



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:55 PM UTC

PDB ID : 3IPD / pdb\_00003ipd  
Title : Helical extension of the neuronal SNARE complex into the membrane, space-group I 21 21 21  
Authors : Stein, A.; Weber, G.; Wahl, M.C.; Jahn, R.  
Deposited on : 2009-08-17  
Resolution : 4.80 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0  
Xtrriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

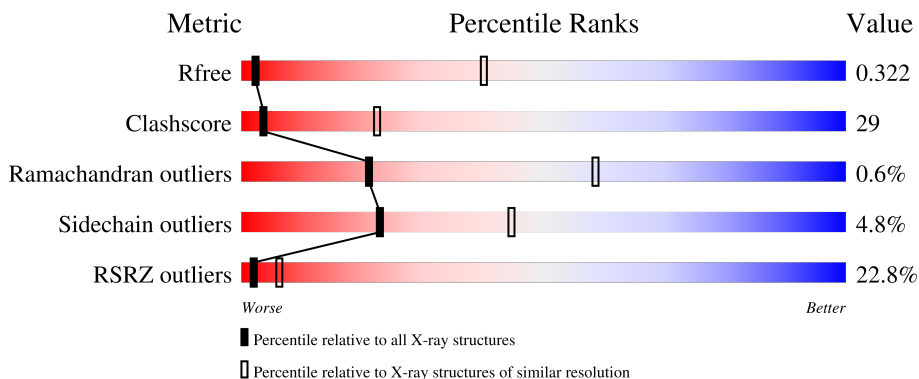
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 4.80 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.




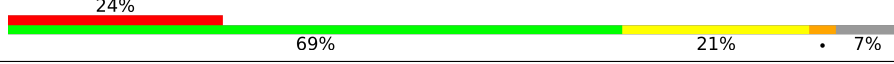
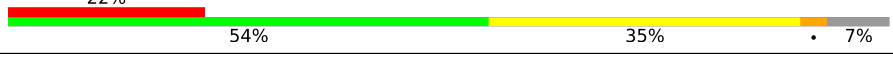
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	1018 (5.66-3.94)
Clashscore	190562	1001 (5.60-3.98)
Ramachandran outliers	187476	1104 (5.70-3.90)
Sidechain outliers	187428	1085 (5.70-3.90)
RSRZ outliers	180081	1013 (5.66-3.94)

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  $\geq 3$ , 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	91	
1	E	91	
2	B	109	
2	F	109	
3	C	80	

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Mol	Chain	Length	Quality of chain
3	G	80	
4	D	68	
4	H	68	

## 2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 5242 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 Vesicle-associated membrane protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	91	733	467	129	132	5	0	0	0
1	E	91	733	467	129	132	5	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	GLY	-	expression tag	UNP P63045
A	27	SER	-	expression tag	UNP P63045
A	28	HIS	-	expression tag	UNP P63045
A	29	MET	-	expression tag	UNP P63045
E	26	GLY	-	expression tag	UNP P63045
E	27	SER	-	expression tag	UNP P63045
E	28	HIS	-	expression tag	UNP P63045
E	29	MET	-	expression tag	UNP P63045

- Molecule 2 is a protein called Syntaxin-1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	98	779	490	135	146	8	0	0	0
2	F	98	779	490	135	146	8	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	180	GLY	-	expression tag	UNP P32851
B	181	SER	-	expression tag	UNP P32851
B	182	HIS	-	expression tag	UNP P32851
F	180	GLY	-	expression tag	UNP P32851

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Chain	Residue	Modelled	Actual	Comment	Reference
F	181	SER	-	expression tag	UNP P32851
F	182	HIS	-	expression tag	UNP P32851

- Molecule 3 is a protein called Synaptosomal-associated protein 25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	75	Total	C	N	O	S	0	0	0
			607	360	112	130	5			
3	G	75	Total	C	N	O	S	0	0	0
			607	360	112	130	5			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	4	GLY	-	expression tag	UNP P60881
C	5	SER	-	expression tag	UNP P60881
C	6	HIS	-	expression tag	UNP P60881
G	4	GLY	-	expression tag	UNP P60881
G	5	SER	-	expression tag	UNP P60881
G	6	HIS	-	expression tag	UNP P60881

- Molecule 4 is a protein called Synaptosomal-associated protein 25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	63	Total	C	N	O	S	0	0	0
			502	293	98	106	5			
4	H	63	Total	C	N	O	S	0	0	0
			502	293	98	106	5			

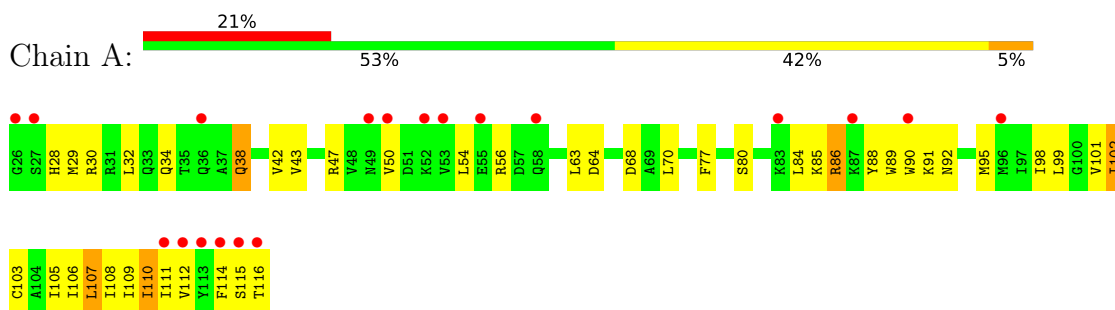
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	137	GLY	-	expression tag	UNP P60881
D	138	SER	-	expression tag	UNP P60881
D	139	HIS	-	expression tag	UNP P60881
D	140	MET	-	expression tag	UNP P60881
H	137	GLY	-	expression tag	UNP P60881
H	138	SER	-	expression tag	UNP P60881
H	139	HIS	-	expression tag	UNP P60881
H	140	MET	-	expression tag	UNP P60881

