



# Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 07:09 am BST

PDB ID : 2OR9  
Title : The structure of the anti-c-myc antibody 9E10 Fab fragment/epitope peptide complex reveals a novel binding mode dominated by the heavy chain hyper-variable loops  
Authors : Krauss, N.; Scheerer, P.; Hoehne, W.  
Deposited on : 2007-02-02  
Resolution : 2.70 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.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

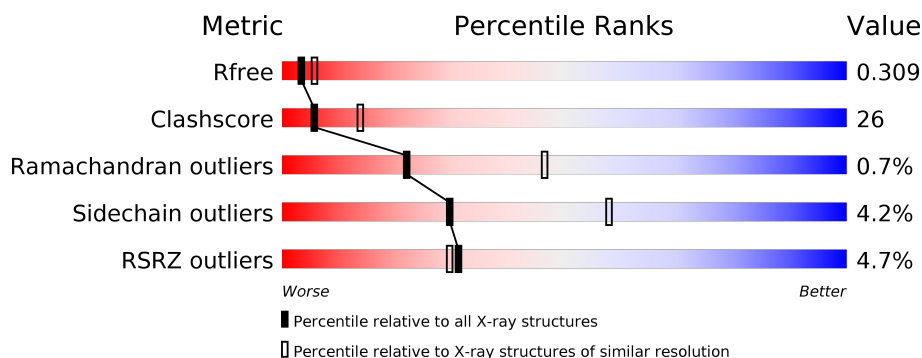
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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))
$R_{free}$	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on 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	L	218	<div> <div>3%</div> <div> <div></div> <div>61%</div> <div>36%</div> <div>..</div> </div> </div>
1	M	218	<div> <div>%</div> <div> <div></div> <div>63%</div> <div>33%</div> <div>.</div> </div> </div>
2	H	228	<div> <div>13%</div> <div> <div></div> <div>50%</div> <div>44%</div> <div>..</div> </div> </div>
2	I	228	<div> <div>2%</div> <div> <div></div> <div>58%</div> <div>37%</div> <div>.</div> </div> </div>
3	P	11	<div> <div></div> <div> <div></div> <div>55%</div> <div>45%</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7182 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 Monoclonal anti-c-myc antibody 9E10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	216	Total	C	N	O	S	0	0	0
			1671	1045	284	335	7			
1	M	218	Total	C	N	O	S	0	0	0
			1687	1053	286	340	8			

- Molecule 2 is a protein called Monoclonal anti-c-myc antibody 9E10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	221	Total	C	N	O	S	0	0	0
			1698	1079	278	332	9			
2	I	228	Total	C	N	O	S	0	0	0
			1752	1108	290	345	9			

- Molecule 3 is a protein called synthetic epitope peptide of 9E10.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	11	Total	C	N	O	0	0	0
			92	55	14	23			

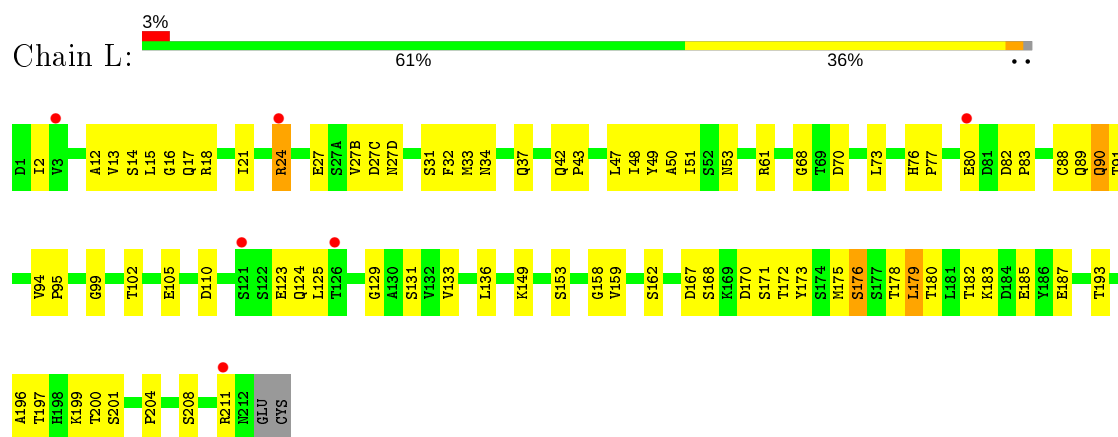
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	48	Total	O	0	0
			48	48		
4	H	35	Total	O	0	0
			35	35		
4	M	91	Total	O	0	0
			91	91		
4	I	94	Total	O	0	0
			94	94		
4	P	14	Total	O	0	0
			14	14		

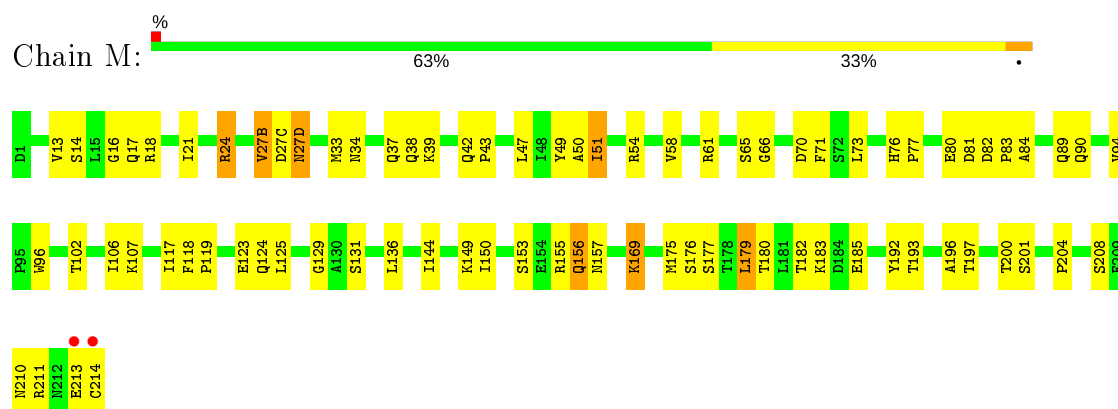
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

