



wwPDB X-ray Structure Validation Summary Report ⓘ

May 21, 2020 – 05:18 am BST

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 wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The 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

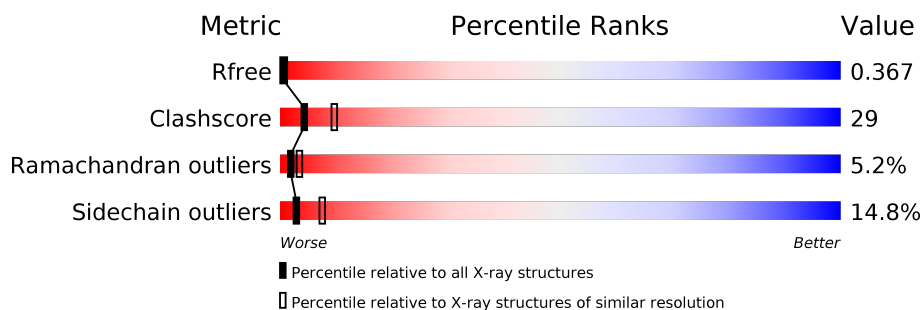
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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}	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)

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 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 $\leq 5\%$

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

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2715 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, Linker, IQ motif of Neuromodulin.

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

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	O	0	0
			40	40		
2	B	29	Total	O	0	0
			29	29		

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



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	79.39 Å 79.28 Å 136.06 Å 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
$\langle I/\sigma(I) \rangle$ ¹	2.44 (at 2.68 Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.276 , 0.318 0.325 , 0.367	Depositor DCC
R_{free} test set	1324 reflections (5.68%)	wwPDB-VP
Wilson B-factor (Å ²)	59.3	Xtriage
Anisotropy	0.097	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 16.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	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
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 ²)	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.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [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 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	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

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.

The worst 5 of 151 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: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:A:100:TYR:HD2	2:A:214:HOH:O	1.56	0.88

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	164/185 (89%)	130 (79%)	25 (15%)	9 (6%)	2	3
1	B	164/185 (89%)	121 (74%)	35 (21%)	8 (5%)	2	4
All	All	328/370 (89%)	251 (76%)	60 (18%)	17 (5%)	2	3

5 of 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

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	142/152 (93%)	118 (83%)	24 (17%)	2	5
1	B	142/152 (93%)	124 (87%)	18 (13%)	4	10
All	All	284/304 (93%)	242 (85%)	42 (15%)	3	7

5 of 42 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	130	ASP
1	A	158	LYS
1	B	147	THR
1	A	131	ILE
1	A	141	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

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

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 ⓘ

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 ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.