



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:42 PM UTC

PDB ID : 4HCW / pdb_00004hcw
Title : Structure of a eukaryotic thiaminase-I
Authors : Kreinbring, C.A.; Hubbard, P.A.; Petsko, G.A.; Ringe, D.
Deposited on : 2012-10-01
Resolution : 2.71 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

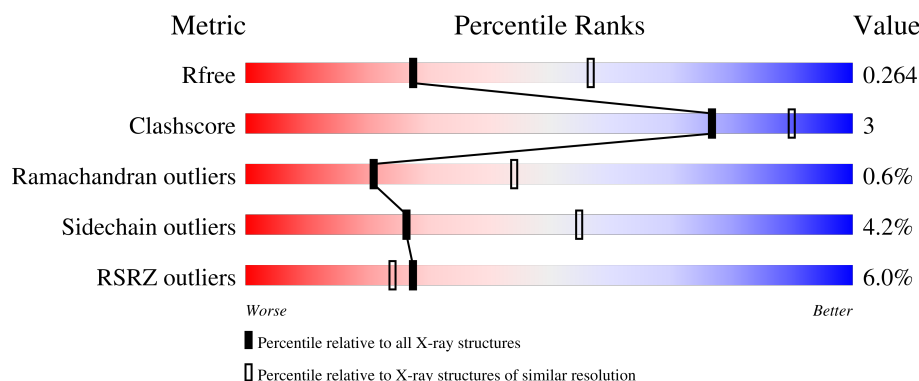
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.71 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	357	8% 87% 11% ..
1	B	357	11% 88% 8% ..
1	C	357	3% 88% 8% ..
1	D	357	5% 87% 10% ..
1	E	357	5% 88% 10% ..

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Mol	Chain	Length	Quality of chain
1	F	357	3% 90% 7% ..

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 16370 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 thiaminase-I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	352	Total	C	N	O	S	79	0	0
			2724	1731	441	545	7			
1	B	352	Total	C	N	O	S	114	0	0
			2724	1731	441	545	7			
1	C	352	Total	C	N	O	S	31	1	0
			2733	1739	441	546	7			
1	D	352	Total	C	N	O	S	44	1	0
			2732	1736	444	545	7			
1	E	352	Total	C	N	O	S	45	1	0
			2733	1739	441	546	7			
1	F	352	Total	C	N	O	S	59	0	0
			2724	1731	441	545	7			

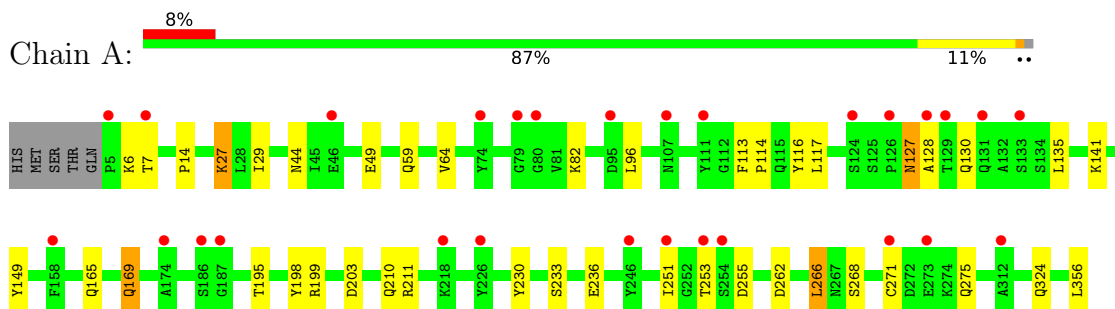
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	HIS	-	expression tag	UNP D2V4Z5
B	0	HIS	-	expression tag	UNP D2V4Z5
C	0	HIS	-	expression tag	UNP D2V4Z5
D	0	HIS	-	expression tag	UNP D2V4Z5
E	0	HIS	-	expression tag	UNP D2V4Z5
F	0	HIS	-	expression tag	UNP D2V4Z5

