	192	PRO
1	C	130	GLY
1	C	186	ASN
1	C	192	PRO
1	C	41	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	276/277 (100%)	260 (94%)	16 (6%)	18	32
1	B	276/277 (100%)	254 (92%)	22 (8%)	11	19
1	C	282/277 (102%)	258 (92%)	24 (8%)	10	16
All	All	834/831 (100%)	772 (93%)	62 (7%)	13	22

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	THR
1	A	32	ARG
1	A	92	LYS
1	A	94	GLU
1	A	99	ILE
1	A	102	SER
1	A	124	VAL
1	A	127	THR
1	A	133	LYS
1	A	142	THR
1	A	169	LYS
1	A	188	THR
1	A	189	GLU
1	A	194	ILE
1	A	265	GLU
1	A	326	ARG
1	B	59	ARG
1	B	64	ASN
1	B	77	VAL
1	B	95	TYR
1	B	97	LYS
1	B	99	ILE
1	B	102	SER
1	B	127	THR
1	B	154	LEU
1	B	158	LEU
1	B	169	LYS
1	B	175	LYS
1	B	185	THR
1	B	190	LEU
1	B	191	LEU
1	B	192	PRO

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Mol	Chain	Res	Type
1	B	194	ILE
1	B	201	GLU
1	B	223	ASN
1	B	320	GLU
1	B	323	LYS
1	B	336	VAL
1	C	2	THR
1	C	3	LEU
1	C	54	GLU
1	C	94	GLU
1	C	100	SER
1	C	102	SER
1	C	105	LEU
1	C	114	SER
1	C	121	ASN
1	C	126	LYS
1	C	128	VAL
1	C	134	VAL
1	C	142	THR
1	C	145	ARG
1	C	169	LYS
1	C	172	LEU
1	C	177	ILE
1	C	181	ARG
1	C	183	ARG
1	C	189	GLU
1	C	190	LEU
1	C	191	LEU
1	C	195	THR
1	C	332	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	ASN
1	A	83	ASN
1	A	155	ASN
1	A	186	ASN
1	A	238	GLN
1	B	16	ASN
1	B	26	GLN
1	B	64	ASN

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Mol	Chain	Res	Type
1	B	83	ASN
1	B	166	ASN
1	B	223	ASN
1	B	244	ASN
1	B	298	GLN
1	C	50	ASN
1	C	79	HIS
1	C	109	ASN
1	C	121	ASN
1	C	166	ASN
1	C	198	GLN
1	C	253	HIS

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](#)

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 [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	326/337 (96%)	-0.38	2 (0%) 85 83	26, 41, 66, 97	0
1	B	325/337 (96%)	-0.34	2 (0%) 85 83	27, 42, 87, 99	0
1	C	332/337 (98%)	-0.11	0 100 100	29, 48, 88, 98	0
All	All	983/1011 (97%)	-0.28	4 (0%) 88 86	26, 43, 86, 99	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	173	ALA	3.2
1	A	176	GLY	2.9
1	B	125	VAL	2.3
1	B	128	VAL	2.2

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.