



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 07:09 PM UTC

PDB ID : 2H66 / pdb_00002h66
Title : The Crystal Structure of Plasmodium Vivax 2-Cys peroxiredoxin
Authors : Wernimont, A.K.; Dong, A.; Zhao, Y.; Lew, J.; Melone, M.; Kozieradzki, I.; Weigelt, J.; Sundstrom, M.; Edwards, A.M.; Arrowsmith, C.H.; Bochkarev, A.; Hui, R.; Artz, J.D.; Structural Genomics Consortium (SGC)
Deposited on : 2006-05-30
Resolution : 2.50 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:

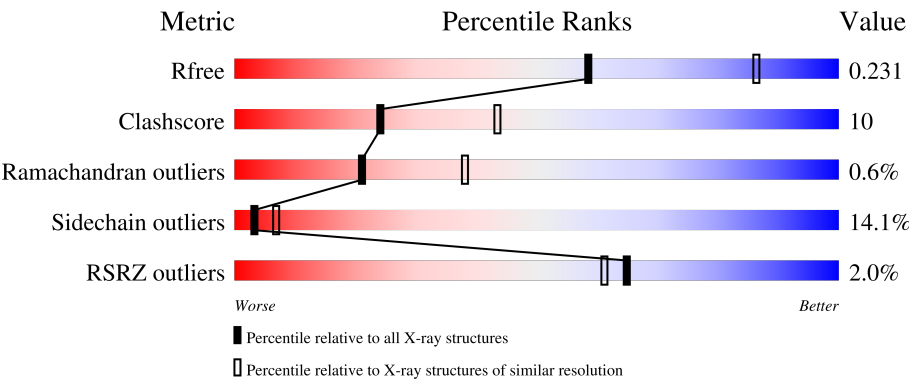
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	213	51%22%5%22%
1	B	213	3% 58%16%5%20%
1	C	213	% 60%15%20%
1	D	213	2% 61%17%21%
1	E	213	% 54%18%24%

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Mol	Chain	Length	Quality of chain
1	F	213	3%54%20%5%21%
1	G	213	57%15%7%20%
1	H	213	%58%18%•22%
1	I	213	2%63%14%6%17%
1	J	213	%53%19%••23%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 13448 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 PV-PF14_0368.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	167	Total	C	N	O	S	0	0	0
			1325	862	216	244	3			
1	B	171	Total	C	N	O	S	0	0	0
			1346	875	219	248	4			
1	C	170	Total	C	N	O	S	0	0	0
			1333	865	217	247	4			
1	D	169	Total	C	N	O	S	0	0	0
			1335	868	218	246	3			
1	E	161	Total	C	N	O	S	0	0	0
			1273	830	207	233	3			
1	F	168	Total	C	N	O	S	0	0	0
			1331	865	217	245	4			
1	G	171	Total	C	N	O	S	0	0	0
			1346	875	219	248	4			
1	H	167	Total	C	N	O	S	0	0	0
			1322	861	216	242	3			
1	I	177	Total	C	N	O	S	0	0	0
			1395	906	229	256	4			
1	J	163	Total	C	N	O	S	0	0	0
			1288	839	211	235	3			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	10	Total	O	0	0
			10	10		
2	B	8	Total	O	0	0
			8	8		
2	C	23	Total	O	0	0
			23	23		
2	D	14	Total	O	0	0
			14	14		

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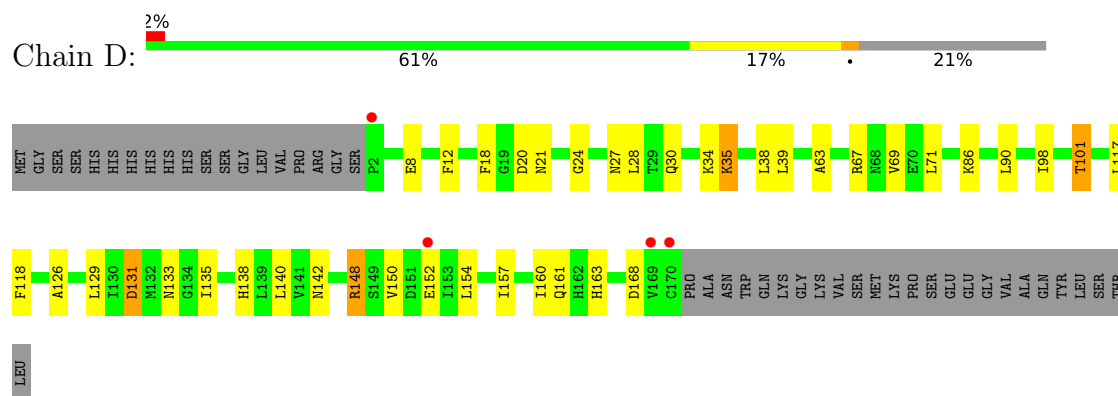
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	E	12	Total 12	O 12	0	0
2	F	12	Total 12	O 12	0	0
2	G	10	Total 10	O 10	0	0
2	H	18	Total 18	O 18	0	0
2	I	23	Total 23	O 23	0	0
2	J	24	Total 24	O 24	0	0

