



# Full wwPDB X-ray Structure Validation Report i

Feb 28, 2014 – 09:18 PM GMT

PDB ID : 4E53  
Title : Calmodulin and Nm peptide complex  
Authors : Kumar, V.; Sivaraman, J.  
Deposited on : 2012-03-13  
Resolution : 2.69 Å(reported)

This is a full wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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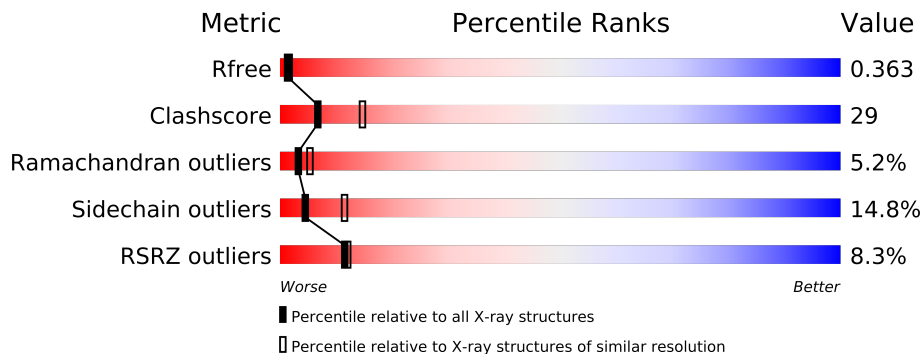
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

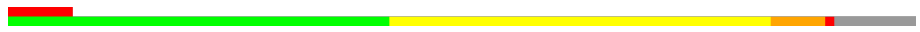

The reported resolution of this entry is 2.69 Å.

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))
$R_{free}$	66092	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	185	
1	B	185	

## 2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 2715 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Calmodulin, Linker, IQ motif of Neuromodulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	168	1323	813	220	280	10	0	0	0
1	B	168	1323	813	220	280	10	0	0	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	-	EXPRESSION TAG	UNP P62204
A	-5	HIS	-	EXPRESSION TAG	UNP P62204
A	-4	HIS	-	EXPRESSION TAG	UNP P62204
A	-3	HIS	-	EXPRESSION TAG	UNP P62204
A	-2	HIS	-	EXPRESSION TAG	UNP P62204
A	-1	HIS	-	EXPRESSION TAG	UNP P62204
A	0	HIS	-	EXPRESSION TAG	UNP P62204
B	-6	MET	-	EXPRESSION TAG	UNP P62204
B	-5	HIS	-	EXPRESSION TAG	UNP P62204
B	-4	HIS	-	EXPRESSION TAG	UNP P62204
B	-3	HIS	-	EXPRESSION TAG	UNP P62204
B	-2	HIS	-	EXPRESSION TAG	UNP P62204
B	-1	HIS	-	EXPRESSION TAG	UNP P62204
B	0	HIS	-	EXPRESSION TAG	UNP P62204

- Molecule 2 is water.

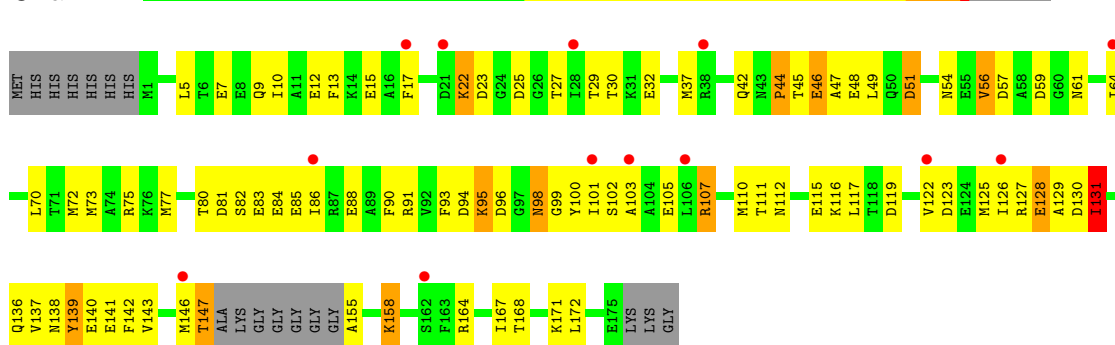
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	40	Total 40	O 40	0	0
2	B	29	Total 29	O 29	0	0

