



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:51 PM UTC

PDB ID : 4TZO / pdb_00004tzo
Title : Structure of *C. elegans* HTP-1 bound to HIM-3 closure motif
Authors : Rosenberg, S.C.; Corbett, K.D.
Deposited on : 2014-07-10
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
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

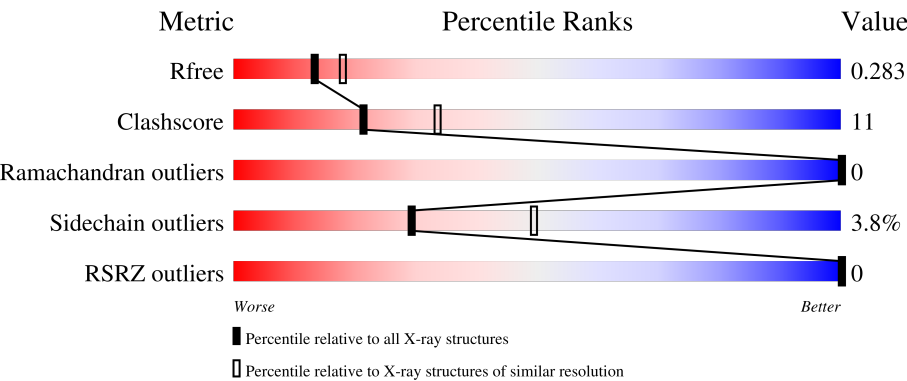
1 Overall quality at a glance i

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	253	69%23%• 7%
1	C	253	72%18%• 7%
1	E	253	71%21%• 7%
1	G	253	70%23%• 7%
2	B	20	40%30%30%

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Mol	Chain	Length	Quality of chain
2	D	20	 60% 10% 30%
2	F	20	 40% 30% 30%
2	H	20	 55% 15% 30%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8235 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 Protein HTP-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	236	Total	C	N	O	S	0	0	0
			1884	1189	321	362	12			
1	C	236	Total	C	N	O	S	0	0	0
			1884	1189	321	362	12			
1	E	236	Total	C	N	O	S	0	0	0
			1884	1189	321	362	12			
1	G	236	Total	C	N	O	S	0	0	0
			1884	1189	321	362	12			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	84	LEU	PRO	engineered mutation	UNP Q20305
C	84	LEU	PRO	engineered mutation	UNP Q20305
E	84	LEU	PRO	engineered mutation	UNP Q20305
G	84	LEU	PRO	engineered mutation	UNP Q20305

- Molecule 2 is a protein called C. elegans HIM-3 closure motif.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	14	Total	C	N	O	0	0	0
			104	64	19	21			
2	D	14	Total	C	N	O	0	0	0
			104	64	19	21			
2	F	14	Total	C	N	O	0	0	0
			104	64	19	21			
2	H	14	Total	C	N	O	0	0	0
			104	64	19	21			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	272	SER	-	expression tag	UNP G5EBG0
B	273	ASN	-	expression tag	UNP G5EBG0
B	274	ALA	-	expression tag	UNP G5EBG0
D	272	SER	-	expression tag	UNP G5EBG0
D	273	ASN	-	expression tag	UNP G5EBG0
D	274	ALA	-	expression tag	UNP G5EBG0
F	272	SER	-	expression tag	UNP G5EBG0
F	273	ASN	-	expression tag	UNP G5EBG0
F	274	ALA	-	expression tag	UNP G5EBG0
H	272	SER	-	expression tag	UNP G5EBG0
H	273	ASN	-	expression tag	UNP G5EBG0
H	274	ALA	-	expression tag	UNP G5EBG0

- Molecule 3 is water.

