Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 01:59 AM GMT

PDB ID	:	2F2O
Title	:	Structure of calmodulin bound to a calcineurin peptide: a new way of making an old binding mode
Authors	:	Ye, Q.; Wong, A.; Jia, Z.
Deposited on	:	2005-11-17
Resolution	:	2.17 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

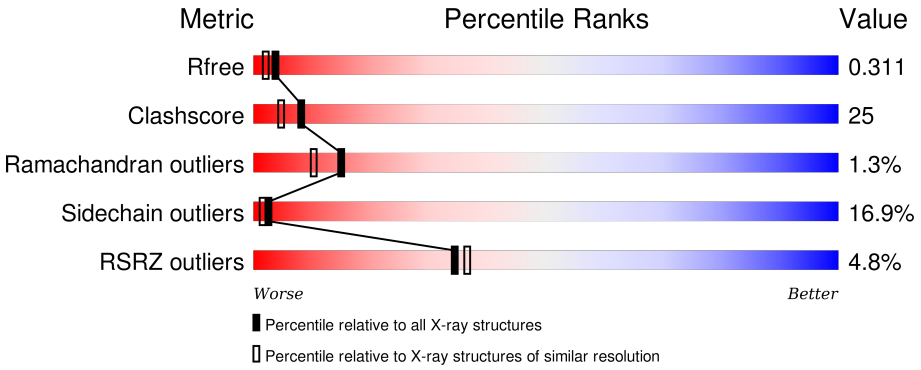
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.17 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}	91344	5130 (2.20-2.16)
Clashscore	102246	5965 (2.20-2.16)
Ramachandran outliers	100387	5863 (2.20-2.16)
Sidechain outliers	100360	5864 (2.20-2.16)
RSRZ outliers	91569	5142 (2.20-2.16)

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 >=3, 2, 1 and 0 types of geometric quality criteria. 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 <=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	179	5% 44% 26% 13% • 13%
1	B	179	3% 41% 35% 9% • 11%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CA	A	1001	-	-	-	X
2	CA	A	1004	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2676 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 Calmodulin fused with calmodulin-binding domain of calcineurin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	156	Total	C	N	O	S	0	0	0
			1236	762	206	258	10			
1	B	159	Total	C	N	O	S	0	0	0
			1254	773	212	259	10			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	149	GLY	-	LINKER	UNP P62157
A	150	GLY	-	LINKER	UNP P62157
A	151	GLY	-	LINKER	UNP P62157
A	152	GLY	-	LINKER	UNP P62157
A	153	GLY	-	LINKER	UNP P62157
B	149	GLY	-	LINKER	UNP P62157
B	150	GLY	-	LINKER	UNP P62157
B	151	GLY	-	LINKER	UNP P62157
B	152	GLY	-	LINKER	UNP P62157
B	153	GLY	-	LINKER	UNP P62157

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	4	Total	Ca	0	0
			4	4		
2	A	4	Total	Ca	0	0
			4	4		

- Molecule 3 is water.

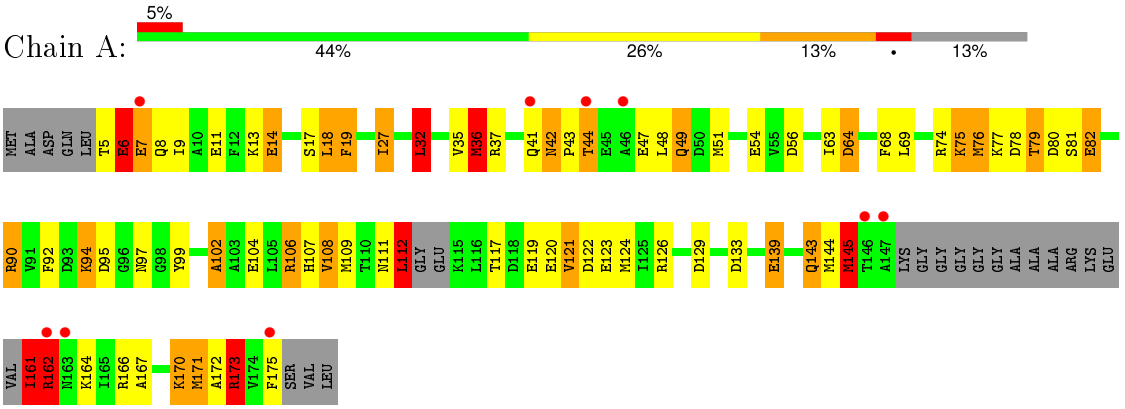
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	85	Total O 85 85	0	0
3	B	93	Total O 93 93	0	0

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 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.