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

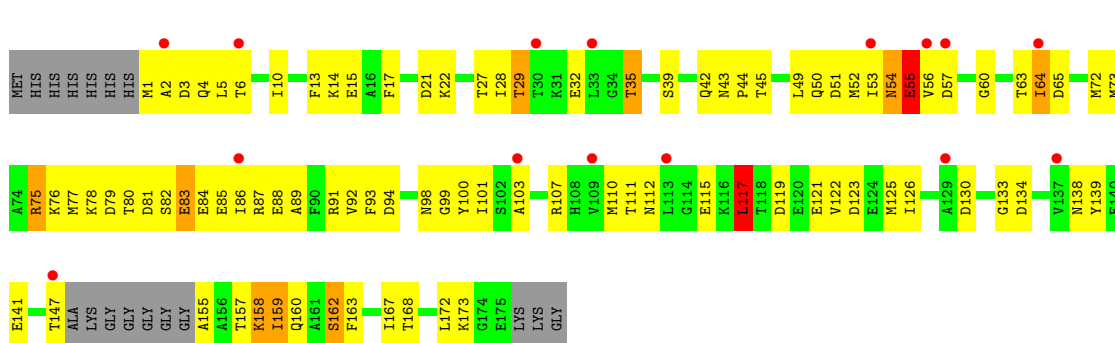
- Molecule 1: Calmodulin, Linker, IQ motif of Neuromodulin

Chain A:



- Molecule 1: Calmodulin, Linker, IQ motif of Neuromodulin

Chain B:



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.39Å 79.28Å 136.06Å 90.00° 90.19° 90.00°	Depositor
Resolution (Å)	30.00 – 2.69 24.51 – 2.69	Depositor EDS
% Data completeness (in resolution range)	94.7 (30.00-2.69) 97.8 (24.51-2.69)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.44 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.276 , 0.318 0.322 , 0.363	Depositor DCC
$R_{free}$ test set	1304 reflections (5.66%)	DCC
Wilson B-factor (Å <sup>2</sup> )	59.3	Xtriage
Anisotropy	0.097	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 22.9	EDS
Estimated twinning fraction	0.255 for H, K, L 0.247 for -K, -H, -L 0.241 for K, H, -L 0.257 for -H, -K, L 0.032 for k,h,-l 0.044 for -k,-h,-l 0.078 for -h,-k,l	Xtriage
Reported twinning fraction	0.255 for H, K, L 0.247 for -K, -H, -L 0.241 for K, H, -L 0.257 for -H, -K, L	Depositor
L-test for twinning	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	0 of 23303 reflections	Xtriage
$F_o, F_c$ correlation	0.76	EDS
Total number of atoms	2715	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 4.81% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 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	0.38	0/1336	0.59	0/1788
1	B	0.38	0/1336	0.60	1/1788 (0.1%)
All	All	0.38	0/2672	0.59	1/3576 (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	B	117	LEU	CA-CB-CG	5.75	128.52	115.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1323	0	1267	76	0
1	B	1323	0	1267	75	0
2	A	40	0	0	18	0
2	B	29	0	0	11	0
All	All	2715	0	2534	151	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 29.

