



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 04:21 pm BST

PDB ID : 4C92  
Title : Crystal structure of the yeast Lsm1-7 complex  
Authors : Sharif, H.; Conti, E.  
Deposited on : 2013-10-02  
Resolution : 2.30 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](mailto: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.

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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  
Xtrriage (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

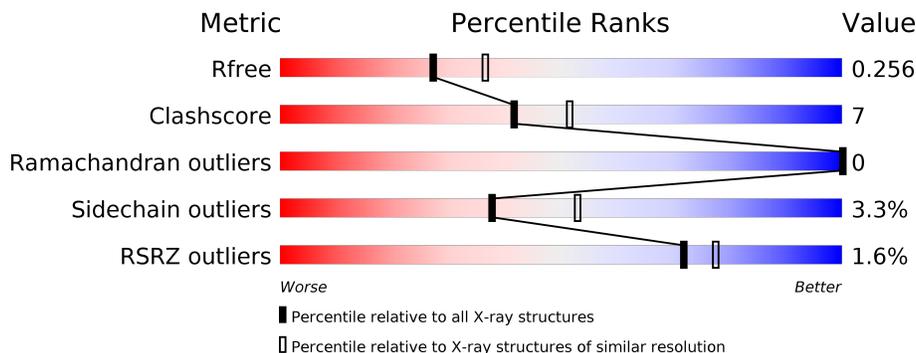
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 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  $\leq 5\%$ . 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	146	 72% 17% 11%
2	B	105	 3% 81% 18%
3	C	89	 3% 71% 15% 11%
4	D	114	 61% 11% 26%
5	E	93	 75% 15% 10%
6	F	86	 78% 12% 10%

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Mol	Chain	Length	Quality of chain
7	G	115	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into three segments: a green segment on the left labeled '56%', a yellow segment in the middle labeled '10%', and a grey segment on the right labeled '35%'.</p>

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 4999 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 SM-LIKE PROTEIN LSM1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	130	1067	678	185	197	7	0	0	0

- Molecule 2 is a protein called U6 SNRNA-ASSOCIATED SM-LIKE PROTEIN LSM2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	105	843	536	143	161	3	0	0	0

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-9	SER	-	expression tag	UNP P38203
B	-8	GLU	-	expression tag	UNP P38203
B	-7	ASN	-	expression tag	UNP P38203
B	-6	LEU	-	expression tag	UNP P38203
B	-5	TYR	-	expression tag	UNP P38203
B	-4	PHE	-	expression tag	UNP P38203
B	-3	GLN	-	expression tag	UNP P38203
B	-2	GLY	-	expression tag	UNP P38203
B	-1	SER	-	expression tag	UNP P38203
B	0	GLY	-	expression tag	UNP P38203
B	1	SER	-	expression tag	UNP P38203

- Molecule 3 is a protein called U6 SNRNA-ASSOCIATED SM-LIKE PROTEIN LSM3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	79	584	366	101	113	4	0	0	0

- Molecule 4 is a protein called U6 SNRNA-ASSOCIATED SM-LIKE PROTEIN LSM4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	84	Total	C	N	O	S	0	0	0
			623	397	104	120	2			

- Molecule 5 is a protein called U6 SNRNA-ASSOCIATED SM-LIKE PROTEIN LSM5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	84	Total	C	N	O	S	0	0	0
			639	407	106	124	2			

- Molecule 6 is a protein called U6 SNRNA-ASSOCIATED SM-LIKE PROTEIN LSM6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	77	Total	C	N	O	S	0	0	0
			587	370	97	118	2			

- Molecule 7 is a protein called U6 SNRNA-ASSOCIATED SM-LIKE PROTEIN LSM7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	75	Total	C	N	O	S	0	0	0
			548	349	92	104	3			

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	20	Total	O	0	0
			20	20		
8	B	21	Total	O	0	0
			21	21		
8	C	9	Total	O	0	0
			9	9		
8	D	17	Total	O	0	0
			17	17		
8	E	20	Total	O	0	0
			20	20		
8	F	11	Total	O	0	0
			11	11		
8	G	10	Total	O	0	0
			10	10		



- Molecule 6: U6 SNRNA-ASSOCIATED SM-LIKE PROTEIN LSM6

Chain F:  78% 12% 10%



- Molecule 7: U6 SNRNA-ASSOCIATED SM-LIKE PROTEIN LSM7

Chain G:  56% 10% 35%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.80Å 90.57Å 68.46Å 90.00° 100.80° 90.00°	Depositor
Resolution (Å)	53.99 – 2.30 67.25 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.5 (53.99-2.30) 99.3 (67.25-2.30)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.89 (at 2.29Å)	Xtrriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.211 , 0.259 0.214 , 0.256	Depositor DCC
$R_{free}$ test set	1641 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	46.2	Xtrriage
Anisotropy	0.652	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 55.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4999	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.70% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical 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	0.56	1/1086 (0.1%)	0.62	0/1455
2	B	0.70	0/855	0.64	0/1153
3	C	0.60	0/588	0.79	3/798 (0.4%)
4	D	0.58	0/629	0.66	1/853 (0.1%)
5	E	0.46	0/648	0.60	0/882
6	F	0.69	0/594	0.66	0/803
7	G	0.48	0/549	0.69	0/737
All	All	0.59	1/4949 (0.0%)	0.66	4/6681 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	82	CYS	CB-SG	-5.35	1.73	1.81