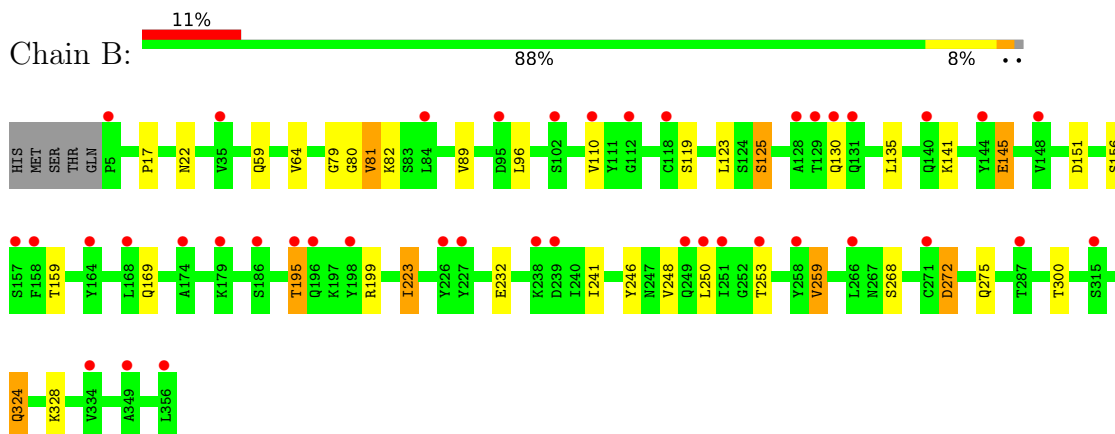
3 Residue-property plots

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.