All (151) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:110:MET:HG2	1:A:115:GLU:HB2	1.38	1.05
1:B:110:MET:HG3	1:B:117:LEU:HG	1.45	0.95
1:A:5:LEU:HB3	1:A:10:ILE:HD11	1.49	0.94
1:B:44:PRO:HB2	2:B:212:HOH:O	1.67	0.93
1:B:80:THR:O	1:B:84:GLU:HB3	1.74	0.88
1:A:100:TYR:HD2	2:A:214:HOH:O	1.56	0.88
1:A:90:PHE:HZ	1:A:99:GLY:HA2	1.38	0.88
1:B:163:PHE:O	1:B:167:ILE:HD12	1.77	0.83
1:A:90:PHE:CZ	1:A:99:GLY:HA2	2.17	0.80
1:A:51:ASP:HA	1:A:54:ASN:ND2	1.98	0.79
1:B:6:THR:O	2:B:221:HOH:O	2.00	0.78
1:B:27:THR:HB	1:B:63:THR:HB	1.65	0.78
1:A:131:ILE:O	2:A:236:HOH:O	2.01	0.77
1:A:102:SER:HA	1:A:136:GLN:HG2	1.67	0.77
1:B:155:ALA:O	1:B:158:LYS:HG2	1.85	0.75
1:A:128:GLU:HB3	2:A:230:HOH:O	1.85	0.74
1:B:65:ASP:HB3	2:B:222:HOH:O	1.86	0.74
1:B:85:GLU:HB3	1:B:167:ILE:HG21	1.72	0.70
1:B:121:GLU:HG3	2:B:211:HOH:O	1.89	0.70
1:B:53:ILE:HG23	1:B:64:ILE:HG12	1.74	0.70
1:A:167:ILE:HA	2:A:222:HOH:O	1.92	0.69
1:A:73:MET:HB2	2:A:233:HOH:O	1.93	0.68
1:A:86:ILE:HG21	1:A:143:VAL:HA	1.74	0.68
1:B:87:ARG:HH22	1:B:91:ARG:NH2	1.92	0.67
1:B:13:PHE:O	1:B:17:PHE:HB2	1.94	0.67
1:A:99:GLY:O	1:A:138:ASN:ND2	2.28	0.66
1:A:46:GLU:O	2:A:234:HOH:O	2.12	0.66
1:A:45:THR:O	1:A:47:ALA:N	2.30	0.64
1:A:88:GLU:HB3	1:A:167:ILE:HG23	1.80	0.64
1:B:51:ASP:HA	1:B:54:ASN:CG	2.18	0.63
1:A:22:LYS:HD3	1:A:32:GLU:HG3	1.79	0.63
1:A:138:ASN:OD1	1:A:139:TYR:N	2.25	0.63
1:B:27:THR:HB	1:B:63:THR:CB	2.29	0.63
1:A:101:ILE:HG23	1:A:105:GLU:HB2	1.81	0.62
1:A:127:ARG:C	1:A:129:ALA:H	2.00	0.62
1:A:168:THR:HB	2:A:213:HOH:O	1.99	0.62
1:B:1:MET:O	1:B:1:MET:HG3	1.98	0.62
1:A:110:MET:HG3	1:A:117:LEU:HD12	1.82	0.61
1:A:10:ILE:HA	1:A:13:PHE:HD1	1.65	0.61
1:B:29:THR:HG22	1:B:63:THR:HG22	1.83	0.61
1:A:56:VAL:HG11	1:A:72:MET:HB2	1.81	0.61
1:A:45:THR:C	1:A:47:ALA:H	2.04	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:9:GLN:O	1:A:12:GLU:HB3	2.00	0.61
1:B:56:VAL:HG21	1:B:72:MET:HB2	1.83	0.61
1:B:10:ILE:O	1:B:14:LYS:HG3	2.01	0.60
1:A:86:ILE:HG23	1:A:146:MET:HE2	1.81	0.60
1:A:98:ASN:HD22	1:A:98:ASN:H	1.48	0.60
1:A:51:ASP:HA	1:A:54:ASN:HD22	1.64	0.59
1:B:28:ILE:HG23	1:B:32:GLU:HB3	1.84	0.59
1:A:44:PRO:HG2	1:A:49:LEU:HD21	1.85	0.59
1:B:42:GLN:HE21	1:B:42:GLN:HA	1.69	0.58