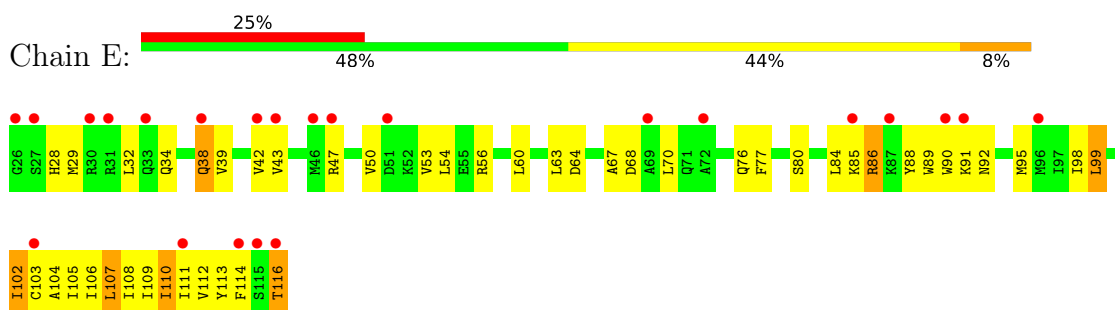
### 3 Residue-property plots [i](#)

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.

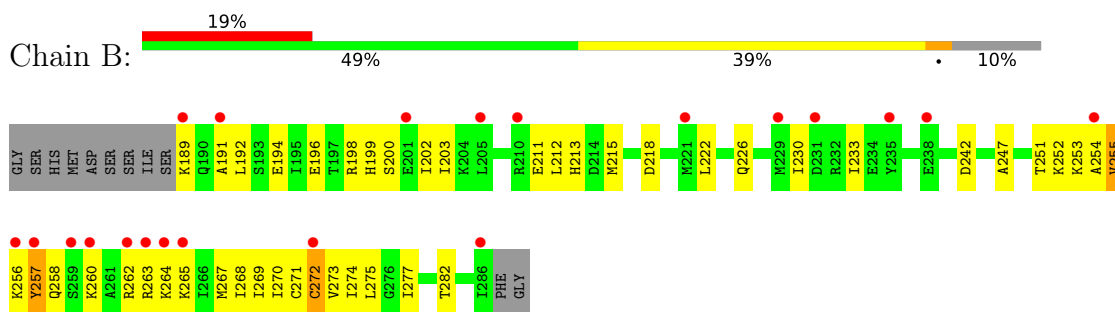
- Molecule 1: Vesicle-associated membrane protein 2



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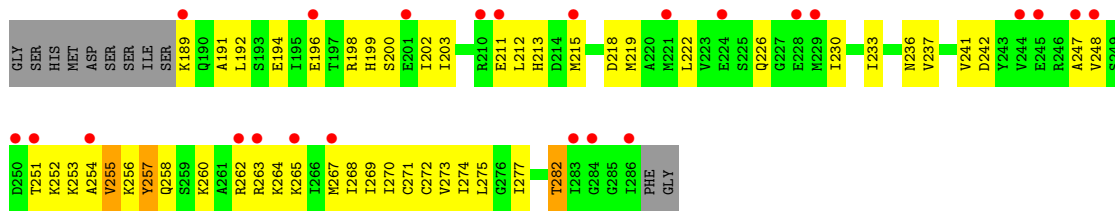


- Molecule 2: Syntaxin-1A

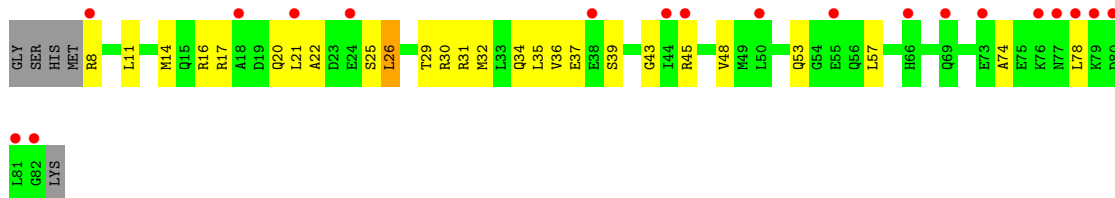


- Molecule 2: Syntaxin-1A

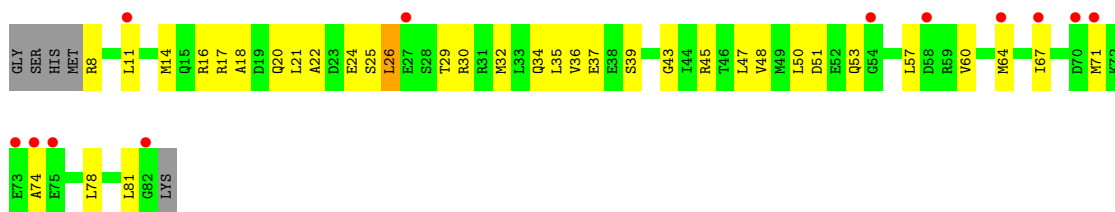




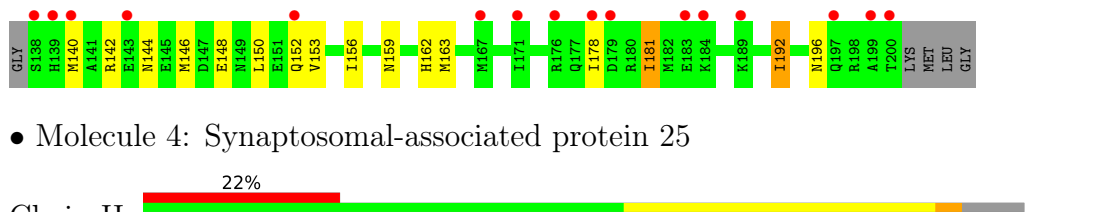
• Molecule 3: Synaptosomal-associated protein 25