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	75	LEU	CA-CB-CG	6.67	130.63	115.30
4	D	56	ASP	N-CA-C	-5.34	96.58	111.00
3	C	43	ASP	CB-CG-OD2	5.20	122.97	118.30
3	C	75	LEU	CB-CG-CD1	5.08	119.64	111.00

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	1067	0	1039	16	0
2	B	843	0	827	11	0
3	C	584	0	575	10	0
4	D	623	0	605	9	0
5	E	639	0	619	11	0
6	F	587	0	564	5	0
7	G	548	0	568	13	0
8	A	20	0	0	0	0
8	B	21	0	0	0	0
8	C	9	0	0	1	0
8	D	17	0	0	1	0
8	E	20	0	0	0	0
8	F	11	0	0	1	0
8	G	10	0	0	0	0
All	All	4999	0	4797	68	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 7.

All (68) 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)
5:E:75:GLY:HA2	5:E:78:ILE:HD12	1.51	0.93
1:A:156:LYS:HE2	3:C:54:GLU:HG3	1.51	0.90
7:G:34:LYS:HE3	7:G:55:TYR:OH	1.80	0.82
7:G:31:ALA:HA	7:G:34:LYS:HE3	1.71	0.71
7:G:34:LYS:HE2	7:G:55:TYR:CE1	2.25	0.71
1:A:52:ARG:HD2	1:A:114:GLU:OE2	1.93	0.69
7:G:34:LYS:CE	7:G:55:TYR:OH	2.41	0.69
2:B:21:ASN:HD21	2:B:23:ILE:CD1	2.06	0.67
2:B:21:ASN:O	2:B:22:ASP:HB2	1.94	0.67
5:E:75:GLY:HA2	5:E:78:ILE:CD1	2.25	0.66
5:E:22:ILE:HG21	5:E:73:LEU:HD21	1.81	0.62
6:F:20:ILE:HD11	6:F:41:ILE:HG22	1.80	0.61
2:B:-5:TYR:HE1	2:B:9:THR:HG21	1.68	0.59
1:A:122:GLN:N	1:A:123:PRO:HD2	2.19	0.58
4:D:29:LEU:HD21	4:D:38:LEU:HD22	1.86	0.57
2:B:84:ASP:OD2	2:B:88:ARG:NH1	2.40	0.55
5:E:57:ASP:N	5:E:57:ASP:OD1	2.38	0.55
2:B:21:ASN:HD21	2:B:23:ILE:HD11	1.72	0.54
4:D:11:LYS:NZ	4:D:31:ASN:OD1	2.41	0.54
7:G:31:ALA:HA	7:G:34:LYS:HG3	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:34:LYS:HE2	7:G:55:TYR:HE1	1.71	0.53
4:D:73:GLY:HA2	4:D:76:ILE:HD12	1.91	0.53
6:F:37:ARG:HB3	6:F:50:SER:OG	2.09	0.53
1:A:59:ARG:NH1	1:A:170:ASP:OD1	2.41	0.52
5:E:23:VAL:HB	5:E:80:ILE:HB	1.91	0.51
1:A:54:ILE:HD11	1:A:79:LEU:HD21	1.93	0.51
3:C:6:ASP:O	3:C:9:LYS:HG2	2.11	0.51
7:G:34:LYS:CE	7:G:55:TYR:CE1	2.94	0.50
1:A:164:TYR:O	1:A:165:ASP:C	2.50	0.50
4:D:1:MET:N	8:D:2001:HOH:O	2.44	0.49
5:E:22:ILE:HG21	5:E:73:LEU:CD2	2.42	0.49
1:A:105:ARG:NH2	1:A:107:GLU:OE2	2.45	0.49
3:C:60:GLU:N	8:C:2007:HOH:O	2.42	0.48
3:C:51:LEU:HD13	3:C:54:GLU:HA	1.97	0.47
2:B:25:ILE:HG21	2:B:60:ILE:HD11	1.97	0.46
7:G:38:ILE:HD11	7:G:63:LEU:HD21	1.97	0.46
4:D:37:ASN:OD1	4:D:72:ARG:HD3	2.17	0.45
2:B:-4:PHE:CD2	2:B:88:ARG:HD3	2.52	0.45
5:E:75:GLY:CA	5:E:78:ILE:HD12	2.35	0.44
6:F:35:SER:O	6:F:52:ALA:HA	2.18	0.44
2:B:58:ARG:HA	2:B:58:ARG:HD3	1.82	0.44
7:G:95:ARG:HG3	7:G:95:ARG:HH11	1.83	0.44
5:E:67:HIS:HD2	8:F:2002:HOH:O	2.00	0.43
1:A:140:GLN:NE2	4:D:49:GLU:OE2	2.41	0.43
3:C:7:LEU:O	3:C:10:LEU:HB2	2.19	0.43
7:G:41:LYS:HB3	7:G:100:VAL:HG23	2.00	0.43
2:B:20:LYS:HE3	3:C:72:THR:HG22	2.00	0.42
1:A:156:LYS:HB3	3:C:51:LEU:HD11	2.01	0.42
3:C:20:LEU:HD11	3:C:26:LEU:HD12	2.00	0.42
5:E:70:ARG:HB2	6:F:15:PHE:HE1	1.85	0.42
7:G:42:LEU:HD22	7:G:98:ILE:HD12	2.01	0.42
4:D:11:LYS:HA	4:D:29:LEU:HD12	2.01	0.42
5:E:82:VAL:HG22	7:G:92:THR:HG22	2.00	0.42
1:A:147:PHE:O	1:A:151:THR:HG23	2.20	0.42
1:A:97:GLU:HG2	1:A:99:ARG:HH12	1.85	0.42
2:B:37:ASN:HB3	2:B:61:PHE:CE1	2.55	0.42
2:B:14:GLU:OE2	2:B:26:LYS:HD3	2.20	0.42
3:C:18:ILE:HD12	3:C:41:LEU:HD11	2.01	0.42
3:C:53:ASN:O	3:C:54:GLU:HB3	2.19	0.42
6:F:38:LEU:HA	6:F:49:LEU:HD23	2.00	0.42
1:A:70:THR:HB	1:A:78:ILE:HD12	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:29:LEU:HD11	4:D:32:VAL:CG2	2.50	0.41
1:A:136:TRP:CE3	4:D:52:ILE:HG21	2.55	0.41
1:A:97:GLU:HG2	1:A:99:ARG:NH1	2.35	0.41
1:A:156:LYS:HD3	1:A:156:LYS:HA	1.84	0.41
7:G:34:LYS:CE	7:G:55:TYR:CZ	3.04	0.41
5:E:44:ILE:CD1	5:E:72:LEU:HD13	2.51	0.40
1:A:122:GLN:N	1:A:123:PRO:CD	2.83	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	128/146 (88%)	127 (99%)	1 (1%)	0	100	100
2	B	103/105 (98%)	101 (98%)	2 (2%)	0	100	100
3	C	77/89 (86%)	76 (99%)	1 (1%)	0	100	100
4	D	82/114 (72%)	82 (100%)	0	0	100	100
5	E	82/93 (88%)	79 (96%)	3 (4%)	0	100	100
6	F	75/86 (87%)	72 (96%)	3 (4%)	0	100	100
7	G	71/115 (62%)	71 (100%)	0	0	100	100
All	All	618/748 (83%)	608 (98%)	10 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	111/130 (85%)	108 (97%)	3 (3%)	44	61
2	B	93/99 (94%)	88 (95%)	5 (5%)	22	30
3	C	62/81 (76%)	58 (94%)	4 (6%)	17	23
4	D	62/104 (60%)	60 (97%)	2 (3%)	39	54
5	E	69/84 (82%)	67 (97%)	2 (3%)	42	58
6	F	64/75 (85%)	63 (98%)	1 (2%)	62	78
7	G	57/103 (55%)	57 (100%)	0	100	100
All	All	518/676 (77%)	501 (97%)	17 (3%)	38	53

