



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:43 AM UTC

PDB ID : 2CUY / pdb_00002cuy
Title : Crystal structure of malonyl CoA-acyl carrier protein transacylase from *Thermophilus thermophilus* HB8
Authors : Misaki, S.; Suzuki, K.; Kunishima, N.; Sugawara, M.; Kuroishi, C.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2005-05-31
Resolution : 2.10 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
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

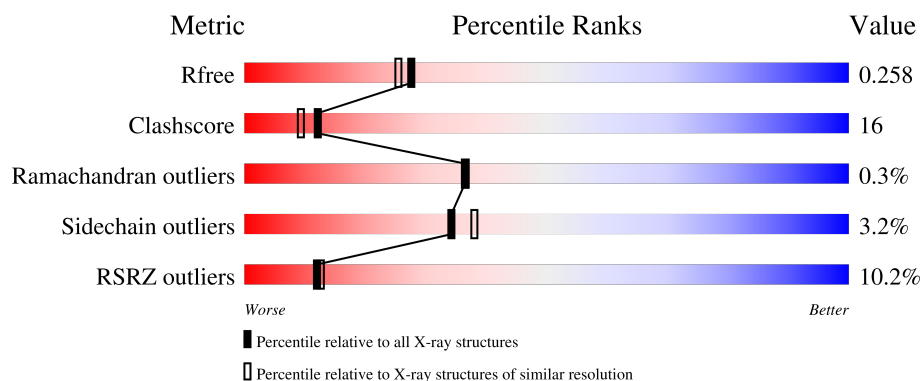
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	305	6% 69% 29% .
1	B	305	14% 68% 30% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5374 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 Malonyl CoA-[acyl carrier protein] transacylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	305	Total	C	N	O	S	0	0	0
			2351	1485	428	431	7			
1	B	305	Total	C	N	O	S	0	0	0
			2351	1485	428	431	7			

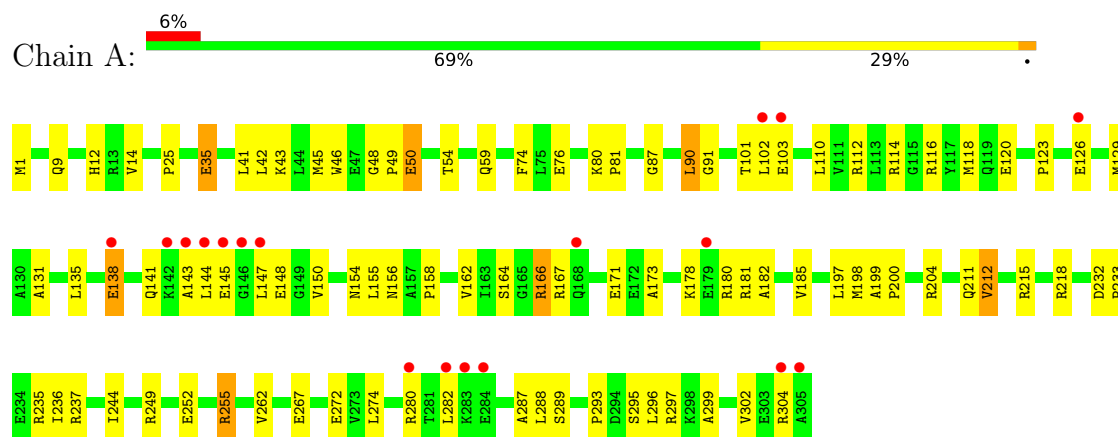
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	357	Total	O	0	0
			357	357		
2	B	315	Total	O	0	0
			315	315		