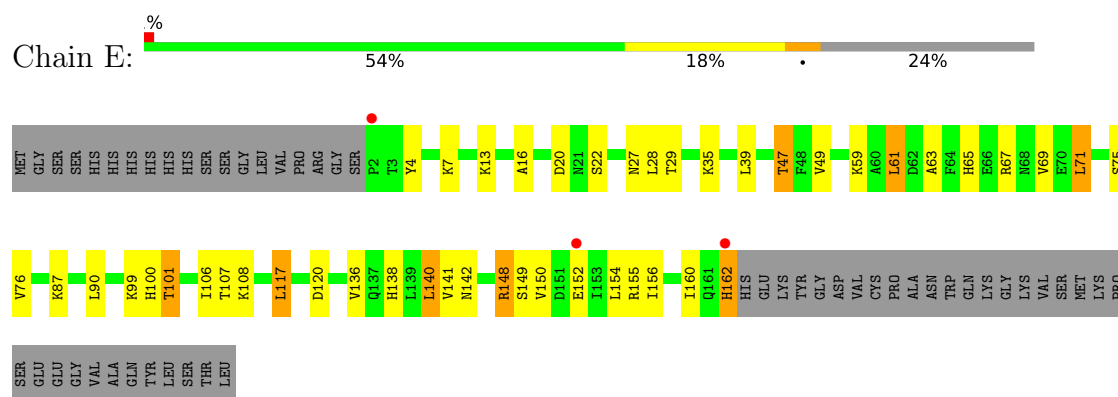
● Molecule 1: PV-PF14 0368



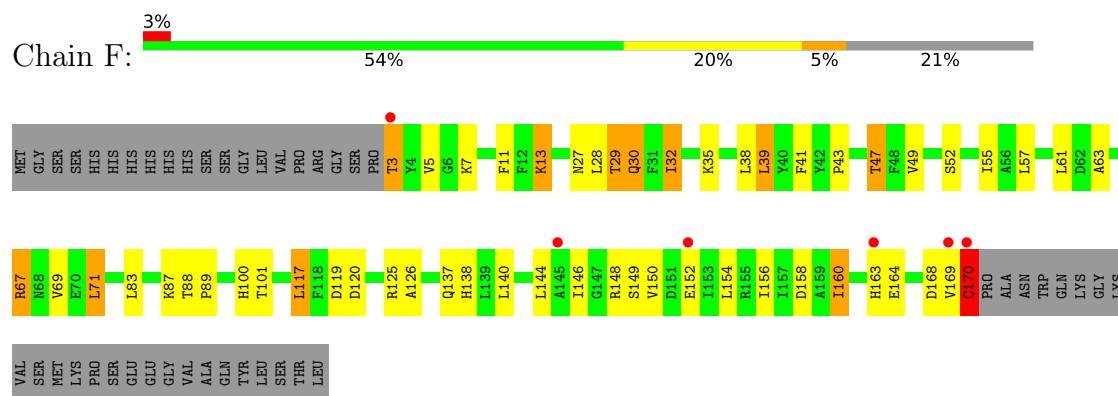
• Molecule 1: PV-PF14_0368



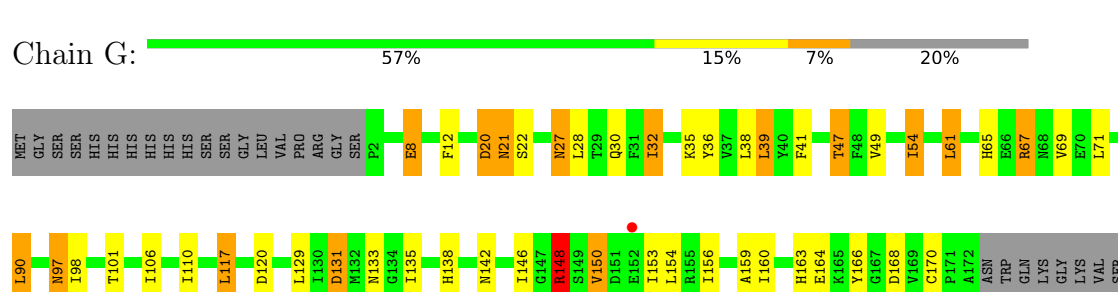
• Molecule 1: PV-PF14_0368



• Molecule 1: PV-PF14_0368



• Molecule 1: PV-PF14_0368



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• Molecule 1: PV-PF14_0368



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• Molecule 1: PV-PF14_0368



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• Molecule 1: PV-PF14_0368



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V69
E70
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V76
K79
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K92
G93
N97

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D105
I106
T107
K108
S109
I110
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R125
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L140
V141
M142
A145
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V150
D151
E152
I153
L154
R155
I156
I157
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.55Å 149.59Å 131.91Å 90.00° 104.88° 90.00°	Depositor
Resolution (Å)	41.56 – 2.50 41.56 – 2.50	Depositor EDS
% Data completeness (in resolution range)	94.7 (41.56-2.50) 94.7 (41.56-2.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.15 (at 2.51Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.194 , 0.232 0.194 , 0.231	Depositor DCC
R_{free} test set	4302 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å ²)	46.1	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 29.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.025 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13448	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% 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	0.82	0/1355	1.06	5/1834 (0.3%)
1	B	0.86	0/1377	1.05	7/1867 (0.4%)
1	C	0.97	3/1363 (0.2%)	1.07	3/1849 (0.2%)
1	D	0.83	0/1365	0.96	2/1848 (0.1%)
1	E	0.89	1/1301 (0.1%)	1.04	1/1762 (0.1%)
1	F	0.92	2/1360 (0.1%)	1.03	0/1841
1	G	0.88	0/1377	1.06	5/1867 (0.3%)
1	H	0.94	1/1352 (0.1%)	1.02	2/1830 (0.1%)
1	I	0.96	1/1428 (0.1%)	1.08	2/1937 (0.1%)
1	J	0.99	0/1317	1.11	6/1784 (0.3%)
All	All	0.91	8/13595 (0.1%)	1.05	33/18419 (0.2%)

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	B	0	3
1	C	0	2
All	All	0	5

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	168	ASP	C-O	11.71	1.47	1.23
1	F	170	CYS	C-O	9.55	1.42	1.23
1	E	162	HIS	C-O	7.31	1.38	1.23
1	C	168	ASP	CG-OD2	6.64	1.38	1.25
1	I	177	GLY	CA-C	6.43	1.63	1.52
1	C	69	VAL	CA-CB	5.25	1.60	1.54
1	F	32	ILE	CA-CB	5.15	1.60	1.54
1	C	135	ILE	CA-CB	5.10	1.60	1.54