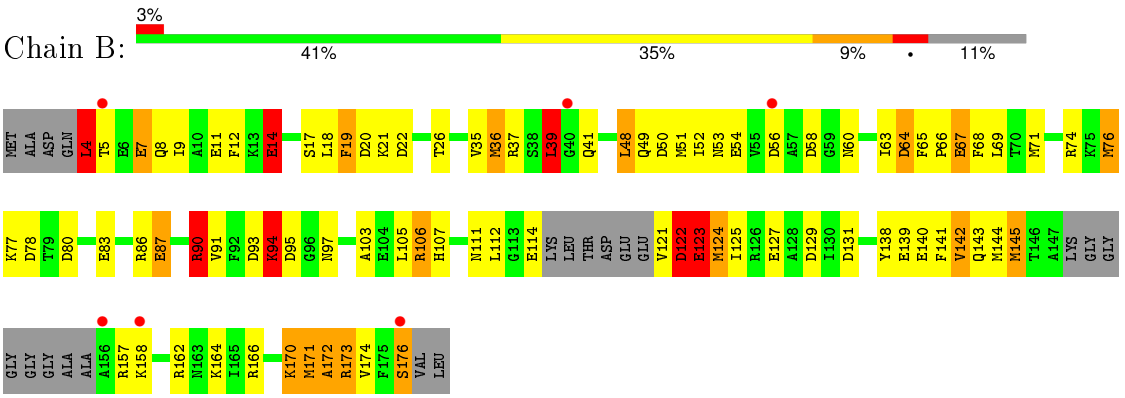
- Molecule 1: Calmodulin fused with calmodulin-binding domain of calcineurin

Chain A:



- Molecule 1: Calmodulin fused with calmodulin-binding domain of calcineurin

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	121.24Å 42.77Å 71.15Å 90.00° 110.39° 90.00°	Depositor
Resolution (Å)	67.42 – 2.17 28.96 – 2.17	Depositor EDS
% Data completeness (in resolution range)	97.9 (67.42-2.17) 97.9 (28.96-2.17)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	6.50 (at 2.18Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.204 , 0.289 0.245 , 0.311	Depositor DCC
R_{free} test set	1777 reflections (9.87%)	DCC
Wilson B-factor (Å^2)	33.1	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 55.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 18124 reflections	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	2676	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.58% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA

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	2.01	29/1247 (2.3%)	1.69	24/1668 (1.4%)
1	B	2.08	37/1265 (2.9%)	1.62	23/1691 (1.4%)
All	All	2.05	66/2512 (2.6%)	1.66	47/3359 (1.4%)