1:A:10:ILE:HB	2:A:237:HOH:O	2.03	0.58
1:A:82:SER:O	1:A:86:ILE:HG13	2.04	0.58
1:A:13:PHE:O	1:A:17:PHE:HB2	2.04	0.58
1:B:126:ILE:O	1:B:130:ASP:HB2	2.04	0.58
1:B:82:SER:O	1:B:86:ILE:HG13	2.04	0.57
1:B:85:GLU:CB	2:B:210:HOH:O	2.52	0.57
1:B:85:GLU:HB2	2:B:210:HOH:O	2.05	0.56
1:B:172:LEU:HD23	1:B:173:LYS:HG2	1.87	0.56
1:A:7:GLU:HA	2:A:237:HOH:O	2.04	0.56
1:A:85:GLU:HA	1:A:167:ILE:HG21	1.88	0.56
1:B:99:GLY:O	1:B:138:ASN:ND2	2.38	0.56
1:A:90:PHE:HZ	1:A:99:GLY:CA	2.16	0.55
1:B:93:PHE:HB3	1:B:101:ILE:HD13	1.87	0.55
1:B:81:ASP:O	1:B:85:GLU:HG3	2.07	0.55
1:B:92:VAL:O	1:B:92:VAL:HG12	2.06	0.55
1:B:45:THR:O	2:B:212:HOH:O	2.18	0.55
1:A:122:VAL:O	1:A:126:ILE:HG13	2.08	0.53
1:B:45:THR:N	2:B:212:HOH:O	2.41	0.53
1:A:91:ARG:HB3	2:A:216:HOH:O	2.08	0.53
1:B:88:GLU:HB3	1:B:167:ILE:HG12	1.89	0.53
1:A:125:MET:HA	2:A:230:HOH:O	2.08	0.53
1:B:79:ASP:O	1:B:83:GLU:N	2.42	0.53
1:B:155:ALA:HA	1:B:158:LYS:HE3	1.91	0.53
1:B:80:THR:O	1:B:84:GLU:CB	2.51	0.53
1:A:127:ARG:C	1:A:129:ALA:N	2.63	0.52
1:B:50:GLN:O	1:B:54:ASN:OD1	2.28	0.52
1:B:77:MET:HA	2:B:217:HOH:O	2.10	0.52
1:A:73:MET:O	1:A:77:MET:HG3	2.10	0.52
1:B:73:MET:HB3	1:B:77:MET:HE3	1.92	0.51
1:A:81:ASP:O	1:A:85:GLU:HG3	2.11	0.51
1:A:94:ASP:HB2	1:A:101:ILE:HG12	1.92	0.51
1:B:52:MET:HG2	1:B:76:LYS:NZ	2.25	0.51
1:A:164:ARG:HA	1:A:167:ILE:HD12	1.92	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:45:THR:OG1	1:A:48:GLU:OE1	2.30	0.50
1:B:138:ASN:HB3	1:B:141:GLU:HG3	1.93	0.50
1:B:122:VAL:HG12	1:B:126:ILE:HD11	1.92	0.50
1:A:61:ASN:HB2	2:A:220:HOH:O	2.11	0.50
1:A:101:ILE:HB	1:A:137:VAL:HB	1.94	0.49
1:B:89:ALA:HB2	1:B:167:ILE:HG13	1.94	0.48
1:B:123:ASP:HA	1:B:126:ILE:HD12	1.96	0.48
1:B:53:ILE:O	1:B:57:ASP:HB2	2.14	0.47
1:A:125:MET:C	2:A:230:HOH:O	2.53	0.47
1:B:29:THR:HG22	1:B:63:THR:CG2	2.44	0.47
1:A:37:MET:CE	1:A:44:PRO:HG3	2.45	0.46
1:A:147:THR:C	2:A:229:HOH:O	2.54	0.46
1:A:100:TYR:CD2	2:A:214:HOH:O	2.42	0.46
1:A:155:ALA:O	1:A:158:LYS:HG3	2.16	0.45
1:B:21:ASP:HB2	1:B:28:ILE:HG13	1.98	0.45
1:B:49:LEU:N	2:B:212:HOH:O	2.49	0.45
1:B:2:ALA:O	1:B:4:GLN:N	2.48	0.45
1:A:10:ILE:HA	1:A:13:PHE:CD1	2.47	0.45
1:B:117:LEU:HD12	1:B:122:VAL:HG22	1.99	0.45
