Full wwPDB X-ray Structure Validation Report (i)

Jan 27, 2024 – 12:41 PM EST

PDB ID	:	1A8Y
Title	:	CRYSTAL STRUCTURE OF CALSEQUESTRIN FROM RABBIT SKELETAL MUSCLE SARCOPLASMIC RETICULUM AT 2.4 A RESOLUTION
Authors	:	Wang, S.; Trumble, W.R.; Liao, H.; Wesson, C.R.; Dunker, A.K.; Kang, C.
Deposited on	:	1998-03-31
Resolution	:	2.40 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 (i) 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 (i)) were used in the production of this report:

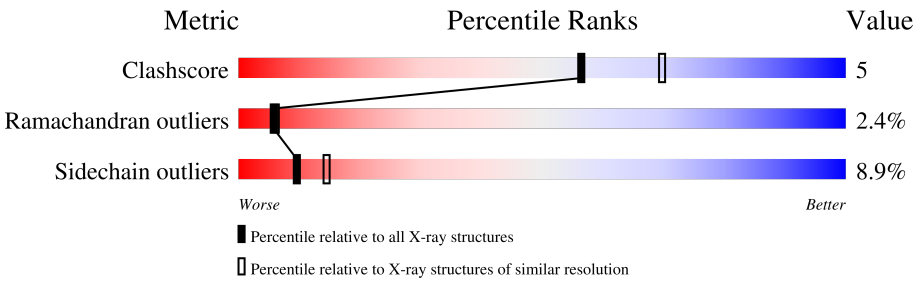
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.40 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	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	367	69% 18% • 8%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 2800 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 CALSEQUESTRIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	338	2760	1775	413	567	5	0	0	0

- Molecule 2 is water.

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

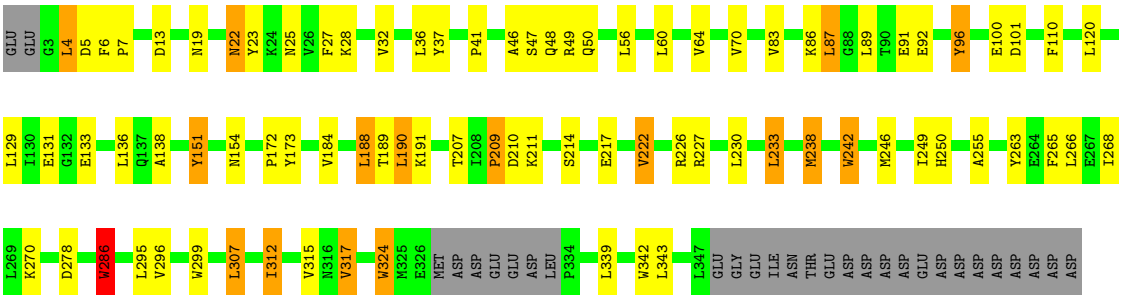
3 Residue-property plots (i)

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

Note EDS was not executed.

• Molecule 1: CALSEQUESTRIN

Chain A: 69% 18% • 8%



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	59.74Å 145.56Å 111.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.40	Depositor
% Data completeness (in resolution range)	78.3 (10.00-2.40)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	0.04	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.188 , 0.242	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2800	wwPDB-VP
Average B, all atoms (Å^2)	17.0	wwPDB-VP

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	0.83	0/2824	1.50	35/3832 (0.9%)

There are no bond length outliers.

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	342	TRP	CD1-CG-CD2	9.23	113.69	106.30
1	A	151	TYR	CB-CG-CD2	-8.86	115.68	121.00
1	A	242	TRP	CD1-CG-CD2	7.98	112.68	106.30
1	A	286	TRP	CE2-CD2-CG	-7.66	101.17	107.30
1	A	342	TRP	CE2-CD2-CG	-7.64	101.19	107.30
1	A	242	TRP	CE2-CD2-CG	-7.60	101.22	107.30
1	A	96	TYR	CB-CG-CD2	-7.37	116.58	121.00
1	A	324	TRP	CE2-CD2-CG	-7.35	101.42	107.30
1	A	210	ASP	CA-C-N	-7.33	101.08	117.20
1	A	286	TRP	CB-CG-CD1	-7.11	117.76	127.00
1	A	37	TYR	CB-CG-CD2	-6.87	116.88	121.00
1	A	324	TRP	CD1-CG-CD2	6.76	111.71	106.30
1	A	317	VAL	CA-CB-CG2	-6.71	100.83	110.90
1	A	286	TRP	NE1-CE2-CZ2	-6.67	123.06	130.40
1	A	64	VAL	CG1-CB-CG2	-6.62	100.31	110.90
1	A	299	TRP	CE2-CD2-CG	-6.25	102.30	107.30
1	A	286	TRP	CA-CB-CG	6.22	125.52	113.70
1	A	295	LEU	CA-CB-CG	5.94	128.96	115.30
1	A	286	TRP	CG-CD2-CE3	5.88	139.19	133.90
1	A	286	TRP	CD1-CG-CD2	5.79	110.93	106.30
1	A	209	PRO	N-CA-C	5.69	126.89	112.10
1	A	324	TRP	CB-CG-CD1	-5.64	119.67	127.00
1	A	242	TRP	CG-CD2-CE3	5.62	138.96	133.90
1	A	13	ASP	CA-C-N	-5.60	104.88	117.20
1	A	342	TRP	CG-CD1-NE1	-5.49	104.61	110.10
1	A	238	MET	N-CA-C	5.45	125.73	111.00
1	A	22	ASN	N-CA-C	5.44	125.69	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	4	LEU	CA-CB-CG	5.40	127.71	115.30
1	A	317	VAL	CA-CB-CG1	5.39	118.99	110.90
1	A	263	TYR	CB-CG-CD2	-5.36	117.78	121.00
1	A	222	VAL	CG1-CB-CG2	-5.29	102.44	110.90
1	A	286	TRP	NE1-CE2-CD2	5.14	112.44	107.30
1	A	173	TYR	CB-CG-CD2	-5.11	117.93	121.00
1	A	188	LEU	CB-CG-CD2	-5.01	102.47	111.00
1	A	49	ARG	NE-CZ-NH2	-5.00	117.80	120.30