• Molecule 3: Synaptosomal-associated protein 25



• Molecule 4: Synaptosomal-associated protein 25



• Molecule 4: Synaptosomal-associated protein 25



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	109.85Å 215.68Å 262.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.88 – 4.80 49.88 – 4.80	Depositor EDS
% Data completeness (in resolution range)	84.2 (49.88-4.80) 99.7 (49.88-4.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.10 (at 4.86Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.4_4)	Depositor
R, $R_{free}$	0.304 , 0.332 0.304 , 0.322	Depositor DCC
$R_{free}$ test set	783 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	202.5	Xtrriage
Anisotropy	0.759	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.26 , 232.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.40$ , $\langle L^2 \rangle = 0.23$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	5242	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	289.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.77% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

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	0.39	0/741	0.81	0/996
1	E	0.37	0/741	0.83	1/996 (0.1%)
2	B	0.36	0/785	0.81	0/1049
2	F	0.36	0/785	0.82	0/1049
3	C	0.28	0/607	0.78	0/807
3	G	0.30	0/607	0.79	0/807
4	D	0.31	0/503	0.79	0/671
4	H	0.32	0/503	0.79	0/671
All	All	0.34	0/5272	0.80	1/7046 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	99	LEU	N-CA-C	-5.09	107.12	113.38

There are no chirality outliers.

There are no planarity outliers.

### 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	A	733	0	768	73	0
1	E	733	0	768	92	0
2	B	779	0	815	69	0
2	F	779	0	815	90	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	607	0	592	23	0
3	G	607	0	592	45	0
4	D	502	0	480	21	0
4	H	502	0	480	41	0
All	All	5242	0	5310	309	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 29.