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	50	CYS	CA-C-N	-7.67	111.89	119.87
1	A	50	CYS	C-N-CA	-7.67	111.89	119.87
1	H	168	ASP	CA-C-O	-7.33	108.35	120.80
1	D	131	ASP	N-CA-C	6.83	119.38	110.43
1	A	50	CYS	N-CA-C	6.59	124.38	109.81
1	A	35	LYS	N-CA-C	6.48	118.92	109.07
1	B	170	CYS	CA-C-N	6.08	141.60	127.00
1	B	170	CYS	C-N-CA	6.08	141.60	127.00
1	G	21	ASN	N-CA-C	6.03	123.65	110.80
1	J	101	THR	CB-CA-C	5.90	119.40	109.48
1	G	170	CYS	CA-C-N	-5.87	113.92	120.14
1	G	170	CYS	C-N-CA	-5.87	113.92	120.14
1	B	170	CYS	C-N-CD	-5.78	107.88	120.60
1	H	97	ASN	N-CA-C	5.78	118.58	109.39
1	J	131	ASP	CB-CA-C	-5.69	99.10	110.42
1	D	35	LYS	N-CA-C	5.66	117.67	109.07
1	I	97	ASN	N-CA-C	5.55	119.80	111.81
1	B	32	ILE	N-CA-C	5.54	115.74	110.42
1	C	168	ASP	OD1-CG-OD2	-5.54	109.60	122.90
1	J	110	ILE	N-CA-CB	5.47	116.95	110.55
1	B	35	LYS	N-CA-C	5.37	116.96	108.96
1	C	101	THR	N-CA-C	5.35	118.82	110.32
1	J	162	HIS	CA-C-N	-5.32	115.24	123.17
1	J	162	HIS	C-N-CA	-5.32	115.24	123.17
1	G	148	ARG	N-CA-C	-5.28	98.30	107.49
1	C	74	CYS	N-CA-C	5.26	116.94	108.32
1	B	170	CYS	N-CA-C	-5.21	98.29	109.81
1	B	131	ASP	N-CA-C	5.20	117.03	110.33
1	J	106	ILE	N-CA-C	5.08	115.80	110.62
1	A	146	ILE	N-CA-C	-5.08	107.89	113.43
1	E	162	HIS	CA-C-O	-5.06	112.19	120.80
1	I	146	ILE	N-CA-CB	5.04	116.21	110.72
1	G	131	ASP	CB-CA-C	-5.01	100.64	109.62

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	169	VAL	Peptide
1	B	170	CYS	Peptide
1	B	33	GLY	Peptide

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Mol	Chain	Res	Type	Group
1	C	168	ASP	Sidechain
1	C	6	GLY	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	1325	0	1327	30	0
1	B	1346	0	1342	25	0
1	C	1333	0	1316	27	0
1	D	1335	0	1331	25	0
1	E	1273	0	1286	25	0
1	F	1331	0	1334	35	0
1	G	1346	0	1342	37	0
1	H	1322	0	1325	18	0
1	I	1395	0	1386	40	0
1	J	1288	0	1294	31	0
2	A	10	0	0	0	0
2	B	8	0	0	0	0
2	C	23	0	0	3	0
2	D	14	0	0	1	0
2	E	12	0	0	3	0
2	F	12	0	0	0	0
2	G	10	0	0	2	0
2	H	18	0	0	0	0
2	I	23	0	0	4	0
2	J	24	0	0	1	0
All	All	13448	0	13283	266	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 10.

All (266) 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:171:PRO:HB2	1:B:172:ALA:O	1.43	1.16

