



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:17 AM UTC

PDB ID : 7DBI / pdb_00007dbi
Title : Crystal structure of the peroxisomal acyl-CoA hydrolase MpaH
Authors : Li, S.Y.; You, C.
Deposited on : 2020-10-20
Resolution : 1.99 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

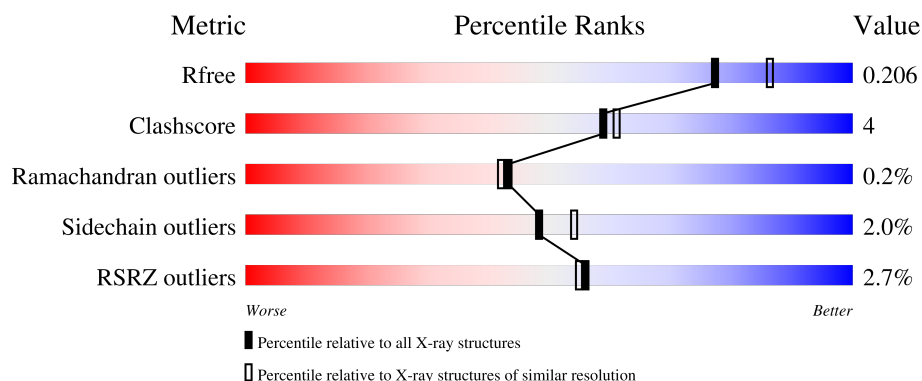
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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	433	4% 85% 11% .
1	B	433	2% 85% 11% .
1	C	433	2% 85% 12% .
1	D	433	2% 87% 8% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14163 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 acyl-CoA hydrolase MpaH'.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	420	Total	C	N	O	S	Se	0	0	0
			3356	2135	591	613	5	12			
1	B	417	Total	C	N	O	S	Se	0	0	0
			3332	2120	587	608	5	12			
1	C	420	Total	C	N	O	S	Se	0	0	0
			3356	2135	591	613	5	12			
1	D	418	Total	C	N	O	S	Se	0	0	0
			3343	2127	589	610	5	12			

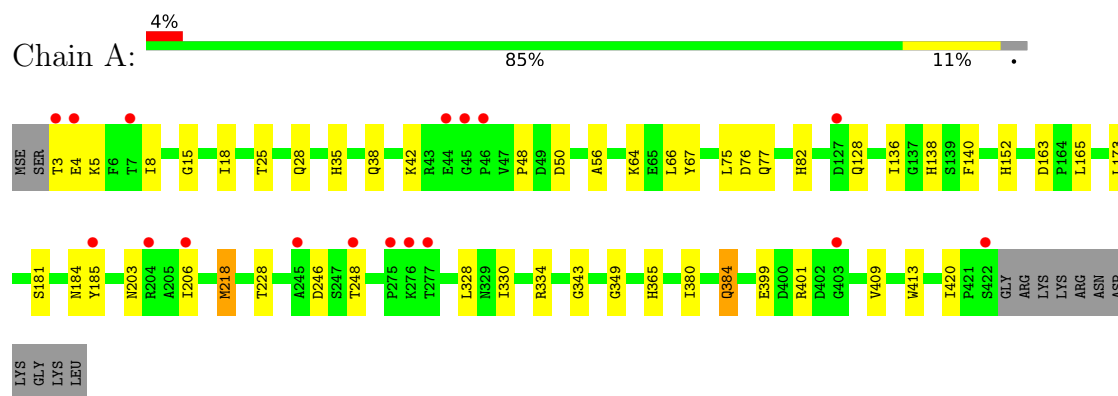
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	204	Total	O	0	0
			204	204		
2	B	210	Total	O	0	0
			210	210		
2	C	188	Total	O	0	0
			188	188		
2	D	174	Total	O	0	0
			174	174		