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	A	0	2
1	B	0	1
All	All	0	3

All (66) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	67	GLU	CD-OE1	15.68	1.42	1.25
1	B	142	VAL	CB-CG1	9.38	1.72	1.52
1	B	53	ASN	CB-CG	8.52	1.70	1.51
1	A	99	TYR	CD2-CE2	8.13	1.51	1.39
1	A	68	PHE	CD2-CE2	7.86	1.54	1.39
1	A	108	VAL	CB-CG2	7.80	1.69	1.52
1	B	58	ASP	CB-CG	7.72	1.68	1.51
1	A	102	ALA	CA-CB	7.41	1.68	1.52
1	B	106	ARG	C-O	7.37	1.37	1.23
1	A	77	LYS	N-CA	7.31	1.60	1.46
1	B	12	PHE	CE1-CZ	7.28	1.51	1.37
1	A	13	LYS	CE-NZ	6.96	1.66	1.49
1	B	78	ASP	CA-CB	-6.96	1.38	1.53
1	B	83	GLU	CD-OE1	6.89	1.33	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	170	LYS	CD-CE	6.86	1.68	1.51
1	A	79	THR	CB-CG2	6.84	1.75	1.52
1	A	94	LYS	CD-CE	6.79	1.68	1.51
1	B	68	PHE	N-CA	-6.71	1.32	1.46
1	A	145	MET	SD-CE	-6.45	1.41	1.77
1	B	67	GLU	CB-CG	6.45	1.64	1.52
1	B	171	MET	SD-CE	6.44	2.13	1.77
1	B	60	ASN	CG-OD1	6.41	1.38	1.24
1	A	121	VAL	CB-CG2	-6.40	1.39	1.52
1	A	80	ASP	C-O	6.39	1.35	1.23
1	B	54	GLU	CD-OE2	6.36	1.32	1.25
1	A	32	LEU	C-O	6.35	1.35	1.23
1	A	139	GLU	CG-CD	6.27	1.61	1.51
1	B	90	ARG	CG-CD	6.26	1.67	1.51
1	B	71	MET	CG-SD	6.25	1.97	1.81
1	B	71	MET	SD-CE	-6.24	1.43	1.77
1	B	145	MET	SD-CE	-6.23	1.43	1.77
1	A	144	MET	CG-SD	-6.22	1.65	1.81
1	B	76	MET	CG-SD	6.19	1.97	1.81
1	A	90	ARG	CB-CG	6.15	1.69	1.52
1	B	19	PHE	CD2-CE2	-6.02	1.27	1.39
1	B	80	ASP	CB-CG	5.95	1.64	1.51
1	A	77	LYS	C-O	-5.90	1.12	1.23
1	B	141	PHE	C-O	5.84	1.34	1.23
1	A	82	GLU	CD-OE1	5.77	1.32	1.25
1	A	90	ARG	CG-CD	5.74	1.66	1.51
1	B	68	PHE	CA-C	5.73	1.67	1.52
1	A	161	ILE	CA-CB	5.72	1.68	1.54
1	A	9	ILE	CB-CG2	5.70	1.70	1.52
1	A	90	ARG	NE-CZ	5.70	1.40	1.33
1	B	41	GLN	CA-CB	-5.63	1.41	1.53
1	B	174	VAL	CB-CG1	-5.61	1.41	1.52
1	A	108	VAL	CB-CG1	5.60	1.64	1.52
1	B	103	ALA	CA-CB	-5.55	1.40	1.52
1	A	171	MET	SD-CE	5.52	2.08	1.77
1	B	12	PHE	CG-CD2	5.52	1.47	1.38
1	B	14	GLU	CD-OE2	5.44	1.31	1.25
1	B	123	GLU	CB-CG	5.41	1.62	1.52
1	B	94	LYS	CB-CG	5.38	1.67	1.52
1	B	172	ALA	CA-CB	5.34	1.63	1.52
1	A	19	PHE	CE1-CZ	-5.33	1.27	1.37
1	B	106	ARG	CG-CD	5.25	1.65	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	20	ASP	CB-CG	5.24	1.62	1.51
1	A	139	GLU	CD-OE2	5.23	1.31	1.25
1	A	123	GLU	CD-OE1	-5.20	1.20	1.25
1	A	27	ILE	CA-CB	-5.18	1.43	1.54
1	B	83	GLU	CG-CD	5.16	1.59	1.51
1	A	68	PHE	CA-CB	-5.13	1.42	1.53
1	B	74	ARG	CG-CD	5.11	1.64	1.51
1	A	104	GLU	CG-CD	5.03	1.59	1.51
1	B	67	GLU	CA-CB	-5.02	1.43	1.53
1	B	94	LYS	CD-CE	5.00	1.63	1.51