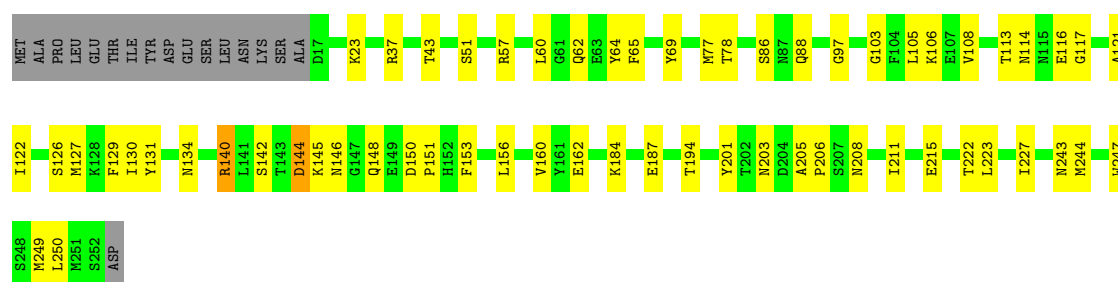
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	60	Total O 60 60	0	0
3	B	4	Total O 4 4	0	0
3	C	69	Total O 69 69	0	0
3	D	7	Total O 7 7	0	0
3	E	70	Total O 70 70	0	0
3	F	6	Total O 6 6	0	0
3	G	62	Total O 62 62	0	0
3	H	5	Total O 5 5	0	0

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.

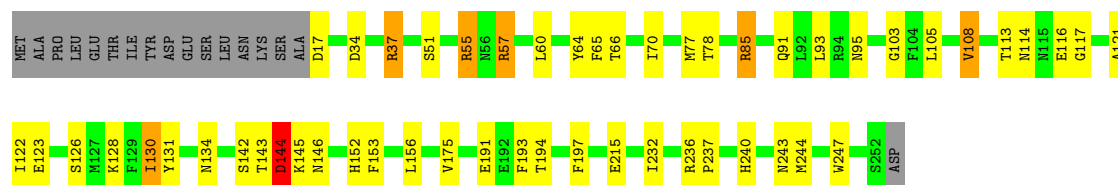
• Molecule 1: Protein HTP-1

Chain A: 



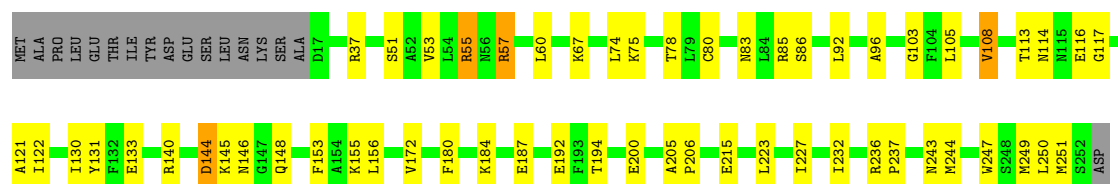
• Molecule 1: Protein HTP-1

Chain C: 



• Molecule 1: Protein HTP-1

Chain E: 



• Molecule 1: Protein HTP-1

Chain G: 

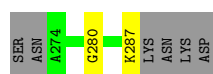




- Molecule 2: C. elegans HIM-3 closure motif



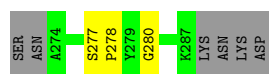
- Molecule 2: C. elegans HIM-3 closure motif



- Molecule 2: C. elegans HIM-3 closure motif



- Molecule 2: C. elegans HIM-3 closure motif



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.90Å 107.87Å 82.97Å 90.00° 90.10° 90.00°	Depositor
Resolution (Å)	45.22 – 2.40 45.22 – 2.40	Depositor EDS
% Data completeness (in resolution range)	94.0 (45.22-2.40) 93.7 (45.22-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 2.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.246 , 0.282 0.251 , 0.283	Depositor DCC
R_{free} test set	2205 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	23.9	Xtriage
Anisotropy	0.587	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 12.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.216 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8235	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.58 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.1945e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.33	0/1921	0.77	2/2595 (0.1%)
1	C	0.35	0/1921	0.76	2/2595 (0.1%)
1	E	0.32	0/1921	0.76	1/2595 (0.0%)
1	G	0.32	0/1921	0.75	1/2595 (0.0%)
2	B	0.33	0/105	0.75	0/140
2	D	0.29	0/105	0.67	0/140
2	F	0.34	0/105	0.66	0/140
2	H	0.38	0/105	0.70	0/140
All	All	0.33	0/8104	0.76	6/10940 (0.1%)