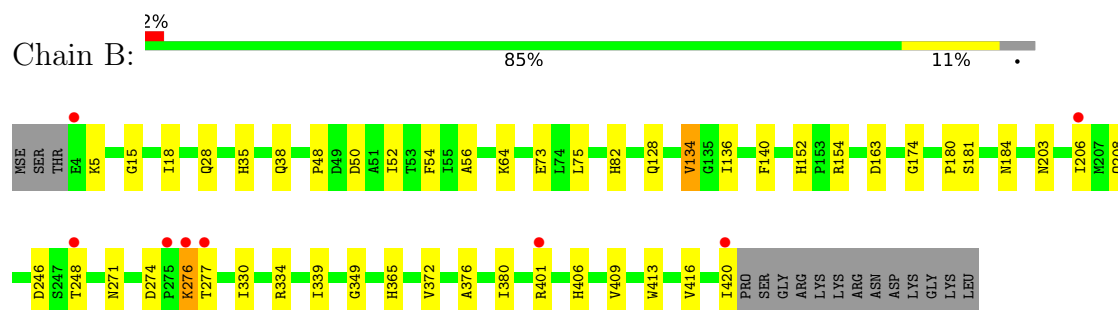
3 Residue-property plots

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.

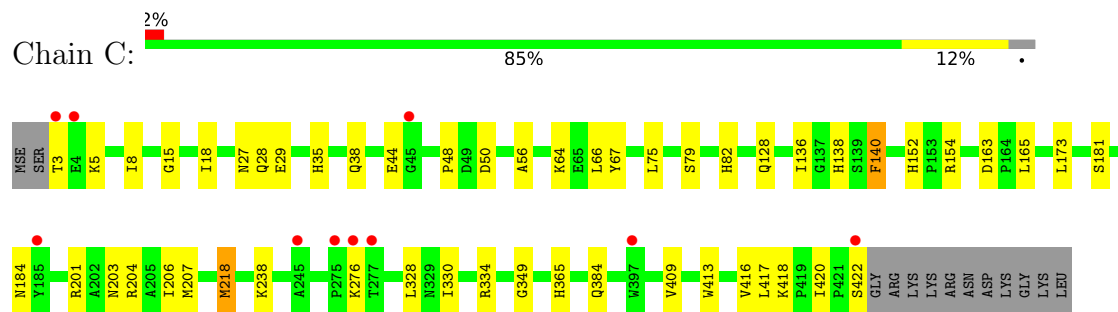
- Molecule 1: acyl-CoA hydrolase MpaH'



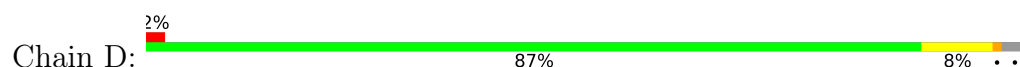
- Molecule 1: acyl-CoA hydrolase MpaH'

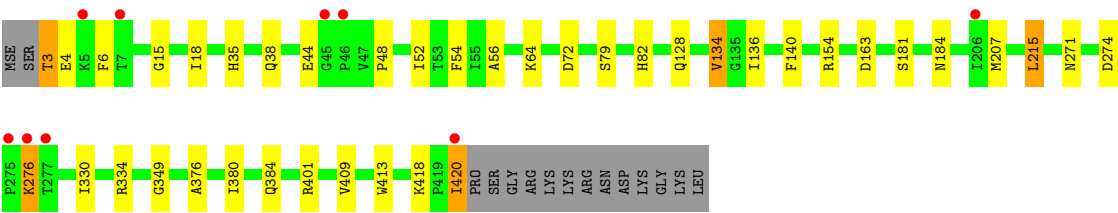


- Molecule 1: acyl-CoA hydrolase MpaH'



- Molecule 1: acyl-CoA hydrolase MpaH'





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.92Å 93.26Å 161.28Å 90.00° 95.12° 90.00°	Depositor
Resolution (Å)	46.63 – 1.99 46.63 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.1 (46.63-1.99) 99.1 (46.63-1.99)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.168 , 0.194 0.182 , 0.206	Depositor DCC
R_{free} test set	6548 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	24.3	Xtriage
Anisotropy	0.756	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 34.5	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	14163	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 5.18% 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	1.02	2/3435 (0.1%)	1.23	6/4657 (0.1%)
1	B	1.02	2/3410 (0.1%)	1.22	4/4623 (0.1%)
1	C	1.01	0/3435	1.22	7/4657 (0.2%)
1	D	1.02	0/3421	1.22	5/4637 (0.1%)
All	All	1.02	4/13701 (0.0%)	1.22	22/18574 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	406	HIS	CE1-NE2	5.87	1.38	1.32
1	A	343	GLY	C-O	5.25	1.28	1.24
1	B	365	HIS	CE1-NE2	5.21	1.37	1.32
1	A	228	THR	N-CA	5.15	1.49	1.46