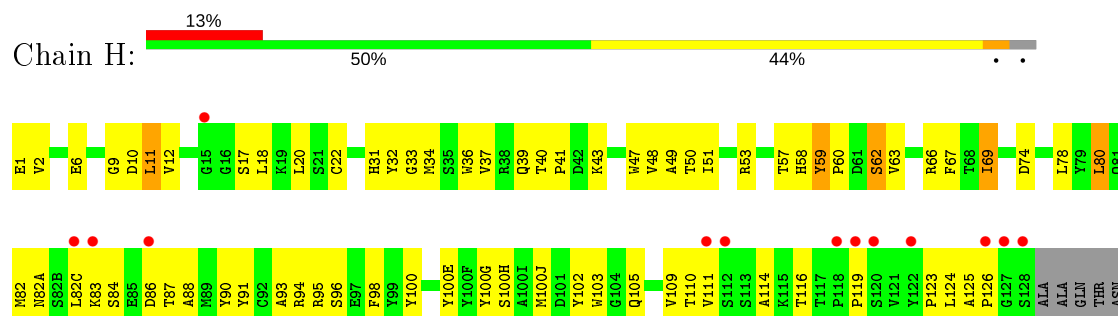
- Molecule 1: Monoclonal anti-c-myc antibody 9E10

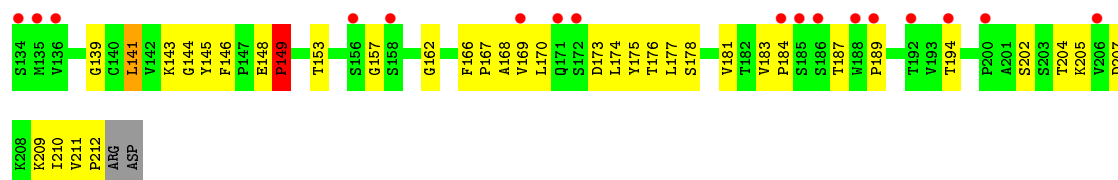


- Molecule 1: Monoclonal anti-c-myc antibody 9E10

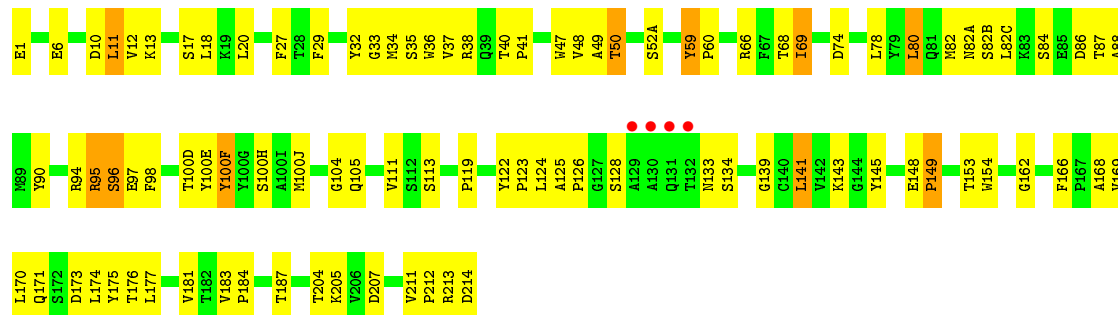


- Molecule 2: Monoclonal anti-c-myc antibody 9E10





- Molecule 2: Monoclonal anti-c-myc antibody 9E10



- Molecule 3: synthetic epitope peptide of 9E10



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	76.07Å 111.88Å 134.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.70 19.36 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.4 (20.00-2.70) 99.4 (19.36-2.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.94 (at 2.70Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.252 , 0.310 0.252 , 0.309	Depositor DCC
$R_{free}$ test set	1626 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	57.0	Xtriage
Anisotropy	0.460	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 49.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	7182	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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

### 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	L	0.41	0/1712	0.65	0/2325
1	M	0.44	0/1728	0.68	0/2345
2	H	0.40	0/1747	0.64	0/2384
2	I	0.41	0/1802	0.67	0/2459
3	P	0.49	0/91	0.60	0/119
All	All	0.41	0/7080	0.66	0/9632

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	1671	0	1598	74	0
1	M	1687	0	1609	75	0
2	H	1698	0	1633	111	0
2	I	1752	0	1682	96	0
3	P	92	0	89	7	0
4	H	35	0	0	5	0
4	I	94	0	0	3	0
4	L	48	0	0	0	0
4	M	91	0	0	0	0
4	P	14	0	0	0	0
All	All	7182	0	6611	345	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 26.