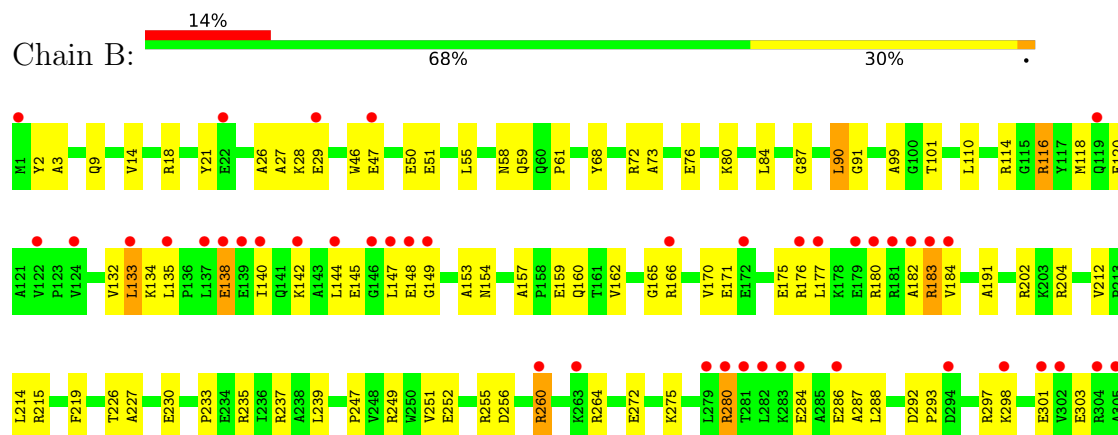
3 Residue-property plots

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: Malonyl CoA-[acyl carrier protein] transacylase



- Molecule 1: Malonyl CoA-[acyl carrier protein] transacylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	121.91Å 121.91Å 114.92Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.10 30.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-2.10) 99.9 (30.00-2.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.226 , 0.258 0.226 , 0.258	Depositor DCC
R_{free} test set	3353 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.513	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 65.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.015 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5374	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.98% 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	0.47	0/2393	0.96	9/3239 (0.3%)
1	B	0.44	0/2393	0.95	4/3239 (0.1%)
All	All	0.46	0/4786	0.95	13/6478 (0.2%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	184	VAL	N-CA-C	6.79	118.09	108.17
1	A	212	VAL	N-CA-C	6.73	115.99	108.05
1	A	235	ARG	N-CA-C	-5.94	104.88	111.36
1	B	116	ARG	CD-NE-CZ	-5.55	116.63	124.40
1	A	112	ARG	N-CA-C	-5.50	105.29	111.28
1	A	156	ASN	N-CA-C	5.47	119.12	112.23
1	B	191	ALA	N-CA-C	-5.25	102.97	109.64
1	A	46	TRP	CA-C-N	-5.17	112.70	121.18
1	A	46	TRP	C-N-CA	-5.17	112.70	121.18
1	B	237	ARG	CD-NE-CZ	-5.13	117.22	124.40
1	A	237	ARG	CD-NE-CZ	-5.09	117.28	124.40
1	A	166	ARG	CD-NE-CZ	-5.06	117.32	124.40
1	A	255	ARG	CD-NE-CZ	-5.05	117.33	124.40

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	2351	0	2411	70	0
1	B	2351	0	2411	84	0
2	A	357	0	0	11	0
2	B	315	0	0	10	0
All	All	5374	0	4822	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 16.