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	384	GLN	CB-CG-CD	-6.69	101.22	112.60
1	B	349	GLY	CA-C-N	5.85	125.01	120.33
1	B	349	GLY	C-N-CA	5.85	125.01	120.33
1	A	384	GLN	CB-CA-C	5.80	120.72	110.85
1	B	15	GLY	CA-C-N	5.66	128.74	120.71
1	B	15	GLY	C-N-CA	5.66	128.74	120.71
1	A	349	GLY	CA-C-N	5.57	124.79	120.33
1	A	349	GLY	C-N-CA	5.57	124.79	120.33
1	D	15	GLY	CA-C-N	5.47	128.48	120.71
1	D	15	GLY	C-N-CA	5.47	128.48	120.71
1	A	218	MSE	CG-SE-CE	5.45	110.91	98.92
1	A	15	GLY	CA-C-N	5.38	128.35	120.71
1	A	15	GLY	C-N-CA	5.38	128.35	120.71
1	C	207	MSE	CG-SE-CE	5.34	110.67	98.92
1	C	140	PHE	CA-CB-CG	5.20	119.00	113.80
1	C	15	GLY	CA-C-N	5.18	128.07	120.71

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	15	GLY	C-N-CA	5.18	128.07	120.71
1	C	349	GLY	CA-C-N	5.15	124.45	120.33
1	C	349	GLY	C-N-CA	5.15	124.45	120.33
1	D	349	GLY	CA-C-N	5.08	124.39	120.33
1	D	349	GLY	C-N-CA	5.08	124.39	120.33
1	C	218	MSE	CG-SE-CE	5.08	110.09	98.92

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	3356	0	3320	32	1
1	B	3332	0	3290	36	0
1	C	3356	0	3320	32	0
1	D	3343	0	3308	30	1
2	A	204	0	0	0	0
2	B	210	0	0	4	0
2	C	188	0	0	1	0
2	D	174	0	0	2	0
All	All	14163	0	13238	117	1

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 4.