All (345) 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)
2:H:87:THR:HG22	2:H:111:VAL:H	0.97	1.12
2:H:87:THR:CG2	2:H:111:VAL:H	1.65	1.10
2:H:87:THR:HG22	2:H:111:VAL:N	1.72	1.02
2:H:49:ALA:HB1	2:H:69:ILE:HD13	1.44	1.00
2:I:87:THR:HG22	2:I:111:VAL:H	1.33	0.94
2:I:34:MET:HE1	2:I:94:ARG:HA	1.52	0.92
1:L:34:ASN:HD22	1:L:49:TYR:HA	1.36	0.91
2:I:49:ALA:HB1	2:I:69:ILE:HD13	1.53	0.90
1:M:51:ILE:HD11	1:M:65:SER:C	1.91	0.89
2:I:119:PRO:HB3	2:I:145:TYR:HB3	1.58	0.84
2:I:34:MET:HE3	2:I:94:ARG:HG3	1.59	0.84
2:H:119:PRO:HB3	2:H:145:TYR:HB3	1.60	0.82
1:L:24:ARG:HH11	1:L:24:ARG:HB3	1.43	0.82
1:M:24:ARG:HB3	1:M:24:ARG:NH1	1.96	0.80
2:H:33:GLY:O	2:H:95:ARG:HB2	1.83	0.78
1:L:33:MET:HE3	1:L:34:ASN:H	1.47	0.77
1:M:24:ARG:HH11	1:M:24:ARG:HB3	1.49	0.76
2:H:84:SER:HA	2:H:111:VAL:HB	1.68	0.76
2:H:40:THR:HB	2:H:41:PRO:HD2	1.68	0.75
2:I:87:THR:CG2	2:I:111:VAL:H	2.00	0.74
2:I:96:SER:O	2:I:100(H):SER:HA	1.88	0.74
2:I:87:THR:HG22	2:I:111:VAL:N	2.04	0.73
1:M:16:GLY:HA2	1:M:77:PRO:HB2	1.69	0.73
2:H:34:MET:HE2	2:H:94:ARG:HA	1.69	0.72
2:I:211:VAL:HG13	2:I:212:PRO:HD2	1.71	0.72
1:L:18:ARG:HB2	1:L:76:HIS:HD2	1.53	0.72
1:L:94:VAL:HG13	2:H:47:TRP:HZ3	1.55	0.71
1:M:193:THR:HG22	1:M:208:SER:CB	2.19	0.71
1:L:133:VAL:HG21	2:H:124:LEU:HD21	1.73	0.71
1:M:210:ASN:HB2	1:M:213:GLU:HB2	1.72	0.70
2:I:40:THR:HB	2:I:41:PRO:HD2	1.73	0.70
1:L:94:VAL:HG13	2:H:47:TRP:CZ3	2.27	0.69
1:M:156:GLN:HA	1:M:156:GLN:HE21	1.57	0.69
2:I:47:TRP:HE1	2:I:50:THR:CG2	2.05	0.69
2:H:34:MET:HE1	2:H:94:ARG:HG3	1.73	0.68
1:L:21:ILE:HG12	1:L:102:THR:HG21	1.73	0.68
2:H:20:LEU:HG	2:H:82:MET:CE	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:34:MET:CE	2:I:94:ARG:HA	2.23	0.68
1:L:33:MET:CE	1:L:88:CYS:HB2	2.24	0.68
1:L:33:MET:HE3	1:L:34:ASN:N	2.09	0.67
2:I:47:TRP:HE1	2:I:50:THR:HG22	1.59	0.67
1:L:193:THR:HG22	1:L:208:SER:CB	2.25	0.66
2:H:189:PRO:HD2	4:H:239:HOH:O	1.96	0.66
2:I:12:VAL:O	2:I:111:VAL:HA	1.95	0.66
1:M:150:ILE:HG23	1:M:192:TYR:CE2	2.31	0.66
1:L:149:LYS:HB2	1:L:193:THR:OG1	1.96	0.66
2:H:20:LEU:HG	2:H:82:MET:HE2	1.78	0.65
2:I:47:TRP:NE1	2:I:50:THR:HG22	2.11	0.65
1:L:197:THR:HG22	1:L:204:PRO:HG3	1.78	0.65
2:I:47:TRP:NE1	2:I:50:THR:CG2	2.60	0.65
1:M:193:THR:HG22	1:M:208:SER:HB3	1.78	0.65
1:M:18:ARG:HB2	1:M:76:HIS:HD2	1.61	0.64
1:L:33:MET:HE2	1:L:88:CYS:HB2	1.79	0.64
1:L:80:GLU:O	1:L:83:PRO:HD2	1.97	0.64
2:H:84:SER:O	2:H:87:THR:HG23	1.97	0.64
2:H:82:MET:HB3	2:H:82(C):LEU:HD21	1.79	0.64
2:I:96:SER:HB3	2:I:98:PHE:HE1	1.62	0.63
1:M:197:THR:HG22	1:M:204:PRO:HG3	1.80	0.63
2:H:34:MET:HB3	2:H:78:LEU:HD22	1.81	0.63
1:L:24:ARG:HG3	1:L:70:ASP:OD1	1.99	0.63
2:H:148:GLU:OE1	2:H:149:PRO:HA	1.98	0.63
1:L:14:SER:HB2	1:L:17:GLN:HG3	1.82	0.62
1:L:31:SER:OG	1:L:51:ILE:HD11	1.99	0.62
1:M:131:SER:OG	1:M:180:THR:HG22	1.98	0.62
1:M:14:SER:HB2	1:M:17:GLN:HG3	1.81	0.62
1:L:94:VAL:HG13	1:L:95:PRO:HA	1.82	0.62
2:I:82:MET:HB3	2:I:82(C):LEU:HD21	1.82	0.62
2:H:6:GLU:OE2	2:H:91:TYR:HA	2.00	0.62
2:I:10:ASP:OD1	2:I:11:LEU:N	2.33	0.62