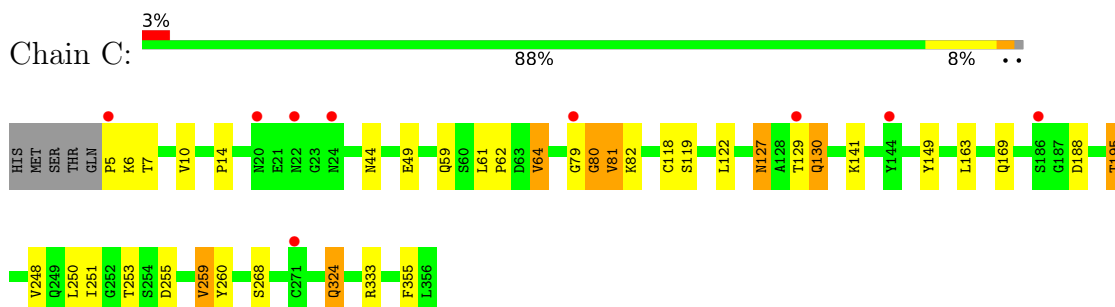
• Molecule 1: thiaminase-I



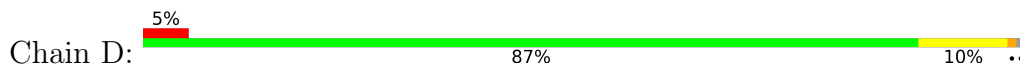
• Molecule 1: thiaminase-I

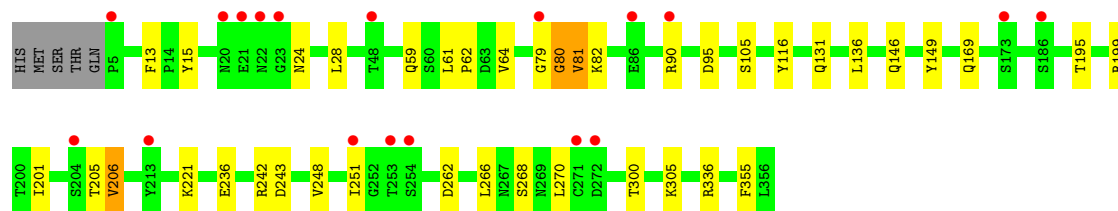


• Molecule 1: thiaminase-I

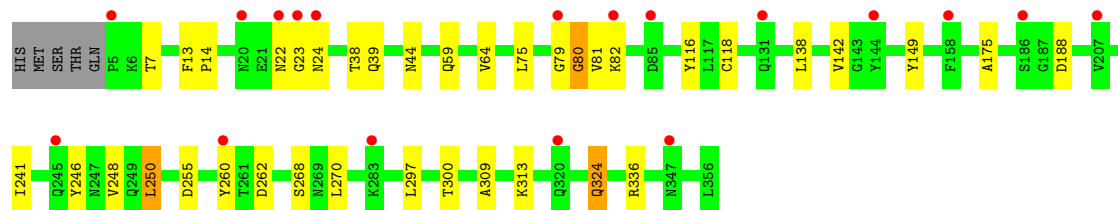
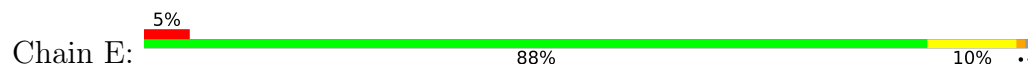


• Molecule 1: thiaminase-I

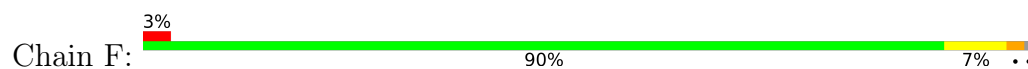




● Molecule 1: thiaminase-I



● Molecule 1: thiaminase-I



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	104.56Å 203.05Å 309.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.30 – 2.71 43.30 – 2.71	Depositor EDS
% Data completeness (in resolution range)	99.5 (43.30-2.71) 99.5 (43.30-2.71)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.21 (at 2.69Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.229 , 0.265 0.228 , 0.264	Depositor DCC
R_{free} test set	4484 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	70.7	Xtriage
Anisotropy	0.314	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 52.0	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	0.95	EDS
Total number of atoms	16370	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.12% 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	0.26	0/2773	0.69	2/3770 (0.1%)
1	B	0.26	0/2773	0.69	2/3770 (0.1%)
1	C	0.26	0/2786	0.68	2/3788 (0.1%)
1	D	0.27	0/2784	0.70	2/3784 (0.1%)
1	E	0.26	0/2786	0.69	2/3788 (0.1%)
1	F	0.26	0/2773	0.68	0/3770
All	All	0.26	0/16675	0.69	10/22670 (0.0%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	149	TYR	CA-C-N	5.83	125.30	119.24
1	D	149	TYR	C-N-CA	5.83	125.30	119.24
1	E	149	TYR	CA-C-N	5.74	125.21	119.24
1	E	149	TYR	C-N-CA	5.74	125.21	119.24
1	B	125	SER	CA-C-N	5.13	125.43	119.47
1	B	125	SER	C-N-CA	5.13	125.43	119.47
1	A	149	TYR	CA-C-N	5.05	125.08	119.32
1	A	149	TYR	C-N-CA	5.05	125.08	119.32
1	C	149	TYR	CA-C-N	5.05	125.08	119.32
1	C	149	TYR	C-N-CA	5.05	125.08	119.32

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	2724	0	2710	16	0
1	B	2724	0	2710	16	0
1	C	2733	0	2719	16	0
1	D	2732	0	2723	13	0
1	E	2733	0	2719	14	0
1	F	2724	0	2710	16	0
All	All	16370	0	16291	91	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 3.