All (117) 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:271:ASN:HB2	2:B:646:HOH:O	1.79	0.82
1:C:201:ARG:NH1	1:C:204:ARG:HD3	1.96	0.80
1:C:5:LYS:HB3	1:C:75:LEU:HD23	1.63	0.79
1:B:276:LYS:HA	1:B:276:LYS:CE	2.18	0.74
1:A:28:GLN:HE22	1:D:154:ARG:HH12	1.38	0.72
1:D:52:ILE:CG2	1:D:134:VAL:HG13	2.21	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:ASP:OD1	1:A:248:THR:HG22	1.91	0.70
1:D:207:MSE:HE2	1:D:215:LEU:HG	1.74	0.70
1:D:3:THR:HG23	1:D:72:ASP:OD1	1.91	0.69
1:D:271:ASN:HB2	2:D:588:HOH:O	1.93	0.69
1:B:28:GLN:HE22	1:C:154:ARG:HH12	1.41	0.68
1:D:35:HIS:HD2	1:D:64:LYS:NZ	1.93	0.66
1:A:35:HIS:HD2	1:A:64:LYS:NZ	1.94	0.66
1:B:52:ILE:CG2	1:B:134:VAL:HG13	2.26	0.66
1:D:3:THR:HG21	1:D:6:PHE:HD1	1.60	0.66
1:B:35:HIS:HD2	1:B:64:LYS:NZ	1.94	0.65
1:C:35:HIS:HD2	1:C:64:LYS:NZ	1.95	0.64
1:D:3:THR:CG2	1:D:72:ASP:OD1	2.45	0.63
1:B:276:LYS:HA	1:B:276:LYS:HE2	1.80	0.62
1:D:52:ILE:HG23	1:D:134:VAL:HG13	1.81	0.62
1:B:339:ILE:CD1	2:B:610:HOH:O	2.49	0.60
1:B:154:ARG:HH12	1:C:28:GLN:HE22	1.48	0.60
1:D:44:GLU:H	1:D:44:GLU:CD	2.09	0.60
1:A:3:THR:HG22	1:A:8:ILE:HD11	1.83	0.60
1:B:52:ILE:HG23	1:B:134:VAL:HG13	1.84	0.58
1:C:3:THR:HG22	1:C:8:ILE:HD11	1.86	0.58
1:C:201:ARG:NH2	1:C:204:ARG:HG3	2.19	0.58
1:C:38:GLN:HE22	1:C:128:GLN:HE22	1.51	0.58
1:B:73:GLU:HG3	1:B:372:VAL:HG12	1.86	0.57
1:A:38:GLN:HE22	1:A:128:GLN:HE22	1.51	0.57
1:A:28:GLN:NE2	1:D:154:ARG:HH12	2.03	0.56
1:B:174:GLY:O	1:C:418:LYS:HE3	2.04	0.56
1:D:181:SER:H	1:D:184:ASN:HD22	1.54	0.55
1:B:181:SER:H	1:B:184:ASN:HD22	1.55	0.55
1:D:52:ILE:HG21	1:D:134:VAL:HG13	1.89	0.54
1:D:181:SER:H	1:D:184:ASN:ND2	2.06	0.54
1:B:181:SER:H	1:B:184:ASN:ND2	2.06	0.54
1:B:38:GLN:HE22	1:B:128:GLN:HE22	1.56	0.53
1:D:38:GLN:HE22	1:D:128:GLN:HE22	1.55	0.53
1:B:5:LYS:CB	1:B:75:LEU:HD23	2.39	0.53
1:A:5:LYS:CB	1:A:75:LEU:HD23	2.39	0.52
1:A:181:SER:H	1:A:184:ASN:ND2	2.08	0.52
1:B:274:ASP:CG	1:B:277:THR:HG22	2.34	0.52
1:A:67:TYR:OH	1:A:138:HIS:HD2	1.92	0.52
1:C:56:ALA:HA	1:C:136:ILE:O	2.10	0.52
1:C:67:TYR:OH	1:C:138:HIS:HD2	1.92	0.52
1:A:5:LYS:HB3	1:A:75:LEU:HD23	1.92	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:GLN:HE21	1:D:154:ARG:HH22	1.58	0.52
1:C:181:SER:H	1:C:184:ASN:ND2	2.08	0.52
1:A:56:ALA:HA	1:A:136:ILE:O	2.10	0.51
1:B:28:GLN:NE2	1:C:154:ARG:HH12	2.08	0.51
1:A:4:GLU:O	1:A:42:LYS:NZ	2.34	0.51
1:D:56:ALA:HA	1:D:136:ILE:O	2.11	0.50
1:A:77:GLN:HB2	1:A:380:ILE:CD1	2.40	0.50
1:A:203:ASN:O	1:A:206:ILE:HG22	2.11	0.50
1:B:56:ALA:HA	1:B:136:ILE:O	2.11	0.50
1:C:181:SER:H	1:C:184:ASN:HD22	1.60	0.50
1:B:35:HIS:HD2	1:B:64:LYS:HZ2	1.57	0.50