All (309) 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:A:109:ILE:HD12	2:B:275:LEU:HD22	1.40	1.03
1:A:28:HIS:HB2	2:B:198:ARG:CZ	1.98	0.94
1:E:106:ILE:HG12	2:F:275:LEU:HD21	1.47	0.93
1:A:91:LYS:O	1:A:95:MET:HG2	1.75	0.86
1:E:91:LYS:O	1:E:95:MET:HG2	1.76	0.86
1:A:95:MET:O	2:B:264:LYS:HB3	1.79	0.82
1:E:43:VAL:HG22	2:F:212:LEU:HD13	1.61	0.81
1:E:106:ILE:HG12	2:F:275:LEU:CD2	2.08	0.81
1:A:88:TYR:HE2	2:B:258:GLN:HB2	1.47	0.80
1:A:28:HIS:HB2	2:B:198:ARG:NH2	1.96	0.80
1:A:88:TYR:CE2	2:B:258:GLN:HB2	2.18	0.79
4:H:146:MET:O	4:H:150:LEU:HB2	1.83	0.79
4:D:146:MET:O	4:D:150:LEU:HB2	1.83	0.79
3:G:22:ALA:HB1	4:H:142:ARG:HD3	1.64	0.78
1:E:88:TYR:OH	2:F:258:GLN:N	2.16	0.78
1:E:109:ILE:HD12	2:F:275:LEU:HD22	1.66	0.78
1:E:39:VAL:CG2	4:H:157:ILE:HD13	2.14	0.76
1:A:34:GLN:HE22	4:H:163:MET:HA	1.51	0.75
1:A:106:ILE:HG12	2:B:275:LEU:HD21	1.71	0.73
1:E:89:TRP:CE3	1:E:90:TRP:HD1	2.08	0.72
1:A:89:TRP:CE3	1:A:90:TRP:HD1	2.06	0.72
3:G:25:SER:HB3	4:H:146:MET:HE1	1.72	0.72
1:A:89:TRP:CE3	1:A:90:TRP:CD1	2.78	0.72
1:E:39:VAL:HG23	4:H:157:ILE:HD13	1.72	0.72
3:G:26:LEU:HD22	4:H:146:MET:HB3	1.72	0.72
1:A:88:TYR:OH	2:B:255:VAL:C	2.32	0.71
1:E:63:LEU:HD23	2:F:233:ILE:HG21	1.70	0.71
1:E:89:TRP:CE3	1:E:90:TRP:CD1	2.78	0.71
2:F:257:TYR:H	2:F:257:TYR:HD2	1.39	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:TYR:OH	2:B:258:GLN:N	2.21	0.71
1:E:95:MET:O	2:F:264:LYS:HB3	1.91	0.71
1:E:108:ILE:HA	1:E:111:ILE:HD12	1.73	0.71
4:D:162:HIS:HE1	1:E:34:GLN:HE22	1.38	0.71
2:B:199:HIS:HE2	2:B:203:ILE:HD11	1.56	0.70
1:A:38:GLN:HG2	4:H:162:HIS:CD2	2.26	0.70
2:F:199:HIS:HE1	3:G:24:GLU:OE1	1.75	0.69
1:E:63:LEU:CD1	4:H:182:MET:HG3	2.23	0.69
2:F:199:HIS:HE2	2:F:203:ILE:HD11	1.58	0.69
1:E:88:TYR:HH	2:F:258:GLN:HB2	1.58	0.68
1:E:63:LEU:HD13	4:H:182:MET:HG3	1.75	0.68
1:A:109:ILE:HD12	2:B:275:LEU:CD2	2.21	0.68
2:F:211:GLU:O	2:F:215:MET:HG2	1.93	0.68
2:B:211:GLU:O	2:B:215:MET:HG2	1.92	0.68
3:G:64:MET:CE	4:H:185:ALA:HA	2.24	0.68
2:B:257:TYR:H	2:B:257:TYR:HD2	1.40	0.67
1:A:108:ILE:HA	1:A:111:ILE:HD12	1.77	0.67
4:D:162:HIS:CE1	1:E:34:GLN:NE2	2.63	0.67
4:D:162:HIS:HE1	1:E:34:GLN:NE2	1.94	0.66
1:A:89:TRP:HE3	1:A:90:TRP:HD1	1.43	0.66
3:C:26:LEU:HD22	4:D:146:MET:HB3	1.76	0.66
1:A:106:ILE:HD11	2:B:271:CYS:HB3	1.77	0.65
2:F:196:GLU:O	2:F:200:SER:HB2	1.95	0.65
3:G:64:MET:HE2	4:H:185:ALA:HA	1.77	0.65
1:A:42:VAL:HG11	2:B:212:LEU:HD21	1.78	0.64
1:A:95:MET:HE2	2:B:265:LYS:HG3	1.79	0.64
2:B:196:GLU:O	2:B:200:SER:HB2	1.98	0.64
1:A:107:LEU:O	1:A:111:ILE:HG13	1.98	0.63
3:C:26:LEU:O	3:C:29:THR:HG22	1.99	0.63
2:F:192:LEU:HD13	3:G:14:MET:HG2	1.81	0.63
2:F:230:ILE:HD13	3:G:57:LEU:HD11	1.80	0.63
1:E:56:ARG:HH21	1:E:56:ARG:HG2	1.64	0.63
1:E:85:LYS:HE2	2:F:253:LYS:CB	2.27	0.63
1:E:88:TYR:OH	2:F:258:GLN:HB2	1.98	0.62
2:F:258:GLN:HG2	2:F:262:ARG:HH11	1.64	0.62
1:A:56:ARG:HG2	1:A:56:ARG:HH21	1.64	0.62
3:G:26:LEU:O	3:G:29:THR:HG22	1.99	0.62
3:G:71:MET:HB3	4:H:191:ARG:HD2	1.81	0.62
4:D:162:HIS:CD2	1:E:38:GLN:OE1	2.53	0.62
1:E:89:TRP:HE3	1:E:90:TRP:HD1	1.44	0.61
1:E:102:ILE:O	1:E:106:ILE:HG13	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:198:ARG:O	2:F:202:ILE:HG13	2.00	0.61
1:A:89:TRP:HE3	1:A:90:TRP:CD1	2.17	0.61
1:E:99:LEU:HD13	2:F:268:ILE:CG1	2.30	0.61
2:F:199:HIS:CE1	3:G:24:GLU:HB2	2.35	0.61
2:B:258:GLN:HG2	2:B:262:ARG:HH11	1.63	0.61
2:B:198:ARG:O	2:B:202:ILE:HG13	2.01	0.60
2:F:213:HIS:HD2	3:G:35:LEU:HD23	1.66	0.60
1:E:105:ILE:O	1:E:109:ILE:HG13	2.01	0.60
1:E:107:LEU:O	1:E:111:ILE:HG13	2.01	0.60
1:E:89:TRP:HE3	1:E:90:TRP:CD1	2.17	0.60
1:A:105:ILE:O	1:A:109:ILE:HG13	2.01	0.60
1:A:106:ILE:HG12	2:B:275:LEU:CD2	2.32	0.59
1:E:67:ALA:HB1	2:F:236:ASN:CB	2.32	0.59
2:F:257:TYR:N	2:F:257:TYR:CD2	2.70	0.59
1:E:103:CYS:O	1:E:107:LEU:HD23	2.02	0.59
1:E:88:TYR:CE2	2:F:258:GLN:HB2	2.38	0.59
2:F:213:HIS:HA	3:G:39:SER:OG	2.03	0.59
2:F:255:VAL:HG12	3:G:81:LEU:CD1	2.32	0.59
2:F:254:ALA:O	3:G:81:LEU:HD21	2.03	0.58
2:F:263:ARG:O	2:F:267:MET:HG2	2.03	0.58
2:F:271:CYS:HA	2:F:274:ILE:HB	1.86	0.58
2:B:263:ARG:O	2:B:267:MET:HG2	2.04	0.57
3:C:74:ALA:O	3:C:78:LEU:HD13	2.04	0.57
2:B:271:CYS:HA	2:B:274:ILE:HB	1.85	0.57
2:B:257:TYR:N	2:B:257:TYR:CD2	2.71	0.57
2:F:248:VAL:HG23	3:G:74:ALA:HB2	1.86	0.57
3:G:22:ALA:O	3:G:26:LEU:HB2	2.04	0.57
3:C:22:ALA:O	3:C:26:LEU:HB2	2.05	0.56
1:A:103:CYS:O	1:A:107:LEU:HD23	2.05	0.56
1:E:99:LEU:HD13	2:F:268:ILE:HG13	1.87	0.56
1:A:109:ILE:CD1	2:B:275:LEU:HD22	2.26	0.56
1:A:109:ILE:O	1:A:109:ILE:HG22	2.06	0.56
2:F:189:LYS:HA	2:F:192:LEU:HD12	1.87	0.56
1:E:53:VAL:HG21	2:F:219:MET:SD	2.46	0.55
1:A:92:ASN:HD21	2:B:264:LYS:NZ	2.04	0.55