2:H:143:LYS:HG3	2:H:176:THR:CG2	2.29	0.62
2:H:211:VAL:HG13	2:H:212:PRO:HD2	1.82	0.61
2:I:124:LEU:HB2	2:I:139:GLY:O	2.01	0.61
1:M:51:ILE:HD12	1:M:71:PHE:HD2	1.64	0.61
1:M:123:GLU:N	1:M:123:GLU:OE1	2.34	0.61
2:I:94:ARG:O	2:I:100(J):MET:HA	2.00	0.61
1:M:211:ARG:HG2	1:M:211:ARG:HH11	1.66	0.61
2:H:39:GLN:O	2:H:88:ALA:HB1	2.01	0.61
2:H:177:LEU:HD12	2:H:177:LEU:C	2.21	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:177:LEU:C	2:I:177:LEU:HD12	2.22	0.60
1:L:34:ASN:ND2	1:L:49:TYR:HA	2.14	0.60
2:H:37:VAL:HG21	2:H:100(J):MET:HE1	1.84	0.59
1:M:214:CYS:SG	2:I:128:SER:HA	2.42	0.59
2:H:125:ALA:HB1	2:H:126:PRO:HD2	1.84	0.59
2:I:84:SER:O	2:I:87:THR:HG23	2.01	0.59
2:I:148:GLU:OE1	2:I:149:PRO:HA	2.03	0.59
2:H:87:THR:HA	2:H:109:VAL:O	2.02	0.59
1:M:94:VAL:HG23	3:P:1:GLU:OE2	2.03	0.59
2:H:184:PRO:HG2	2:H:187:THR:HG23	1.84	0.59
1:M:27(B):VAL:HG22	1:M:90:GLN:HG2	1.85	0.59
1:M:33:MET:HE3	1:M:89:GLN:O	2.03	0.58
1:L:89:GLN:HG2	1:L:90:GLN:N	2.17	0.58
2:H:124:LEU:HD11	2:H:141:LEU:HB2	1.84	0.58
2:H:51:ILE:HG13	2:H:57:THR:HG22	1.85	0.58
2:H:100:TYR:HB3	2:H:100(E):TYR:CE1	2.39	0.58
2:H:170:LEU:HD13	2:H:175:TYR:CE1	2.39	0.58
2:I:125:ALA:HB1	2:I:126:PRO:HD2	1.85	0.58
1:L:123:GLU:OE1	1:L:123:GLU:N	2.36	0.58
1:L:48:ILE:HA	1:L:53:ASN:O	2.04	0.58
2:H:58:HIS:CD2	2:H:59:TYR:H	2.22	0.57
1:L:193:THR:HG22	1:L:208:SER:HB3	1.86	0.57
1:M:34:ASN:OD1	1:M:49:TYR:HA	2.04	0.57
2:H:37:VAL:HG22	2:H:47:TRP:HA	1.86	0.57
1:L:2:ILE:HG12	1:L:27:GLU:OE2	2.04	0.57
2:I:170:LEU:HD13	2:I:175:TYR:CE1	2.39	0.57
1:M:136:LEU:HD21	1:M:196:ALA:HB2	1.87	0.57
1:M:50:ALA:O	1:M:51:ILE:HG22	2.05	0.57
1:M:51:ILE:HD11	1:M:66:GLY:N	2.19	0.57
2:H:168:ALA:HA	2:H:177:LEU:HB3	1.86	0.57
2:I:143:LYS:HG3	2:I:176:THR:CG2	2.35	0.56
2:H:50:THR:C	2:H:69:ILE:HD12	2.26	0.56
2:H:210:ILE:HG13	4:H:234:HOH:O	2.05	0.56
2:I:119:PRO:CB	2:I:145:TYR:HB3	2.32	0.56
2:I:36:TRP:CE2	2:I:80:LEU:HB2	2.41	0.56
2:H:12:VAL:O	2:H:111:VAL:HA	2.06	0.56
2:H:63:VAL:HG11	2:H:67:PHE:CE2	2.40	0.56
2:I:34:MET:HB3	2:I:78:LEU:HD22	1.88	0.56
2:H:10:ASP:OD1	2:H:11:LEU:N	2.38	0.55
2:H:169:VAL:O	2:H:175:TYR:HA	2.06	0.55
1:L:131:SER:OG	1:L:180:THR:HG22	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:211:ARG:HH11	1:L:211:ARG:HG2	1.70	0.55
1:M:124:GLN:HE22	1:M:131:SER:CB	2.20	0.55
2:H:31:HIS:CD2	3:P:8:GLU:HG3	2.42	0.54
2:H:96:SER:HB3	3:P:4:LEU:HB3	1.88	0.54
2:H:183:VAL:HB	2:H:184:PRO:HD2	1.88	0.54
2:H:31:HIS:CG	3:P:8:GLU:HG3	2.43	0.54
1:M:117:ILE:HD13	1:M:208:SER:HA	1.89	0.54
2:H:168:ALA:HB2	2:H:177:LEU:HD23	1.90	0.54
1:M:182:THR:HG23	1:M:185:GLU:OE2	2.08	0.54
1:L:170:ASP:O	1:L:172:THR:HG23	2.07	0.53
2:I:184:PRO:HG2	2:I:187:THR:HG23	1.90	0.53
1:M:27(C):ASP:O	1:M:27(D):ASN:HB2	2.07	0.53
2:H:116:THR:HG22	2:H:202:SER:HB3	1.90	0.53
1:M:38:GLN:O	1:M:84:ALA:HB1	2.09	0.53
2:H:119:PRO:HB3	2:H:145:TYR:CB	2.37	0.53
1:L:162:SER:OG	2:H:166:PHE:HB3	2.08	0.53
1:L:110:ASP:OD2	1:L:199:LYS:HE3	2.07	0.53
1:L:105:GLU:CD	1:L:173:TYR:HH	2.12	0.53
1:M:96:TRP:CZ2	2:I:100(H):SER:HB2	2.44	0.53