1:A:181:SER:H	1:A:184:ASN:HD22	1.59	0.49
1:B:274:ASP:HB3	1:B:277:THR:CG2	2.43	0.49
1:B:52:ILE:HG21	1:B:134:VAL:HG13	1.95	0.48
1:B:180:PRO:HB2	1:C:420:ILE:HD13	1.95	0.48
1:C:203:ASN:O	1:C:206:ILE:HG22	2.12	0.48
1:B:154:ARG:HH12	1:C:28:GLN:NE2	2.11	0.48
1:A:35:HIS:HD2	1:A:64:LYS:HZ2	1.61	0.48
1:B:274:ASP:HB3	1:B:277:THR:HG23	1.97	0.47
1:A:5:LYS:HE3	1:A:76:ASP:HB2	1.97	0.47
1:C:138:HIS:HE1	1:C:365:HIS:O	1.97	0.47
1:A:25:THR:O	1:D:401:ARG:NH2	2.47	0.47
1:D:3:THR:HG21	1:D:6:PHE:CD1	2.46	0.47
1:B:246:ASP:OD1	1:B:248:THR:HG22	2.14	0.47
1:C:201:ARG:NH1	1:C:204:ARG:CD	2.75	0.47
1:D:330:ILE:O	1:D:334:ARG:HG3	2.15	0.47
1:A:138:HIS:HE1	1:A:365:HIS:O	1.98	0.46
1:A:330:ILE:O	1:A:334:ARG:HG3	2.16	0.46
1:A:66:LEU:HD21	1:A:218:MSE:HG2	1.98	0.46
1:A:165:LEU:HD12	1:A:328:LEU:HD21	1.98	0.46
1:C:330:ILE:O	1:C:334:ARG:HG3	2.16	0.45
1:A:50:ASP:OD1	1:A:50:ASP:N	2.42	0.45
1:A:77:GLN:HB2	1:A:380:ILE:HD11	1.98	0.45
1:A:185:TYR:CE1	1:D:420:ILE:HG13	2.52	0.45
1:C:409:VAL:HB	1:C:413:TRP:CD1	2.52	0.45
1:C:50:ASP:OD1	1:C:50:ASP:N	2.41	0.45
1:B:54:PHE:CD1	1:B:134:VAL:HG22	2.52	0.44
1:D:54:PHE:CD1	1:D:134:VAL:HG22	2.52	0.44
1:B:203:ASN:O	1:B:206:ILE:HG22	2.17	0.44
1:A:409:VAL:HB	1:A:413:TRP:CD1	2.52	0.44
1:B:330:ILE:O	1:B:334:ARG:HG3	2.17	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:48:PRO:HG2	1:C:82:HIS:CG	2.52	0.44
1:A:173:LEU:HA	1:A:184:ASN:HD21	1.83	0.43
1:C:35:HIS:HD2	1:C:64:LYS:HZ1	1.65	0.43
1:C:173:LEU:HA	1:C:184:ASN:HD21	1.83	0.43
1:D:274:ASP:OD1	1:D:276:LYS:HG2	2.18	0.43
1:C:79:SER:OG	1:C:384:GLN:NE2	2.52	0.43
1:D:3:THR:HG23	1:D:72:ASP:CG	2.44	0.43
1:B:376:ALA:O	1:B:380:ILE:HG12	2.19	0.42
1:B:409:VAL:HB	1:B:413:TRP:CD1	2.55	0.42
1:D:409:VAL:HB	1:D:413:TRP:CD1	2.55	0.42
1:B:50:ASP:OD1	1:B:50:ASP:N	2.42	0.42
1:C:66:LEU:HD21	1:C:218:MSE:HG2	2.02	0.42
1:B:339:ILE:HD12	2:B:610:HOH:O	2.14	0.41
1:B:28:GLN:HE21	1:C:154:ARG:HH22	1.68	0.41
1:B:48:PRO:HG2	1:B:82:HIS:CG	2.55	0.41
1:A:48:PRO:HG2	1:A:82:HIS:CG	2.54	0.41
1:D:376:ALA:O	1:D:380:ILE:HG12	2.20	0.41
2:B:625:HOH:O	1:C:152:HIS:HD2	2.02	0.41
1:C:27:ASN:ND2	1:C:29:GLU:H	2.19	0.41
1:C:38:GLN:HE22	1:C:128:GLN:NE2	2.18	0.41
1:A:152:HIS:HD2	2:D:619:HOH:O	2.02	0.41
1:D:35:HIS:HD2	1:D:64:LYS:HZ1	1.63	0.41
1:A:28:GLN:NE2	1:D:154:ARG:HH22	2.17	0.41
1:C:165:LEU:HD12	1:C:328:LEU:HD21	2.03	0.41
1:D:48:PRO:HG2	1:D:82:HIS:CG	2.56	0.41
1:A:5:LYS:HB2	1:A:75:LEU:HD23	2.03	0.40
1:B:271:ASN:O	1:B:271:ASN:CG	2.64	0.40
1:D:35:HIS:HD2	1:D:64:LYS:HZ2	1.68	0.40
1:B:152:HIS:HD2	2:C:602:HOH:O	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:399:GLU:OE2	1:D:79:SER:OG[1_455]	2.11	0.09