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:149:SER:HB3	1:F:152:GLU:HG3	1.30	1.08
1:E:16:ALA:HB3	2:E:203:HOH:O	1.60	1.00
1:B:32:ILE:HD13	1:B:33:GLY:H	1.31	0.95
1:I:67:ARG:CZ	2:I:204:HOH:O	2.16	0.93
1:G:47:THR:HB	2:G:197:HOH:O	1.67	0.92
1:I:138:HIS:HE1	1:I:148:ARG:NH2	1.69	0.90
1:C:6:GLY:HA3	1:D:118:PHE:HE2	1.35	0.90
1:I:101:THR:HG22	2:I:202:HOH:O	1.71	0.89
1:E:160:ILE:HG12	1:F:144:LEU:HD12	1.56	0.87
1:D:101:THR:HG22	2:D:204:HOH:O	1.76	0.84
1:J:148:ARG:NE	1:J:152:GLU:OE2	2.09	0.84
1:I:138:HIS:CE1	1:I:148:ARG:NH2	2.46	0.83
1:G:90:LEU:H	1:G:90:LEU:HD12	1.45	0.82
1:J:12:PHE:HB2	1:J:110:ILE:HD12	1.60	0.82
1:B:171:PRO:HB2	1:B:172:ALA:C	2.05	0.80
1:B:90:LEU:H	1:B:90:LEU:HD12	1.46	0.79
1:A:13:LYS:HE3	1:A:25:GLU:OE2	1.83	0.79
1:B:32:ILE:HD13	1:B:33:GLY:N	1.99	0.77
1:H:27:ASN:HB2	1:H:30:GLN:HE21	1.50	0.77
1:I:142:ASN:OD1	1:J:138:HIS:HD2	1.68	0.77
1:F:27:ASN:H	1:F:30:GLN:NE2	1.84	0.75
1:I:35:LYS:HE2	1:I:68:ASN:ND2	2.02	0.75
1:G:39:LEU:HD13	1:G:41:PHE:CE1	2.24	0.73
1:F:149:SER:HB3	1:F:152:GLU:CG	2.14	0.71
1:E:47:THR:HB	2:E:199:HOH:O	1.90	0.70
1:E:149:SER:HB3	1:E:152:GLU:HG3	1.74	0.69
1:C:6:GLY:HA3	1:D:118:PHE:CE2	2.26	0.68
1:D:131:ASP:HB3	1:D:133:ASN:H	1.59	0.68
1:D:90:LEU:HD12	1:D:90:LEU:H	1.59	0.67
1:I:29:THR:O	1:I:32:ILE:HG22	1.94	0.67
1:G:67:ARG:HG3	1:G:67:ARG:HH11	1.58	0.67
1:B:171:PRO:CB	1:B:172:ALA:O	2.33	0.67
1:D:148:ARG:HE	1:D:152:GLU:CD	2.03	0.66
1:I:144:LEU:O	1:I:146:ILE:N	2.29	0.66
1:G:117:LEU:HD13	1:G:120:ASP:HA	1.77	0.66
1:C:101:THR:HG21	2:C:210:HOH:O	1.96	0.66
1:I:175:GLN:H	1:I:175:GLN:HE21	1.44	0.66
1:I:144:LEU:HB2	1:I:146:ILE:HD13	1.79	0.65
1:C:138:HIS:HD2	1:D:142:ASN:OD1	1.80	0.64
1:H:129:LEU:HD23	1:H:160:ILE:HD13	1.78	0.64
1:E:20:ASP:OD1	1:E:22:SER:HB2	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:20:ASP:O	1:G:22:SER:N	2.30	0.63
1:I:144:LEU:C	1:I:146:ILE:H	2.06	0.63
1:C:101:THR:HG22	2:C:198:HOH:O	1.99	0.63
1:I:138:HIS:CE1	1:I:148:ARG:HH21	2.15	0.63
1:E:117:LEU:CD1	1:E:120:ASP:HA	2.29	0.62
1:C:27:ASN:N	1:C:27:ASN:HD22	1.96	0.62
1:B:54:ILE:HD11	1:B:98:ILE:HD13	1.81	0.61
1:I:138:HIS:HD2	1:J:142:ASN:OD1	1.83	0.61
1:C:142:ASN:OD1	1:D:138:HIS:HD2	1.83	0.61
1:C:144:LEU:HD13	1:D:163:HIS:CD2	2.36	0.61
1:J:3:THR:HB	1:J:5:VAL:H	1.65	0.61
1:I:138:HIS:HE1	1:I:148:ARG:HH22	1.46	0.61
1:G:159:ALA:HB1	1:H:146:ILE:HG13	1.83	0.60
1:J:152:GLU:O	1:J:156:ILE:HG12	2.02	0.60
1:G:138:HIS:HD2	1:H:142:ASN:OD1	1.85	0.60
1:F:169:VAL:HG12	1:F:170:CYS:H	1.66	0.60
1:E:152:GLU:O	1:E:156:ILE:HG12	2.02	0.59
1:A:138:HIS:CE1	1:A:148:ARG:HH22	2.21	0.59
1:D:157:ILE:O	1:D:161:GLN:HG3	2.03	0.59
1:J:131:ASP:HB3	1:J:133:ASN:H	1.68	0.59
1:B:32:ILE:O	1:B:132:MET:O	2.21	0.59
1:I:138:HIS:CE1	1:I:148:ARG:HH22	2.17	0.59
1:F:47:THR:HG22	1:F:49:VAL:H	1.67	0.59
1:H:76:VAL:HG22	1:H:108:LYS:HG2	1.84	0.59
1:C:152:GLU:HG2	1:C:155:ARG:NH1	2.18	0.58
1:I:117:LEU:HD13	1:I:120:ASP:HA	1.85	0.58
1:G:27:ASN:HB3	1:G:30:GLN:HE21	1.68	0.58
1:A:12:PHE:HB2	1:A:110:ILE:HD12	1.86	0.58
1:F:27:ASN:HB2	1:F:30:GLN:HE21	1.69	0.57
1:E:138:HIS:CE1	1:F:140:LEU:HD11	2.40	0.57
1:I:112:LYS:HG2	1:I:117:LEU:HD12	1.86	0.57
1:C:26:VAL:C	1:C:27:ASN:HD22	2.13	0.57
1:J:76:VAL:HG22	1:J:108:LYS:HG2	1.85	0.57
1:A:63:ALA:O	1:A:67:ARG:HG2	2.05	0.56
1:F:156:ILE:O	1:F:160:ILE:HG23	2.04	0.56
1:I:21:ASN:HD21	1:I:86:LYS:HE2	1.71	0.56
1:I:138:HIS:HE1	1:I:148:ARG:HH21	1.50	0.56
1:A:155:ARG:HH12	1:B:149:SER:H	1.53	0.56
1:B:90:LEU:H	1:B:90:LEU:CD1	2.18	0.56
1:F:11:PHE:HE1	1:F:29:THR:HG22	1.71	0.56
1:G:27:ASN:HB3	1:G:30:GLN:NE2	2.19	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:20:ASP:O	1:C:21:ASN:HB2	2.05	0.56
1:A:107:THR:OG1	1:A:109:SER:HB2	2.06	0.56
1:F:39:LEU:HD13	1:F:41:PHE:CE1	2.41	0.56
1:A:35:LYS:HD2	1:A:70:GLU:HB2	1.87	0.55
1:F:43:PRO:HD3	1:F:125:ARG:HG3	1.87	0.55
1:E:61:LEU:HD13	1:E:65:HIS:CE1	2.41	0.55
1:I:21:ASN:HD21	1:I:86:LYS:CE	2.20	0.55
1:C:152:GLU:HG2	1:C:155:ARG:HH12	1.72	0.55
1:B:50:CYS:C	1:B:52:SER:H	2.14	0.55
1:I:142:ASN:OD1	1:J:138:HIS:CD2	2.56	0.54