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	36	MET	CG-SD-CE	-12.25	80.59	100.20
1	B	71	MET	CG-SD-CE	-11.58	81.67	100.20
1	B	58	ASP	CB-CG-OD2	11.07	128.26	118.30
1	A	124	MET	CG-SD-CE	-10.73	83.03	100.20
1	A	173	ARG	NE-CZ-NH2	-9.71	115.45	120.30
1	B	36	MET	CG-SD-CE	-9.23	85.43	100.20
1	A	76	MET	CG-SD-CE	9.23	114.97	100.20
1	B	95	ASP	CB-CG-OD2	9.10	126.49	118.30
1	A	173	ARG	NE-CZ-NH1	8.85	124.72	120.30
1	A	18	LEU	CA-CB-CG	8.29	134.36	115.30
1	B	80	ASP	CB-CG-OD2	8.07	125.57	118.30
1	A	129	ASP	CB-CG-OD2	8.01	125.51	118.30
1	A	106	ARG	NE-CZ-NH2	7.76	124.18	120.30
1	A	78	ASP	CB-CG-OD2	7.22	124.79	118.30
1	B	129	ASP	CB-CG-OD2	6.92	124.53	118.30
1	B	22	ASP	CB-CG-OD2	6.80	124.42	118.30
1	B	64	ASP	CB-CG-OD2	6.51	124.16	118.30
1	A	112	LEU	CA-CB-CG	6.51	130.27	115.30
1	B	93	ASP	CB-CG-OD2	6.49	124.14	118.30
1	A	175	PHE	CA-C-O	6.31	133.36	120.10
1	A	122	ASP	CB-CG-OD2	6.25	123.92	118.30
1	A	133	ASP	CB-CG-OD1	6.17	123.85	118.30
1	A	42	ASN	N-CA-C	-6.10	94.53	111.00
1	A	90	ARG	NE-CZ-NH1	6.05	123.32	120.30
1	A	56	ASP	CB-CG-OD2	5.81	123.53	118.30
1	A	162	ARG	NE-CZ-NH2	5.64	123.12	120.30
1	B	51	MET	CG-SD-CE	5.62	109.19	100.20
1	B	112	LEU	CB-CG-CD2	-5.60	101.49	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	14	GLU	OE1-CD-OE2	5.56	129.97	123.30
1	A	99	TYR	CZ-CE2-CD2	-5.49	114.86	119.80
1	A	162	ARG	NE-CZ-NH1	-5.45	117.57	120.30
1	B	122	ASP	CB-CG-OD2	5.43	123.19	118.30
1	A	49	GLN	CB-CA-C	-5.40	99.59	110.40
1	B	93	ASP	CB-CG-OD1	-5.40	113.44	118.30
1	B	140	GLU	OE1-CD-OE2	-5.38	116.85	123.30
1	A	126	ARG	NE-CZ-NH2	5.32	122.96	120.30
1	B	131	ASP	CB-CG-OD1	5.27	123.04	118.30
1	B	50	ASP	CB-CG-OD2	5.25	123.02	118.30
1	A	64	ASP	CB-CG-OD2	5.22	123.00	118.30
1	B	4	LEU	CA-CB-CG	5.21	127.30	115.30
1	B	56	ASP	CB-CG-OD1	-5.17	113.64	118.30
1	B	48	LEU	CB-CG-CD1	-5.16	102.22	111.00
1	B	39	LEU	CB-CG-CD2	5.13	119.73	111.00
1	B	37	ARG	NE-CZ-NH1	5.13	122.86	120.30
1	A	48	LEU	CB-CG-CD2	5.09	119.66	111.00
1	B	74	ARG	N-CA-CB	-5.09	101.43	110.60
1	B	94	LYS	CD-CE-NZ	5.06	123.34	111.70

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	161	ILE	Peptide
1	A	7	GLU	Peptide
1	B	122	ASP	Peptide

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	1236	0	1185	63	0
1	B	1254	0	1209	59	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
3	A	85	0	0	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	93	0	0	22	0
All	All	2676	0	2394	120	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 25.