All (91) 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:130:GLN:HG2	1:C:141:LYS:HB3	1.65	0.78
1:E:250:LEU:HB2	1:E:324:GLN:HG2	1.65	0.77
1:F:136:LEU:HD11	1:F:197:LYS:HD2	1.76	0.68
1:F:250:LEU:H	1:F:324:GLN:HG2	1.59	0.68
1:C:163:LEU:HD22	1:C:253:THR:HG21	1.77	0.67
1:D:82:LYS:HB2	1:D:268:SER:HA	1.77	0.66
1:B:272:ASP:N	1:B:272:ASP:OD1	2.33	0.61
1:A:82:LYS:HB2	1:A:268:SER:HA	1.83	0.60
1:D:95:ASP:HB3	1:D:305:LYS:HD2	1.85	0.59
1:B:82:LYS:HB2	1:B:268:SER:HA	1.85	0.58
1:B:156:SER:HB2	1:B:159:THR:HG22	1.85	0.58
1:E:7:THR:HG22	1:E:44:ASN:HB2	1.86	0.58
1:A:135:LEU:HB2	1:A:253:THR:HA	1.86	0.57
1:A:199:ARG:NH2	1:A:203:ASP:OD1	2.38	0.57
1:F:82:LYS:HB2	1:F:268:SER:HA	1.86	0.57
1:C:82:LYS:HB2	1:C:268:SER:HA	1.87	0.56
1:D:146:GLN:HB3	1:D:206:VAL:HG22	1.87	0.55
1:A:14:PRO:HB3	1:A:49:GLU:HB3	1.87	0.55
1:F:17:PRO:HG3	1:F:232:GLU:HB3	1.89	0.55
1:B:250:LEU:HB2	1:B:324:GLN:HG2	1.88	0.55
1:B:79:GLY:O	1:B:81:VAL:N	2.38	0.54
1:A:7:THR:HG22	1:A:44:ASN:HB3	1.89	0.54
1:C:127:ASN:N	1:C:127:ASN:OD1	2.39	0.54
1:F:186:SER:OG	1:F:188:ASP:OD1	2.26	0.53
1:B:130:GLN:HG2	1:B:141:LYS:HD3	1.91	0.52
1:E:118:CYS:HB2	1:E:260:TYR:HB2	1.91	0.52
1:D:105:SER:O	1:D:336:ARG:NH2	2.42	0.52
1:D:199:ARG:NH1	1:D:355:PHE:O	2.41	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:241:ILE:HG23	1:B:246:TYR:HB2	1.90	0.52
1:D:79:GLY:O	1:D:81:VAL:N	2.37	0.52
1:A:165:GLN:OE1	1:A:169:GLN:NE2	2.43	0.51
1:F:79:GLY:O	1:F:81:VAL:N	2.40	0.51
1:D:266:LEU:HB3	1:D:270:LEU:HD12	1.92	0.51
1:A:116:TYR:HB2	1:A:262:ASP:HB2	1.93	0.51
1:C:250:LEU:H	1:C:324:GLN:HG2	1.76	0.51
1:A:135:LEU:HG	1:A:251:ILE:HD12	1.92	0.51
1:F:8:LEU:HD23	1:F:274:LYS:HE2	1.93	0.51
1:E:79:GLY:O	1:E:81:VAL:N	2.42	0.50
1:F:119:SER:HB2	1:F:259:VAL:HG23	1.95	0.49
1:E:116:TYR:HB2	1:E:262:ASP:HB2	1.95	0.48
1:A:130:GLN:HG2	1:A:141:LYS:HB3	1.95	0.48
1:B:17:PRO:HD3	1:B:232:GLU:HG2	1.95	0.48
1:E:22:ASN:O	1:E:24:ASN:N	2.47	0.47
1:C:119:SER:HB2	1:C:259:VAL:HG23	1.97	0.47
1:B:135:LEU:HB2	1:B:253:THR:HA	1.97	0.47
1:B:250:LEU:O	1:B:328:LYS:NZ	2.47	0.47
1:F:13:PHE:HA	1:F:14:PRO:HD3	1.81	0.46