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	C	0	1
1	E	0	1
All	All	0	2

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	144	ASP	CA-C-N	6.64	133.65	121.70
1	A	144	ASP	C-N-CA	6.64	133.65	121.70
1	C	144	ASP	CA-C-N	6.25	132.95	121.70
1	C	144	ASP	C-N-CA	6.25	132.95	121.70
1	G	145	LYS	N-CA-C	-5.85	106.10	113.18
1	E	133	GLU	N-CA-C	5.03	116.76	111.28

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	144	ASP	Peptide
1	E	144	ASP	Peptide

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	1884	0	1840	44	0
1	C	1884	0	1840	37	0
1	E	1884	0	1840	43	0
1	G	1884	0	1840	46	0
2	B	104	0	103	3	0
2	D	104	0	103	3	0
2	F	104	0	103	5	0
2	H	104	0	103	5	0
3	A	60	0	0	0	0
3	B	4	0	0	0	0
3	C	69	0	0	1	0
3	D	7	0	0	0	0
3	E	70	0	0	2	0
3	F	6	0	0	0	0
3	G	62	0	0	1	0
3	H	5	0	0	0	0
All	All	8235	0	7772	171	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 (171) 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:E:144:ASP:HB3	1:E:146:ASN:H	0.94	1.09
1:E:144:ASP:HA	1:E:145:LYS:HB2	1.43	0.97
1:E:144:ASP:HB3	1:E:146:ASN:N	1.79	0.96
1:G:130:ILE:HD12	1:G:156:LEU:HB2	1.47	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:ILE:HD11	1:A:153:PHE:HA	1.49	0.94
1:A:103:GLY:O	1:A:140:ARG:NH2	2.03	0.90
1:G:130:ILE:HD11	1:G:153:PHE:HA	1.53	0.90
1:E:144:ASP:CA	1:E:145:LYS:HB2	2.01	0.90
1:C:144:ASP:HB3	1:C:145:LYS:HB3	1.54	0.87
1:E:144:ASP:CB	1:E:146:ASN:H	1.87	0.84
1:C:144:ASP:HB3	1:C:146:ASN:H	1.45	0.81
1:E:130:ILE:HD11	1:E:153:PHE:HA	1.64	0.80
1:A:144:ASP:HB3	1:A:146:ASN:H	1.46	0.79
1:A:227:ILE:HA	1:A:249:MET:HE1	1.65	0.77
1:G:142:SER:OG	1:G:150:ASP:OD2	2.02	0.77
1:A:116:GLU:HB2	1:A:194:THR:HG21	1.68	0.75
1:C:105:LEU:HD11	1:C:108:VAL:HG22	1.68	0.74
1:G:134:ASN:O	3:G:301:HOH:O	2.06	0.73
1:C:144:ASP:CB	1:C:146:ASN:H	2.02	0.72
1:G:116:GLU:HB2	1:G:194:THR:HG21	1.72	0.72
1:C:215:GLU:HG2	3:C:369:HOH:O	1.90	0.71
1:C:65:PHE:HB2	1:C:77:MET:HE3	1.73	0.71
1:A:144:ASP:HB3	1:A:145:LYS:HB3	1.75	0.68
1:E:130:ILE:HD12	1:E:156:LEU:HB2	1.75	0.68
1:G:156:LEU:HD21	1:G:243:ASN:HB2	1.76	0.67
1:G:130:ILE:CD1	1:G:153:PHE:HA	2.24	0.67
1:G:130:ILE:HD12	1:G:156:LEU:CB	2.24	0.66
1:E:55:ARG:NH1	1:E:60:LEU:O	2.29	0.66
1:C:34:ASP:OD1	1:C:37:ARG:HD3	1.96	0.65
1:A:144:ASP:HB3	1:A:146:ASN:N	2.10	0.65
1:A:144:ASP:CB	1:A:146:ASN:H	2.11	0.64
1:E:116:GLU:HB2	1:E:194:THR:HG21	1.79	0.64