All (120) 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:79:THR:CG2	1:A:79:THR:CB	1.74	1.56
1:A:171:MET:CE	1:A:171:MET:SD	2.08	1.42
1:B:171:MET:CE	1:B:171:MET:SD	2.13	1.35
1:B:76:MET:SD	3:B:1091:HOH:O	2.15	1.05
1:B:4:LEU:O	1:B:8:GLN:HG3	1.61	0.98
1:B:121:VAL:O	1:B:122:ASP:C	2.00	0.96
1:B:7:GLU:O	1:B:11:GLU:HG3	1.69	0.92
1:A:106:ARG:HB3	3:A:1078:HOH:O	1.72	0.89
1:A:95:ASP:HA	3:A:1043:HOH:O	1.73	0.88
1:A:117:THR:HG21	3:A:1061:HOH:O	1.76	0.85
1:A:64:ASP:HB2	3:A:1073:HOH:O	1.78	0.83
1:B:121:VAL:HG23	1:B:124:MET:HE3	1.60	0.83
1:B:4:LEU:N	1:B:4:LEU:HD23	1.94	0.80
1:A:172:ALA:HB2	1:B:39:LEU:HD22	1.68	0.76
1:A:108:VAL:HG22	3:B:1017:HOH:O	1.85	0.76
1:A:117:THR:CG2	1:A:120:GLU:H	2.00	0.74
1:A:117:THR:HG22	1:A:120:GLU:H	1.52	0.73
1:A:139:GLU:CG	3:A:1010:HOH:O	2.36	0.72
1:B:176:SER:CB	3:B:1095:HOH:O	2.38	0.72
1:B:121:VAL:O	1:B:123:GLU:N	2.24	0.70
1:B:14:GLU:HG3	3:B:1060:HOH:O	1.92	0.69
1:B:19:PHE:HA	1:B:35:VAL:HG21	1.74	0.69
1:B:123:GLU:HA	3:B:1073:HOH:O	1.94	0.68
1:B:106:ARG:NH1	3:B:1049:HOH:O	2.26	0.68
1:B:4:LEU:CD2	1:B:9:ILE:HD11	2.25	0.66
1:A:107:HIS:O	1:A:111:ASN:ND2	2.28	0.66
1:B:176:SER:HB2	3:B:1095:HOH:O	1.96	0.64
1:B:4:LEU:N	1:B:4:LEU:CD2	2.60	0.64
1:A:5:THR:O	1:A:6:GLU:CG	2.46	0.64
1:A:112:LEU:HB2	3:A:1047:HOH:O	1.97	0.64
1:B:76:MET:HB3	3:B:1033:HOH:O	1.97	0.63
1:B:4:LEU:HD21	1:B:9:ILE:HD11	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:139:GLU:OE1	3:B:1029:HOH:O	2.15	0.63
1:B:121:VAL:HG23	1:B:124:MET:CE	2.29	0.62
1:B:124:MET:HB2	3:B:1020:HOH:O	1.97	0.62
1:A:139:GLU:HG3	3:A:1010:HOH:O	1.96	0.62
1:A:7:GLU:HG3	3:A:1040:HOH:O	2.00	0.61
1:A:79:THR:CA	1:A:79:THR:CG2	2.73	0.60
1:B:106:ARG:CZ	3:B:1049:HOH:O	2.49	0.60
1:A:121:VAL:HG23	3:A:1066:HOH:O	2.02	0.60
1:A:161:ILE:C	3:A:1022:HOH:O	2.41	0.59
1:A:79:THR:CG2	1:A:79:THR:OG1	2.45	0.59
1:B:121:VAL:C	1:B:123:GLU:N	2.58	0.57
1:A:112:LEU:CB	3:A:1047:HOH:O	2.53	0.56
1:A:36:MET:HG2	1:A:43:PRO:HG3	1.85	0.56
1:B:138:TYR:O	1:B:142:VAL:HG23	2.07	0.55
1:A:42:ASN:ND2	3:A:1068:HOH:O	2.40	0.55
1:B:173:ARG:HD2	3:B:1057:HOH:O	2.07	0.54
1:A:6:GLU:CG	1:A:8:GLN:HG3	2.37	0.54
1:A:121:VAL:C	3:A:1066:HOH:O	2.46	0.54
1:B:124:MET:HG3	3:B:1054:HOH:O	2.08	0.53
1:A:139:GLU:HG2	3:A:1010:HOH:O	2.04	0.52
1:A:5:THR:N	3:A:1025:HOH:O	2.43	0.52
1:A:7:GLU:O	1:A:11:GLU:HG3	2.09	0.51
1:A:42:ASN:CG	3:A:1068:HOH:O	2.47	0.51
1:A:167:ALA:O	1:A:171:MET:HG2	2.10	0.51
1:A:121:VAL:HB	3:A:1066:HOH:O	2.09	0.51
1:A:121:VAL:CB	3:A:1066:HOH:O	2.58	0.51
1:B:121:VAL:N	1:B:124:MET:HB3	2.26	0.51
1:B:87:GLU:HG3	1:B:172:ALA:HB1	1.93	0.51
1:B:87:GLU:HG3	1:B:90:ARG:HH11	1.76	0.51
1:B:21:LYS:HD3	3:B:1035:HOH:O	2.10	0.51