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:A:41:LEU:HG	1:A:45:MET:HE2	1.10	1.09
1:A:41:LEU:HG	1:A:45:MET:CE	1.91	0.99
1:B:272:GLU:HB3	1:B:275:LYS:HE3	1.50	0.92
1:A:129:MET:HG2	1:A:164:SER:HB2	1.57	0.84
1:A:35:GLU:HG2	1:A:42:LEU:HG	1.60	0.81
1:A:135:LEU:HD22	1:A:182:ALA:HB2	1.61	0.81
1:B:147:LEU:HD11	1:B:176:ARG:CZ	2.10	0.81
1:A:41:LEU:CG	1:A:45:MET:HE2	2.03	0.80
1:B:26:ALA:CB	1:B:76:GLU:HG3	2.14	0.78
1:A:204:ARG:HD2	2:A:651:HOH:O	1.85	0.76
1:A:42:LEU:HA	1:A:45:MET:HE3	1.68	0.76
1:B:26:ALA:HB2	1:B:76:GLU:HG3	1.71	0.73
1:A:129:MET:HG2	1:A:164:SER:CB	2.20	0.72
1:A:118:MET:HE2	2:A:445:HOH:O	1.88	0.72
1:B:292:ASP:HB2	1:B:293:PRO:HD2	1.71	0.72
1:B:147:LEU:HD11	1:B:176:ARG:NH1	2.05	0.72
1:A:154:ASN:HB2	1:A:162:VAL:HB	1.71	0.71
1:A:118:MET:HE3	1:A:244:ILE:HD12	1.71	0.71
1:B:58:ASN:O	1:B:61:PRO:HD2	1.90	0.71
1:A:14:VAL:HG23	1:A:50:GLU:OE2	1.91	0.70
1:A:131:ALA:CB	1:B:51:GLU:HG2	2.22	0.69
1:A:185:VAL:HG22	1:B:55:LEU:HD21	1.75	0.69
1:B:50:GLU:HG3	2:B:601:HOH:O	1.92	0.69
1:A:138:GLU:H	1:A:138:GLU:CD	1.98	0.68
1:A:287:ALA:C	1:A:288:LEU:HD12	2.20	0.67
1:B:204:ARG:HD2	2:B:548:HOH:O	1.95	0.67
1:B:87:GLY:HA3	1:B:91:GLY:C	2.21	0.66
1:A:43:LYS:NZ	2:A:598:HOH:O	2.27	0.66
1:B:235:ARG:HG2	1:B:235:ARG:HH11	1.61	0.65
1:B:120:GLU:HG3	1:B:204:ARG:NH2	2.12	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:LEU:HD12	1:B:183:ARG:HB3	1.78	0.65
1:B:157:ALA:HB3	1:B:160:GLN:HB2	1.78	0.65
1:B:90:LEU:C	1:B:90:LEU:HD13	2.22	0.65
1:A:35:GLU:HG2	1:A:42:LEU:CG	2.27	0.65
1:B:249:ARG:HD2	1:B:252:GLU:OE1	1.97	0.64
1:B:214:LEU:C	1:B:215:ARG:HD2	2.22	0.63
1:B:9:GLN:HG3	2:B:565:HOH:O	1.98	0.63
1:A:116:ARG:O	1:A:120:GLU:HG2	1.98	0.62
1:B:58:ASN:C	1:B:61:PRO:HD2	2.24	0.62
1:B:135:LEU:CD1	1:B:182:ALA:HB2	2.29	0.62
1:A:42:LEU:HD23	1:A:45:MET:HE3	1.82	0.62
1:B:260:ARG:HG3	1:B:260:ARG:HH11	1.64	0.61
1:B:280:ARG:HA	1:B:280:ARG:HE	1.65	0.61
1:A:42:LEU:HD23	1:A:45:MET:CE	2.32	0.60
1:A:249:ARG:HD2	1:A:252:GLU:OE2	2.01	0.60
1:B:287:ALA:C	1:B:288:LEU:HD12	2.26	0.60
1:B:297:ARG:O	1:B:301:GLU:HG3	2.02	0.59
1:B:298:LYS:HE2	2:B:592:HOH:O	2.03	0.59
1:B:297:ARG:HG2	1:B:297:ARG:HH11	1.67	0.59
1:A:232:ASP:O	1:A:236:ILE:HG12	2.02	0.59
1:B:233:PRO:HB2	2:B:572:HOH:O	2.03	0.58
