



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:07 PM UTC

PDB ID : 3CLN / pdb_00003cln
Title : STRUCTURE OF CALMODULIN REFINED AT 2.2 ANGSTROMS RESOLUTION
Authors : Babu, Y.S.; Bugg, C.E.; Cook, W.J.
Deposited on : 1988-05-11
Resolution : 2.20 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

<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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

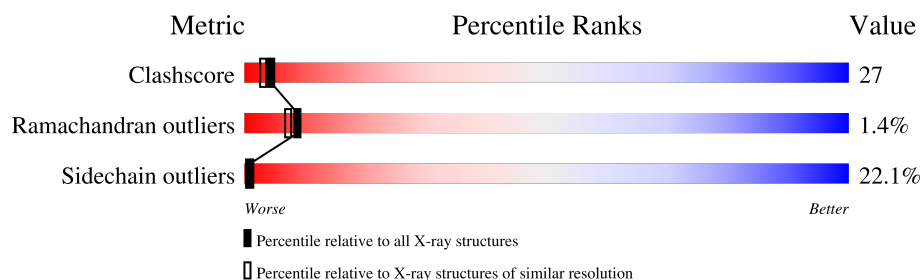
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)

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 ≥ 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	148	 46% 30% 16% 5% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1199 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	143	Total	C	N	O	S	0	0	0
			1126	690	182	245	9			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

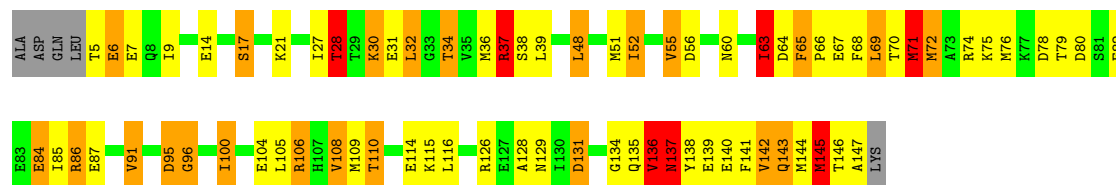
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	Ca	0	0
			4	4		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	69	Total	O	0	0
			69	69		

Note EDS was not executed.

Chain A: 46% 30% 16% 5%



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	29.71Å 53.79Å 24.99Å 94.13° 97.57° 89.46°	Depositor
Resolution (Å)	(Not available) – 2.20	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.175 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1199	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	1.24	1/1138 (0.1%)	2.07	35/1528 (2.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	52	ILE	C-O	5.64	1.30	1.24

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	17	SER	CA-C-O	-7.86	111.22	120.10
1	A	28	THR	N-CA-CB	-7.61	98.69	110.56
1	A	65	PHE	CA-CB-CG	-7.39	106.41	113.80
1	A	32	LEU	CA-C-N	6.74	127.42	119.94
1	A	32	LEU	C-N-CA	6.74	127.42	119.94
1	A	115	LYS	CA-C-O	6.74	128.11	120.84
1	A	34	THR	CA-C-N	6.46	128.84	120.56
1	A	34	THR	C-N-CA	6.46	128.84	120.56
1	A	37	ARG	CD-NE-CZ	-6.35	115.50	124.40
1	A	27	ILE	CA-C-O	6.34	127.05	120.39
1	A	71	MET	CA-C-N	6.19	128.49	120.44
1	A	71	MET	C-N-CA	6.19	128.49	120.44
1	A	60	ASN	N-CA-C	-5.96	105.96	113.18
1	A	52	ILE	CB-CG1-CD1	5.89	126.18	113.80
1	A	21	LYS	N-CA-C	5.86	117.67	111.28
1	A	131	ASP	CA-C-N	5.77	128.82	120.11
1	A	131	ASP	C-N-CA	5.77	128.82	120.11
1	A	63	ILE	CA-CB-CG2	5.73	120.23	110.50
1	A	6	GLU	CB-CG-CD	5.72	122.33	112.60
1	A	28	THR	CA-CB-CG2	5.71	120.21	110.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	38	SER	N-CA-C	-5.62	105.53	112.38
1	A	95	ASP	CA-CB-CG	5.55	118.15	112.60
1	A	48	LEU	N-CA-C	-5.46	105.23	111.07
1	A	145	MET	CA-C-N	5.37	127.75	120.44
1	A	145	MET	C-N-CA	5.37	127.75	120.44
1	A	136	VAL	CB-CA-C	5.33	118.34	110.77
1	A	86	ARG	CD-NE-CZ	5.32	131.85	124.40
1	A	115	LYS	CA-C-N	5.30	130.54	122.65
1	A	115	LYS	C-N-CA	5.30	130.54	122.65
1	A	85	ILE	CA-C-O	-5.22	115.31	120.85
1	A	100	ILE	CA-C-O	-5.20	114.41	120.75
1	A	28	THR	OG1-CB-CG2	5.15	119.60	109.30
1	A	7	GLU	CA-CB-CG	5.11	124.32	114.10
1	A	106	ARG	CA-C-O	-5.04	115.51	120.70
1	A	137	ASN	CB-CA-C	5.01	119.12	110.45

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	A	1126	0	1054	58	2
2	A	4	0	0	0	0
3	A	69	0	0	0	0
All	All	1199	0	1054	58	2

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 27.