1:L:15:LEU:HD21	1:L:80:GLU:CD	2.28	0.53
1:L:27(B):VAL:HG12	1:L:27(B):VAL:O	2.08	0.53
1:L:124:GLN:HE22	1:L:131:SER:CB	2.23	0.52
1:L:199:LYS:HD2	1:L:199:LYS:O	2.09	0.52
2:I:169:VAL:O	2:I:175:TYR:HA	2.10	0.52
2:H:53:ARG:HD3	4:H:246:HOH:O	2.09	0.52
2:I:168:ALA:HA	2:I:177:LEU:HB3	1.91	0.52
2:H:143:LYS:HG3	2:H:176:THR:HG21	1.90	0.52
2:H:119:PRO:CB	2:H:145:TYR:HB3	2.35	0.52
2:H:60:PRO:HA	4:H:245:HOH:O	2.09	0.52
2:I:32:TYR:O	2:I:52(A):SER:HA	2.08	0.52
1:L:80:GLU:O	1:L:83:PRO:CD	2.58	0.52
1:M:150:ILE:HD11	1:M:155:ARG:HG2	1.91	0.52
2:H:173:ASP:O	2:H:174:LEU:HG	2.10	0.52
1:L:61:ARG:NH2	1:L:82:ASP:OD2	2.43	0.52
1:L:82:ASP:N	1:L:83:PRO:HD2	2.25	0.52
2:I:183:VAL:HB	2:I:184:PRO:HD2	1.91	0.52
1:M:18:ARG:HB2	1:M:76:HIS:CD2	2.42	0.52
1:L:27(B):VAL:HG12	1:L:32:PHE:H	1.74	0.51
1:L:50:ALA:HB3	1:L:53:ASN:HD22	1.75	0.51
2:H:124:LEU:HB2	2:H:139:GLY:O	2.10	0.51
2:I:211:VAL:CG1	2:I:212:PRO:HD2	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:87:THR:HG22	2:H:110:THR:HA	1.92	0.51
2:I:34:MET:CE	2:I:94:ARG:HG3	2.36	0.51
1:M:33:MET:HE3	1:M:89:GLN:C	2.31	0.51
1:L:105:GLU:OE2	1:L:173:TYR:OH	2.25	0.50
2:I:33:GLY:O	2:I:34:MET:HE1	2.12	0.50
1:L:13:VAL:CG1	1:L:17:GLN:HB2	2.41	0.50
1:L:18:ARG:HB2	1:L:76:HIS:CD2	2.42	0.50
2:I:124:LEU:HD11	2:I:141:LEU:HB2	1.92	0.50
1:L:136:LEU:HD21	1:L:196:ALA:HB2	1.94	0.50
1:M:156:GLN:HA	1:M:156:GLN:NE2	2.26	0.50
1:M:80:GLU:O	1:M:83:PRO:HD2	2.11	0.50
2:H:17:SER:HB2	2:H:82(A):ASN:HD22	1.77	0.50
2:I:84:SER:HA	2:I:111:VAL:HB	1.93	0.50
1:L:33:MET:HE3	1:L:89:GLN:O	2.11	0.50
1:M:169:LYS:NZ	1:M:169:LYS:HB3	2.27	0.50
2:H:34:MET:CE	2:H:94:ARG:HA	2.40	0.50
2:I:50:THR:O	2:I:69:ILE:CD1	2.61	0.49
2:H:103:TRP:N	2:H:103:TRP:CD1	2.80	0.49
2:H:123:PRO:O	2:H:124:LEU:HD12	2.13	0.49
2:I:13:LYS:HG2	2:I:113:SER:HA	1.94	0.49
1:M:156:GLN:CA	1:M:156:GLN:HE21	2.19	0.49
2:I:105:GLN:N	2:I:105:GLN:OE1	2.44	0.49
2:I:6:GLU:H	2:I:105:GLN:HE22	1.60	0.49
1:M:21:ILE:HG12	1:M:102:THR:HG21	1.95	0.49
1:M:51:ILE:HD12	1:M:71:PHE:CD2	2.47	0.49
2:I:27:PHE:CE1	2:I:29:PHE:HA	2.48	0.49
2:I:33:GLY:O	2:I:34:MET:CE	2.61	0.49
1:M:61:ARG:NH2	1:M:82:ASP:OD1	2.40	0.49
1:M:42:GLN:HB3	1:M:43:PRO:HD2	1.95	0.48
2:I:47:TRP:NE1	2:I:50:THR:HG23	2.28	0.48
1:M:54:ARG:HD2	1:M:58:VAL:O	2.14	0.48
2:H:83:LYS:O	2:H:111:VAL:HG21	2.13	0.48
1:M:24:ARG:HG3	1:M:70:ASP:OD1	2.13	0.48
1:M:61:ARG:NH2	1:M:82:ASP:OD2	2.46	0.48
2:H:48:VAL:O	2:H:60:PRO:HD2	2.14	0.48
2:H:66:ARG:NH2	2:H:86:ASP:OD2	2.47	0.48
2:I:20:LEU:HG	2:I:82:MET:CE	2.42	0.48
2:I:97:GLU:HG3	2:I:100(H):SER:HB3	1.95	0.48
1:L:133:VAL:HG11	2:H:124:LEU:HD23	1.96	0.48
2:H:9:GLY:O	2:H:10:ASP:HB2	2.13	0.48
1:M:82:ASP:N	1:M:83:PRO:HD2	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:36:TRP:CE2	2:H:80:LEU:HB2	2.48	0.48
2:H:32:TYR:HD2	2:H:95:ARG:O	1.96	0.47
1:M:149:LYS:HA	1:M:153:SER:O	2.14	0.47
3:P:10:LEU:HD23	3:P:10:LEU:C	2.34	0.47
2:I:166:PHE:CD1	2:I:166:PHE:N	2.82	0.47
2:H:204:THR:HG22	2:H:205:LYS:N	2.29	0.47
2:I:173:ASP:O	2:I:174:LEU:HG	2.14	0.47
1:L:175:MET:HG2	1:L:176:SER:N	2.29	0.47