2:B:189:LYS:HA	2:B:192:LEU:HD12	1.87	0.55
3:C:17:ARG:HA	3:C:20:GLN:HB3	1.88	0.55
3:G:8:ARG:HG2	3:G:11:LEU:HD12	1.89	0.55
3:G:74:ALA:O	3:G:78:LEU:HD13	2.07	0.55
2:B:192:LEU:HD13	3:C:14:MET:HE2	1.88	0.55
3:G:17:ARG:HA	3:G:20:GLN:HB3	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:ILE:O	1:A:106:ILE:HG13	2.07	0.54
3:C:43:GLY:HA3	4:D:163:MET:HE1	1.90	0.54
3:G:36:VAL:HG11	4:H:156:ILE:HB	1.89	0.54
1:A:115:SER:O	1:A:116:THR:CB	2.55	0.54
2:B:213:HIS:HA	3:C:39:SER:OG	2.07	0.54
1:E:109:ILE:HG22	1:E:109:ILE:O	2.06	0.54
1:A:30:ARG:HH11	3:G:51:ASP:HB2	1.71	0.54
3:C:8:ARG:HG2	3:C:11:LEU:HD12	1.88	0.54
2:B:258:GLN:HG2	2:B:262:ARG:NH1	2.23	0.53
2:B:213:HIS:HD2	3:C:35:LEU:HD23	1.74	0.53
2:F:273:VAL:O	2:F:277:ILE:HG13	2.07	0.53
2:F:241:VAL:HA	3:G:67:ILE:HD11	1.91	0.53
1:A:102:ILE:HD12	2:B:268:ILE:HG23	1.90	0.53
3:G:32:MET:O	3:G:36:VAL:HG23	2.09	0.53
2:F:255:VAL:HG12	3:G:81:LEU:HD13	1.89	0.53
2:B:273:VAL:O	2:B:277:ILE:HG13	2.09	0.52
2:F:199:HIS:CE1	3:G:24:GLU:OE1	2.61	0.52
4:D:152:GLN:O	4:D:156:ILE:HG13	2.10	0.52
2:F:258:GLN:HG2	2:F:262:ARG:NH1	2.23	0.52
1:A:56:ARG:HD3	2:B:226:GLN:OE1	2.10	0.52
1:E:67:ALA:HB1	2:F:236:ASN:HB2	1.90	0.52
1:E:106:ILE:HG12	2:F:275:LEU:CG	2.39	0.52
3:C:32:MET:O	3:C:36:VAL:HG23	2.09	0.52
4:H:152:GLN:O	4:H:156:ILE:HG13	2.10	0.52
3:G:26:LEU:HD23	4:H:142:ARG:HG3	1.92	0.52
1:E:28:HIS:HB2	2:F:198:ARG:NH2	2.25	0.52
2:F:213:HIS:HD1	2:F:213:HIS:C	2.18	0.52
3:G:43:GLY:HA3	4:H:163:MET:HE1	1.92	0.52
1:A:98:ILE:O	1:A:98:ILE:HG22	2.09	0.51
1:E:85:LYS:HE2	2:F:253:LYS:HB3	1.91	0.51
1:E:60:LEU:HD21	2:F:230:ILE:HG13	1.92	0.51
1:E:98:ILE:HG22	1:E:98:ILE:O	2.10	0.51
1:E:108:ILE:C	1:E:110:ILE:N	2.68	0.51
2:F:257:TYR:HA	2:F:260:LYS:HB2	1.92	0.51
1:E:56:ARG:HD3	2:F:226:GLN:OE1	2.10	0.51
1:A:64:ASP:HB2	2:B:233:ILE:HD11	1.93	0.51
2:B:213:HIS:C	2:B:213:HIS:HD1	2.18	0.50
1:A:88:TYR:CZ	2:B:258:GLN:N	2.79	0.50
2:B:265:LYS:O	2:B:269:ILE:HG13	2.12	0.50
1:A:95:MET:HE3	2:B:268:ILE:HD12	1.94	0.50
1:A:108:ILE:C	1:A:110:ILE:H	2.20	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:LEU:HD13	2:B:268:ILE:HG13	1.92	0.49
3:G:16:ARG:O	3:G:20:GLN:HB2	2.13	0.49
4:D:150:LEU:HD12	4:D:153:VAL:HG21	1.94	0.49
1:E:106:ILE:HD11	2:F:271:CYS:HB3	1.94	0.49
3:G:29:THR:CG2	4:H:149:ASN:HD21	2.26	0.49
1:A:88:TYR:HH	2:B:258:GLN:HB2	1.77	0.49
1:A:88:TYR:HH	2:B:255:VAL:C	2.20	0.49
1:A:92:ASN:ND2	2:B:260:LYS:HB3	2.28	0.49
3:G:45:ARG:HA	3:G:48:VAL:HG12	1.94	0.49
1:E:108:ILE:C	1:E:110:ILE:H	2.21	0.49
1:E:113:TYR:CE1	2:F:282:THR:HG21	2.48	0.49
2:B:230:ILE:HD13	3:C:57:LEU:HD11	1.93	0.49
1:A:63:LEU:HD11	4:D:181:ILE:HG22	1.94	0.48
1:E:88:TYR:CE1	2:F:257:TYR:HB2	2.48	0.48
1:A:108:ILE:C	1:A:110:ILE:N	2.67	0.48
1:A:43:VAL:HG22	2:B:212:LEU:HD13	1.95	0.48
1:E:92:ASN:HD21	2:F:264:LYS:NZ	2.11	0.48
3:C:45:ARG:HA	3:C:48:VAL:HG12	1.95	0.48
1:E:77:PHE:HA	4:H:196:ASN:OD1	2.13	0.48
1:E:99:LEU:HD13	2:F:268:ILE:HA	1.96	0.48
1:E:92:ASN:OD1	2:F:264:LYS:NZ	2.46	0.48
3:G:64:MET:HE2	4:H:185:ALA:CA	2.44	0.48
2:B:247:ALA:O	2:B:251:THR:HG23	2.13	0.48
3:C:16:ARG:O	3:C:20:GLN:HB2	2.13	0.48
2:F:192:LEU:HD22	3:G:18:ALA:HB2	1.95	0.48
1:A:85:LYS:HE2	2:B:253:LYS:CB	2.43	0.48
2:F:269:ILE:O	2:F:273:VAL:HG23	2.14	0.48
4:H:150:LEU:HD12	4:H:153:VAL:HG21	1.95	0.48
1:A:38:GLN:O	1:A:42:VAL:HG23	2.14	0.48
2:F:265:LYS:O	2:F:269:ILE:HG13	2.14	0.48
2:B:269:ILE:O	2:B:273:VAL:HG23	2.13	0.47
1:E:38:GLN:O	1:E:42:VAL:HG23	2.13	0.47
2:F:252:LYS:O	2:F:255:VAL:HG22	2.14	0.47
1:E:63:LEU:HD12	4:H:182:MET:CG	2.45	0.47
2:B:257:TYR:HA	2:B:260:LYS:HB2	1.97	0.47
1:E:88:TYR:CZ	2:F:258:GLN:HB2	2.49	0.47
3:G:64:MET:HE1	4:H:185:ALA:HA	1.96	0.47
2:F:265:LYS:O	2:F:268:ILE:HB	2.15	0.47
2:B:252:LYS:O	2:B:255:VAL:HG22	2.15	0.47
3:C:25:SER:HB3	4:D:146:MET:HE1	1.97	0.47
1:E:89:TRP:CD1	2:F:257:TYR:CE1	3.02	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:56:ARG:HH21	1:E:56:ARG:CG	2.26	0.46
1:E:63:LEU:HD12	4:H:182:MET:HG3	1.97	0.46
4:H:144:ASN:O	4:H:148:GLU:HB2	2.16	0.46
1:E:60:LEU:HD11	2:F:230:ILE:HG12	1.98	0.46
1:E:88:TYR:OH	2:F:258:GLN:CB	2.64	0.46
1:E:113:TYR:CZ	2:F:282:THR:HG21	2.50	0.46
3:G:60:VAL:HG12	4:H:181:ILE:HD13	1.96	0.46
3:C:17:ARG:O	3:C:21:LEU:HG	2.15	0.46
1:E:88:TYR:OH	2:F:255:VAL:C	2.58	0.46
1:A:86:ARG:O	1:A:89:TRP:N	2.49	0.46
1:E:92:ASN:ND2	2:F:260:LYS:HB3	2.30	0.46
1:E:88:TYR:HE2	2:F:258:GLN:HB2	1.80	0.45
1:A:56:ARG:HH21	1:A:56:ARG:CG	2.26	0.45
1:E:67:ALA:HB1	2:F:236:ASN:HB3	1.97	0.45
1:E:95:MET:HE3	2:F:268:ILE:HD12	1.98	0.45
2:F:247:ALA:O	2:F:251:THR:HG23	2.16	0.45
2:B:199:HIS:HD2	2:B:199:HIS:O	2.00	0.45
3:G:17:ARG:O	3:G:21:LEU:HG	2.16	0.45
1:E:86:ARG:O	1:E:89:TRP:N	2.49	0.45