1:A:75:LYS:HD3	3:A:1021:HOH:O	2.11	0.50
1:A:112:LEU:HD21	1:A:164:LYS:HE3	1.93	0.50
1:B:4:LEU:O	1:B:5:THR:OG1	2.26	0.50
1:A:117:THR:HG23	1:A:119:GLU:N	2.27	0.50
1:A:161:ILE:HD13	1:A:161:ILE:N	2.27	0.49
1:A:121:VAL:CG2	3:A:1066:HOH:O	2.59	0.49
1:A:5:THR:O	1:A:6:GLU:HG3	2.12	0.49
1:A:37:ARG:HA	1:A:41:GLN:O	2.12	0.49
1:A:44:THR:CG2	1:A:47:GLU:H	2.25	0.49
1:A:170:LYS:O	1:A:173:ARG:HB3	2.13	0.49
1:B:111:ASN:C	1:B:111:ASN:OD1	2.52	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:65:PHE:HB2	3:B:1064:HOH:O	2.13	0.48
1:A:14:GLU:OE1	1:B:164:LYS:HD3	2.14	0.47
1:B:162:ARG:HE	1:B:166:ARG:NH2	2.12	0.47
1:A:27:ILE:HB	1:A:63:ILE:HB	1.95	0.47
1:B:17:SER:HB2	3:B:1094:HOH:O	2.15	0.47
1:B:121:VAL:N	1:B:124:MET:H	2.14	0.46
1:A:44:THR:HG22	1:A:47:GLU:HB2	1.98	0.46
1:A:92:PHE:CE1	1:A:108:VAL:HG11	2.51	0.45
1:B:65:PHE:N	1:B:66:PRO:HD2	2.31	0.45
1:A:6:GLU:HG3	1:A:8:GLN:H	1.82	0.45
1:A:145:MET:CE	1:A:162:ARG:HG3	2.46	0.45
1:A:54:GLU:O	1:A:74:ARG:NH1	2.51	0.44
1:A:109:MET:HA	1:A:109:MET:HE2	1.98	0.44
1:B:52:ILE:HG12	1:B:63:ILE:HG13	2.00	0.44
1:B:52:ILE:HG12	1:B:63:ILE:CD1	2.47	0.44
1:A:92:PHE:CE1	1:A:108:VAL:CG1	3.01	0.44
1:A:112:LEU:O	1:A:112:LEU:HD22	2.18	0.44
1:A:49:GLN:HG3	3:A:1081:HOH:O	2.17	0.44
1:A:63:ILE:HG21	1:A:63:ILE:HD13	1.75	0.43
1:B:107:HIS:CD2	3:B:1043:HOH:O	2.69	0.43
1:B:64:ASP:OD1	1:B:67:GLU:HG3	2.19	0.43
1:A:5:THR:O	1:A:6:GLU:CB	2.66	0.43
1:B:123:GLU:O	1:B:127:GLU:HG3	2.19	0.42
1:A:117:THR:HG23	1:A:119:GLU:H	1.84	0.42
1:B:7:GLU:HG3	1:B:11:GLU:OE1	2.20	0.42
1:B:4:LEU:HG	1:B:9:ILE:HD13	2.01	0.42
1:B:63:ILE:HA	1:B:67:GLU:OE2	2.19	0.42
1:B:105:LEU:HD22	1:B:125:ILE:HD11	2.01	0.42
1:B:76:MET:HG2	3:B:1033:HOH:O	2.20	0.42
1:A:170:LYS:HE3	1:A:170:LYS:HB2	1.94	0.42
1:A:32:LEU:HD11	1:A:51:MET:HE2	2.02	0.42
1:A:94:LYS:HG3	3:A:1008:HOH:O	2.19	0.42
1:B:145:MET:HA	1:B:145:MET:CE	2.49	0.42
1:B:4:LEU:HG	1:B:9:ILE:CD1	2.50	0.41
1:B:144:MET:HE2	3:B:1054:HOH:O	2.19	0.41
1:B:106:ARG:HB2	3:B:1050:HOH:O	2.20	0.41
1:A:19:PHE:HA	1:A:35:VAL:HG21	2.03	0.41
1:A:64:ASP:OD1	1:A:64:ASP:C	2.59	0.41
1:B:76:MET:CG	3:B:1033:HOH:O	2.68	0.40
1:A:102:ALA:HB1	3:A:1066:HOH:O	2.20	0.40
1:B:76:MET:CB	3:B:1033:HOH:O	2.61	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:114:GLU:OE2	1:B:164:LYS:CE	2.70	0.40
1:B:91:VAL:O	1:B:94:LYS:NZ	2.46	0.40
1:A:143:GLN:HB3	1:A:143:GLN:HE21	1.59	0.40
1:B:106:ARG:HG3	1:B:121:VAL:CG1	2.52	0.40
1:A:117:THR:HG23	1:A:120:GLU:H	1.82	0.40
1:B:48:LEU:HD23	1:B:48:LEU:HA	1.90	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	150/179 (84%)	142 (95%)	6 (4%)	2 (1%)	15	10
1	B	153/179 (86%)	144 (94%)	7 (5%)	2 (1%)	15	10
All	All	303/358 (85%)	286 (94%)	13 (4%)	4 (1%)	15	10