1:G:107:GLU:OE2	1:G:109:SER:OG	2.16	0.64
1:A:227:ILE:HD11	1:A:247:TRP:HB3	1.80	0.63
1:G:127:MET:HE2	1:G:129:PHE:CZ	2.33	0.63
1:A:130:ILE:HD12	1:A:156:LEU:HB2	1.81	0.62
1:G:105:LEU:HD21	1:G:108:VAL:HG23	1.80	0.62
1:C:116:GLU:HB2	1:C:194:THR:HG21	1.82	0.61
1:C:144:ASP:CB	1:C:145:LYS:HB3	2.29	0.61
1:G:130:ILE:CD1	1:G:156:LEU:HB2	2.28	0.61
1:A:144:ASP:CB	1:A:145:LYS:HB3	2.30	0.61
1:E:105:LEU:HD11	1:E:108:VAL:HG22	1.83	0.60
1:E:140:ARG:NH2	3:E:302:HOH:O	2.35	0.60
1:C:51:SER:HB3	1:C:78:THR:HA	1.84	0.59
1:E:92:LEU:HD21	2:F:282:SER:HB2	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:PHE:HB2	1:A:77:MET:HE3	1.84	0.59
2:F:275:ARG:NH1	2:F:277:SER:O	2.35	0.58
1:E:227:ILE:HA	1:E:249:MET:HE1	1.85	0.58
1:C:130:ILE:HD11	1:C:153:PHE:HA	1.85	0.57
1:E:144:ASP:CB	1:E:145:LYS:HB2	2.34	0.57
1:A:227:ILE:CA	1:A:249:MET:HE1	2.34	0.57
1:C:126:SER:HG	1:C:247:TRP:HZ3	1.52	0.57
1:A:156:LEU:HD21	1:A:243:ASN:HB2	1.87	0.56
1:C:64:TYR:OH	1:C:85:ARG:NH1	2.37	0.56
1:E:187:GLU:HB2	1:E:250:LEU:HB3	1.86	0.56
1:C:130:ILE:HD12	1:C:156:LEU:HD12	1.87	0.56
1:G:57:ARG:NH2	1:G:123:GLU:OE1	2.39	0.55
1:G:113:THR:HG22	1:G:121:ALA:HA	1.88	0.55
1:E:223:LEU:HB3	1:E:227:ILE:HD13	1.87	0.55
1:A:223:LEU:HB3	1:A:227:ILE:HD13	1.89	0.55
1:G:145:LYS:HZ3	2:H:278:PRO:HG3	1.72	0.55
1:G:51:SER:HB3	1:G:78:THR:HA	1.89	0.54
1:E:53:VAL:O	1:E:57:ARG:HB2	2.07	0.54
1:A:105:LEU:HD21	1:A:108:VAL:HG23	1.89	0.54
1:C:17:ASP:OD1	1:C:17:ASP:N	2.40	0.54
1:A:23:LYS:HB3	1:A:69:TYR:HB3	1.89	0.54
1:C:113:THR:HG22	1:C:121:ALA:HA	1.90	0.53
1:A:103:GLY:HA2	1:A:131:TYR:CZ	2.43	0.53
1:A:227:ILE:HG12	1:A:247:TRP:HD1	1.74	0.52
1:E:236:ARG:NH1	1:E:237:PRO:O	2.43	0.52
1:G:65:PHE:HB2	1:G:77:MET:HE3	1.89	0.52
1:A:144:ASP:HB2	1:A:148:GLN:H	1.75	0.52
1:E:130:ILE:CD1	1:E:153:PHE:HA	2.35	0.52
1:A:142:SER:OG	1:A:150:ASP:OD2	2.26	0.52
1:C:130:ILE:CD1	1:C:153:PHE:HA	2.40	0.52
1:E:74:LEU:HD22	1:E:172:VAL:HG13	1.92	0.52
1:C:126:SER:OG	1:C:247:TRP:HZ3	1.94	0.51
1:G:145:LYS:NZ	2:H:278:PRO:HG3	2.25	0.51
1:C:215:GLU:O	2:D:280:GLY:HA2	2.11	0.51
1:G:17:ASP:OD2	1:G:161:TYR:OH	2.24	0.50
1:E:103:GLY:HA2	1:E:131:TYR:CZ	2.45	0.50
1:G:127:MET:CE	1:G:129:PHE:CZ	2.95	0.50
1:G:103:GLY:HA2	1:G:131:TYR:CZ	2.46	0.50
1:A:106:LYS:HE2	1:A:150:ASP:OD2	2.12	0.50
1:G:183:THR:HG22	1:G:184:LYS:HG2	1.93	0.49