1:A:29:MET:SD	1:A:32:LEU:HD12	2.57	0.45
4:D:144:ASN:O	4:D:148:GLU:HB2	2.16	0.45
1:E:85:LYS:HE2	2:F:253:LYS:HB2	1.96	0.45
2:F:218:ASP:O	2:F:222:LEU:HD13	2.16	0.45
1:E:64:ASP:HB2	2:F:233:ILE:HD11	2.00	0.44
4:D:192:ILE:O	4:D:196:ASN:HB2	2.17	0.44
3:C:36:VAL:HG11	4:D:156:ILE:HB	2.00	0.44
1:E:77:PHE:CD2	1:E:77:PHE:O	2.70	0.44
3:G:71:MET:CB	4:H:191:ARG:HD2	2.47	0.44
4:H:161:ARG:O	4:H:164:ALA:HB3	2.17	0.44
4:D:162:HIS:O	4:D:163:MET:C	2.61	0.44
2:B:265:LYS:O	2:B:268:ILE:HB	2.17	0.44
1:E:89:TRP:CE3	1:E:90:TRP:HA	2.52	0.44
1:A:88:TYR:CE1	2:B:257:TYR:HB2	2.53	0.44
1:A:99:LEU:HD13	2:B:268:ILE:CG1	2.47	0.44
3:G:67:ILE:HG22	4:H:188:ASN:ND2	2.33	0.44
4:H:162:HIS:O	4:H:163:MET:C	2.61	0.43
4:H:192:ILE:O	4:H:196:ASN:HB2	2.18	0.43
1:A:89:TRP:CE3	1:A:90:TRP:HA	2.52	0.43
1:E:88:TYR:CZ	2:F:258:GLN:N	2.86	0.43
1:E:99:LEU:HD13	2:F:268:ILE:CA	2.48	0.43
1:A:34:GLN:HE22	4:H:163:MET:CA	2.25	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:253:LYS:O	2:B:257:TYR:CD2	2.71	0.43
2:F:256:LYS:HE2	2:F:260:LYS:NZ	2.33	0.43
2:F:273:VAL:HG12	2:F:277:ILE:HD11	2.00	0.43
1:A:80:SER:O	1:A:84:LEU:HG	2.18	0.43
1:A:68:ASP:C	1:A:70:LEU:H	2.27	0.43
3:C:26:LEU:HD23	4:D:142:ARG:HG3	2.01	0.43
1:A:98:ILE:HG22	2:B:268:ILE:HD13	2.01	0.43
2:B:256:LYS:HE2	2:B:260:LYS:NZ	2.33	0.43
2:B:213:HIS:HD2	3:C:35:LEU:CD2	2.31	0.43
1:E:68:ASP:C	1:E:70:LEU:H	2.26	0.43
1:E:113:TYR:OH	2:F:282:THR:HG21	2.19	0.43
1:A:34:GLN:NE2	4:H:163:MET:HA	2.27	0.43
1:A:112:VAL:O	1:A:112:VAL:HG22	2.19	0.43
4:D:159:ASN:H	4:D:159:ASN:HD22	1.65	0.43
1:A:88:TYR:CD1	2:B:254:ALA:HA	2.54	0.42
2:B:213:HIS:C	2:B:213:HIS:ND1	2.77	0.42
1:E:76:GLN:HG2	4:H:196:ASN:HD21	1.84	0.42
2:F:191:ALA:HA	2:F:194:GLU:CD	2.44	0.42
1:A:43:VAL:O	1:A:47:ARG:HG3	2.19	0.42
4:D:159:ASN:HD21	4:H:155:GLY:HA3	1.83	0.42
2:B:191:ALA:HA	2:B:194:GLU:CD	2.45	0.42
2:B:218:ASP:O	2:B:222:LEU:HD13	2.19	0.42
1:A:102:ILE:HD11	2:B:272:CYS:SG	2.59	0.42
2:F:213:HIS:C	2:F:213:HIS:ND1	2.77	0.42
3:G:43:GLY:O	3:G:47:LEU:HG	2.20	0.42
2:B:273:VAL:HG12	2:B:277:ILE:HD11	2.01	0.42
3:G:34:GLN:HA	3:G:37:GLU:OE1	2.20	0.42
1:E:102:ILE:HD12	2:F:268:ILE:HG23	2.01	0.42
2:F:237:VAL:O	2:F:241:VAL:N	2.50	0.42
4:H:159:ASN:H	4:H:159:ASN:HD22	1.65	0.42
1:A:92:ASN:HD21	2:B:264:LYS:HZ1	1.67	0.42
1:A:98:ILE:HB	2:B:268:ILE:CD1	2.50	0.42
1:E:43:VAL:O	1:E:47:ARG:HG3	2.20	0.42
3:G:29:THR:HG21	4:H:149:ASN:HD21	1.84	0.42
1:A:29:MET:HG3	1:A:29:MET:O	2.20	0.42
1:A:77:PHE:O	1:A:77:PHE:CD2	2.72	0.42
1:E:43:VAL:HG13	2:F:215:MET:HG3	2.02	0.42
1:E:80:SER:O	1:E:84:LEU:HG	2.19	0.42
2:B:222:LEU:O	2:B:226:GLN:HG3	2.20	0.41
1:E:29:MET:SD	1:E:32:LEU:HD12	2.60	0.41
2:F:258:GLN:OE1	3:G:81:LEU:CD1	2.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:50:VAL:O	1:E:54:LEU:HD13	2.20	0.41
1:E:56:ARG:NH1	4:H:174:GLN:OE1	2.53	0.41
3:G:53:GLN:O	3:G:57:LEU:HD13	2.20	0.41
1:E:113:TYR:HA	1:E:116:THR:OG1	2.20	0.41
2:F:199:HIS:HD2	2:F:199:HIS:O	2.02	0.41
2:F:222:LEU:O	2:F:226:GLN:HG3	2.20	0.41
1:A:92:ASN:HD21	2:B:260:LYS:HB3	1.86	0.41
1:E:70:LEU:HD12	1:E:70:LEU:HA	1.82	0.41
1:E:112:VAL:O	1:E:112:VAL:HG22	2.19	0.41
1:A:98:ILE:CG2	2:B:268:ILE:HD13	2.50	0.41
4:D:150:LEU:HD12	4:D:150:LEU:HA	1.88	0.41
1:E:43:VAL:HG22	2:F:212:LEU:CD1	2.42	0.41
1:E:34:GLN:O	1:E:38:GLN:HG3	2.21	0.41
1:E:104:ALA:HA	1:E:107:LEU:HB2	2.03	0.41
1:A:50:VAL:O	1:A:54:LEU:HD13	2.21	0.41
3:C:31:ARG:O	3:C:35:LEU:HG	2.21	0.41
1:E:42:VAL:HG11	2:F:212:LEU:HD21	2.02	0.41
1:A:101:VAL:HG23	1:A:102:ILE:N	2.36	0.41
3:C:26:LEU:O	3:C:30:ARG:HG3	2.20	0.41
1:E:95:MET:N	1:E:95:MET:SD	2.94	0.41
4:D:140:MET:O	4:D:144:ASN:HB2	2.21	0.41
4:D:156:ILE:O	4:D:159:ASN:HB2	2.21	0.41
1:E:43:VAL:HG22	2:F:212:LEU:HB2	2.03	0.41
3:C:34:GLN:HA	3:C:37:GLU:OE1	2.22	0.40
3:C:53:GLN:O	3:C:57:LEU:HD13	2.20	0.40
4:H:156:ILE:O	4:H:159:ASN:HB2	2.22	0.40
1:A:34:GLN:O	1:A:38:GLN:HG3	2.21	0.40
3:G:50:LEU:HD23	3:G:50:LEU:HA	1.89	0.40
1:A:88:TYR:CZ	2:B:258:GLN:HB2	2.55	0.40
3:G:26:LEU:O	3:G:30:ARG:HG3	2.21	0.40
4:H:157:ILE:O	4:H:157:ILE:HG22	2.22	0.40
1:E:88:TYR:HH	2:F:258:GLN:CB	2.31	0.40
2:F:253:LYS:O	2:F:257:TYR:CD2	2.74	0.40
4:H:177:GLN:O	4:H:180:ARG:HB3	2.22	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	89/91 (98%)	83 (93%)	5 (6%)	1 (1%)	11	45
1	E	89/91 (98%)	84 (94%)	4 (4%)	1 (1%)	11	45
2	B	96/109 (88%)	88 (92%)	7 (7%)	1 (1%)	12	47
2	F	96/109 (88%)	88 (92%)	7 (7%)	1 (1%)	12	47
3	C	73/80 (91%)	72 (99%)	1 (1%)	0	100	100
3	G	73/80 (91%)	72 (99%)	1 (1%)	0	100	100
4	D	61/68 (90%)	59 (97%)	2 (3%)	0	100	100
4	H	61/68 (90%)	59 (97%)	2 (3%)	0	100	100
All	All	638/696 (92%)	605 (95%)	29 (4%)	4 (1%)	21	58