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	2760	0	2611	27	0
2	A	40	0	0	2	0
All	All	2800	0	2611	27	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 5.

All (27) 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:255:ALA:O	1:A:286:TRP:HZ3	1.77	0.67
1:A:87:LEU:HB3	1:A:89:LEU:HD13	1.80	0.64
1:A:296:VAL:HB	1:A:307:LEU:HD23	1.80	0.63
1:A:41:PRO:HB2	1:A:48:GLN:HG2	1.81	0.63
1:A:214:SER:OG	1:A:217:GLU:HG3	2.01	0.60
1:A:46:ALA:O	1:A:50:GLN:HG3	2.03	0.58
1:A:227:ARG:HH12	1:A:246:MET:HB3	1.71	0.56
1:A:242:TRP:O	1:A:250:HIS:HE1	1.90	0.55
1:A:19:ASN:H	1:A:22:ASN:HB2	1.72	0.54
1:A:154:ASN:HB2	2:A:519:HOH:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:LEU:O	1:A:270:LYS:HG3	2.14	0.48
1:A:25:ASN:HA	1:A:28:LYS:HG2	1.98	0.46
1:A:151:TYR:HB2	1:A:190:LEU:HD23	1.98	0.45
1:A:136:LEU:HD12	1:A:184:VAL:HG22	1.98	0.45
1:A:91:GLU:HG2	1:A:96:TYR:OH	2.18	0.44
1:A:191:LYS:HD2	1:A:211:LYS:NZ	2.33	0.43
1:A:246:MET:O	1:A:249:ILE:HG22	2.19	0.42
1:A:32:VAL:O	1:A:70:VAL:HA	2.20	0.42
1:A:83:VAL:O	1:A:87:LEU:HB2	2.20	0.42
1:A:265:PHE:HA	1:A:268:ILE:HD12	2.02	0.42
1:A:6:PHE:HB3	2:A:501:HOH:O	2.20	0.41
1:A:222:VAL:O	1:A:226:ARG:HG3	2.20	0.41
1:A:23:TYR:O	1:A:27:PHE:HD2	2.03	0.41
1:A:233:LEU:HD23	1:A:233:LEU:HA	1.79	0.41
1:A:6:PHE:HA	1:A:7:PRO:HD3	1.90	0.40
1:A:266:LEU:HD12	1:A:266:LEU:HA	1.89	0.40
1:A:265:PHE:HZ	1:A:312:ILE:HB	1.87	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	A	334/367 (91%)	313 (94%)	13 (4%)	8 (2%)	6 6

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	92	GLU
1	A	133	GLU
1	A	189	THR
1	A	190	LEU

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Mol	Chain	Res	Type
1	A	100	GLU
1	A	238	MET
1	A	138	ALA
1	A	209	PRO

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	303/331 (92%)	276 (91%)	27 (9%)	9 14

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

Mol	Chain	Res	Type
1	A	4	LEU
1	A	5	ASP
1	A	36	LEU
1	A	47	SER
1	A	56	LEU
1	A	60	LEU
1	A	86	LYS
1	A	87	LEU
1	A	101	ASP
1	A	110	PHE
1	A	120	LEU
1	A	129	LEU
1	A	131	GLU
1	A	172	PRO
1	A	188	LEU
1	A	207	THR
1	A	230	LEU
1	A	233	LEU
1	A	278	ASP
1	A	286	TRP
1	A	307	LEU
1	A	312	ILE

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Mol	Chain	Res	Type
1	A	315	VAL
1	A	317	VAL
1	A	324	TRP
1	A	339	LEU
1	A	343	LEU

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	50	GLN
1	A	171	HIS
1	A	193	ASN
1	A	220	ASN
1	A	250	HIS
1	A	311	GLN

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 monosaccharides 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 (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.