1:A:113:THR:HG22	1:A:121:ALA:HA	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:156:LEU:HD21	1:E:243:ASN:HB2	1.95	0.48
1:E:232:ILE:HD12	1:E:244:MET:HG2	1.95	0.48
1:A:206:PRO:HB2	1:A:208:ASN:OD1	2.13	0.48
1:A:103:GLY:HA2	1:A:131:TYR:CE2	2.49	0.48
2:B:275:ARG:NH1	2:B:277:SER:O	2.37	0.48
1:C:103:GLY:HA2	1:C:131:TYR:CE1	2.49	0.48
1:G:206:PRO:HB2	1:G:208:ASN:OD1	2.13	0.48
1:C:156:LEU:HD21	1:C:243:ASN:HB2	1.95	0.48
1:C:232:ILE:HD12	1:C:244:MET:HG2	1.95	0.48
1:G:17:ASP:OD1	1:G:17:ASP:N	2.46	0.48
1:G:130:ILE:HD12	1:G:156:LEU:HD12	1.96	0.48
1:C:116:GLU:HA	1:C:117:GLY:HA2	1.52	0.47
1:E:215:GLU:O	2:F:280:GLY:HA2	2.14	0.47
1:C:70:ILE:HD13	1:C:175:VAL:HG21	1.96	0.47
1:G:127:MET:HE2	1:G:129:PHE:HZ	1.79	0.47
1:G:223:LEU:HD22	1:G:227:ILE:HD13	1.97	0.47
1:E:80:CYS:O	1:E:86:SER:OG	2.21	0.47
1:E:144:ASP:HB2	1:E:148:GLN:N	2.30	0.47
1:C:128:LYS:HB3	1:C:152:HIS:CD2	2.51	0.46
1:C:236:ARG:NH1	1:C:237:PRO:O	2.45	0.46
1:E:113:THR:HG22	1:E:121:ALA:HA	1.95	0.46
1:C:237:PRO:HD2	1:C:240:HIS:O	2.15	0.46
1:A:227:ILE:HA	1:A:249:MET:CE	2.39	0.46
1:C:103:GLY:HA2	1:C:131:TYR:CZ	2.50	0.46
1:E:103:GLY:HA2	1:E:131:TYR:CE2	2.51	0.46
1:E:227:ILE:HD11	1:E:247:TRP:HB3	1.98	0.46
1:E:51:SER:HB3	1:E:78:THR:HA	1.98	0.46
1:C:55:ARG:NH1	1:C:60:LEU:O	2.24	0.46
1:G:96:ALA:HB1	1:G:108:VAL:HG21	1.97	0.46
1:E:144:ASP:HB2	1:E:148:GLN:H	1.80	0.45
1:E:145:LYS:HE3	1:E:200:GLU:OE2	2.16	0.45
1:A:114:ASN:CG	1:A:122:ILE:HD11	2.41	0.45
1:A:126:SER:O	1:A:244:MET:HA	2.16	0.45
1:G:187:GLU:HB2	1:G:250:LEU:HB3	1.98	0.45
1:G:236:ARG:NH1	1:G:237:PRO:O	2.45	0.45
1:E:57:ARG:HA	1:E:57:ARG:HD2	1.61	0.45
1:E:205:ALA:HA	1:E:206:PRO:HD3	1.85	0.45
1:G:128:LYS:HB3	1:G:152:HIS:CD2	2.52	0.45
1:E:180:PHE:CE1	1:E:184:LYS:HD2	2.52	0.44
1:G:217:SER:O	2:H:277:SER:HB3	2.18	0.44
1:C:114:ASN:CG	1:C:122:ILE:HD11	2.41	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:ALA:HA	1:A:206:PRO:HD3	1.80	0.44
1:G:227:ILE:HA	1:G:249:MET:HE2	1.99	0.44
1:G:23:LYS:HB3	1:G:69:TYR:HB3	1.98	0.44
1:G:116:GLU:HA	1:G:117:GLY:HA2	1.57	0.44
1:A:222:THR:HG21	1:G:75:LYS:HE2	2.00	0.44
1:G:127:MET:HE2	1:G:127:MET:HB3	1.79	0.43
1:A:64:TYR:HB3	1:A:86:SER:HB3	2.00	0.43
1:A:187:GLU:HB2	1:A:250:LEU:HB3	1.99	0.43
1:E:114:ASN:CG	1:E:122:ILE:HD11	2.42	0.43
1:E:155:LYS:NZ	3:E:340:HOH:O	2.51	0.43