1:G:129:LEU:HD23	1:G:160:ILE:HD13	1.90	0.54
1:G:97:ASN:C	1:G:97:ASN:HD22	2.16	0.54
1:D:90:LEU:HD12	1:D:90:LEU:N	2.21	0.54
1:A:108:LYS:HE3	1:A:121:SER:O	2.08	0.54
1:A:135:ILE:HG22	1:A:137:GLN:NE2	2.23	0.54
1:G:117:LEU:CD1	1:G:120:ASP:HA	2.37	0.53
1:G:131:ASP:HB3	1:G:133:ASN:H	1.73	0.53
1:A:142:ASN:OD1	1:B:138:HIS:HD2	1.90	0.53
1:F:71:LEU:HD13	1:F:100:HIS:CD2	2.43	0.53
1:A:117:LEU:CD1	1:A:120:ASP:HA	2.38	0.53
1:D:126:ALA:HA	1:D:140:LEU:O	2.08	0.53
1:F:57:LEU:HD23	1:F:150:VAL:HG23	1.91	0.52
1:G:32:ILE:HG12	2:G:202:HOH:O	2.09	0.52
1:G:61:LEU:HD13	1:G:65:HIS:CE1	2.45	0.52
1:C:39:LEU:HD13	1:C:41:PHE:CE1	2.45	0.52
1:E:117:LEU:HD11	1:E:120:ASP:HA	1.91	0.52
1:C:168:ASP:C	1:C:168:ASP:OD1	2.52	0.51
1:C:112:LYS:HG3	1:C:117:LEU:HD12	1.92	0.51
1:I:115:ASN:O	1:J:3:THR:HG23	2.10	0.51
1:I:144:LEU:C	1:I:146:ILE:N	2.68	0.51
1:H:140:LEU:HD23	1:H:140:LEU:C	2.35	0.51
1:C:110:ILE:HD13	1:C:110:ILE:N	2.26	0.51
1:G:67:ARG:HG3	1:G:67:ARG:NH1	2.25	0.51
1:F:47:THR:CG2	1:F:49:VAL:H	2.23	0.51
1:B:90:LEU:HD12	1:B:90:LEU:N	2.22	0.51
1:F:137:GLN:HG3	1:F:160:ILE:HD13	1.92	0.51
1:J:11:PHE:HE1	1:J:29:THR:HG23	1.76	0.51
1:J:20:ASP:OD1	1:J:20:ASP:C	2.54	0.51
1:A:15:GLU:OE2	1:A:79:LYS:HE2	2.11	0.50
1:A:54:ILE:HD11	1:A:98:ILE:HD13	1.93	0.50
1:F:117:LEU:HD13	1:F:120:ASP:HA	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:152:GLU:O	1:H:156:ILE:HG12	2.11	0.50
1:I:67:ARG:NH2	2:I:204:HOH:O	2.35	0.50
1:J:131:ASP:HB2	1:J:135:ILE:H	1.76	0.50
1:C:144:LEU:HD13	1:D:163:HIS:CG	2.46	0.50
1:A:138:HIS:HE1	1:A:148:ARG:HH22	1.59	0.50
1:G:160:ILE:O	1:G:164:GLU:HG3	2.11	0.50
1:I:63:ALA:O	1:I:67:ARG:HG3	2.11	0.50
1:H:18:PHE:HE1	1:H:24:GLY:HA3	1.76	0.50
1:I:163:HIS:HD2	2:I:206:HOH:O	1.95	0.50
1:C:27:ASN:H	1:C:30:GLN:NE2	2.10	0.49
1:B:64:PHE:HB3	1:B:69:VAL:HG22	1.95	0.49
1:B:131:ASP:HB3	1:B:133:ASN:H	1.77	0.49
1:J:163:HIS:O	1:J:164:GLU:CB	2.61	0.49
1:D:27:ASN:HB3	1:D:30:GLN:HE21	1.77	0.49
1:G:54:ILE:HD11	1:G:98:ILE:HG21	1.94	0.49
1:C:93:GLY:O	2:C:202:HOH:O	2.19	0.49
1:H:20:ASP:O	1:H:21:ASN:HB2	2.12	0.49
1:A:55:ILE:HD13	1:A:56:ALA:N	2.26	0.49
1:G:54:ILE:CD1	1:G:98:ILE:HD13	2.42	0.49
1:A:15:GLU:OE2	1:A:79:LYS:CE	2.61	0.48
1:I:163:HIS:HE1	1:I:168:ASP:OD1	1.96	0.48
1:E:27:ASN:HD22	1:E:29:THR:H	1.62	0.48
1:F:3:THR:O	1:F:7:LYS:HE2	2.13	0.48
1:J:117:LEU:HD13	1:J:120:ASP:HA	1.94	0.48
1:F:163:HIS:CE1	1:F:168:ASP:HB3	2.49	0.48
1:A:131:ASP:HB3	1:A:133:ASN:H	1.77	0.48
1:J:70:GLU:OE2	1:J:70:GLU:HA	2.13	0.48
1:C:27:ASN:H	1:C:30:GLN:HE21	1.62	0.48
1:I:8:GLU:H	1:I:8:GLU:CD	2.22	0.48
1:I:38:LEU:HD11	1:I:153:ILE:CD1	2.44	0.48
1:D:90:LEU:H	1:D:90:LEU:CD1	2.26	0.47
1:I:39:LEU:HD13	1:I:41:PHE:CE1	2.49	0.47
1:C:27:ASN:N	1:C:27:ASN:ND2	2.62	0.47
1:F:160:ILE:O	1:F:164:GLU:HB2	2.14	0.47
1:G:39:LEU:CD1	1:G:41:PHE:CE1	2.97	0.47
1:A:110:ILE:N	1:A:110:ILE:HD13	2.30	0.47
1:E:160:ILE:HG12	1:F:144:LEU:CD1	2.38	0.47
1:D:63:ALA:O	1:D:67:ARG:HG2	2.15	0.47
1:E:138:HIS:CE1	1:F:140:LEU:CD1	2.97	0.47
1:E:141:VAL:HG12	1:F:5:VAL:HG11	1.97	0.47
1:E:152:GLU:HG2	1:F:152:GLU:HG2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:21:ASN:ND2	1:I:86:LYS:NZ	2.62	0.47
1:D:12:PHE:CE1	1:D:28:LEU:HD12	2.49	0.47
1:F:13:LYS:HG3	1:F:27:ASN:ND2	2.30	0.47
1:D:86:LYS:HE3	1:D:98:ILE:HG13	1.96	0.46
1:G:61:LEU:CD1	1:G:65:HIS:CE1	2.98	0.46
1:A:117:LEU:HD13	1:A:120:ASP:HA	1.96	0.46
1:D:163:HIS:CE1	1:D:168:ASP:OD1	2.68	0.46
1:E:138:HIS:HE1	1:F:140:LEU:HD11	1.78	0.46
1:H:64:PHE:HB3	1:H:69:VAL:HG22	1.98	0.46
1:J:140:LEU:HD23	1:J:141:VAL:N	2.30	0.46
1:C:126:ALA:HA	1:C:140:LEU:O	2.15	0.46
1:A:149:SER:HB2	1:A:152:GLU:HG3	1.98	0.46
1:A:163:HIS:O	1:A:167:GLY:HA2	2.15	0.46
1:F:13:LYS:HG3	1:F:27:ASN:HD21	1.79	0.46
1:F:67:ARG:NH1	1:F:158:ASP:OD2	2.48	0.46
1:G:163:HIS:O	1:G:166:TYR:O	2.34	0.46
1:A:18:PHE:HE1	1:A:24:GLY:HA3	1.80	0.45
1:A:138:HIS:HD2	1:B:142:ASN:OD1	1.99	0.45
1:H:62:ASP:O	1:H:66:GLU:HG3	2.16	0.45
1:G:47:THR:HG22	1:G:49:VAL:H	1.81	0.45
1:H:140:LEU:HD23	1:H:141:VAL:N	2.31	0.45
1:F:88:THR:HA	1:F:89:PRO:HD2	1.84	0.45