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	255	VAL
1	A	86	ARG
1	E	86	ARG
2	F	255	VAL

### 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	80/80 (100%)	75 (94%)	5 (6%)	16	38

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	80/80 (100%)	74 (92%)	6 (8%)	12	33
2	B	87/96 (91%)	82 (94%)	5 (6%)	18	40
2	F	87/96 (91%)	82 (94%)	5 (6%)	18	40
3	C	67/71 (94%)	66 (98%)	1 (2%)	57	71
3	G	67/71 (94%)	66 (98%)	1 (2%)	57	71
4	D	55/58 (95%)	52 (94%)	3 (6%)	19	42
4	H	55/58 (95%)	53 (96%)	2 (4%)	31	52
All	All	578/610 (95%)	550 (95%)	28 (5%)	23	45

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	GLN
1	A	102	ILE
1	A	107	LEU
1	A	110	ILE
1	A	114	PHE
2	B	242	ASP
2	B	257	TYR
2	B	270	ILE
2	B	272	CYS
2	B	282	THR
3	C	26	LEU
4	D	178	ILE
4	D	181	ILE
4	D	192	ILE
1	E	38	GLN
1	E	102	ILE
1	E	107	LEU
1	E	110	ILE
1	E	114	PHE
1	E	116	THR
2	F	242	ASP
2	F	257	TYR
2	F	270	ILE
2	F	272	CYS
2	F	282	THR
3	G	26	LEU
4	H	181	ILE
4	H	192	ILE

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	34	GLN
1	A	92	ASN
4	D	159	ASN
4	D	162	HIS
1	E	34	GLN
2	F	199	HIS
2	F	213	HIS
3	G	56	GLN
3	G	65	ASN
4	H	149	ASN
4	H	162	HIS
4	H	175	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 5.7 Other polymers [i](#)

There are no such residues in this entry.