1:C:14:PRO:HB3	1:C:49:GLU:HB3	1.98	0.46
1:B:123:LEU:HD11	1:B:246:TYR:HB3	1.98	0.46
1:B:89:VAL:HG21	1:B:110:VAL:HG11	1.97	0.46
1:F:259:VAL:HG12	1:F:334:VAL:HG22	1.98	0.45
1:D:201:ILE:O	1:D:205:THR:HG23	2.16	0.45
1:D:242:ARG:NE	1:D:243:ASP:OD1	2.39	0.45
1:F:80:GLY:O	1:F:268:SER:N	2.46	0.45
1:A:127:ASN:N	1:A:127:ASN:OD1	2.49	0.45
1:B:195:THR:O	1:B:199:ARG:HB2	2.16	0.45
1:D:80:GLY:O	1:D:268:SER:N	2.42	0.45
1:C:195:THR:HG21	1:C:355:PHE:HB3	1.99	0.45
1:A:96:LEU:HD13	1:A:114:PRO:HG2	1.98	0.45
1:C:7:THR:HG22	1:C:44:ASN:HB3	1.98	0.45
1:C:10:VAL:HG13	1:C:64:VAL:HG22	1.99	0.45
1:E:82:LYS:HB2	1:E:268:SER:HA	1.98	0.45
1:F:123:LEU:HB3	1:F:227:TYR:HB3	1.99	0.45
1:B:151:ASP:HB3	1:B:159:THR:HG21	1.98	0.44
1:D:13:PHE:CE2	1:D:15:TYR:HB2	2.53	0.44
1:A:230:TYR:O	1:A:233:SER:OG	2.33	0.44
1:A:128:ALA:O	1:A:130:GLN:N	2.46	0.43
1:E:309:ALA:O	1:E:313:LYS:HG3	2.18	0.43
1:E:175:ALA:HB1	1:E:336:ARG:HH11	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:123:LEU:HD23	1:F:248:VAL:HG13	2.01	0.43
1:B:119:SER:HB2	1:B:259:VAL:HG22	2.01	0.42
1:C:118:CYS:HB3	1:C:260:TYR:HB2	2.01	0.42
1:E:241:ILE:HA	1:E:246:TYR:HB2	2.00	0.42
1:A:27:LYS:HB3	1:A:29:ILE:HG22	2.01	0.42
1:C:80:GLY:O	1:C:268:SER:N	2.50	0.41
1:E:13:PHE:HA	1:E:14:PRO:HD3	1.79	0.41
1:F:116:TYR:HB2	1:F:262:ASP:HB2	2.03	0.41
1:C:61:LEU:HA	1:C:62:PRO:HD3	1.85	0.41
1:C:79:GLY:O	1:C:81:VAL:N	2.42	0.41
1:D:116:TYR:HB2	1:D:262:ASP:HB2	2.02	0.41
1:E:80:GLY:O	1:E:268:SER:N	2.44	0.41
1:C:5:PRO:HB2	1:C:6:LYS:H	1.76	0.41
1:A:198:TYR:CE1	1:A:356:LEU:HD13	2.56	0.41
1:E:75:LEU:HD12	1:E:75:LEU:HA	1.94	0.41
1:C:122:LEU:HB2	1:C:251:ILE:HG21	2.03	0.40
1:D:61:LEU:HA	1:D:62:PRO:HD3	1.90	0.40
1:F:167:LEU:HD13	1:F:197:LYS:HB2	2.02	0.40
1:F:188:ASP:OD1	1:F:188:ASP:N	2.54	0.40
1:A:113:PHE:HE1	1:A:266:LEU:HD21	1.86	0.40
1:B:223:ILE:H	1:B:223:ILE:HG12	1.63	0.40
1:E:138:LEU:O	1:E:142:VAL:HG22	2.22	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	350/357 (98%)	334 (95%)	14 (4%)	2 (1%)	21	44
1	B	350/357 (98%)	327 (93%)	20 (6%)	3 (1%)	14	35
1	C	351/357 (98%)	338 (96%)	11 (3%)	2 (1%)	21	44