1:A:114:ARG:O	1:A:118:MET:HG3	2.03	0.58
1:B:133:LEU:O	1:B:134:LYS:HB2	2.04	0.58
1:A:289:SER:O	1:A:295:SER:HB2	2.04	0.58
1:A:87:GLY:HA3	1:A:91:GLY:C	2.29	0.57
1:A:90:LEU:C	1:A:90:LEU:HD13	2.31	0.56
1:A:148:GLU:HG2	2:A:629:HOH:O	2.06	0.56
1:A:74:PHE:HA	1:A:296:LEU:HD11	1.88	0.55
1:B:26:ALA:HB1	1:B:76:GLU:HG3	1.86	0.55
1:B:133:LEU:CD1	1:B:183:ARG:HB3	2.37	0.55
1:A:180:ARG:O	1:A:181:ARG:HB2	2.06	0.54
1:B:176:ARG:O	1:B:180:ARG:HG3	2.08	0.53
1:A:178:LYS:HE2	2:A:610:HOH:O	2.08	0.53
1:B:230:GLU:HG3	1:B:239:LEU:HD12	1.89	0.53
1:B:202:ARG:NH1	2:B:337:HOH:O	2.42	0.53
1:B:80:LYS:HD2	1:B:303:GLU:OE2	2.07	0.53
1:B:116:ARG:O	1:B:120:GLU:HG2	2.08	0.52
1:B:120:GLU:HG3	1:B:204:ARG:HH22	1.74	0.52
1:A:143:ALA:HB2	1:A:180:ARG:NE	2.24	0.52
1:B:142:LYS:O	1:B:145:GLU:HG3	2.10	0.52
1:A:299:ALA:O	1:A:302:VAL:HG22	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:LEU:HD23	1:A:150:VAL:HG11	1.92	0.52
1:B:149:GLY:O	1:B:165:GLY:HA3	2.11	0.51
1:B:166:ARG:O	1:B:170:VAL:HG23	2.10	0.51
1:B:132:VAL:HG21	1:B:177:LEU:HD13	1.92	0.51
1:A:103:GLU:HG2	2:A:373:HOH:O	2.10	0.51
1:B:110:LEU:HD11	1:B:212:VAL:HG21	1.93	0.50
1:A:131:ALA:HB3	1:B:51:GLU:HG2	1.94	0.50
1:B:21:TYR:O	1:B:28:LYS:HB2	2.11	0.50
1:B:18:ARG:HH11	1:B:18:ARG:HG2	1.78	0.49
1:A:35:GLU:HG2	1:A:42:LEU:CD1	2.43	0.49
1:A:233:PRO:HG2	2:A:602:HOH:O	2.12	0.49
1:B:14:VAL:HG12	1:B:46:TRP:O	2.13	0.48
1:B:51:GLU:H	1:B:51:GLU:CD	2.21	0.48
1:B:264:ARG:NH1	1:B:286:GLU:OE2	2.47	0.48
1:B:80:LYS:HG2	1:B:303:GLU:O	2.13	0.48
1:A:101:THR:HG23	1:A:215:ARG:O	2.13	0.48
1:B:9:GLN:HA	1:B:59:GLN:OE1	2.14	0.48
1:B:280:ARG:HG2	2:B:322:HOH:O	2.14	0.48
1:B:2:TYR:CD2	1:B:303:GLU:HG2	2.49	0.48
1:B:171:GLU:O	1:B:175:GLU:HG3	2.13	0.48
1:A:167:ARG:HH11	1:A:167:ARG:HG2	1.79	0.48
1:B:140:ILE:O	1:B:144:LEU:HG	2.13	0.48
1:B:230:GLU:HG3	1:B:239:LEU:CD1	2.44	0.48
1:B:297:ARG:HD3	2:B:460:HOH:O	2.15	0.47
1:A:80:LYS:HG3	1:A:81:PRO:HD2	1.97	0.47
1:B:215:ARG:HD2	1:B:215:ARG:N	2.30	0.46
1:A:48:GLY:CA	1:A:49:PRO:C	2.88	0.46
1:A:199:ALA:HB3	1:A:200:PRO:HD3	1.97	0.46
1:A:144:LEU:HD21	1:A:173:ALA:HB1	1.97	0.46
1:A:1:MET:HE2	1:A:262:VAL:HG22	1.96	0.46
1:A:166:ARG:HG3	1:A:166:ARG:HH11	1.81	0.45
1:B:235:ARG:HG2	1:B:235:ARG:NH1	2.28	0.45