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	6	GLU
1	A	145	MET
1	B	122	ASP
1	B	123	GLU

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	133/146 (91%)	112 (84%)	21 (16%)	3 2
1	B	134/146 (92%)	110 (82%)	24 (18%)	2 1
All	All	267/292 (91%)	222 (83%)	45 (17%)	2 1

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLU
1	A	17	SER
1	A	18	LEU
1	A	32	LEU
1	A	36	MET
1	A	44	THR
1	A	69	LEU
1	A	75	LYS
1	A	76	MET
1	A	81	SER
1	A	82	GLU
1	A	90	ARG
1	A	97	ASN
1	A	112	LEU
1	A	143	GLN
1	A	145	MET
1	A	161	ILE
1	A	162	ARG
1	A	166	ARG
1	A	170	LYS
1	A	173	ARG
1	B	4	LEU
1	B	7	GLU
1	B	14	GLU
1	B	18	LEU
1	B	26	THR
1	B	36	MET
1	B	39	LEU
1	B	49	GLN
1	B	69	LEU
1	B	77	LYS
1	B	86	ARG
1	B	87	GLU
1	B	90	ARG

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Mol	Chain	Res	Type
1	B	94	LYS
1	B	97	ASN
1	B	122	ASP
1	B	123	GLU
1	B	124	MET
1	B	143	GLN
1	B	157	ARG
1	B	158	LYS
1	B	170	LYS
1	B	173	ARG
1	B	176	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	ASN
1	A	49	GLN
1	A	97	ASN
1	A	111	ASN
1	A	135	GLN
1	A	143	GLN
1	B	49	GLN
1	B	53	ASN
1	B	97	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	156/179 (87%)	0.64	9 (5%) 26 29	27, 32, 43, 50	0
1	B	159/179 (88%)	0.54	6 (3%) 44 46	21, 33, 45, 62	0
All	All	315/358 (87%)	0.59	15 (4%) 34 36	21, 33, 45, 62	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	156	ALA	4.7
1	A	46	ALA	4.5
1	B	40	GLY	4.4
1	A	147	ALA	4.4
1	B	158	LYS	3.2
1	A	175	PHE	2.8
1	A	162	ARG	2.7
1	A	7	GLU	2.6
1	A	146	THR	2.5
1	B	5	THR	2.4
1	A	163	ASN	2.2
1	A	44	THR	2.2
1	A	41	GLN	2.1
1	B	56	ASP	2.1
1	B	176	SER	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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	CA	A	1004	1/1	0.98	0.24	4.49	33,33,33,33	0
2	CA	A	1001	1/1	0.96	0.21	2.44	36,36,36,36	0
2	CA	B	1001	1/1	0.96	0.23	1.84	62,62,62,62	0
2	CA	B	1004	1/1	0.99	0.19	1.41	34,34,34,34	0
2	CA	B	1002	1/1	0.94	0.18	0.55	48,48,48,48	0
2	CA	A	1002	1/1	0.99	0.14	0.24	44,44,44,44	0
2	CA	B	1003	1/1	0.99	0.12	-1.05	39,39,39,39	0
2	CA	A	1003	1/1	0.99	0.11	-1.32	32,32,32,32	0

6.5 Other polymers

There are no such residues in this entry.