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	ASP
1	A	98	ASP
1	A	115	VAL
2	B	9	THR
2	B	41	ASP
2	B	48	GLU
2	B	72	LEU
2	B	89	GLU
3	C	26	LEU
3	C	37	CYS
3	C	51	LEU
3	C	75	LEU
4	D	43	VAL
4	D	52	ILE
5	E	74	SER
5	E	76	ASN
6	F	17	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	21	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 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 [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	130/146 (89%)	0.13	2 (1%) 73 79	19, 39, 85, 112	0
2	B	105/105 (100%)	0.11	3 (2%) 51 58	21, 42, 94, 114	0
3	C	79/89 (88%)	0.21	3 (3%) 40 47	21, 42, 85, 132	0
4	D	84/114 (73%)	0.12	1 (1%) 79 83	16, 38, 92, 127	0
5	E	84/93 (90%)	0.04	1 (1%) 79 83	19, 41, 87, 138	0
6	F	77/86 (89%)	-0.02	0 100 100	20, 38, 64, 77	0
7	G	75/115 (65%)	-0.04	0 100 100	12, 33, 73, 123	0
All	All	634/748 (84%)	0.09	10 (1%) 72 77	12, 40, 88, 138	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	163	VAL	3.6
3	C	54	GLU	3.4
3	C	53	ASN	3.3
3	C	56	LEU	3.2
2	B	49	LYS	3.2
4	D	61	SER	3.1
5	E	56	GLU	2.9
2	B	-1	SER	2.4
2	B	95	LYS	2.3
1	A	144	GLU	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.