1:G:20:ASP:O	1:G:20:ASP:CG	2.60	0.45
1:G:148:ARG:NH2	1:G:156:ILE:HD13	2.31	0.45
1:F:163:HIS:C	1:F:163:HIS:CD2	2.93	0.45
1:G:12:PHE:HB2	1:G:110:ILE:HG12	1.98	0.45
1:I:61:LEU:HD12	1:I:99:LYS:HB2	1.98	0.45
1:G:36:TYR:OH	1:G:164:GLU:OE1	2.32	0.45
1:B:61:LEU:CD1	1:B:65:HIS:CE1	3.00	0.44
1:D:20:ASP:O	1:D:21:ASN:HB2	2.17	0.44
1:J:8:GLU:CD	1:J:8:GLU:H	2.25	0.44
1:E:63:ALA:O	1:E:67:ARG:HG2	2.18	0.44
1:I:57:LEU:HB3	1:I:71:LEU:HD11	1.99	0.44
1:B:12:PHE:HB2	1:B:110:ILE:HG12	1.99	0.44
1:B:171:PRO:HB2	1:B:172:ALA:CA	2.47	0.44
1:E:142:ASN:OD1	1:F:138:HIS:HD2	2.01	0.44
1:J:76:VAL:CG2	1:J:108:LYS:HG2	2.48	0.44
1:C:71:LEU:HD13	1:C:100:HIS:CG	2.52	0.44
1:J:23:PHE:CD2	1:J:79:LYS:HD2	2.53	0.44
1:I:35:LYS:HE2	1:I:68:ASN:HD22	1.80	0.44
1:D:131:ASP:HB2	1:D:135:ILE:H	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:150:VAL:HA	1:G:153:ILE:HD12	2.00	0.43
1:J:67:ARG:NH1	1:J:158:ASP:OD1	2.50	0.43
1:A:16:ALA:HB2	1:A:103:LEU:HD23	2.00	0.43
1:C:76:VAL:HG22	1:C:108:LYS:HG2	2.00	0.43
1:D:129:LEU:HD23	1:D:160:ILE:HD12	1.99	0.43
1:J:18:PHE:HA	1:J:101:THR:HB	1.98	0.43
1:B:39:LEU:HD13	1:B:41:PHE:CE1	2.53	0.43
1:B:153:ILE:O	1:B:157:ILE:HG13	2.18	0.43
1:E:101:THR:HG22	2:E:197:HOH:O	2.18	0.43
1:F:63:ALA:O	1:F:67:ARG:HG2	2.18	0.43
1:H:107:THR:O	1:H:108:LYS:HB2	2.18	0.43
1:A:148:ARG:HD3	1:A:153:ILE:HG13	2.00	0.43
1:J:57:LEU:HB3	1:J:71:LEU:HD11	2.00	0.43
1:B:61:LEU:HD13	1:B:65:HIS:CE1	2.53	0.43
1:E:47:THR:HG23	1:E:49:VAL:H	1.83	0.43
1:I:76:VAL:HG22	1:I:108:LYS:HG2	1.99	0.43
1:H:76:VAL:CG2	1:H:108:LYS:HG2	2.49	0.43
1:J:11:PHE:CE1	1:J:29:THR:HG23	2.53	0.43
1:D:131:ASP:HB3	1:D:133:ASN:N	2.30	0.43
1:G:47:THR:CG2	1:G:49:VAL:H	2.32	0.43
1:C:140:LEU:HD12	1:D:140:LEU:HD12	2.01	0.42
1:F:29:THR:HA	1:F:32:ILE:HG12	2.00	0.42
1:F:126:ALA:HA	1:F:140:LEU:O	2.19	0.42
1:G:67:ARG:HD3	1:G:67:ARG:N	2.34	0.42
1:J:105:ASP:OD2	1:J:110:ILE:HG12	2.19	0.42
1:I:5:VAL:HG11	1:J:141:VAL:HG12	2.00	0.42
1:G:131:ASP:HB2	1:G:135:ILE:H	1.84	0.42
1:J:12:PHE:HB2	1:J:110:ILE:CD1	2.40	0.42
1:E:4:TYR:CD2	1:E:136:VAL:HG21	2.55	0.42
1:I:53:GLU:OE1	1:I:148:ARG:O	2.37	0.42
1:H:45:ASP:O	1:H:46:PHE:HB2	2.20	0.42
1:A:140:LEU:HD23	1:A:140:LEU:C	2.43	0.42
1:A:67:ARG:HB2	1:A:69:VAL:CG1	2.51	0.41
1:B:59:LYS:HD3	1:B:59:LYS:HA	1.94	0.41
1:F:83:LEU:HD11	1:F:87:LYS:HE3	2.02	0.41
1:H:27:ASN:H	1:H:30:GLN:NE2	2.17	0.41
1:A:36:TYR:OH	1:A:164:GLU:OE1	2.37	0.41
1:B:50:CYS:C	1:B:52:SER:N	2.78	0.41
1:G:27:ASN:O	1:G:30:GLN:HG2	2.20	0.41
1:G:142:ASN:OD1	1:H:138:HIS:HD2	2.03	0.41
1:E:71:LEU:HD13	1:E:100:HIS:CG	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:129:LEU:HD23	1:C:160:ILE:HD12	2.03	0.41
1:E:107:THR:O	1:E:108:LYS:HB2	2.21	0.41
1:G:97:ASN:C	1:G:97:ASN:ND2	2.78	0.41
1:I:21:ASN:HD21	1:I:86:LYS:NZ	2.19	0.41
1:I:144:LEU:HD12	1:J:160:ILE:HG12	2.03	0.41
1:J:132:MET:CE	2:J:203:HOH:O	2.68	0.41
1:A:107:THR:O	1:A:108:LYS:HB2	2.20	0.41
1:D:18:PHE:CE1	1:D:24:GLY:HA3	2.56	0.41
1:E:20:ASP:OD1	1:E:22:SER:CB	2.68	0.41
1:G:8:GLU:H	1:G:8:GLU:HG2	1.52	0.41
1:H:138:HIS:HE1	1:H:148:ARG:HH12	1.68	0.41
1:E:140:LEU:CD2	1:E:148:ARG:HG3	2.51	0.40
1:G:27:ASN:OD1	1:G:27:ASN:C	2.65	0.40
1:B:58:ASP:O	1:B:61:LEU:HB2	2.21	0.40
1:I:114:TYR:O	1:I:115:ASN:HB2	2.22	0.40
1:A:4:TYR:CD2	1:A:136:VAL:HG21	2.57	0.40
1:C:51:PRO:O	1:C:55:ILE:HG12	2.21	0.40
1:J:18:PHE:HE1	1:J:24:GLY:HA3	1.87	0.40
1:J:125:ARG:CZ	1:J:145:ALA:HB2	2.52	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	165/213 (78%)	153 (93%)	10 (6%)	2 (1%)	10	20
1	B	169/213 (79%)	158 (94%)	8 (5%)	3 (2%)	6	12
1	C	168/213 (79%)	157 (94%)	9 (5%)	2 (1%)	10	20
1	D	167/213 (78%)	160 (96%)	7 (4%)	0	100	100
1	E	159/213 (75%)	153 (96%)	6 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	166/213 (78%)	156 (94%)	10 (6%)	0	100	100
1	G	169/213 (79%)	162 (96%)	6 (4%)	1 (1%)	21	38
1	H	165/213 (78%)	160 (97%)	5 (3%)	0	100	100
1	I	175/213 (82%)	166 (95%)	8 (5%)	1 (1%)	21	38
1	J	161/213 (76%)	155 (96%)	5 (3%)	1 (1%)	21	38
All	All	1664/2130 (78%)	1580 (95%)	74 (4%)	10 (1%)	21	38