1:G:215:GLU:O	2:H:280:GLY:HA2	2.19	0.43
1:A:223:LEU:HD13	1:A:227:ILE:HD13	2.02	0.42
1:G:43:THR:HG23	1:G:97:GLY:HA3	2.01	0.42
1:G:145:LYS:NZ	2:H:278:PRO:CG	2.82	0.42
1:A:144:ASP:CA	1:A:145:LYS:HB3	2.50	0.42
1:G:103:GLY:HA2	1:G:131:TYR:CE2	2.54	0.42
1:A:150:ASP:HA	1:A:151:PRO:HD3	1.93	0.42
1:C:105:LEU:HD21	1:C:108:VAL:CG2	2.50	0.42
1:E:96:ALA:O	1:E:105:LEU:HD22	2.19	0.42
1:E:116:GLU:HA	1:E:117:GLY:HA2	1.58	0.42
1:A:60:LEU:HD21	2:B:285:ILE:HD13	2.01	0.42
1:A:43:THR:HG23	1:A:97:GLY:HA3	2.02	0.42
1:E:85:ARG:NH2	2:F:285:ILE:HG23	2.35	0.41
1:A:51:SER:HB3	1:A:78:THR:HA	2.02	0.41
1:C:191:GLU:O	2:D:287:LYS:HG2	2.20	0.41
1:E:83:ASN:HB3	1:E:86:SER:OG	2.20	0.41
1:G:57:ARG:HD2	1:G:57:ARG:HA	1.78	0.41
1:C:55:ARG:HD2	1:C:65:PHE:CD1	2.55	0.41
1:C:193:PHE:CE2	2:D:287:LYS:HD3	2.55	0.41
1:A:57:ARG:HA	1:A:57:ARG:HD2	1.75	0.41
1:G:17:ASP:HB2	1:G:164:GLY:H	1.85	0.41
1:C:93:LEU:HD21	1:C:197:PHE:CZ	2.55	0.41
1:A:215:GLU:O	2:B:280:GLY:HA2	2.21	0.41
1:C:57:ARG:NH2	1:C:123:GLU:OE1	2.47	0.41
1:E:67:LYS:HD2	1:E:75:LYS:HD2	2.03	0.41
1:A:160:VAL:O	1:A:162:GLU:HG3	2.21	0.40
1:A:116:GLU:HA	1:A:117:GLY:HA2	1.60	0.40
1:A:127:MET:HE2	1:A:129:PHE:CZ	2.56	0.40
1:E:60:LEU:HD21	2:F:285:ILE:HD13	2.04	0.40
1:A:201:TYR:HE1	1:A:211:ILE:HG13	1.86	0.40
1:G:105:LEU:HD11	1:G:108:VAL:CG2	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:127:MET:CE	1:G:129:PHE:HZ	2.34	0.40
1:C:91:GLN:NE2	1:C:95:ASN:OD1	2.54	0.40
1:G:144:ASP:HB3	1:G:146:ASN:H	1.86	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	234/253 (92%)	229 (98%)	5 (2%)	0	100	100
1	C	234/253 (92%)	228 (97%)	6 (3%)	0	100	100
1	E	234/253 (92%)	230 (98%)	4 (2%)	0	100	100
1	G	234/253 (92%)	229 (98%)	5 (2%)	0	100	100
2	B	12/20 (60%)	12 (100%)	0	0	100	100
2	D	12/20 (60%)	12 (100%)	0	0	100	100
2	F	12/20 (60%)	11 (92%)	1 (8%)	0	100	100
2	H	12/20 (60%)	11 (92%)	1 (8%)	0	100	100
All	All	984/1092 (90%)	962 (98%)	22 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	209/224 (93%)	202 (97%)	7 (3%)	33	55
1	C	209/224 (93%)	199 (95%)	10 (5%)	23	40
1	E	209/224 (93%)	203 (97%)	6 (3%)	37	60
1	G	209/224 (93%)	202 (97%)	7 (3%)	33	55
2	B	11/17 (65%)	9 (82%)	2 (18%)	2	2
2	D	11/17 (65%)	11 (100%)	0	100	100
2	F	11/17 (65%)	10 (91%)	1 (9%)	9	14
2	H	11/17 (65%)	11 (100%)	0	100	100
All	All	880/964 (91%)	847 (96%)	33 (4%)	29	49

