



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:08 PM UTC

PDB ID : 4G80 / pdb_00004g80
Title : Crystal structure of voltage sensing domain of Ci-VSP with fragment antibody (WT, 3.8 Å)
Authors : Li, Q.
Deposited on : 2012-07-20
Resolution : 3.58 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

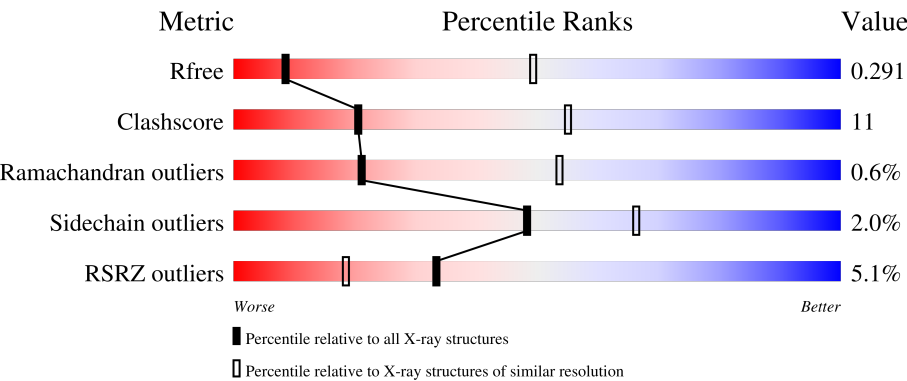
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.58 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	1521 (3.66-3.50)
Clashscore	190562	1595 (3.66-3.50)
Ramachandran outliers	187476	1551 (3.66-3.50)
Sidechain outliers	187428	1551 (3.66-3.50)
RSRZ outliers	180081	1520 (3.66-3.50)

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	I	155	9% 68% 21% 10%
1	J	155	6% 70% 19% 10%
1	S	155	7% 64% 19% 17%
1	T	155	6% 69% 19% 10%
2	A	224	5% 69% 28% ..

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Mol	Chain	Length	Quality of chain
2	C	224	2%71%26%.
2	E	224	4%73%25%..
2	G	224	9%71%27%.
3	B	211	%77%22%
3	D	211	4%79%20%
3	F	211	5%76%23%
3	H	211	3%80%20%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17446 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 Voltage-sensor containing phosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	S	128	Total	C	N	O	S	0	0	0
			1020	677	165	173	5			
1	T	139	Total	C	N	O	S	0	0	0
			1111	734	181	190	6			
1	I	139	Total	C	N	O	S	0	0	0
			1111	734	181	190	6			
1	J	139	Total	C	N	O	S	0	0	0
			1111	734	181	190	6			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
S	90	MET	-	expression tag	UNP Q4W8A1
S	91	ARG	-	expression tag	UNP Q4W8A1
S	92	GLY	-	expression tag	UNP Q4W8A1
S	93	SER	-	expression tag	UNP Q4W8A1
S	94	HIS	-	expression tag	UNP Q4W8A1
S	95	HIS	-	expression tag	UNP Q4W8A1
S	96	HIS	-	expression tag	UNP Q4W8A1
S	97	HIS	-	expression tag	UNP Q4W8A1
S	98	HIS	-	expression tag	UNP Q4W8A1
S	99	HIS	-	expression tag	UNP Q4W8A1
S	100	GLY	-	expression tag	UNP Q4W8A1
S	101	GLU	-	expression tag	UNP Q4W8A1
S	102	ASN	-	expression tag	UNP Q4W8A1
S	103	LEU	-	expression tag	UNP Q4W8A1
S	104	TYR	-	expression tag	UNP Q4W8A1
S	105	PHE	-	expression tag	UNP Q4W8A1
T	90	MET	-	expression tag	UNP Q4W8A1
T	91	ARG	-	expression tag	UNP Q4W8A1
T	92	GLY	-	expression tag	UNP Q4W8A1
T	93	SER	-	expression tag	UNP Q4W8A1
T	94	HIS	-	expression tag	UNP Q4W8A1

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Chain	Residue	Modelled	Actual	Comment	Reference
T	95	HIS	-	expression tag	UNP Q4W8A1
T	96	HIS	-	expression tag	UNP Q4W8A1
T	97	HIS	-	expression tag	UNP Q4W8A1
T	98	HIS	-	expression tag	UNP Q4W8A1
T	99	HIS	-	expression tag	UNP Q4W8A1
T	100	GLY	-	expression tag	UNP Q4W8A1
T	101	GLU	-	expression tag	UNP Q4W8A1
T	102	ASN	-	expression tag	UNP Q4W8A1
T	103	LEU	-	expression tag	UNP Q4W8A1
T	104	TYR	-	expression tag	UNP Q4W8A1
T	105	PHE	-	expression tag	UNP Q4W8A1
I	90	MET	-	expression tag	UNP Q4W8A1
I	91	ARG	-	expression tag	UNP Q4W8A1
I	92	GLY	-	expression tag	UNP Q4W8A1
I	93	SER	-	expression tag	UNP Q4W8A1
I	94	HIS	-	expression tag	UNP Q4W8A1
I	95	HIS	-	expression tag	UNP Q4W8A1
I	96	HIS	-	expression tag	UNP Q4W8A1
I	97	HIS	-	expression tag	UNP Q4W8A1
I	98	HIS	-	expression tag	UNP Q4W8A1
I	99	HIS	-	expression tag	UNP Q4W8A1
I	100	GLY	-	expression tag	UNP Q4W8A1
I	101	GLU	-	expression tag	UNP Q4W8A1
I	102	ASN	-	expression tag	UNP Q4W8A1
I	103	LEU	-	expression tag	UNP Q4W8A1
I	104	TYR	-	expression tag	UNP Q4W8A1
I	105	PHE	-	expression tag	UNP Q4W8A1
J	90	MET	-	expression tag	UNP Q4W8A1
J	91	ARG	-	expression tag	UNP Q4W8A1
J	92	GLY	-	expression tag	UNP Q4W8A1
J	93	SER	-	expression tag	UNP Q4W8A1
J	94	HIS	-	expression tag	UNP Q4W8A1
J	95	HIS	-	expression tag	UNP Q4W8A1
J	96	HIS	-	expression tag	UNP Q4W8A1
J	97	HIS	-	expression tag	UNP Q4W8A1
J	98	HIS	-	expression tag	UNP Q4W8A1
J	99	HIS	-	expression tag	UNP Q4W8A1
J	100	GLY	-	expression tag	UNP Q4W8A1
J	101	GLU	-	expression tag	UNP Q4W8A1
J	102	ASN	-	expression tag	UNP Q4W8A1
J	103	LEU	-	expression tag	UNP Q4W8A1
J	104	TYR	-	expression tag	UNP Q4W8A1

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Chain	Residue	Modelled	Actual	Comment	Reference
J	105	PHE	-	expression tag	UNP Q4W8A1

- Molecule 2 is a protein called fragment antibody heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	221	Total	C	N	O	S	0	0	0
			1661	1052	276	327	6			
2	C	219	Total	C	N	O	S	0	0	0
			1646	1044	274	322	6			
2	E	221	Total	C	N	O	S	0	0	0
			1661	1052	276	327	6			
2	G	221	Total	C	N	O	S	0	0	0
			1661	1052	276	327	6			

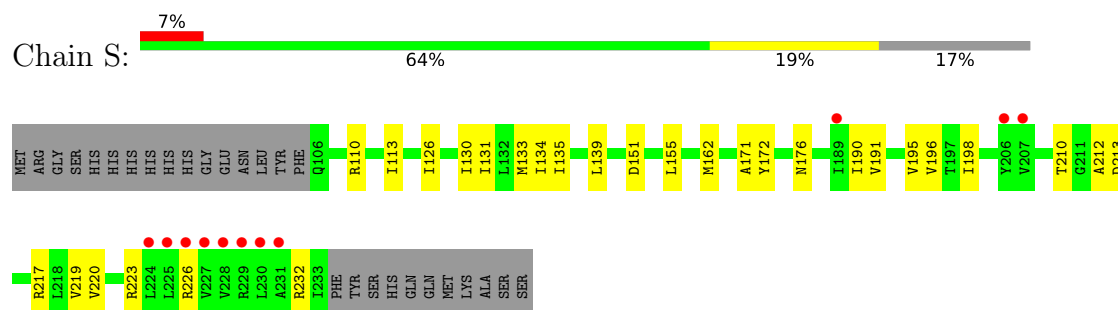
- Molecule 3 is a protein called fragment antibody light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	B	211	Total	C	N	O	S	0	0	0
			1616	1011	272	328	5			
3	D	211	Total	C	N	O	S	0	0	0
			1616	1011	272	328	5			
3	F	211	Total	C	N	O	S	0	0	0
			1616	1011	272	328	5			
3	H	211	Total	C	N	O	S	0	0	0
			1616	1011	272	328	5			