1:B:107:ARG:HE	1:B:119:ASP:CG	2.21	0.45
1:B:85:GLU:O	1:B:167:ILE:HD13	2.17	0.45
1:A:110:MET:SD	1:A:111:THR:N	2.89	0.45
1:B:55:GLU:OE1	1:B:72:MET:HG3	2.17	0.44
1:B:93:PHE:HD1	2:B:209:HOH:O	1.99	0.44
1:B:89:ALA:O	1:B:93:PHE:N	2.50	0.44
1:A:45:THR:C	1:A:47:ALA:N	2.69	0.44
1:A:168:THR:O	1:A:172:LEU:HB3	2.18	0.44
1:A:46:GLU:HG2	1:A:46:GLU:H	1.66	0.44
1:B:126:ILE:O	1:B:130:ASP:CB	2.65	0.44
1:B:54:ASN:O	1:B:55:GLU:C	2.57	0.43
1:B:103:ALA:HB2	1:B:126:ILE:HD13	2.01	0.43
1:B:93:PHE:CB	1:B:101:ILE:HD13	2.49	0.43
1:A:85:GLU:CA	1:A:167:ILE:HG21	2.48	0.43
1:A:98:ASN:ND2	1:A:98:ASN:H	2.16	0.43
1:B:4:GLN:O	1:B:6:THR:N	2.50	0.43
1:A:110:MET:HG2	1:A:115:GLU:CB	2.28	0.43
1:B:21:ASP:HB2	1:B:28:ILE:CG1	2.49	0.43
1:B:75:ARG:HG2	1:B:78:LYS:HD3	2.01	0.42
1:A:15:GLU:OE2	1:A:15:GLU:N	2.51	0.42
1:B:35:THR:O	1:B:39:SER:OG	2.37	0.42
1:A:98:ASN:HD22	1:A:99:GLY:N	2.16	0.42
1:B:54:ASN:O	1:B:56:VAL:N	2.52	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:155:ALA:O	1:B:158:LYS:HE3	2.20	0.42
1:A:95:LYS:HZ3	1:A:96:ASP:N	2.17	0.42
1:B:94:ASP:OD2	1:B:100:TYR:O	2.38	0.42
1:B:125:MET:SD	1:B:159:ILE:CD1	3.08	0.42
1:A:168:THR:HA	1:A:171:LYS:HG2	2.02	0.42
1:A:42:GLN:HE21	1:A:42:GLN:HA	1.84	0.42
1:A:98:ASN:N	1:A:98:ASN:ND2	2.68	0.41
1:B:121:GLU:O	1:B:125:MET:HG2	2.20	0.41
1:A:142:PHE:HB2	2:A:205:HOH:O	2.20	0.41
1:A:98:ASN:HD22	1:A:98:ASN:N	2.10	0.41
1:A:93:PHE:O	1:A:105:GLU:HB3	2.20	0.41
1:A:137:VAL:HG11	1:A:142:PHE:HD1	1.85	0.41
1:B:51:ASP:OD1	1:B:54:ASN:ND2	2.54	0.41
1:A:9:GLN:NE2	2:A:224:HOH:O	2.53	0.41
1:A:88:GLU:OE2	1:A:91:ARG:NH1	2.51	0.41
1:B:51:ASP:HA	1:B:54:ASN:ND2	2.35	0.41
1:B:73:MET:O	1:B:77:MET:HG3	2.20	0.40
1:B:160:GLN:HA	1:B:163:PHE:HB2	2.04	0.40
1:A:107:ARG:NH2	2:A:202:HOH:O	2.52	0.40
1:A:86:ILE:HG13	1:A:86:ILE:H	1.76	0.40
1:A:94:ASP:HB2	1:A:101:ILE:CG1	2.51	0.40
1:B:52:MET:HG2	1:B:76:LYS:HZ1	1.86	0.40
1:B:115:GLU:OE2	1:B:162:SER:HB3	2.21	0.40
1:A:119:ASP:O	1:A:123:ASP:HB3	2.21	0.40
1:B:138:ASN:CG	1:B:139:TYR:H	2.24	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	164/185 (89%)	130 (79%)	25 (15%)	9 (6%)	3	4
1	B	164/185 (89%)	121 (74%)	35 (21%)	8 (5%)	3	6
All	All	328/370 (89%)	251 (76%)	60 (18%)	17 (5%)	3	5