1:M:33:MET:CE	1:M:89:GLN:C	2.83	0.47
2:H:50:THR:H	2:H:69:ILE:CD1	2.28	0.47
1:M:13:VAL:CG1	1:M:17:GLN:HB2	2.44	0.47
1:M:200:THR:O	1:M:201:SER:HB2	2.14	0.47
2:I:119:PRO:HB3	2:I:145:TYR:CB	2.37	0.47
2:I:143:LYS:HG3	2:I:176:THR:HG21	1.97	0.47
2:I:36:TRP:CD1	2:I:69:ILE:HG22	2.49	0.47
2:H:6:GLU:OE1	2:H:90:TYR:O	2.32	0.47
3:P:7:GLU:O	3:P:10:LEU:HB3	2.14	0.47
2:I:100(E):TYR:O	2:I:100(F):TYR:C	2.54	0.47
2:H:114:ALA:HB3	2:H:146:PHE:CE2	2.50	0.47
1:L:61:ARG:NH2	1:L:82:ASP:OD1	2.45	0.47
1:M:193:THR:HG22	1:M:208:SER:HB2	1.95	0.47
2:I:33:GLY:O	2:I:95:ARG:N	2.39	0.46
1:L:50:ALA:O	1:L:51:ILE:HB	2.15	0.46
1:L:27(C):ASP:O	1:L:27(D):ASN:HB2	2.15	0.46
2:H:37:VAL:CG2	2:H:100(J):MET:HE1	2.46	0.46
2:I:66:ARG:HG2	2:I:82(B):SER:HB2	1.97	0.46
1:M:136:LEU:HD23	1:M:144:ILE:CD1	2.46	0.46
2:I:37:VAL:HG22	2:I:47:TRP:HA	1.97	0.46
2:I:100(D):THR:HG22	2:I:100(E):TYR:O	2.16	0.45
2:I:168:ALA:HB2	2:I:177:LEU:HD23	1.98	0.45
2:I:66:ARG:NH2	2:I:86:ASP:OD2	2.49	0.45
1:L:158:GLY:O	1:L:179:LEU:HA	2.17	0.45
2:H:143:LYS:HG3	2:H:176:THR:HG22	1.99	0.45
2:I:171:GLN:HG2	4:I:229:HOH:O	2.17	0.45
1:L:32:PHE:HA	2:H:100(G):TYR:OH	2.16	0.45
2:H:11:LEU:HD23	2:H:12:VAL:H	1.82	0.45
2:H:59:TYR:CE2	2:H:69:ILE:HG12	2.51	0.45
1:L:162:SER:OG	2:H:167:PRO:HD2	2.17	0.45
2:I:204:THR:HG22	2:I:205:LYS:N	2.31	0.45
1:L:200:THR:O	1:L:201:SER:HB2	2.17	0.45
2:H:60:PRO:HG2	2:H:62:SER:HB3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:38:ARG:NH1	2:I:86:ASP:HA	2.32	0.45
1:L:182:THR:HG23	1:L:185:GLU:OE2	2.16	0.45
1:L:88:CYS:O	1:L:99:GLY:N	2.47	0.45
2:H:98:PHE:HB2	2:H:100(G):TYR:HB2	1.98	0.45
2:H:94:ARG:O	2:H:100(J):MET:HA	2.16	0.45
2:H:178:SER:HB3	4:H:247:HOH:O	2.17	0.45
2:I:40:THR:CB	2:I:41:PRO:HD2	2.44	0.45
2:I:47:TRP:HE1	2:I:50:THR:HG23	1.81	0.44
1:L:33:MET:HE1	1:L:88:CYS:HB2	1.95	0.44
2:H:166:PHE:N	2:H:166:PHE:CD1	2.85	0.44
1:M:27(B):VAL:HG22	1:M:90:GLN:CG	2.47	0.44
2:H:31:HIS:HA	3:P:8:GLU:CG	2.47	0.44
2:H:177:LEU:HD12	2:H:177:LEU:O	2.18	0.44
2:I:6:GLU:OE1	2:I:90:TYR:O	2.34	0.44
1:M:37:GLN:HB2	1:M:47:LEU:HD11	1.99	0.44
2:H:40:THR:CB	2:H:41:PRO:HD2	2.42	0.44
2:I:35:SER:HA	2:I:50:THR:HA	2.00	0.44
1:M:211:ARG:HG2	1:M:211:ARG:NH1	2.32	0.44
1:M:80:GLU:O	1:M:83:PRO:CD	2.65	0.44
1:M:118:PHE:HA	1:M:119:PRO:HD3	1.82	0.44
1:L:2:ILE:HG23	1:L:27:GLU:H	1.82	0.44
2:H:162:GLY:O	2:H:181:VAL:HA	2.18	0.44
2:H:22:CYS:HB3	2:H:78:LEU:HB3	1.99	0.44
2:H:83:LYS:O	2:H:111:VAL:HG11	2.18	0.44
1:M:43:PRO:HB3	2:I:104:GLY:O	2.17	0.44
2:I:48:VAL:O	2:I:60:PRO:HD2	2.17	0.44
2:I:96:SER:HB3	2:I:98:PHE:CE1	2.49	0.44
2:H:211:VAL:CG1	2:H:212:PRO:HD2	2.48	0.44
1:M:125:LEU:O	1:M:183:LYS:HD2	2.17	0.44
1:M:39:LYS:NZ	1:M:81:ASP:OD1	2.51	0.44
2:H:12:VAL:HG11	2:H:82(C):LEU:HD12	2.00	0.43
2:I:50:THR:C	2:I:69:ILE:HD12	2.39	0.43
2:H:144:GLY:C	2:H:174:LEU:HD22	2.38	0.43
1:M:51:ILE:CD1	1:M:66:GLY:O	2.66	0.43
1:M:51:ILE:HD13	1:M:66:GLY:O	2.18	0.43
2:H:20:LEU:HG	2:H:82:MET:HE1	1.97	0.43
2:I:13:LYS:HE3	2:I:113:SER:O	2.18	0.43
2:I:87:THR:O	2:I:88:ALA:HB2	2.18	0.43