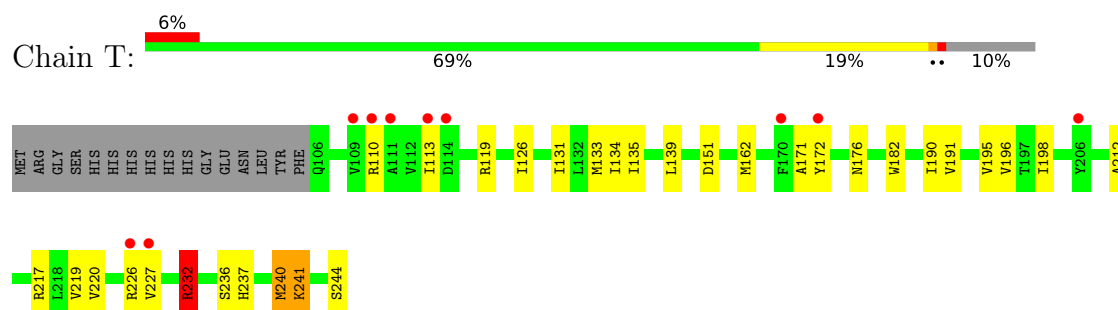
3 Residue-property plots [i](#)

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.

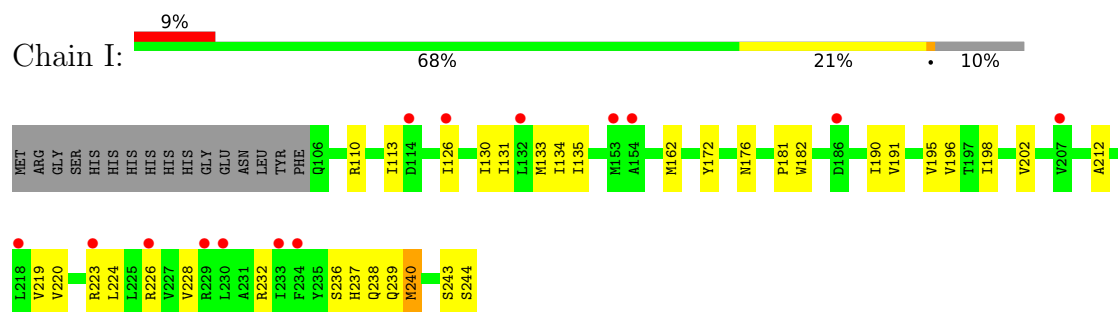
- Molecule 1: Voltage-sensor containing phosphatase



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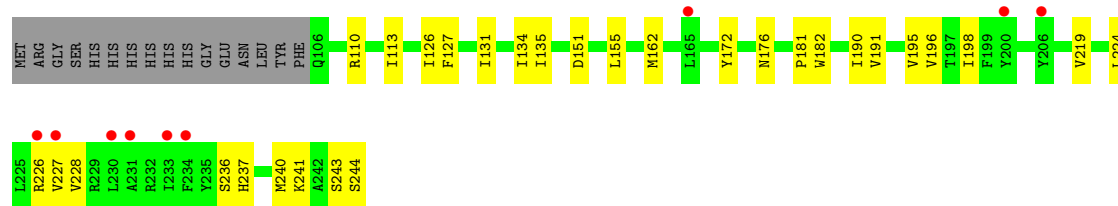


- Molecule 1: Voltage-sensor containing phosphatase

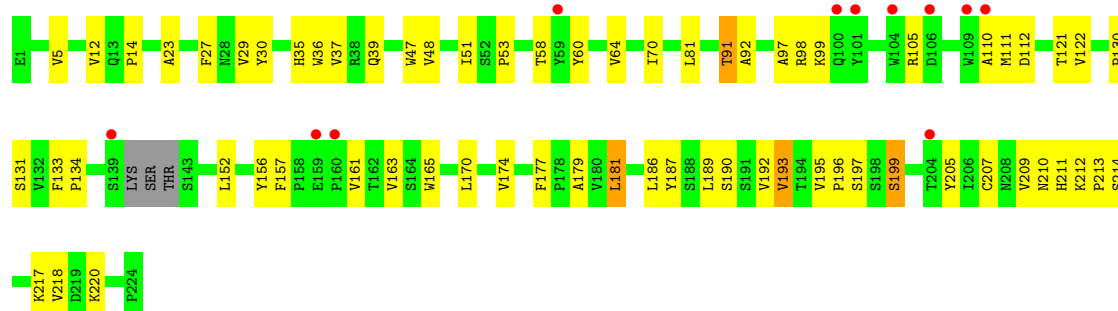


- Molecule 1: Voltage-sensor containing phosphatase

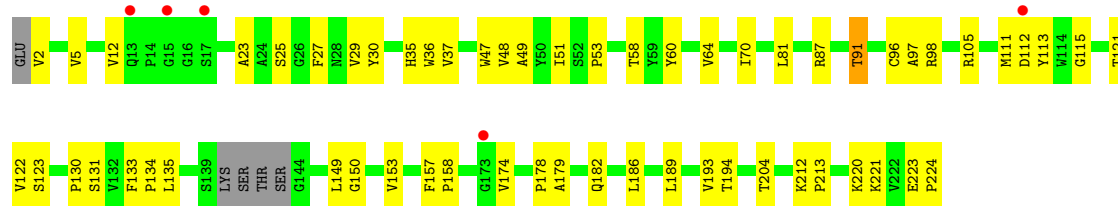




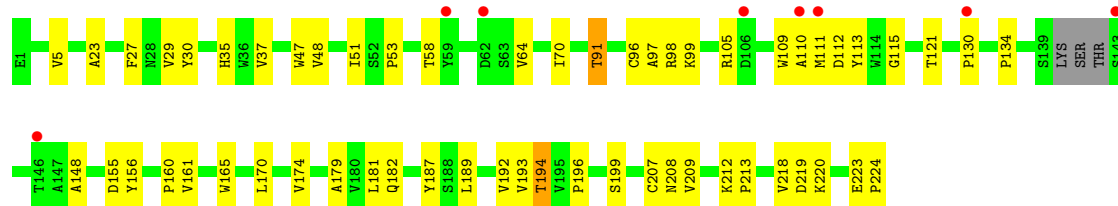
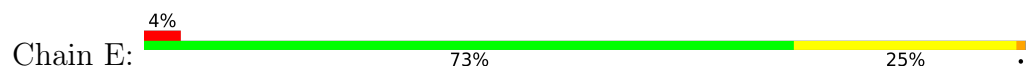
• Molecule 2: fragment antibody heavy chain



• Molecule 2: fragment antibody heavy chain

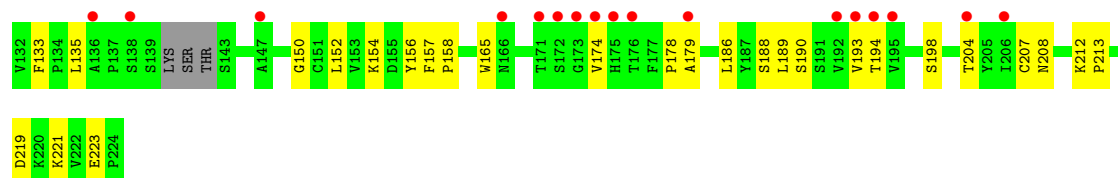


• Molecule 2: fragment antibody heavy chain

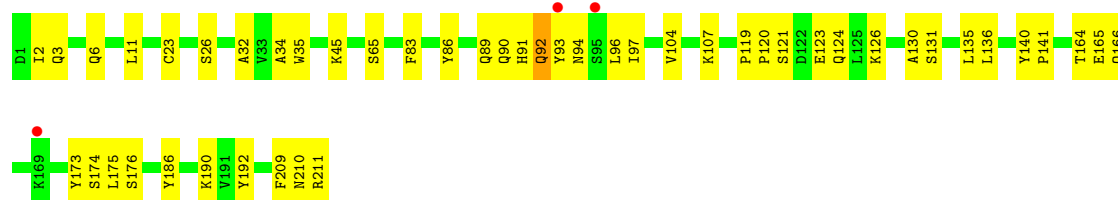
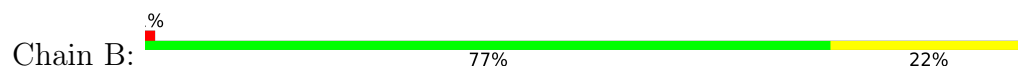