1:A:218:ARG:HH11	1:A:218:ARG:HG2	1.81	0.45
1:B:133:LEU:H	1:B:133:LEU:HG	1.66	0.45
1:A:141:GLN:O	1:A:145:GLU:HG3	2.16	0.45
1:B:247:PRO:HG3	2:B:436:HOH:O	2.17	0.45
1:B:256:ASP:O	1:B:260:ARG:HD3	2.16	0.45
1:B:260:ARG:HH11	1:B:260:ARG:CG	2.28	0.45
1:A:249:ARG:HD3	2:A:561:HOH:O	2.17	0.45
1:B:297:ARG:HG2	1:B:297:ARG:NH1	2.30	0.45
1:A:155:LEU:N	1:A:155:LEU:HD22	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:GLN:HA	1:A:59:GLN:OE1	2.18	0.44
1:B:138:GLU:O	1:B:142:LYS:HG3	2.18	0.44
1:A:211:GLN:NE2	2:A:521:HOH:O	2.50	0.44
1:B:99:ALA:HB2	1:B:219:PHE:CE1	2.53	0.44
1:B:133:LEU:HA	1:B:159:GLU:O	2.17	0.43
1:A:267:GLU:CD	1:A:274:LEU:HB2	2.43	0.43
1:B:68:TYR:CZ	1:B:72:ARG:HD3	2.54	0.43
1:A:110:LEU:HD11	1:A:212:VAL:HG21	2.00	0.43
1:A:131:ALA:HB2	1:B:51:GLU:HG2	1.96	0.43
1:B:27:ALA:HB2	1:B:73:ALA:HB2	1.98	0.43
1:A:118:MET:HB3	1:A:198:MET:HE1	2.01	0.43
1:A:293:PRO:O	1:A:297:ARG:HG2	2.18	0.43
1:B:288:LEU:HD12	1:B:288:LEU:N	2.33	0.43
1:B:101:THR:HG23	1:B:215:ARG:O	2.19	0.43
1:A:171:GLU:OE1	2:A:408:HOH:O	2.21	0.42
1:B:154:ASN:HB2	1:B:162:VAL:HB	2.01	0.42
1:A:181:ARG:HH11	1:A:181:ARG:HG2	1.85	0.42
1:B:3:ALA:HA	1:B:84:LEU:O	2.18	0.42
1:A:135:LEU:HD22	1:A:182:ALA:CB	2.40	0.42
1:B:114:ARG:O	1:B:118:MET:HG3	2.20	0.42
1:B:226:THR:O	1:B:227:ALA:HB3	2.20	0.42
1:A:197:LEU:HD12	1:A:197:LEU:N	2.35	0.41
1:A:158:PRO:O	2:A:584:HOH:O	2.22	0.41
1:A:25:PRO:HD2	1:A:76:GLU:OE2	2.20	0.41
1:A:166:ARG:HG3	1:A:166:ARG:NH1	2.35	0.41
1:B:51:GLU:CD	1:B:51:GLU:N	2.78	0.41
1:B:183:ARG:HH11	1:B:183:ARG:HG3	1.86	0.41
1:B:153:ALA:O	1:B:154:ASN:ND2	2.53	0.41
1:A:12:HIS:CG	1:A:59:GLN:HG3	2.56	0.41
1:B:132:VAL:HG12	2:B:488:HOH:O	2.20	0.41
1:A:54:THR:O	1:A:54:THR:HG22	2.20	0.41
1:A:280:ARG:NH2	1:B:47:GLU:O	2.53	0.41
1:B:90:LEU:C	1:B:90:LEU:CD1	2.92	0.40
1:A:123:PRO:HD2	1:A:126:GLU:OE2	2.21	0.40
1:B:132:VAL:CG2	1:B:177:LEU:HD13	2.51	0.40
1:A:272:GLU:HG2	1:A:289:SER:OG	2.21	0.40
1:B:251:VAL:O	1:B:255:ARG:HG2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	303/305 (99%)	294 (97%)	8 (3%)	1 (0%)	36	36
1	B	303/305 (99%)	291 (96%)	11 (4%)	1 (0%)	36	36
All	All	606/610 (99%)	585 (96%)	19 (3%)	2 (0%)	36	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	148	GLU
1	A	304	ARG