2:H:40:THR:HB	2:H:41:PRO:CD	2.44	0.43
2:I:153:THR:HG22	2:I:154:TRP:N	2.34	0.43
1:L:159:VAL:HA	1:L:178:THR:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:102:TYR:C	2:H:103:TRP:CD1	2.92	0.43
2:I:50:THR:O	2:I:69:ILE:HD12	2.18	0.43
1:L:16:GLY:HA2	1:L:77:PRO:HB2	2.01	0.43
1:M:51:ILE:HD11	1:M:65:SER:CA	2.47	0.43
2:I:20:LEU:HG	2:I:82:MET:HE2	1.99	0.43
1:L:105:GLU:CD	1:L:173:TYR:OH	2.55	0.43
1:L:12:ALA:HA	1:L:105:GLU:O	2.18	0.43
2:I:59:TYR:CD1	2:I:59:TYR:N	2.87	0.43
1:L:167:ASP:OD2	1:L:168:SER:N	2.52	0.43
2:I:11:LEU:HD23	2:I:12:VAL:H	1.84	0.42
1:M:118:PHE:CD2	2:I:124:LEU:HB3	2.54	0.42
1:M:24:ARG:HH11	1:M:24:ARG:CB	2.25	0.42
1:L:32:PHE:HB3	1:L:91:THR:OG1	2.18	0.42
2:I:17:SER:HB2	2:I:82(A):ASN:HD22	1.84	0.42
1:L:37:GLN:HB2	1:L:47:LEU:HD11	2.01	0.42
2:H:100(G):TYR:O	2:H:100(H):SER:HB3	2.20	0.42
2:H:39:GLN:NE2	2:H:43:LYS:O	2.45	0.42
2:I:36:TRP:HD1	2:I:69:ILE:CG2	2.33	0.42
2:I:97:GLU:HG2	4:I:258:HOH:O	2.19	0.42
2:H:1:GLU:HG3	2:H:2:VAL:N	2.35	0.42
1:M:131:SER:HA	1:M:179:LEU:O	2.19	0.42
1:M:96:TRP:CH2	2:I:100(H):SER:HB2	2.54	0.42
2:I:123:PRO:O	2:I:124:LEU:HD12	2.20	0.42
2:I:162:GLY:O	2:I:181:VAL:HA	2.19	0.42
1:M:106:ILE:HG22	1:M:107:LYS:N	2.35	0.42
2:H:105:GLN:OE1	2:H:105:GLN:N	2.53	0.42
1:L:193:THR:HG22	1:L:208:SER:HB2	1.99	0.42
1:M:123:GLU:O	1:M:124:GLN:C	2.58	0.42
1:M:124:GLN:HG2	1:M:129:GLY:C	2.40	0.42
1:M:124:GLN:HB2	2:I:122:TYR:CE2	2.55	0.42
2:H:11:LEU:HD23	2:H:12:VAL:N	2.35	0.41
2:H:194:THR:HG23	2:H:209:LYS:N	2.35	0.41
1:L:27(C):ASP:OD1	1:L:68:GLY:N	2.52	0.41
1:L:89:GLN:CG	1:L:90:GLN:N	2.82	0.41
2:I:6:GLU:H	2:I:105:GLN:NE2	2.18	0.41
1:L:124:GLN:HG2	1:L:129:GLY:C	2.40	0.41
1:M:81:ASP:C	1:M:83:PRO:HD2	2.40	0.41
1:L:149:LYS:HA	1:L:153:SER:O	2.20	0.41
2:H:59:TYR:CD1	2:H:59:TYR:N	2.88	0.41
2:I:36:TRP:HD1	2:I:69:ILE:HG22	1.85	0.41
2:I:59:TYR:CE2	2:I:68:THR:HA	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:42:GLN:HB3	1:L:43:PRO:HD2	2.02	0.41
1:M:124:GLN:HE22	1:M:131:SER:HB2	1.86	0.41
2:H:194:THR:HG23	2:H:209:LYS:HA	2.03	0.41
2:H:39:GLN:C	2:H:88:ALA:HB1	2.41	0.41
2:H:87:THR:O	2:H:88:ALA:HB2	2.21	0.41
2:I:213:ARG:O	2:I:214:ASP:HB2	2.21	0.41
1:M:175:MET:HG2	1:M:176:SER:N	2.36	0.41
2:I:133:ASN:OD1	2:I:134:SER:N	2.48	0.41
1:M:76:HIS:HA	1:M:77:PRO:HA	1.94	0.41
2:H:83:LYS:C	2:H:111:VAL:HG11	2.41	0.41
1:L:125:LEU:HD12	1:L:125:LEU:HA	1.91	0.41
1:L:136:LEU:N	1:L:136:LEU:HD12	2.36	0.40
2:H:63:VAL:HB	2:H:67:PHE:CG	2.57	0.40
2:H:100(E):TYR:CD1	2:H:100(E):TYR:N	2.90	0.40
2:H:153:THR:CG2	2:H:157:GLY:N	2.85	0.40
2:H:177:LEU:CD1	2:H:177:LEU:C	2.89	0.40
1:M:175:MET:CE	1:M:177:SER:HB2	2.51	0.40
2:I:1:GLU:HB2	4:I:303:HOH:O	2.21	0.40
2:H:93:ALA:HB3	2:H:100(J):MET:CE	2.52	0.40
2:I:40:THR:HB	2:I:41:PRO:CD	2.48	0.40
1:L:183:LYS:HG2	1:L:187:GLU:OE2	2.21	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	L	214/218 (98%)	190 (89%)	23 (11%)	1 (0%)	29	54
1	M	216/218 (99%)	199 (92%)	15 (7%)	2 (1%)	17	40
2	H	217/228 (95%)	198 (91%)	17 (8%)	2 (1%)	17	40
2	I	226/228 (99%)	211 (93%)	14 (6%)	1 (0%)	34	60