• Molecule 2: fragment antibody heavy chain

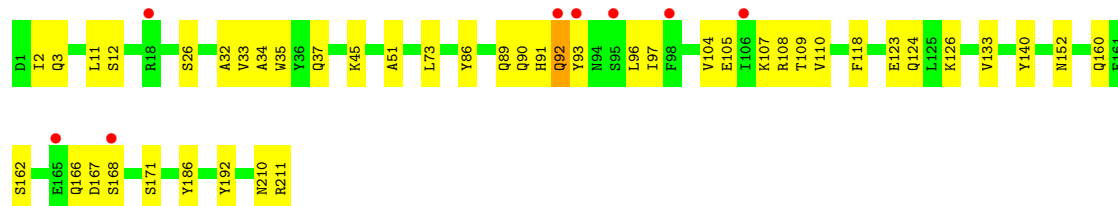
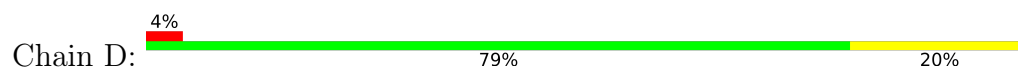




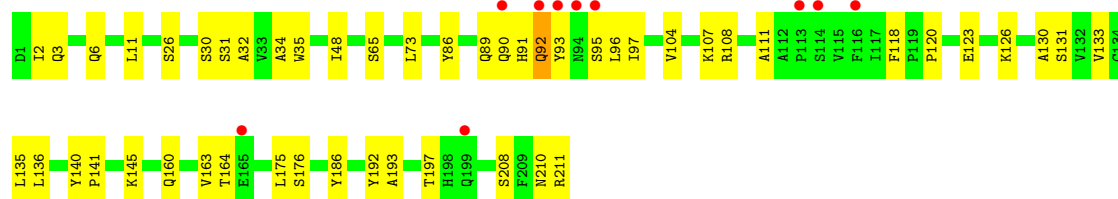
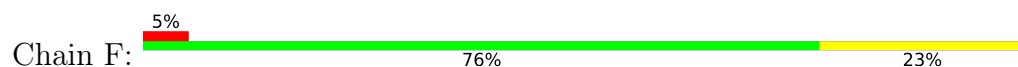
- Molecule 3: fragment antibody light chain



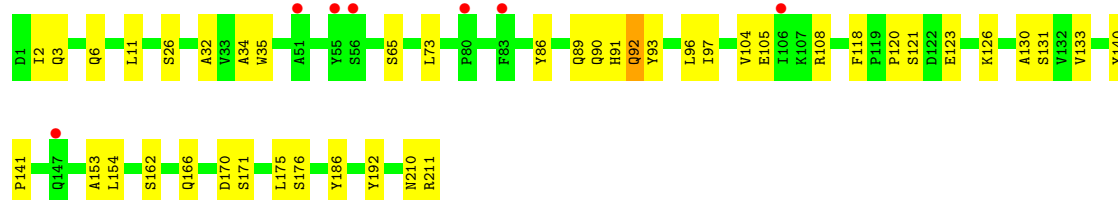
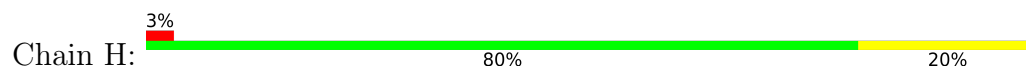
- Molecule 3: fragment antibody light chain



- Molecule 3: fragment antibody light chain



- Molecule 3: fragment antibody light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	77.29Å 94.24Å 193.95Å 102.63° 93.45° 105.25°	Depositor
Resolution (Å)	50.00 – 3.58 50.00 – 3.58	Depositor EDS
% Data completeness (in resolution range)	92.6 (50.00-3.58) 92.6 (50.00-3.58)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.52 (at 3.57Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.248 , 0.292 0.245 , 0.291	Depositor DCC
R_{free} test set	2803 reflections (4.66%)	wwPDB-VP
Wilson B-factor (Å ²)	130.0	Xtriage
Anisotropy	0.142	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 75.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	17446	wwPDB-VP
Average B, all atoms (Å ²)	149.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% 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	I	0.42	0/1135	0.66	0/1536
1	J	0.43	0/1135	0.67	0/1536
1	S	0.45	1/1041 (0.1%)	0.68	0/1411
1	T	1.01	3/1135 (0.3%)	0.72	1/1536 (0.1%)
2	A	0.50	0/1705	0.67	0/2326
2	C	0.49	0/1690	0.68	0/2306
2	E	0.50	0/1705	0.68	0/2326
2	G	0.49	0/1705	0.65	0/2326
3	B	0.44	0/1650	0.65	0/2240
3	D	0.43	0/1650	0.64	0/2240
3	F	0.44	0/1650	0.66	0/2240
3	H	0.44	0/1650	0.66	0/2240
All	All	0.51	4/17851 (0.0%)	0.67	1/24263 (0.0%)

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	T	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	T	232	ARG	NE-CZ	21.03	1.56	1.33
1	T	232	ARG	CZ-NH2	18.24	1.57	1.33
1	T	232	ARG	CD-NE	13.88	1.65	1.46
1	S	232	ARG	CZ-NH2	7.24	1.42	1.33

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	T	232	ARG	NE-CZ-NH2	9.38	127.64	119.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	T	232	ARG	Sidechain