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	418/433 (96%)	406 (97%)	11 (3%)	1 (0%)	43	42
1	B	415/433 (96%)	406 (98%)	8 (2%)	1 (0%)	43	42
1	C	418/433 (96%)	410 (98%)	7 (2%)	1 (0%)	43	42
1	D	416/433 (96%)	405 (97%)	10 (2%)	1 (0%)	43	42
All	All	1667/1732 (96%)	1627 (98%)	36 (2%)	4 (0%)	43	42

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	18	Ile
1	A	18	Ile
1	B	18	Ile
1	D	18	Ile

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	368/366 (100%)	363 (99%)	5 (1%)	59	66
1	B	364/366 (100%)	356 (98%)	8 (2%)	45	50
1	C	368/366 (100%)	360 (98%)	8 (2%)	45	50
1	D	366/366 (100%)	357 (98%)	9 (2%)	42	45
All	All	1466/1464 (100%)	1436 (98%)	30 (2%)	48	54

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	140	PHE
1	A	163	ASP
1	A	384	GLN
1	A	401	ARG
1	A	420	ILE
1	B	134	VAL
1	B	140	PHE
1	B	163	ASP
1	B	208	GLN
1	B	276	LYS
1	B	401	ARG
1	B	416	VAL
1	B	420	ILE
1	C	44	GLU
1	C	140	PHE
1	C	163	ASP
1	C	238	LYS
1	C	276	LYS
1	C	416	VAL
1	C	417	LEU
1	C	422	SER
1	D	3	THR
1	D	4	GLU
1	D	134	VAL
1	D	140	PHE
1	D	163	ASP
1	D	215	LEU
1	D	276	LYS
1	D	418	LYS
1	D	420	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	28	GLN
1	A	35	HIS
1	A	77	GLN
1	A	82	HIS
1	A	128	GLN
1	A	138	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	147	ASN
1	A	152	HIS
1	A	184	ASN
1	A	384	GLN
1	B	27	ASN
1	B	28	GLN
1	B	35	HIS
1	B	77	GLN
1	B	96	GLN
1	B	100	HIS
1	B	128	GLN
1	B	147	ASN
1	B	152	HIS
1	B	184	ASN
1	B	265	GLN
1	B	384	GLN
1	B	396	GLN
1	C	28	GLN
1	C	35	HIS
1	C	77	GLN
1	C	128	GLN
1	C	138	HIS
1	C	147	ASN
1	C	152	HIS
1	C	184	ASN
1	C	384	GLN
1	D	27	ASN
1	D	35	HIS
1	D	77	GLN
1	D	82	HIS
1	D	95	ASN
1	D	128	GLN
1	D	147	ASN
1	D	152	HIS
1	D	184	ASN
1	D	265	GLN

5.3.3 RNA ⓘ

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

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	408/433 (94%)	0.03	17 (4%) 40 39	19, 26, 47, 82	0
1	B	405/433 (93%)	0.07	8 (1%) 65 64	18, 27, 49, 88	0
1	C	408/433 (94%)	0.07	10 (2%) 58 58	19, 26, 48, 76	0
1	D	406/433 (93%)	0.02	9 (2%) 62 61	19, 26, 49, 77	0
All	All	1627/1732 (93%)	0.05	44 (2%) 56 55	18, 26, 49, 88	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	420	ILE	6.3
1	D	206	ILE	5.3
1	A	248	THR	4.8
1	C	3	THR	4.3
1	A	3	THR	4.0
1	B	277	THR	3.9
1	D	420	ILE	3.9
1	B	248	THR	3.2
1	C	275	PRO	3.2
1	C	422	SER	3.2
1	D	277	THR	3.2
1	B	275	PRO	3.1
1	B	4	GLU	3.0
1	B	276	LYS	3.0
1	C	185	TYR	2.9
1	A	45	GLY	2.8
1	C	45	GLY	2.7
1	A	4	GLU	2.7
1	D	5	LYS	2.6
1	A	277	THR	2.6
1	C	277	THR	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	46	PRO	2.6
1	A	403	GLY	2.6
1	D	275	PRO	2.5
1	A	204	ARG	2.5
1	A	422	SER	2.5
1	C	397	TRP	2.5
1	D	46	PRO	2.4
1	B	401	ARG	2.4
1	D	45	GLY	2.4
1	A	7	THR	2.3
1	C	245	ALA	2.3
1	A	276	LYS	2.3
1	C	276	LYS	2.3
1	B	206	ILE	2.2
1	A	127	ASP	2.2
1	A	275	PRO	2.1
1	C	4	GLU	2.1
1	A	245	ALA	2.1
1	A	185	TYR	2.1
1	D	276	LYS	2.0
1	A	44	GLU	2.0
1	A	206	ILE	2.0
1	D	7	THR	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 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.