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	351/357 (98%)	338 (96%)	11 (3%)	2 (1%)	21	44
1	E	351/357 (98%)	334 (95%)	14 (4%)	3 (1%)	14	35
1	F	350/357 (98%)	333 (95%)	16 (5%)	1 (0%)	36	60
All	All	2103/2142 (98%)	2004 (95%)	86 (4%)	13 (1%)	21	44

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	145	GLU
1	A	271	CYS
1	B	80	GLY
1	D	80	GLY
1	E	23	GLY
1	C	255	ASP
1	D	24	ASN
1	E	255	ASP
1	F	80	GLY
1	E	80	GLY
1	A	255	ASP
1	B	22	ASN
1	C	80	GLY

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	314/319 (98%)	300 (96%)	14 (4%)	24	52
1	B	314/319 (98%)	299 (95%)	15 (5%)	23	50
1	C	315/319 (99%)	302 (96%)	13 (4%)	27	56
1	D	315/319 (99%)	299 (95%)	16 (5%)	21	48
1	E	315/319 (99%)	304 (96%)	11 (4%)	32	61
1	F	314/319 (98%)	302 (96%)	12 (4%)	29	58

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1887/1914 (99%)	1806 (96%)	81 (4%)	26 54

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LYS
1	A	27	LYS
1	A	59	GLN
1	A	64	VAL
1	A	117	LEU
1	A	127	ASN
1	A	169	GLN
1	A	195	THR
1	A	210	GLN
1	A	211	ARG
1	A	236	GLU
1	A	266	LEU
1	A	275	GLN
1	A	324	GLN
1	B	59	GLN
1	B	64	VAL
1	B	81	VAL
1	B	96	LEU
1	B	125	SER
1	B	145	GLU
1	B	169	GLN
1	B	195	THR
1	B	223	ILE
1	B	248	VAL
1	B	259	VAL
1	B	272	ASP
1	B	275	GLN
1	B	300	THR
1	B	324	GLN
1	C	59	GLN
1	C	64	VAL
1	C	81	VAL
1	C	127	ASN
1	C	129	THR
1	C	130	GLN
1	C	169	GLN
1	C	188	ASP

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Mol	Chain	Res	Type
1	C	195	THR
1	C	248	VAL
1	C	259	VAL
1	C	324	GLN
1	C	333	ARG
1	D	28	LEU
1	D	59	GLN
1	D	64	VAL
1	D	81	VAL
1	D	90[A]	ARG
1	D	90[B]	ARG
1	D	131	GLN
1	D	136	LEU
1	D	169	GLN
1	D	195	THR
1	D	206	VAL
1	D	221	LYS
1	D	236	GLU
1	D	248	VAL
1	D	251	ILE
1	D	300	THR
1	E	38	THR
1	E	39	GLN
1	E	59	GLN
1	E	64	VAL
1	E	188	ASP
1	E	248	VAL
1	E	250	LEU
1	E	270	LEU
1	E	297	LEU
1	E	300	THR
1	E	324	GLN
1	F	8	LEU
1	F	59	GLN
1	F	64	VAL
1	F	81	VAL
1	F	85	ASP
1	F	123	LEU
1	F	136	LEU
1	F	146	GLN
1	F	169	GLN
1	F	259	VAL

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Mol	Chain	Res	Type
1	F	275	GLN
1	F	324	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	169	GLN
1	A	249	GLN
1	A	324	GLN
1	B	24	ASN
1	B	30	ASN
1	C	120	ASN
1	C	169	GLN
1	C	249	GLN
1	D	24	ASN
1	D	97	HIS
1	D	120	ASN
1	D	146	GLN
1	D	169	GLN
1	D	275	GLN
1	E	59	GLN
1	E	120	ASN
1	E	247	ASN
1	F	97	HIS
1	F	247	ASN
1	F	316	ASN
1	F	320	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	352/357 (98%)	0.78	28 (7%) 18 15	43, 82, 125, 207	23 (6%)
1	B	352/357 (98%)	1.01	41 (11%) 9 8	56, 94, 141, 182	32 (9%)
1	C	352/357 (98%)	0.20	9 (2%) 57 54	30, 65, 101, 183	10 (2%)
1	D	352/357 (98%)	0.34	18 (5%) 33 29	35, 67, 103, 183	19 (5%)
1	E	351/357 (98%)	0.35	18 (5%) 33 29	30, 67, 106, 148	14 (3%)
1	F	352/357 (98%)	0.47	12 (3%) 48 44	40, 76, 115, 157	18 (5%)
All	All	2111/2142 (98%)	0.53	126 (5%) 27 24	30, 74, 119, 207	116 (5%)