5.2 Too-close contacts [i](#)

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	I	1111	0	1140	26	0
1	J	1111	0	1140	17	0
1	S	1020	0	1057	17	0
1	T	1111	0	1140	38	0
2	A	1661	0	1598	59	0
2	C	1646	0	1584	42	0
2	E	1661	0	1598	38	0
2	G	1661	0	1598	42	0
3	B	1616	0	1576	41	0
3	D	1616	0	1576	37	0
3	F	1616	0	1576	37	0
3	H	1616	0	1576	33	0
All	All	17446	0	17159	388	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 (388) 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:T:182:TRP:CH2	1:T:240:MET:HA	1.82	1.12
1:T:182:TRP:CZ3	1:T:240:MET:HA	1.89	1.07
1:I:182:TRP:HZ3	1:I:243:SER:HG	1.01	0.99
1:J:237:HIS:O	1:J:240:MET:HG3	1.65	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:182:TRP:CE3	1:T:240:MET:HB2	2.11	0.85
2:E:182:GLN:HA	3:F:160:GLN:HE22	1.42	0.83
1:T:232:ARG:NH1	1:T:236:SER:HB2	1.96	0.81
3:D:108:ARG:HG2	3:D:109:THR:H	1.46	0.80
2:E:170:LEU:HD21	2:E:193:VAL:HG21	1.63	0.79
1:T:182:TRP:CZ3	1:T:240:MET:CA	2.65	0.79
3:D:210:ASN:O	3:D:211:ARG:HB2	1.83	0.78
3:D:108:ARG:CG	3:D:109:THR:H	2.00	0.75
1:T:182:TRP:CE3	1:T:240:MET:CB	2.72	0.73
3:B:175:LEU:HD23	3:B:176:SER:N	2.05	0.72
1:T:232:ARG:NH1	1:T:236:SER:CB	2.52	0.71
1:J:182:TRP:CE3	1:J:240:MET:HB2	2.26	0.70
3:B:107:LYS:HA	3:B:140:TYR:OH	1.92	0.69
2:A:181:LEU:HD12	2:A:187:TYR:CZ	2.27	0.69
3:B:210:ASN:O	3:B:211:ARG:HB2	1.93	0.69
2:E:97:ALA:HB3	2:E:111:MET:HE2	1.75	0.68
1:I:237:HIS:HA	1:I:240:MET:CG	2.23	0.68
3:F:210:ASN:O	3:F:211:ARG:HB2	1.92	0.68
3:B:96:LEU:HD23	3:B:96:LEU:H	1.59	0.68
2:G:97:ALA:HB3	2:G:111:MET:HE2	1.76	0.68
2:A:97:ALA:HB3	2:A:111:MET:HE2	1.76	0.67
3:D:89:GLN:HG2	3:D:90:GLN:H	1.59	0.67
3:D:12:SER:OG	3:D:107:LYS:HB2	1.95	0.67
3:F:96:LEU:HD23	3:F:96:LEU:H	1.61	0.66
1:T:232:ARG:HG3	1:T:232:ARG:HH11	1.61	0.65
3:H:96:LEU:H	3:H:96:LEU:HD23	1.62	0.65
1:I:182:TRP:CZ3	1:I:240:MET:HA	2.31	0.64
2:A:134:PRO:HD3	2:A:220:LYS:HE2	1.78	0.64
1:T:237:HIS:O	1:T:240:MET:HG3	1.98	0.63
3:D:89:GLN:NE2	3:D:96:LEU:HD12	2.13	0.63
1:S:162:MET:HE1	1:S:191:VAL:HG22	1.79	0.63
2:A:110:ALA:HB2	3:B:91:HIS:CE1	2.33	0.63
2:A:152:LEU:HD12	2:A:190:SER:HB3	1.81	0.62
3:B:89:GLN:HG2	3:B:90:GLN:H	1.63	0.62
3:D:96:LEU:HD23	3:D:96:LEU:H	1.64	0.62
3:D:108:ARG:HG2	3:D:109:THR:N	2.14	0.62
1:T:237:HIS:O	1:T:240:MET:CG	2.47	0.62
1:T:237:HIS:HA	1:T:240:MET:HG2	1.82	0.61
3:F:89:GLN:HG2	3:F:90:GLN:H	1.64	0.61
2:C:97:ALA:HB3	2:C:111:MET:HE2	1.80	0.61
2:C:178:PRO:HG2	3:D:162:SER:OG	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:11:LEU:HD11	3:B:104:VAL:HG22	1.83	0.60
2:A:47:TRP:CD1	3:B:96:LEU:HD11	2.36	0.60
2:A:27:PHE:CZ	2:A:98:ARG:HD3	2.36	0.60
1:T:162:MET:HE1	1:T:191:VAL:HG22	1.84	0.60
1:T:182:TRP:CH2	1:T:240:MET:CA	2.74	0.60
2:A:98:ARG:NH2	2:A:112:ASP:OD1	2.35	0.59
2:A:170:LEU:HD21	2:A:193:VAL:HG21	1.84	0.59
2:A:218:VAL:HG22	3:H:154:LEU:HB3	1.82	0.59
1:I:172:TYR:O	1:I:176:ASN:ND2	2.36	0.59
1:I:236:SER:O	1:I:240:MET:HG2	2.02	0.59
2:G:174:VAL:HG22	2:G:193:VAL:HB	1.84	0.58
1:I:237:HIS:O	1:I:240:MET:HG3	2.03	0.58
1:S:172:TYR:O	1:S:176:ASN:ND2	2.36	0.58
2:A:211:HIS:ND1	2:A:214:SER:OG	2.16	0.58
2:C:204:THR:HG23	2:C:221:LYS:HE3	1.86	0.58
3:F:107:LYS:HA	3:F:140:TYR:OH	2.04	0.58
2:A:212:LYS:N	2:A:213:PRO:CD	2.67	0.58
3:H:32:ALA:HB1	3:H:91:HIS:HB2	1.86	0.58
1:T:172:TYR:O	1:T:176:ASN:ND2	2.37	0.57
2:A:157:PHE:HB2	2:A:186:LEU:HD22	1.85	0.57
3:F:34:ALA:HB2	3:F:91:HIS:HE1	1.68	0.57
1:J:162:MET:HE1	1:J:191:VAL:HG22	1.84	0.57
2:A:91:THR:HG23	2:A:121:THR:HA	1.86	0.57
3:H:210:ASN:O	3:H:211:ARG:HB2	2.03	0.57
3:H:89:GLN:HG2	3:H:90:GLN:H	1.69	0.57
1:J:236:SER:O	1:J:240:MET:HB3	2.04	0.57
2:C:134:PRO:HD3	2:C:220:LYS:HE2	1.87	0.56
1:T:195:VAL:O	1:T:198:ILE:HG22	2.06	0.56
3:D:108:ARG:HD3	3:D:109:THR:O	2.05	0.56
1:T:182:TRP:CD2	1:T:240:MET:HB3	2.41	0.56
2:G:221:LYS:HE2	2:G:223:GLU:OE1	2.05	0.56
1:J:172:TYR:O	1:J:176:ASN:ND2	2.38	0.56
3:B:6:GLN:NE2	3:B:86:TYR:O	2.34	0.56
2:C:212:LYS:N	2:C:213:PRO:CD	2.68	0.56
2:C:182:GLN:HA	3:D:160:GLN:HE22	1.69	0.56
1:I:162:MET:HE1	1:I:191:VAL:HG22	1.88	0.56
3:B:34:ALA:HB2	3:B:91:HIS:HE1	1.70	0.55
1:T:232:ARG:CZ	1:T:236:SER:HB2	2.36	0.55
1:I:237:HIS:HA	1:I:240:MET:SD	2.46	0.55
1:T:240:MET:HG3	1:T:241:LYS:H	1.72	0.55
3:D:108:ARG:CG	3:D:109:THR:N	2.70	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:27:PHE:CZ	2:E:98:ARG:HD3	2.41	0.55
2:G:179:ALA:HA	2:G:189:LEU:HB3	1.88	0.55
2:G:179:ALA:HB2	2:G:189:LEU:HD23	1.88	0.55
2:G:157:PHE:HB2	2:G:186:LEU:HD22	1.88	0.55