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

Mol	Chain	Res	Type
1	A	37	ARG
1	A	62	GLN
1	A	88	GLN
1	A	134	ASN
1	A	140	ARG
1	A	184	LYS
1	A	203	ASN
2	B	276	ASP
2	B	282	SER
1	C	37	ARG
1	C	55	ARG
1	C	57	ARG
1	C	66	THR
1	C	85	ARG
1	C	108	VAL
1	C	130	ILE
1	C	134	ASN
1	C	142	SER
1	C	143	THR
1	E	37	ARG
1	E	55	ARG
1	E	57	ARG
1	E	108	VAL
1	E	192	GLU
1	E	251	MET
2	F	287	LYS
1	G	22	LYS

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Mol	Chain	Res	Type
1	G	37	ARG
1	G	40	ASN
1	G	116	GLU
1	G	138	VAL
1	G	203	ASN
1	G	218	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	ASN
1	A	87	ASN
1	A	134	ASN
1	C	68	ASN
1	C	87	ASN
1	C	91	GLN
1	C	95	ASN
1	C	134	ASN
2	D	283	GLN
1	E	87	ASN
1	G	87	ASN
2	H	283	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 [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	236/253 (93%)	-1.60	0 100 100	25, 38, 54, 76	0
1	C	236/253 (93%)	-1.67	0 100 100	15, 28, 46, 59	0
1	E	236/253 (93%)	-1.64	0 100 100	22, 32, 49, 63	0
1	G	236/253 (93%)	-1.66	0 100 100	17, 30, 47, 57	0
2	B	14/20 (70%)	-1.41	0 100 100	34, 46, 58, 60	0
2	D	14/20 (70%)	-1.48	0 100 100	20, 31, 52, 55	0
2	F	14/20 (70%)	-1.42	0 100 100	29, 42, 54, 57	0
2	H	14/20 (70%)	-1.39	0 100 100	25, 42, 56, 56	0
All	All	1000/1092 (91%)	-1.63	0 100 100	15, 32, 52, 76	0

There are no RSRZ outliers to report.

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.