All (126) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	271	CYS	5.4
1	B	5	PRO	5.3
1	F	255	ASP	5.0
1	B	186	SER	4.8
1	B	95	ASP	4.4
1	A	186	SER	4.4
1	E	5	PRO	4.3
1	C	129	THR	4.3
1	D	23	GLY	3.8
1	B	118	CYS	3.8
1	B	130	GLN	3.7
1	C	24	ASN	3.7
1	D	5	PRO	3.7
1	A	129	THR	3.7
1	B	164	TYR	3.6
1	B	249	GLN	3.5
1	A	253	THR	3.4
1	B	253	THR	3.3
1	C	5	PRO	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	174	ALA	3.3
1	B	195	THR	3.3
1	A	126	PRO	3.3
1	D	86	GLU	3.2
1	B	196	GLN	3.2
1	A	128	ALA	3.1
1	A	5	PRO	3.1
1	E	320	GLN	3.1
1	C	271	CYS	3.1
1	E	347	ASN	3.1
1	E	144[A]	TYR	3.0
1	B	174	ALA	3.0
1	D	90[A]	ARG	3.0
1	D	272	ASP	3.0
1	D	173	SER	2.9
1	A	107	ASN	2.9
1	E	24	ASN	2.9
1	E	85	ASP	2.9
1	D	253	THR	2.9
1	A	187	GLY	2.9
1	A	7	THR	2.9
1	C	79	GLY	2.9
1	E	207	VAL	2.9
1	B	84	LEU	2.8
1	E	20	ASN	2.8
1	B	144	TYR	2.8
1	D	213	TYR	2.8
1	B	140	GLN	2.8
1	F	5	PRO	2.7
1	A	46	GLU	2.6
1	E	23	GLY	2.6
1	A	246	TYR	2.6
1	B	226	TYR	2.6
1	A	79	GLY	2.6
1	A	80	GLY	2.6
1	B	227	TYR	2.6
1	B	128	ALA	2.6
1	D	22	ASN	2.6
1	E	186	SER	2.5
1	A	218	LYS	2.5
1	C	22	ASN	2.5
1	B	157	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	74	TYR	2.5
1	F	187	GLY	2.5
1	B	131	GLN	2.5
1	B	349	ALA	2.5
1	F	22	ASN	2.5
1	F	82	LYS	2.5
1	E	82	LYS	2.4
1	B	179	LYS	2.4
1	A	226	TYR	2.4
1	B	334	VAL	2.4
1	B	251	ILE	2.4
1	B	129	THR	2.4
1	C	186	SER	2.4
1	D	186	SER	2.4
1	E	245	GLN	2.4
1	A	111	TYR	2.3
1	B	168	LEU	2.3
1	F	253	THR	2.3
1	F	254	SER	2.3
1	A	312	ALA	2.3
1	C	20	ASN	2.3
1	D	204	SER	2.3
1	F	79	GLY	2.3
1	A	254	SER	2.3
1	B	112	GLY	2.3
1	B	250	LEU	2.3
1	B	271	CYS	2.2
1	A	251	ILE	2.2
1	E	260	TYR	2.2
1	B	158	PHE	2.2
1	B	238	LYS	2.2
1	E	79	GLY	2.2
1	B	315	SER	2.2
1	D	254	SER	2.2
1	F	186	SER	2.2
1	F	248	VAL	2.2
1	D	251	ILE	2.2
1	F	48	THR	2.2
1	B	239	ASP	2.2
1	A	133	SER	2.2
1	B	148	VAL	2.2
1	D	48	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	271	CYS	2.1
1	D	20	ASN	2.1
1	E	22	ASN	2.1
1	B	198	TYR	2.1
1	E	131	GLN	2.1
1	B	356	LEU	2.1
1	B	266	LEU	2.1
1	D	79	GLY	2.1
1	A	95	ASP	2.1
1	D	21	GLU	2.1
1	F	157	SER	2.1
1	B	110	VAL	2.1
1	A	158	PHE	2.1
1	A	273	GLU	2.0
1	A	124	SER	2.0
1	B	102	SER	2.0
1	E	283	LYS	2.0
1	A	131	GLN	2.0
1	C	144[A]	TYR	2.0
1	E	158	PHE	2.0
1	B	287	THR	2.0
1	B	35	VAL	2.0
1	B	258	TYR	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.