2:E:110:ALA:HB2	3:F:91:HIS:CE1	2.42	0.55
3:D:32:ALA:HB1	3:D:91:HIS:HB2	1.88	0.55
1:S:217:ARG:HH11	1:T:139:LEU:HD21	1.72	0.54
2:E:165:TRP:CH2	2:E:207:CYS:HB3	2.43	0.54
3:F:123:GLU:HA	3:F:126:LYS:HD3	1.90	0.54
2:A:217:LYS:HE3	3:H:153:ALA:HB2	1.88	0.54
2:C:37:VAL:HG22	2:C:47:TRP:HA	1.89	0.54
2:C:91:THR:HG23	2:C:121:THR:HA	1.89	0.54
3:F:11:LEU:HD11	3:F:104:VAL:HG22	1.88	0.54
2:E:29:VAL:HG12	2:E:53:PRO:HG3	1.90	0.54
2:A:196:PRO:O	2:A:199:SER:OG	2.23	0.54
2:A:133:PHE:CE1	3:B:124:GLN:HG3	2.43	0.54
3:H:34:ALA:HB2	3:H:91:HIS:HE1	1.72	0.54
1:I:110:ARG:HA	1:I:113:ILE:HG22	1.90	0.54
2:A:212:LYS:N	2:A:213:PRO:HD2	2.23	0.53
3:H:11:LEU:HD11	3:H:104:VAL:HG22	1.89	0.53
1:I:182:TRP:CD2	1:I:240:MET:HB3	2.42	0.53
1:S:110:ARG:HA	1:S:113:ILE:HG22	1.89	0.53
2:G:98:ARG:HH21	2:G:113:TYR:HD2	1.56	0.53
2:A:60:TYR:CE1	2:A:70:ILE:HG22	2.43	0.53
2:C:60:TYR:CE1	2:C:70:ILE:HG22	2.43	0.53
3:H:105:GLU:OE1	3:H:166:GLN:NE2	2.41	0.53
2:A:35:HIS:CE1	3:B:93:TYR:HD2	2.27	0.53
2:A:163:VAL:HG22	2:A:209:VAL:HG22	1.90	0.53
3:D:89:GLN:HE21	3:D:96:LEU:HD12	1.74	0.53
3:F:6:GLN:NE2	3:F:86:TYR:O	2.34	0.53
2:E:98:ARG:HH21	2:E:113:TYR:HD2	1.57	0.53
3:H:108:ARG:HD2	3:H:170:ASP:O	2.08	0.53
1:T:110:ARG:HA	1:T:113:ILE:HG22	1.90	0.53
2:E:91:THR:HG23	2:E:121:THR:HA	1.90	0.53
3:F:34:ALA:HB2	3:F:91:HIS:CE1	2.43	0.53
2:C:123:SER:HB3	2:C:157:PHE:CZ	2.43	0.53
1:T:240:MET:HG3	1:T:241:LYS:N	2.23	0.53
2:A:30:TYR:HA	2:A:53:PRO:HG2	1.91	0.53
2:E:181:LEU:HD12	2:E:187:TYR:CZ	2.44	0.52
1:J:110:ARG:HA	1:J:113:ILE:HG22	1.90	0.52
2:C:179:ALA:HB2	2:C:189:LEU:HD23	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:212:LYS:N	2:C:213:PRO:HD2	2.24	0.52
1:T:182:TRP:CZ3	1:T:240:MET:CB	2.91	0.52
2:G:37:VAL:HG22	2:G:47:TRP:HA	1.91	0.52
2:G:135:LEU:HB2	2:G:150:GLY:O	2.10	0.52
1:T:133:MET:SD	1:T:220:VAL:HG13	2.49	0.52
3:H:175:LEU:HD23	3:H:176:SER:N	2.24	0.52
3:D:105:GLU:OE1	3:D:140:TYR:HE2	1.92	0.52
2:G:204:THR:HG23	2:G:221:LYS:HE3	1.91	0.52
1:J:181:PRO:HG3	1:J:244:SER:HB2	1.92	0.52
2:C:27:PHE:CZ	2:C:98:ARG:HD3	2.45	0.52
3:D:34:ALA:HB2	3:D:91:HIS:HE1	1.75	0.52
3:D:118:PHE:HB2	3:D:133:VAL:HB	1.91	0.52
2:G:29:VAL:HG12	2:G:53:PRO:HG3	1.92	0.52
2:G:36:TRP:CG	2:G:81:LEU:HD22	2.45	0.52
2:G:91:THR:HG23	2:G:121:THR:HA	1.90	0.52
1:I:237:HIS:C	1:I:240:MET:HG3	2.35	0.52
3:B:186:TYR:HA	3:B:192:TYR:OH	2.10	0.51
2:G:165:TRP:CH2	2:G:207:CYS:HB3	2.45	0.51
1:J:182:TRP:CZ3	1:J:240:MET:HB2	2.45	0.51
2:G:36:TRP:HB2	2:G:49:ALA:HB3	1.92	0.51
2:C:30:TYR:HA	2:C:53:PRO:HG2	1.91	0.51
2:C:29:VAL:HG12	2:C:53:PRO:HG3	1.93	0.51
2:E:47:TRP:CD1	3:F:96:LEU:HD11	2.46	0.51
2:E:5:VAL:HG23	2:E:23:ALA:HB3	1.93	0.51
1:S:195:VAL:O	1:S:198:ILE:HG22	2.11	0.51
2:E:182:GLN:HA	3:F:160:GLN:NE2	2.20	0.51
3:D:11:LEU:HD11	3:D:104:VAL:HG22	1.92	0.50
2:E:208:ASN:HD22	2:E:219:ASP:CG	2.19	0.50
1:J:195:VAL:O	1:J:198:ILE:HG22	2.11	0.50
2:C:36:TRP:HD1	2:C:70:ILE:HD12	1.77	0.50
1:J:196:VAL:HB	1:J:219:VAL:HG21	1.94	0.50
2:A:177:PHE:CD1	3:B:164:THR:HG23	2.46	0.50
3:B:123:GLU:HA	3:B:126:LYS:HD3	1.94	0.50
3:H:34:ALA:HB2	3:H:91:HIS:CE1	2.47	0.50
2:E:196:PRO:HG2	2:E:199:SER:OG	2.12	0.50
3:F:186:TYR:HA	3:F:192:TYR:OH	2.12	0.50
1:J:131:ILE:O	1:J:135:ILE:HG13	2.12	0.50
3:F:32:ALA:HB1	3:F:91:HIS:HB2	1.94	0.50
2:C:133:PHE:CE1	3:D:124:GLN:HG3	2.47	0.50
3:D:34:ALA:HB2	3:D:91:HIS:CE1	2.47	0.50
2:E:212:LYS:N	2:E:213:PRO:CD	2.73	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:151:ASP:OD2	2:A:105:ARG:HD2	2.11	0.49
2:E:30:TYR:HA	2:E:53:PRO:HG2	1.94	0.49
2:E:165:TRP:CZ3	2:E:207:CYS:HB3	2.47	0.49
2:G:30:TYR:HA	2:G:53:PRO:HG2	1.94	0.49
2:C:48:VAL:HG13	2:C:64:VAL:HG21	1.95	0.49
3:H:91:HIS:O	3:H:92:GLN:HB3	2.13	0.49
2:A:156:TYR:CE2	2:A:161:VAL:HG23	2.47	0.49
2:A:210:ASN:HD21	2:A:212:LYS:HE2	1.77	0.49
3:B:34:ALA:HB2	3:B:91:HIS:CE1	2.48	0.49
1:T:182:TRP:CD2	1:T:240:MET:CB	2.94	0.49
2:G:135:LEU:N	2:G:150:GLY:O	2.46	0.49
1:I:196:VAL:HB	1:I:219:VAL:HG21	1.93	0.49
2:A:189:LEU:C	2:A:189:LEU:HD12	2.38	0.49
2:C:36:TRP:HB2	2:C:49:ALA:HB3	1.95	0.48
1:I:237:HIS:ND1	1:I:240:MET:SD	2.86	0.48
2:C:179:ALA:HA	2:C:189:LEU:HB3	1.95	0.48
1:T:196:VAL:HB	1:T:219:VAL:HG21	1.95	0.48
2:G:178:PRO:HG2	3:H:162:SER:OG	2.14	0.48
3:B:83:PHE:HE1	3:B:165:GLU:HB3	1.79	0.48
2:G:212:LYS:HB2	2:G:213:PRO:HD3	1.96	0.48
3:H:6:GLN:NE2	3:H:86:TYR:O	2.37	0.48
2:A:165:TRP:CH2	2:A:207:CYS:HB3	2.49	0.48
3:D:186:TYR:HA	3:D:192:TYR:OH	2.13	0.48
2:A:177:PHE:HE1	3:B:174:SER:O	1.96	0.48
3:D:35:TRP:CD2	3:D:73:LEU:HB2	2.49	0.48