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	P	9/11 (82%)	8 (89%)	1 (11%)	0	100	100
All	All	882/903 (98%)	806 (91%)	70 (8%)	6 (1%)	22	46

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	171	SER
2	H	62	SER
1	M	51	ILE
1	M	27(D)	ASN
2	I	100(F)	TYR
2	H	149	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	L	189/191 (99%)	184 (97%)	5 (3%)	46	75
1	M	191/191 (100%)	184 (96%)	7 (4%)	34	63
2	H	191/196 (97%)	182 (95%)	9 (5%)	26	54
2	I	196/196 (100%)	184 (94%)	12 (6%)	18	41
3	P	11/11 (100%)	11 (100%)	0	100	100
All	All	778/785 (99%)	745 (96%)	33 (4%)	30	58

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

Mol	Chain	Res	Type
1	L	24	ARG
1	L	73	LEU
1	L	90	GLN
1	L	176	SER
1	L	179	LEU
2	H	11	LEU

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Mol	Chain	Res	Type
2	H	18	LEU
2	H	59	TYR
2	H	69	ILE
2	H	74	ASP
2	H	80	LEU
2	H	141	LEU
2	H	149	PRO
2	H	207	ASP
1	M	24	ARG
1	M	27(B)	VAL
1	M	73	LEU
1	M	156	GLN
1	M	157	ASN
1	M	169	LYS
1	M	179	LEU
2	I	11	LEU
2	I	18	LEU
2	I	50	THR
2	I	59	TYR
2	I	69	ILE
2	I	74	ASP
2	I	80	LEU
2	I	95	ARG
2	I	96	SER
2	I	141	LEU
2	I	149	PRO
2	I	207	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	34	ASN
1	L	53	ASN
1	L	76	HIS
1	L	89	GLN
1	L	90	GLN
1	L	137	ASN
2	H	31	HIS
2	H	58	HIS
2	H	81	GLN
2	H	82(A)	ASN
1	M	53	ASN

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Mol	Chain	Res	Type
1	M	76	HIS
1	M	137	ASN
1	M	156	GLN
2	I	31	HIS
2	I	81	GLN
2	I	82(A)	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 carbohydrates 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	L	216/218 (99%)	0.18	6 (2%) 53 54	35, 61, 82, 99	0
1	M	218/218 (100%)	-0.15	2 (0%) 84 85	25, 45, 75, 99	0
2	H	221/228 (96%)	0.85	30 (13%) 3 2	23, 72, 92, 102	0
2	I	228/228 (100%)	-0.01	4 (1%) 68 70	23, 55, 82, 111	0
3	P	11/11 (100%)	-0.45	0 100 100	35, 40, 59, 61	0
All	All	894/903 (99%)	0.21	42 (4%) 31 30	23, 58, 86, 111	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	206	VAL	5.4
2	I	131	GLN	5.4
2	I	130	ALA	4.8
2	H	134	SER	4.4
1	M	214	CYS	3.9
2	H	136	VAL	3.5
2	H	200	PRO	3.5
2	H	126	PRO	3.4
2	H	184	PRO	3.3
2	H	186	SER	3.2
1	L	3	VAL	3.1
1	L	121	SER	3.0
2	H	128	SER	3.0
2	H	15	GLY	2.9
2	H	172	SER	2.9
2	H	185	SER	2.9
2	H	158	SER	2.9
2	H	188	TRP	2.8
2	H	192	THR	2.8
2	H	156	SER	2.7

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Mol	Chain	Res	Type	RSRZ
2	H	127	GLY	2.7
2	I	129	ALA	2.5
1	L	126	THR	2.4
2	H	112	SER	2.4
2	H	83	LYS	2.4
2	H	111	VAL	2.4
2	H	86	ASP	2.4
2	H	122	TYR	2.3
2	H	171	GLN	2.3
2	H	118	PRO	2.2
2	H	119	PRO	2.2
2	H	82(C)	LEU	2.2
2	H	169	VAL	2.2
2	H	135	MET	2.2
2	H	189	PRO	2.1
1	L	80	GLU	2.1
2	H	194	THR	2.1
1	L	211	ARG	2.1
1	M	213	GLU	2.1
2	H	120	SER	2.1
1	L	24	ARG	2.0
2	I	132	THR	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

There are no ligands in this entry.

## 6.5 Other polymers ⓘ

There are no such residues in this entry.