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	235/235 (100%)	228 (97%)	7 (3%)	36	41
1	B	235/235 (100%)	227 (97%)	8 (3%)	32	35
All	All	470/470 (100%)	455 (97%)	15 (3%)	34	38

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	GLU
1	A	50	GLU
1	A	90	LEU
1	A	102	LEU

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Mol	Chain	Res	Type
1	A	138	GLU
1	A	255	ARG
1	A	282	LEU
1	B	29	GLU
1	B	90	LEU
1	B	133	LEU
1	B	138	GLU
1	B	183	ARG
1	B	260	ARG
1	B	280	ARG
1	B	284	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	211	GLN
1	B	154	ASN
1	B	160	GLN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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 ⓘ

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	305/305 (100%)	0.41	18 (5%)	28 30	16, 28, 47, 75	0
1	B	305/305 (100%)	0.72	44 (14%)	6 6	17, 31, 56, 67	0
All	All	610/610 (100%)	0.57	62 (10%)	12 12	16, 30, 53, 75	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	305	ALA	8.1
1	B	147	LEU	4.6
1	A	304	ARG	4.4
1	B	284	GLU	4.4
1	B	282	LEU	4.2
1	B	138	GLU	3.6
1	A	284	GLU	3.5
1	A	147	LEU	3.4
1	B	148	GLU	3.3
1	B	263	LYS	3.2
1	A	283	LYS	3.2
1	B	133	LEU	3.2
1	B	305	ALA	3.1
1	A	143	ALA	3.1
1	A	179	GLU	3.1
1	B	179	GLU	3.1
1	A	146	GLY	3.0
1	B	283	LYS	3.0
1	A	282	LEU	3.0
1	A	144	LEU	3.0
1	B	172	GLU	3.0
1	B	22	GLU	3.0
1	B	298	LYS	2.9
1	B	146	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	280	ARG	2.8
1	B	260	ARG	2.8
1	B	180	ARG	2.7
1	B	182	ALA	2.7
1	B	183	ARG	2.7
1	B	135	LEU	2.6
1	A	168	GLN	2.6
1	B	1	MET	2.5
1	A	102	LEU	2.5
1	B	140	ILE	2.5
1	B	304	ARG	2.5
1	A	280	ARG	2.4
1	B	29	GLU	2.4
1	B	294	ASP	2.4
1	B	302	VAL	2.4
1	B	279	LEU	2.3
1	B	301	GLU	2.3
1	B	144	LEU	2.3
1	B	122	VAL	2.3
1	A	138	GLU	2.3
1	B	139	GLU	2.3
1	B	124	VAL	2.3
1	B	166	ARG	2.3
1	B	181	ARG	2.3
1	A	142	LYS	2.2
1	B	176	ARG	2.2
1	B	47	GLU	2.2
1	A	126	GLU	2.1
1	B	177	LEU	2.1
1	A	145	GLU	2.1
1	B	184	VAL	2.1
1	B	137	LEU	2.1
1	B	286	GLU	2.1
1	B	149	GLY	2.1
1	B	119	GLN	2.0
1	B	281	THR	2.0
1	B	142	LYS	2.0
1	A	103	GLU	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 [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.