2:G:27:PHE:CZ	2:G:98:ARG:HD3	2.49	0.48
1:J:190:ILE:HA	1:J:226:ARG:NH2	2.29	0.48
2:C:87:ARG:O	2:C:122:VAL:HG11	2.14	0.48
2:E:196:PRO:O	2:E:199:SER:OG	2.32	0.48
1:T:110:ARG:HD3	1:T:171:ALA:O	2.14	0.48
3:B:3:GLN:HB2	3:B:26:SER:HB3	1.95	0.47
2:E:98:ARG:NH2	2:E:112:ASP:OD1	2.47	0.47
2:A:211:HIS:CE1	2:A:214:SER:HG	2.30	0.47
2:G:35:HIS:CE1	3:H:93:TYR:HD2	2.32	0.47
3:H:89:GLN:NE2	3:H:96:LEU:HD12	2.29	0.47
3:B:96:LEU:H	3:B:96:LEU:CD2	2.26	0.47
2:E:99:LYS:HZ1	3:F:91:HIS:HB3	1.79	0.47
1:I:182:TRP:CH2	1:I:240:MET:HA	2.49	0.47
2:A:179:ALA:HB2	2:A:189:LEU:HD23	1.97	0.47
3:B:32:ALA:HB1	3:B:91:HIS:HB2	1.95	0.47
2:A:181:LEU:HD12	2:A:187:TYR:CE1	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:47:TRP:CD1	3:D:96:LEU:HD11	2.50	0.47
3:F:89:GLN:NE2	3:F:96:LEU:HD12	2.28	0.47
2:G:130:PRO:HB3	2:G:156:TYR:HB3	1.97	0.47
3:F:91:HIS:O	3:F:92:GLN:HB3	2.15	0.47
2:A:14:PRO:HG2	3:D:152:ASN:HD21	1.79	0.47
1:T:190:ILE:HA	1:T:226:ARG:NH2	2.30	0.47
1:T:232:ARG:HH12	1:T:236:SER:HB2	1.78	0.47
3:F:175:LEU:HD23	3:F:176:SER:N	2.29	0.47
3:H:3:GLN:HB2	3:H:26:SER:HB3	1.97	0.47
1:I:195:VAL:O	1:I:198:ILE:HG22	2.14	0.47
1:S:131:ILE:O	1:S:135:ILE:HG13	2.14	0.47
1:S:196:VAL:HB	1:S:219:VAL:HG21	1.97	0.47
2:A:192:VAL:HG11	3:B:135:LEU:HD22	1.96	0.47
2:C:51:ILE:HG13	2:C:58:THR:HG22	1.95	0.47
2:E:134:PRO:HD3	2:E:220:LYS:HE2	1.95	0.47
2:A:36:TRP:CG	2:A:81:LEU:HD22	2.50	0.47
3:D:123:GLU:HA	3:D:126:LYS:HD3	1.96	0.47
3:F:136:LEU:HB2	3:F:175:LEU:HB3	1.96	0.47
2:A:133:PHE:HB3	3:B:121:SER:OG	2.15	0.46
3:B:91:HIS:O	3:B:92:GLN:HB3	2.15	0.46
3:D:89:GLN:HG2	3:D:90:GLN:N	2.28	0.46
2:E:35:HIS:CE1	3:F:93:TYR:HD2	2.33	0.46
2:E:179:ALA:HB2	2:E:189:LEU:HD23	1.97	0.46
3:B:166:GLN:HB2	3:B:173:TYR:CE2	2.50	0.46
1:T:151:ASP:OD2	2:E:105:ARG:HD2	2.14	0.46
3:H:123:GLU:HA	3:H:126:LYS:HD3	1.98	0.46
2:A:195:VAL:HG21	2:A:205:TYR:CE2	2.51	0.46
3:H:35:TRP:CD2	3:H:73:LEU:HB2	2.51	0.46
2:A:130:PRO:HB3	2:A:156:TYR:HB3	1.96	0.46
2:A:161:VAL:HG11	2:A:209:VAL:HG13	1.98	0.46
2:C:98:ARG:HH21	2:C:113:TYR:HD2	1.64	0.46
2:C:133:PHE:CG	3:D:124:GLN:HB2	2.51	0.46
3:D:37:GLN:HB2	3:D:86:TYR:CE1	2.51	0.46
2:E:192:VAL:HG11	3:F:135:LEU:HD22	1.97	0.46
1:S:155:LEU:HD11	1:S:198:ILE:HD12	1.98	0.46
2:A:99:LYS:HD3	3:B:93:TYR:CE2	2.50	0.46
3:B:140:TYR:CG	3:B:141:PRO:HA	2.50	0.46
2:G:36:TRP:HD1	2:G:70:ILE:HD12	1.80	0.46
2:A:165:TRP:CZ3	2:A:207:CYS:HB3	2.50	0.46
2:G:154:LYS:HA	2:G:188:SER:OG	2.15	0.46
3:H:105:GLU:OE1	3:H:140:TYR:HE2	1.99	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:91:HIS:O	3:D:92:GLN:HB3	2.15	0.46
3:H:118:PHE:HB2	3:H:133:VAL:HB	1.97	0.46
2:A:174:VAL:HG22	2:A:193:VAL:HB	1.97	0.46
3:H:166:GLN:NE2	3:H:171:SER:HB3	2.31	0.46
2:A:48:VAL:HG13	2:A:64:VAL:HG21	1.98	0.45
3:F:34:ALA:HB3	3:F:89:GLN:HB3	1.98	0.45
3:H:140:TYR:CG	3:H:141:PRO:HA	2.51	0.45
1:I:130:ILE:HD11	1:I:223:ARG:O	2.16	0.45
2:A:210:ASN:ND2	2:A:212:LYS:HE2	2.31	0.45
2:C:5:VAL:HG23	2:C:23:ALA:HB3	1.98	0.45
1:T:237:HIS:O	1:T:240:MET:HG2	2.15	0.45
1:I:181:PRO:HG3	1:I:244:SER:OG	2.16	0.45
2:C:157:PHE:HB2	2:C:186:LEU:HD22	1.98	0.45
1:T:131:ILE:O	1:T:135:ILE:HG13	2.16	0.45
1:I:236:SER:HB2	1:I:239:GLN:HG3	1.98	0.45
2:C:35:HIS:CE1	3:D:93:TYR:HD2	2.35	0.45
2:G:96:CYS:O	2:G:115:GLY:N	2.50	0.45
3:F:89:GLN:HG2	3:F:90:GLN:N	2.31	0.45
3:H:2:ILE:HB	3:H:97:ILE:HD12	1.99	0.45
2:E:51:ILE:HG13	2:E:58:THR:HG22	1.99	0.45
1:J:127:PHE:O	1:J:131:ILE:HG13	2.16	0.45
1:S:190:ILE:HA	1:S:226:ARG:NH2	2.32	0.45
2:C:60:TYR:CZ	2:C:70:ILE:HG22	2.51	0.45
2:G:135:LEU:HB2	2:G:150:GLY:C	2.42	0.45
3:H:120:PRO:HG3	3:H:130:ALA:HB1	1.99	0.45
2:C:96:CYS:O	2:C:115:GLY:N	2.50	0.45
2:C:123:SER:HB3	2:C:157:PHE:HZ	1.82	0.45
2:C:135:LEU:N	2:C:150:GLY:O	2.49	0.45
1:I:133:MET:SD	1:I:220:VAL:HG13	2.57	0.45
2:A:37:VAL:HG22	2:A:47:TRP:HA	1.99	0.44
2:G:5:VAL:HG23	2:G:23:ALA:HB3	2.00	0.44
2:G:51:ILE:HG13	2:G:58:THR:HG22	1.99	0.44
3:D:166:GLN:NE2	3:D:171:SER:HB3	2.32	0.44
3:D:3:GLN:HB2	3:D:26:SER:HB3	1.98	0.44
3:F:140:TYR:CG	3:F:141:PRO:HA	2.52	0.44
2:G:152:LEU:CD1	2:G:190:SER:HB3	2.47	0.44
3:H:34:ALA:HB3	3:H:89:GLN:HB3	1.99	0.44
3:F:120:PRO:HB3	3:F:131:SER:H	1.81	0.44
2:A:5:VAL:HG23	2:A:23:ALA:HB3	2.00	0.44
1:I:190:ILE:HA	1:I:226:ARG:NH2	2.32	0.44
2:G:60:TYR:CE1	2:G:70:ILE:HG22	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:120:PRO:HB3	3:B:131:SER:H	1.83	0.44
2:G:208:ASN:HD22	2:G:219:ASP:CG	2.26	0.44
1:S:130:ILE:HD11	1:S:223:ARG:O	2.18	0.44
2:G:48:VAL:HG13	2:G:64:VAL:HG21	2.00	0.44