All (58) 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:28:THR:HG23	1:A:30:LYS:HD2	1.20	1.13
1:A:100:ILE:HB	1:A:136:VAL:HG13	1.24	1.10

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:PRO:O	1:A:70:THR:HG23	1.66	0.93
1:A:63:ILE:HD11	1:A:68:PHE:HB2	1.52	0.92
1:A:28:THR:CG2	1:A:30:LYS:HD2	2.00	0.91
1:A:129:ASN:HD21	1:A:134:GLY:H	1.19	0.90
1:A:28:THR:HG22	1:A:31:GLU:H	1.38	0.89
1:A:75:LYS:O	1:A:75:LYS:HG2	1.80	0.81
1:A:129:ASN:ND2	1:A:134:GLY:H	1.86	0.73
1:A:82:GLU:OE1	1:A:86:ARG:NH1	2.22	0.73
1:A:105:LEU:O	1:A:109:MET:HB2	1.94	0.68
1:A:63:ILE:HD11	1:A:68:PHE:CB	2.24	0.67
1:A:100:ILE:HB	1:A:136:VAL:CG1	2.16	0.66
1:A:75:LYS:O	1:A:79:THR:HG22	1.95	0.66
1:A:63:ILE:CD1	1:A:68:PHE:N	2.62	0.62
1:A:137:ASN:ND2	1:A:140:GLU:H	1.97	0.61
1:A:71:MET:SD	1:A:72:MET:SD	2.99	0.61
1:A:138:TYR:O	1:A:142:VAL:HG13	2.03	0.59
1:A:100:ILE:CB	1:A:136:VAL:HG13	2.17	0.57
1:A:63:ILE:HD12	1:A:67:GLU:HB2	1.85	0.57
1:A:28:THR:HG23	1:A:30:LYS:H	1.70	0.56
1:A:74:ARG:HG3	1:A:74:ARG:NH1	2.19	0.56
1:A:28:THR:HG23	1:A:30:LYS:CD	2.14	0.55
1:A:137:ASN:HD22	1:A:140:GLU:H	1.54	0.55
1:A:75:LYS:O	1:A:75:LYS:CG	2.53	0.54
1:A:28:THR:CG2	1:A:30:LYS:CD	2.84	0.52
1:A:74:ARG:HG3	1:A:74:ARG:HH11	1.75	0.52
1:A:63:ILE:HD11	1:A:68:PHE:N	2.24	0.51
1:A:63:ILE:HD11	1:A:68:PHE:CA	2.40	0.51
1:A:84:GLU:O	1:A:87:GLU:HB2	2.11	0.51
1:A:146:THR:OG1	1:A:147:ALA:N	2.45	0.50
1:A:63:ILE:HA	1:A:67:GLU:OE1	2.12	0.49
1:A:131:ASP:C	1:A:131:ASP:OD2	2.53	0.49
1:A:65:PHE:HB3	1:A:66:PRO:HD3	1.93	0.49
1:A:95:ASP:O	1:A:96:GLY:C	2.56	0.49
1:A:78:ASP:OD2	1:A:78:ASP:N	2.47	0.47
1:A:36:MET:HE2	1:A:51:MET:SD	2.55	0.46
1:A:30:LYS:HD2	1:A:30:LYS:H	1.80	0.46
1:A:137:ASN:HD22	1:A:137:ASN:C	2.24	0.46
1:A:100:ILE:O	1:A:136:VAL:HG12	2.17	0.45
1:A:80:ASP:O	1:A:84:GLU:OE1	2.35	0.45
1:A:32:LEU:HD22	1:A:52:ILE:HD12	1.99	0.44
1:A:128:ALA:HB2	1:A:144:MET:HG2	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:ILE:HD12	1:A:64:ASP:O	2.17	0.43
1:A:136:VAL:HA	1:A:140:GLU:OE1	2.18	0.43
1:A:63:ILE:CD1	1:A:68:PHE:H	2.30	0.43
1:A:104:GLU:O	1:A:108:VAL:HG13	2.18	0.43
1:A:55:VAL:CG2	1:A:63:ILE:HG21	2.48	0.43
1:A:65:PHE:CZ	1:A:69:LEU:HD23	2.54	0.43
1:A:55:VAL:HG21	1:A:63:ILE:HG21	2.01	0.42
1:A:106:ARG:O	1:A:110:THR:CG2	2.67	0.42
1:A:82:GLU:CG	1:A:86:ARG:HH11	2.33	0.42
1:A:141:PHE:CZ	1:A:145:MET:SD	3.13	0.41
1:A:37:ARG:C	1:A:39:LEU:N	2.79	0.41
1:A:68:PHE:O	1:A:72:MET:HB2	2.21	0.40
1:A:139:GLU:O	1:A:143:GLN:CG	2.69	0.40
1:A:71:MET:CE	1:A:72:MET:SD	3.10	0.40
1:A:128:ALA:HA	1:A:144:MET:HE3	2.02	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:VAL:O	1:A:126:ARG:NH1[1_556]	2.01	0.19
1:A:37:ARG:NH2	1:A:114:GLU:OE1[1_566]	2.11	0.09

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	141/148 (95%)	129 (92%)	10 (7%)	2 (1%)	9 7

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	56	ASP
1	A	96	GLY

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	122/126 (97%)	95 (78%)	27 (22%)	1 1

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	A	6	GLU
1	A	9	ILE
1	A	14	GLU
1	A	17	SER
1	A	28	THR
1	A	30	LYS
1	A	34	THR
1	A	37	ARG
1	A	48	LEU
1	A	55	VAL
1	A	63	ILE
1	A	69	LEU
1	A	71	MET
1	A	72	MET
1	A	76	MET
1	A	84	GLU
1	A	91	VAL
1	A	108	VAL
1	A	110	THR
1	A	116	LEU
1	A	135	GLN
1	A	136	VAL
1	A	137	ASN
1	A	142	VAL

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Mol	Chain	Res	Type
1	A	143	GLN
1	A	145	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	41	GLN
1	A	49	GLN
1	A	53	ASN
1	A	111	ASN
1	A	137	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](#)

Of 4 ligands modelled in this entry, 4 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 [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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.