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	167	GLY
1	C	168	ASP
1	G	21	ASN
1	I	145	ALA
1	C	167	GLY
1	J	131	ASP
1	A	50	CYS
1	B	43	PRO
1	B	51	PRO
1	B	171	PRO

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	145/184 (79%)	123 (85%)	22 (15%)	3	5
1	B	147/184 (80%)	122 (83%)	25 (17%)	2	4
1	C	144/184 (78%)	124 (86%)	20 (14%)	3	7
1	D	145/184 (79%)	133 (92%)	12 (8%)	10	22
1	E	140/184 (76%)	116 (83%)	24 (17%)	2	4
1	F	146/184 (79%)	123 (84%)	23 (16%)	2	5
1	G	147/184 (80%)	123 (84%)	24 (16%)	2	4

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	144/184 (78%)	123 (85%)	21 (15%)	3	6
1	I	151/184 (82%)	134 (89%)	17 (11%)	5	12
1	J	141/184 (77%)	125 (89%)	16 (11%)	5	12
All	All	1450/1840 (79%)	1246 (86%)	204 (14%)	3	7

All (204) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	29	THR
1	A	32	ILE
1	A	34	LYS
1	A	35	LYS
1	A	38	LEU
1	A	39	LEU
1	A	54	ILE
1	A	55	ILE
1	A	61	LEU
1	A	69	VAL
1	A	71	LEU
1	A	76	VAL
1	A	87	LYS
1	A	92	LYS
1	A	109	SER
1	A	110	ILE
1	A	117	LEU
1	A	131	ASP
1	A	135	ILE
1	A	150	VAL
1	A	154	LEU
1	B	27	ASN
1	B	28	LEU
1	B	32	ILE
1	B	35	LYS
1	B	38	LEU
1	B	39	LEU
1	B	54	ILE
1	B	59	LYS
1	B	61	LEU
1	B	69	VAL
1	B	71	LEU

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Mol	Chain	Res	Type
1	B	90	LEU
1	B	92	LYS
1	B	97	ASN
1	B	106	ILE
1	B	119	ASP
1	B	120	ASP
1	B	139	LEU
1	B	140	LEU
1	B	146	ILE
1	B	148	ARG
1	B	150	VAL
1	B	154	LEU
1	B	169	VAL
1	B	170	CYS
1	C	8	GLU
1	C	15	GLU
1	C	27	ASN
1	C	29	THR
1	C	34	LYS
1	C	38	LEU
1	C	39	LEU
1	C	61	LEU
1	C	67	ARG
1	C	69	VAL
1	C	71	LEU
1	C	76	VAL
1	C	87	LYS
1	C	101	THR
1	C	110	ILE
1	C	117	LEU
1	C	139	LEU
1	C	152	GLU
1	C	154	LEU
1	C	168	ASP
1	D	8	GLU
1	D	34	LYS
1	D	35	LYS
1	D	38	LEU
1	D	39	LEU
1	D	69	VAL
1	D	71	LEU
1	D	101	THR

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Mol	Chain	Res	Type
1	D	117	LEU
1	D	148	ARG
1	D	150	VAL
1	D	154	LEU
1	E	7	LYS
1	E	13	LYS
1	E	28	LEU
1	E	35	LYS
1	E	39	LEU
1	E	47	THR
1	E	59	LYS
1	E	61	LEU
1	E	69	VAL
1	E	71	LEU
1	E	75	SER
1	E	76	VAL
1	E	87	LYS
1	E	90	LEU
1	E	99	LYS
1	E	101	THR
1	E	106	ILE
1	E	117	LEU
1	E	140	LEU
1	E	148	ARG
1	E	150	VAL
1	E	154	LEU
1	E	155	ARG
1	E	162	HIS
1	F	3	THR
1	F	13	LYS
1	F	28	LEU
1	F	29	THR
1	F	30	GLN
1	F	35	LYS
1	F	38	LEU
1	F	39	LEU
1	F	47	THR
1	F	52	SER
1	F	55	ILE
1	F	61	LEU
1	F	67	ARG
1	F	69	VAL