3:H:166:GLN:HG2	3:H:171:SER:HA	1.99	0.44
2:C:130:PRO:HB2	2:C:153:VAL:HG13	2.00	0.44
2:G:112:ASP:OD2	2:G:112:ASP:N	2.46	0.44
2:G:152:LEU:HD12	2:G:190:SER:HB3	2.00	0.44
2:G:103:TYR:HB3	1:I:202:VAL:HG13	2.00	0.43
1:S:133:MET:SD	1:S:220:VAL:HG13	2.58	0.43
2:A:99:LYS:O	3:B:93:TYR:OH	2.31	0.43
2:G:212:LYS:N	2:G:213:PRO:CD	2.81	0.43
2:G:157:PHE:HA	2:G:158:PRO:HA	1.79	0.43
3:B:190:LYS:O	3:B:210:ASN:HA	2.18	0.43
2:E:27:PHE:CE1	2:E:98:ARG:HD3	2.53	0.43
2:E:37:VAL:HG22	2:E:47:TRP:HA	2.01	0.43
1:S:139:LEU:HD21	1:T:217:ARG:HH11	1.83	0.43
1:I:232:ARG:HB3	1:I:236:SER:OG	2.18	0.43
2:A:12:VAL:HG13	2:A:122:VAL:HG22	2.00	0.43
2:A:51:ILE:HG13	2:A:58:THR:HG22	2.01	0.43
2:C:189:LEU:C	2:C:189:LEU:HD12	2.44	0.43
2:E:148:ALA:HB2	2:E:194:THR:HG22	2.00	0.43
2:C:36:TRP:CG	2:C:81:LEU:HD22	2.53	0.43
2:E:130:PRO:HB3	2:E:156:TYR:HB3	2.01	0.43
1:J:155:LEU:HD11	1:J:198:ILE:HD12	2.00	0.43
3:D:2:ILE:HB	3:D:97:ILE:HD12	2.00	0.42
3:F:3:GLN:HB2	3:F:26:SER:HB3	2.01	0.42
1:I:131:ILE:O	1:I:135:ILE:HG13	2.19	0.42
2:C:174:VAL:HG22	2:C:193:VAL:HB	2.01	0.42
2:G:47:TRP:CD1	3:H:96:LEU:HD11	2.54	0.42
3:B:23:CYS:HB2	3:B:35:TRP:CH2	2.54	0.42
1:T:232:ARG:HD2	1:T:232:ARG:HA	1.73	0.42
1:I:224:LEU:O	1:I:228:VAL:HG23	2.19	0.42
2:C:2:VAL:HA	2:C:25:SER:O	2.19	0.42
3:B:119:PRO:HB3	3:B:209:PHE:CE1	2.55	0.42
2:C:112:ASP:OD2	2:C:112:ASP:N	2.48	0.42
3:D:33:VAL:O	3:D:51:ALA:N	2.52	0.42
3:D:45:LYS:HD2	3:D:45:LYS:HA	1.87	0.42
1:T:151:ASP:OD1	1:T:151:ASP:N	2.52	0.42
1:S:210:THR:OG1	1:S:213:ASP:HB3	2.20	0.42
3:F:35:TRP:CD2	3:F:73:LEU:HB2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:145:LYS:HE2	3:F:197:THR:HB	2.02	0.42
1:J:190:ILE:HA	1:J:226:ARG:HH21	1.85	0.42
2:A:179:ALA:HA	2:A:189:LEU:HB3	2.02	0.42
1:I:182:TRP:CE3	1:I:240:MET:HB3	2.55	0.42
3:F:108:ARG:HH21	3:F:111:ALA:HB2	1.85	0.42
2:G:133:PHE:HB3	3:H:121:SER:OG	2.19	0.42
1:J:224:LEU:O	1:J:228:VAL:HG23	2.19	0.42
2:A:99:LYS:HZ1	3:B:92:GLN:H	1.67	0.41
3:F:2:ILE:HB	3:F:97:ILE:HD12	2.01	0.41
2:E:48:VAL:HG13	2:E:64:VAL:HG21	2.02	0.41
3:F:35:TRP:HB2	3:F:48:ILE:HB	2.01	0.41
3:H:186:TYR:HA	3:H:192:TYR:OH	2.19	0.41
2:A:27:PHE:CE1	2:A:98:ARG:HD3	2.55	0.41
2:A:36:TRP:NE1	2:A:81:LEU:HB2	2.36	0.41
2:C:223:GLU:HA	2:C:224:PRO:HD3	1.90	0.41
1:T:119:ARG:NH1	1:T:119:ARG:HA	2.35	0.41
3:F:118:PHE:HB2	3:F:133:VAL:HB	2.03	0.41
2:A:29:VAL:HG12	2:A:53:PRO:HG3	2.02	0.41
3:B:136:LEU:HB2	3:B:175:LEU:HB3	2.01	0.41
3:D:34:ALA:HB3	3:D:89:GLN:HB3	2.02	0.41
1:T:190:ILE:HA	1:T:226:ARG:HH21	1.84	0.41
3:F:163:VAL:HG12	3:F:164:THR:O	2.19	0.41
2:G:2:VAL:HA	2:G:25:SER:O	2.19	0.41
2:C:221:LYS:HE2	2:C:223:GLU:OE1	2.21	0.41
3:H:120:PRO:HB3	3:H:131:SER:H	1.85	0.41
2:A:99:LYS:HZ1	3:B:91:HIS:HB3	1.85	0.41
3:F:120:PRO:HG3	3:F:130:ALA:HB1	2.03	0.41
3:F:193:ALA:HB2	3:F:208:SER:HB3	2.03	0.41
2:A:36:TRP:HD1	2:A:70:ILE:HD12	1.86	0.41
3:H:96:LEU:H	3:H:96:LEU:CD2	2.32	0.41
1:S:110:ARG:HB2	1:S:171:ALA:HB1	2.03	0.41
2:A:97:ALA:CB	2:A:111:MET:HE2	2.49	0.41
2:A:122:VAL:O	2:A:122:VAL:HG12	2.20	0.41
3:B:107:LYS:HA	3:B:140:TYR:HH	1.85	0.41
3:B:175:LEU:HD23	3:B:175:LEU:C	2.45	0.41
1:T:196:VAL:CB	1:T:219:VAL:HG21	2.51	0.41
2:E:99:LYS:HE3	2:E:109:TRP:O	2.20	0.41
2:E:174:VAL:HG22	2:E:193:VAL:CG2	2.51	0.41
2:E:223:GLU:HA	2:E:224:PRO:HD3	1.94	0.41
2:G:104:TRP:HB3	2:G:105:ARG:H	1.67	0.41
2:G:189:LEU:C	2:G:189:LEU:HD12	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:217:ARG:NH1	1:T:139:LEU:HD21	2.36	0.41
2:A:39:GLN:C	2:A:92:ALA:HB1	2.46	0.41
2:C:157:PHE:HA	2:C:158:PRO:HA	1.83	0.41
2:E:96:CYS:O	2:E:115:GLY:N	2.54	0.41
3:F:30:SER:OG	3:F:31:SER:N	2.54	0.41
2:E:161:VAL:HG11	2:E:209:VAL:HG13	2.03	0.40
3:B:89:GLN:HG2	3:B:90:GLN:N	2.32	0.40
2:E:70:ILE:HG23	2:E:70:ILE:O	2.21	0.40
3:B:45:LYS:HD2	3:B:45:LYS:HA	1.89	0.40
3:B:120:PRO:HG3	3:B:130:ALA:HB1	2.02	0.40
2:C:105:ARG:HD2	1:J:151:ASP:OD2	2.22	0.40
1:I:237:HIS:HA	1:I:240:MET:HG2	2.02	0.40
3:D:167:ASP:OD1	3:D:168:SER:N	2.54	0.40
1:S:151:ASP:OD1	1:S:151:ASP:N	2.53	0.40
3:B:2:ILE:HB	3:B:97:ILE:HD12	2.04	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	I	137/155 (88%)	126 (92%)	10 (7%)	1 (1%)	18	51
1	J	137/155 (88%)	126 (92%)	11 (8%)	0	100	100
1	S	126/155 (81%)	115 (91%)	10 (8%)	1 (1%)	16	49
1	T	137/155 (88%)	126 (92%)	10 (7%)	1 (1%)	18	51
2	A	217/224 (97%)	205 (94%)	11 (5%)	1 (0%)	24	57
2	C	215/224 (96%)	200 (93%)	14 (6%)	1 (0%)	24	57
2	E	217/224 (97%)	209 (96%)	5 (2%)	3 (1%)	9	38
2	G	217/224 (97%)	208 (96%)	8 (4%)	1 (0%)	24	57