### 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ > 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	91/91 (100%)	1.27	19 (20%) 2 7	175, 264, 308, 361	0
1	E	91/91 (100%)	1.43	23 (25%) 1 5	187, 264, 353, 387	0
2	B	98/109 (89%)	1.05	21 (21%) 2 7	181, 286, 355, 422	0
2	F	98/109 (89%)	1.20	24 (24%) 2 6	191, 281, 364, 503	0
3	C	75/80 (93%)	1.34	19 (25%) 1 5	187, 314, 388, 395	0
3	G	75/80 (93%)	1.07	12 (16%) 5 10	205, 304, 377, 441	0
4	D	63/68 (92%)	1.38	16 (25%) 1 5	222, 295, 380, 411	0
4	H	63/68 (92%)	1.30	15 (23%) 2 6	193, 290, 362, 426	0
All	All	654/696 (93%)	1.25	149 (22%) 2 6	175, 284, 373, 503	0

All (149) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	114	PHE	9.4
1	A	116	THR	9.4
4	D	138	SER	8.9
3	G	75	GLU	8.8
4	H	183	GLU	8.7
1	E	33	GLN	8.1
1	E	114	PHE	8.0
3	C	80	ASP	6.5
1	A	115	SER	6.2
3	C	81	LEU	5.9
3	G	71	MET	5.8
1	E	30	ARG	5.4
1	E	87	LYS	5.4
1	E	27	SER	5.2
4	D	140	MET	5.2
1	A	113	TYR	5.1

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
2	F	228	GLU	5.1
3	C	73	GLU	4.8
4	H	191	ARG	4.8
2	F	248	VAL	4.5
2	B	260	LYS	4.4
3	C	55	GLU	4.4
4	H	198	ARG	4.4
2	F	245	GLU	4.4
2	B	259	SER	4.3
2	F	221	MET	4.3
1	E	96	MET	4.3
3	G	82	GLY	4.2
2	F	247	ALA	4.2
4	D	183	GLU	4.2
1	A	26	GLY	4.2
3	C	82	GLY	4.2
1	A	90	TRP	4.1
2	B	286	ILE	4.0
1	A	87	LYS	4.0
4	H	200	THR	4.0
3	C	77	ASN	3.8
3	G	58	ASP	3.7
1	E	69	ALA	3.7
2	B	231	ASP	3.7
2	F	286	ILE	3.7
4	D	139	HIS	3.6
3	G	74	ALA	3.6
3	C	8	ARG	3.6
1	E	42	VAL	3.6
2	F	262	ARG	3.6
2	B	254	ALA	3.6
1	E	46	MET	3.6
1	E	111	ILE	3.5
2	F	251	THR	3.5
4	D	200	THR	3.5
1	A	27	SER	3.5
2	B	238	GLU	3.5
2	F	250	ASP	3.4
2	B	201	GLU	3.4
1	E	90	TRP	3.4
1	E	103	CYS	3.3
1	A	36	GLN	3.2

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
3	G	64	MET	3.2
1	E	47	ARG	3.2
4	H	157	ILE	3.1
4	D	171	ILE	3.1
2	B	265	LYS	3.1
1	E	72	ALA	3.1
4	H	139	HIS	3.0
2	F	210	ARG	3.0
4	H	186	ASP	3.0
1	E	51	ASP	3.0
4	D	189	LYS	2.9
1	A	112	VAL	2.9
2	F	284	GLY	2.9
1	A	49	ASN	2.9
2	B	263	ARG	2.9
2	F	244	VAL	2.9
4	D	184	LYS	2.9
4	D	199	ALA	2.9
2	F	211	GLU	2.9
2	B	205	LEU	2.8
4	H	160	LEU	2.8
2	B	264	LYS	2.8
4	D	152	GLN	2.8
2	B	272	CYS	2.7
1	E	85	LYS	2.7
3	G	11	LEU	2.7
1	A	50	VAL	2.7
3	C	21	LEU	2.6
1	A	55	GLU	2.6
2	F	196	GLU	2.6
2	F	254	ALA	2.6
1	E	31	ARG	2.6
1	A	111	ILE	2.6
1	A	58	GLN	2.6
1	A	53	VAL	2.6
4	D	143	GLU	2.6
1	E	115	SER	2.5
4	H	140	MET	2.5
1	A	96	MET	2.5
2	F	263	ARG	2.5
4	H	153	VAL	2.5
2	F	215	MET	2.5

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
2	F	267	MET	2.5
3	C	66	HIS	2.5
3	C	44	ILE	2.5
1	E	91	LYS	2.5
4	H	193	ASP	2.5
1	E	43	VAL	2.5
4	H	161	ARG	2.5
1	E	26	GLY	2.5
4	D	197	GLN	2.4
2	F	283	ILE	2.4
4	H	189	LYS	2.4
4	D	178	ILE	2.4
3	C	79	LYS	2.4
4	D	179	ASP	2.4
2	F	201	GLU	2.4
4	H	187	SER	2.4
2	F	265	LYS	2.4
4	D	167	MET	2.4
3	C	69	GLN	2.3
2	B	189	LYS	2.3
2	B	262	ARG	2.3
2	B	229	MET	2.3
3	C	38	GLU	2.3
1	E	116	THR	2.3
3	C	76	LYS	2.3
2	B	256	LYS	2.3
3	C	78	LEU	2.2
2	F	224	GLU	2.2
2	B	235	TYR	2.2
3	G	70	ASP	2.2
4	D	176	ARG	2.2
2	B	191	ALA	2.1
3	C	18	ALA	2.1
3	G	54	GLY	2.1
1	A	83	LYS	2.1
2	B	257	TYR	2.1
3	G	27	GLU	2.1
2	B	221	MET	2.1
3	G	73	GLU	2.1
3	C	50	LEU	2.1
3	C	45	ARG	2.1
3	G	67	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	38	GLN	2.1
2	B	210	ARG	2.0
2	F	229	MET	2.0
1	A	52	LYS	2.0
2	F	189	LYS	2.0
4	H	188	ASN	2.0
3	C	24	GLU	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.