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Mol	Chain	Res	Type
1	F	71	LEU
1	F	101	THR
1	F	117	LEU
1	F	119	ASP
1	F	146	ILE
1	F	148	ARG
1	F	154	LEU
1	F	160	ILE
1	F	170	CYS
1	G	8	GLU
1	G	20	ASP
1	G	27	ASN
1	G	28	LEU
1	G	32	ILE
1	G	35	LYS
1	G	38	LEU
1	G	39	LEU
1	G	47	THR
1	G	54	ILE
1	G	61	LEU
1	G	67	ARG
1	G	69	VAL
1	G	71	LEU
1	G	90	LEU
1	G	97	ASN
1	G	101	THR
1	G	106	ILE
1	G	117	LEU
1	G	146	ILE
1	G	148	ARG
1	G	150	VAL
1	G	154	LEU
1	G	168	ASP
1	H	13	LYS
1	H	25	GLU
1	H	27	ASN
1	H	28	LEU
1	H	29	THR
1	H	34	LYS
1	H	35	LYS
1	H	38	LEU
1	H	39	LEU

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Mol	Chain	Res	Type
1	H	61	LEU
1	H	67	ARG
1	H	69	VAL
1	H	71	LEU
1	H	76	VAL
1	H	97	ASN
1	H	99	LYS
1	H	101	THR
1	H	117	LEU
1	H	146	ILE
1	H	150	VAL
1	H	154	LEU
1	I	8	GLU
1	I	28	LEU
1	I	38	LEU
1	I	39	LEU
1	I	61	LEU
1	I	69	VAL
1	I	71	LEU
1	I	76	VAL
1	I	101	THR
1	I	117	LEU
1	I	120	ASP
1	I	146	ILE
1	I	148	ARG
1	I	153	ILE
1	I	154	LEU
1	I	164	GLU
1	I	175	GLN
1	J	3	THR
1	J	8	GLU
1	J	38	LEU
1	J	39	LEU
1	J	59	LYS
1	J	67	ARG
1	J	69	VAL
1	J	71	LEU
1	J	76	VAL
1	J	97	ASN
1	J	101	THR
1	J	117	LEU
1	J	121	SER

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Mol	Chain	Res	Type
1	J	150	VAL
1	J	154	LEU
1	J	155	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (60) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	HIS
1	A	68	ASN
1	A	115	ASN
1	A	137	GLN
1	A	138	HIS
1	A	143	ASN
1	A	163	HIS
1	B	27	ASN
1	B	65	HIS
1	B	68	ASN
1	B	137	GLN
1	B	138	HIS
1	B	143	ASN
1	C	27	ASN
1	C	30	GLN
1	C	137	GLN
1	C	138	HIS
1	C	143	ASN
1	C	162	HIS
1	D	30	GLN
1	D	115	ASN
1	D	138	HIS
1	D	143	ASN
1	D	161	GLN
1	D	163	HIS
1	E	27	ASN
1	E	65	HIS
1	E	68	ASN
1	E	143	ASN
1	F	27	ASN
1	F	30	GLN
1	F	68	ASN
1	F	97	ASN
1	F	138	HIS
1	F	161	GLN

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Mol	Chain	Res	Type
1	F	163	HIS
1	G	30	GLN
1	G	65	HIS
1	G	68	ASN
1	G	133	ASN
1	G	138	HIS
1	G	143	ASN
1	H	27	ASN
1	H	30	GLN
1	H	65	HIS
1	H	68	ASN
1	H	137	GLN
1	H	138	HIS
1	I	21	ASN
1	I	27	ASN
1	I	68	ASN
1	I	137	GLN
1	I	138	HIS
1	I	143	ASN
1	I	163	HIS
1	I	175	GLN
1	J	68	ASN
1	J	138	HIS
1	J	143	ASN
1	J	162	HIS

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	167/213 (78%)	0.09	0 100 100	34, 49, 67, 84	0
1	B	171/213 (80%)	0.09	7 (4%) 41 37	36, 47, 67, 85	0
1	C	170/213 (79%)	-0.10	3 (1%) 67 64	30, 41, 59, 73	0
1	D	169/213 (79%)	-0.01	4 (2%) 59 55	32, 48, 61, 70	0
1	E	161/213 (75%)	-0.09	3 (1%) 66 62	33, 45, 61, 76	0
1	F	168/213 (78%)	0.06	6 (3%) 46 41	33, 46, 65, 86	0
1	G	171/213 (80%)	0.00	1 (0%) 85 83	33, 47, 62, 67	0
1	H	167/213 (78%)	-0.14	2 (1%) 76 73	26, 41, 61, 84	0
1	I	177/213 (83%)	-0.11	4 (2%) 61 57	30, 43, 58, 65	0
1	J	163/213 (76%)	-0.31	3 (1%) 67 64	27, 37, 51, 64	0
All	All	1684/2130 (79%)	-0.05	33 (1%) 65 61	26, 44, 62, 86	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	172	ALA	6.4
1	D	170	CYS	5.2
1	F	145	ALA	3.9
1	F	170	CYS	3.8
1	I	177	GLY	3.8
1	E	2	PRO	3.8
1	C	170	CYS	3.7
1	B	171	PRO	3.7
1	F	3	THR	3.7
1	B	170	CYS	3.4
1	F	169	VAL	3.3
1	C	1	SER	3.3
1	H	168	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	G	152	GLU	2.9
1	B	169	VAL	2.9
1	E	152	GLU	2.9
1	I	1	SER	2.6
1	I	152	GLU	2.6
1	J	164	GLU	2.5
1	D	152	GLU	2.5
1	F	152	GLU	2.5
1	E	162	HIS	2.4
1	H	152	GLU	2.3
1	D	169	VAL	2.3
1	J	93	GLY	2.3
1	B	152	GLU	2.2
1	I	146	ILE	2.2
1	B	163	HIS	2.2
1	J	91	ALA	2.2
1	F	163	HIS	2.1
1	B	32	ILE	2.1
1	D	2	PRO	2.1
1	C	168	ASP	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.