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	B	209/211 (99%)	194 (93%)	14 (7%)	1 (0%)	24	57
3	D	209/211 (99%)	195 (93%)	13 (6%)	1 (0%)	24	57
3	F	209/211 (99%)	195 (93%)	12 (6%)	2 (1%)	12	44
3	H	209/211 (99%)	195 (93%)	13 (6%)	1 (0%)	24	57
All	All	2239/2360 (95%)	2094 (94%)	131 (6%)	14 (1%)	21	54

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	F	95	SER
1	S	212	ALA
2	E	155	ASP
3	F	92	GLN
2	A	91	THR
3	B	92	GLN
2	C	91	THR
3	D	92	GLN
1	T	212	ALA
2	E	91	THR
2	G	91	THR
3	H	92	GLN
1	I	212	ALA
2	E	160	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	I	120/134 (90%)	116 (97%)	4 (3%)	33	58
1	J	120/134 (90%)	115 (96%)	5 (4%)	26	52
1	S	110/134 (82%)	108 (98%)	2 (2%)	51	68
1	T	120/134 (90%)	113 (94%)	7 (6%)	18	45
2	A	182/187 (97%)	177 (97%)	5 (3%)	39	61

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	180/187 (96%)	176 (98%)	4 (2%)	45	65
2	E	182/187 (97%)	180 (99%)	2 (1%)	65	74
2	G	182/187 (97%)	178 (98%)	4 (2%)	45	65
3	B	185/185 (100%)	183 (99%)	2 (1%)	65	74
3	D	185/185 (100%)	184 (100%)	1 (0%)	81	80
3	F	185/185 (100%)	184 (100%)	1 (0%)	81	80
3	H	185/185 (100%)	184 (100%)	1 (0%)	81	80
All	All	1936/2024 (96%)	1898 (98%)	38 (2%)	48	66

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	S	126	ILE
1	S	134	ILE
2	A	131	SER
2	A	181	LEU
2	A	193	VAL
2	A	197	SER
2	A	199	SER
3	B	65	SER
3	B	94	ASN
2	C	12	VAL
2	C	131	SER
2	C	149	LEU
2	C	194	THR
3	D	110	VAL
1	T	126	ILE
1	T	134	ILE
1	T	227	VAL
1	T	232	ARG
1	T	240	MET
1	T	241	LYS
1	T	244	SER
2	E	194	THR
2	E	218	VAL
3	F	65	SER
2	G	12	VAL
2	G	131	SER
2	G	194	THR
2	G	198	SER

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Mol	Chain	Res	Type
3	H	65	SER
1	I	126	ILE
1	I	134	ILE
1	I	238	GLN
1	I	240	MET
1	J	126	ILE
1	J	134	ILE
1	J	227	VAL
1	J	241	LYS
1	J	243	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	182	GLN
2	A	210	ASN
3	B	91	HIS
3	B	160	GLN
2	C	182	GLN
3	D	160	GLN
2	E	182	GLN
3	F	37	GLN
3	F	94	ASN
3	F	138	ASN
3	F	160	GLN
2	G	182	GLN
1	J	237	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	I	139/155 (89%)	0.67	14 (10%) 12 8	160, 217, 278, 289	0
1	J	139/155 (89%)	0.53	9 (6%) 25 15	146, 202, 271, 292	0
1	S	128/155 (82%)	0.33	11 (8%) 16 11	95, 150, 248, 278	0
1	T	139/155 (89%)	0.36	10 (7%) 21 13	106, 157, 255, 293	0
2	A	221/224 (98%)	0.06	11 (4%) 34 19	70, 104, 159, 214	0
2	C	219/224 (97%)	0.14	5 (2%) 61 34	66, 135, 208, 261	0
2	E	221/224 (98%)	0.05	8 (3%) 46 25	68, 113, 171, 223	0
2	G	221/224 (98%)	0.40	20 (9%) 15 10	104, 170, 228, 263	0
3	B	211/211 (100%)	-0.04	3 (1%) 73 45	67, 112, 146, 185	0
3	D	211/211 (100%)	0.08	8 (3%) 44 24	70, 125, 186, 237	0
3	F	211/211 (100%)	0.16	10 (4%) 36 19	90, 135, 178, 201	0
3	H	211/211 (100%)	0.14	7 (3%) 49 27	91, 157, 212, 269	0
All	All	2271/2360 (96%)	0.21	116 (5%) 33 18	66, 140, 239, 293	0

All (116) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	G	174	VAL	7.2
3	F	95	SER	6.9
1	S	228	VAL	6.3
2	C	17	SER	6.3
1	T	114	ASP	5.9
1	I	226	ARG	5.8
1	S	206	TYR	5.0
1	I	153	MET	4.9
2	E	62	ASP	4.8
1	T	172	TYR	4.8
1	J	226	ARG	4.8

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Mol	Chain	Res	Type	RSRZ
3	F	114	SER	4.5
2	G	194	THR	4.4
1	T	206	TYR	4.4
1	I	233	ILE	4.4
1	T	170	PHE	4.4
3	D	95	SER	4.3
1	J	230	LEU	4.1
3	H	55	TYR	4.0
3	F	199	GLN	4.0
2	C	15	GLY	4.0
2	G	175	HIS	4.0
1	J	227	VAL	4.0
1	S	229	ARG	3.9
2	E	143	SER	3.8
2	A	59	TYR	3.7
2	A	160	PRO	3.6
1	I	234	PHE	3.6
3	D	18	ARG	3.6
2	G	195	VAL	3.5
1	S	225	LEU	3.5
2	A	110	ALA	3.5
3	D	92	GLN	3.5
3	D	165	GLU	3.3
1	S	226	ARG	3.3
1	T	110	ARG	3.3
1	I	154	ALA	3.3
1	I	230	LEU	3.2
1	J	234	PHE	3.1
2	C	13	GLN	3.1
3	F	92	GLN	3.1
3	D	168	SER	3.0
2	A	109	TRP	3.0
3	H	80	PRO	3.0
2	A	139	SER	3.0
3	F	165	GLU	2.9
2	G	173	GLY	2.9
2	G	193	VAL	2.9
2	E	111	MET	2.8
1	I	126	ILE	2.7
1	T	109	VAL	2.7
1	S	224	LEU	2.7
1	J	231	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	J	233	ILE	2.6
2	G	136	ALA	2.6
2	A	204	THR	2.6
2	G	111	MET	2.6
2	G	147	ALA	2.6
3	H	83	PHE	2.6
2	A	106	ASP	2.6
3	F	94	ASN	2.6
2	G	171	THR	2.5
1	J	206	TYR	2.5
2	C	173	GLY	2.5
2	A	159	GLU	2.5
2	E	110	ALA	2.5
3	F	113	PRO	2.5
1	I	186	ASP	2.5
2	G	112	ASP	2.5
2	G	179	ALA	2.5
2	G	138	SER	2.5
1	T	113	ILE	2.5
1	I	207	VAL	2.5
2	E	106	ASP	2.4
3	D	98	PHE	2.4
2	A	104	TRP	2.4
1	T	111	ALA	2.4
3	B	93	TYR	2.4
3	D	93	TYR	2.4
1	S	231	ALA	2.4
2	G	166	ASN	2.4
3	B	169	LYS	2.4
2	G	176	THR	2.3
3	H	106	ILE	2.3
1	S	227	VAL	2.3
1	I	132	LEU	2.3
3	F	90	GLN	2.3
3	B	95	SER	2.3
1	S	230	LEU	2.2
2	E	59	TYR	2.2
1	J	165	LEU	2.2
2	G	206	ILE	2.2
1	T	226	ARG	2.2
2	E	146	THR	2.2
2	C	112	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
3	H	56	SER	2.2
3	D	106	ILE	2.2
3	F	93	TYR	2.2
1	I	114	ASP	2.1
1	T	227	VAL	2.1
1	J	200	TYR	2.1
2	G	192	VAL	2.1
2	G	172	SER	2.1
1	S	207	VAL	2.1
3	F	116	PHE	2.1
1	I	223	ARG	2.1
2	G	204	THR	2.1
1	I	229	ARG	2.1
2	E	130	PRO	2.0
2	G	73	ASP	2.0
1	S	189	ILE	2.0
1	I	218	LEU	2.0
2	A	100	GLN	2.0
3	H	147	GLN	2.0
2	A	101	TYR	2.0
3	H	51	ALA	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.