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	54	ASN
1	B	55	GLU
1	B	60	GLY
1	A	44	PRO
1	A	46	GLU
1	A	70	LEU
1	A	103	ALA
1	A	139	TYR
1	B	3	ASP
1	B	133	GLY
1	A	128	GLU
1	B	5	LEU
1	A	23	ASP
1	A	25	ASP
1	B	158	LYS
1	B	159	ILE
1	A	131	ILE

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	142/152 (93%)	118 (83%)	24 (17%)	<b>3</b> <b>8</b>
1	B	142/152 (93%)	124 (87%)	18 (13%)	<b>6</b> <b>15</b>
All	All	284/304 (93%)	242 (85%)	42 (15%)	<b>4</b> <b>11</b>

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LYS
1	A	27	THR
1	A	29	THR
1	A	30	THR
1	A	51	ASP
1	A	56	VAL
1	A	57	ASP

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Mol	Chain	Res	Type
1	A	59	ASP
1	A	64	ILE
1	A	75	ARG
1	A	80	THR
1	A	83	GLU
1	A	84	GLU
1	A	95	LYS
1	A	98	ASN
1	A	107	ARG
1	A	112	ASN
1	A	116	LYS
1	A	130	ASP
1	A	131	ILE
1	A	140	GLU
1	A	141	GLU
1	A	147	THR
1	A	158	LYS
1	B	15	GLU
1	B	22	LYS
1	B	29	THR
1	B	35	THR
1	B	43	ASN
1	B	55	GLU
1	B	64	ILE
1	B	75	ARG
1	B	83	GLU
1	B	98	ASN
1	B	111	THR
1	B	112	ASN
1	B	117	LEU
1	B	134	ASP
1	B	147	THR
1	B	157	THR
1	B	162	SER
1	B	168	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	A	54	ASN
1	A	61	ASN

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Mol	Chain	Res	Type
1	A	98	ASN
1	A	112	ASN
1	A	160	GLN
1	B	4	GLN
1	B	42	GLN
1	B	98	ASN
1	B	112	ASN

### 5.3.3 RNA [i](#)

There are no RNA chains 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

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	168/185 (90%)	0.62	13 (7%) <b>13</b> <b>14</b>	50, 58, 64, 64	0
1	B	168/185 (90%)	0.56	15 (8%) <b>10</b> <b>10</b>	49, 57, 61, 63	0
All	All	336/370 (90%)	0.59	28 (8%) <b>11</b> <b>12</b>	49, 57, 63, 64	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	56	VAL	7.9
1	A	64	ILE	6.4
1	A	103	ALA	5.4
1	B	109	VAL	4.4
1	B	57	ASP	4.1
1	A	106	LEU	3.9
1	A	122	VAL	3.8
1	B	137	VAL	3.7
1	A	146	MET	3.4
1	B	113	LEU	3.3
1	B	86	ILE	3.3
1	B	147	THR	3.2
1	B	30	THR	3.0
1	B	64	ILE	2.8
1	B	129	ALA	2.8
1	B	33	LEU	2.6
1	A	21	ASP	2.6
1	B	103	ALA	2.5
1	A	38	ARG	2.4
1	A	86	ILE	2.3
1	B	53	ILE	2.3
1	B	2	ALA	2.3
1	A	28	ILE	2.2
1	A	162	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	126	ILE	2.1
1	B	6	THR	2.1
1	A	101	ILE	2.0
1	A	17	PHE	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 carbohydrates 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.