



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2018 – 10:00 pm GMT

PDB ID : 4C8Q  
Title : Crystal structure of the yeast Lsm1-7-Pat1 complex  
Authors : Sharif, H.; Conti, E.  
Deposited on : 2013-10-01  
Resolution : 3.70 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  
Xtriage (Phenix) : 1.13  
EDS : trunk30967  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk30967

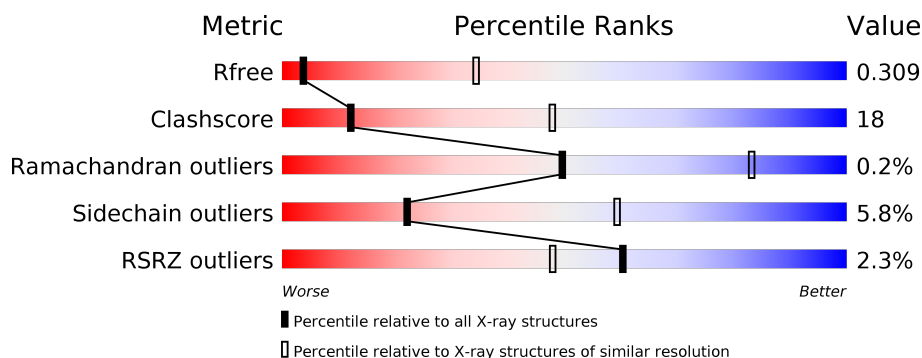
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.70 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))
$R_{free}$	111664	1413 (3.90-3.50)
Clashscore	122126	1524 (3.90-3.50)
Ramachandran outliers	120053	1470 (3.90-3.50)
Sidechain outliers	120020	1467 (3.90-3.50)
RSRZ outliers	108989	1298 (3.90-3.50)

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. 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	101	<div> <div>8%</div> <div>68%</div> <div>20%</div> <div>7%</div> </div>
2	B	105	<div> <div>%</div> <div>69%</div> <div>24%</div> <div>7%</div> </div>
3	C	89	<div> <div>62%</div> <div>22%</div> <div>11%</div> </div>
4	D	114	<div> <div>7%</div> <div>58%</div> <div>11%</div> <div>29%</div> </div>
5	E	93	<div> <div>65%</div> <div>23%</div> <div>11%</div> </div>
6	F	86	<div> <div>65%</div> <div>23%</div> <div>10%</div> </div>

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Mol	Chain	Length	Quality of chain
7	G	115	
8	H	328	

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
9	CO	H	1080	-	-	-	X

## 2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 6280 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
1	A	94	Total	C	N	O	S	0	0	0
			673	428	115	125	5			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	104	Total	C	N	O	S	0	0	0
			796	505	134	154	3			

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
3	C	79	Total	C	N	O	S	0	0	0
			589	369	105	112	3			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	81	Total	C	N	O	S	0	0	0
			504	321	85	97	1			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	83	Total	C	N	O	S	0	0	0
			600	383	104	111	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
			571	361	95	113	2			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	69	Total	C	N	O	S	0	0	0
			456	289	79	87	1			

- Molecule 8 is a protein called DNA TOPOISOMERASE 2-ASSOCIATED PROTEIN PAT1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	279	Total	C	N	O	S	0	0	0
			2089	1366	332	384	7			

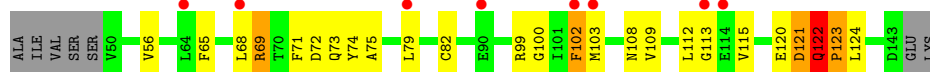
- Molecule 9 is COBALT (II) ION (three-letter code: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	H	1	Total	Co	0	0
			1	1		
9	C	1	Total	Co	0	0
			1	1		

### 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 the various outlier classes 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: SM-LIKE PROTEIN LSM1

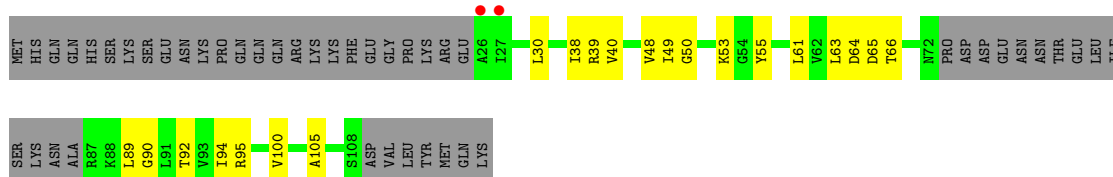


Chain F:  65% 23% 10%



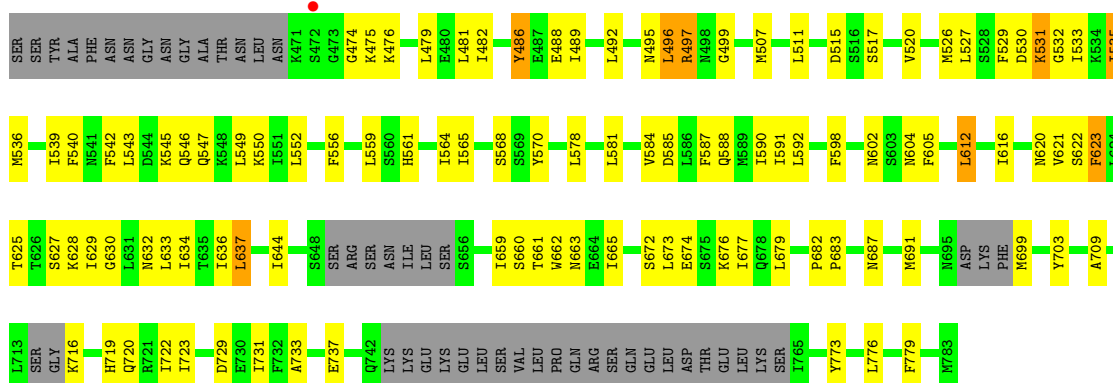
• Molecule 7: U6 SNRNA-ASSOCIATED SM-LIKE PROTEIN LSM7

Chain G:  2% 42% 18% 40%



• Molecule 8: DNA TOPOISOMERASE 2-ASSOCIATED PROTEIN PAT1

Chain H:  53% 30% 15%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	258.69Å 258.69Å 47.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.42 – 3.70 84.68 – 3.70	Depositor EDS
% Data completeness (in resolution range)	96.6 (43.42-3.70) 99.3 (84.68-3.70)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.49 (at 3.67Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.249 , 0.295 0.272 , 0.309	Depositor DCC
$R_{free}$ test set	990 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	135.7	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 122.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.064 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	6280	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	128.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% 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

### 5.1 Standard geometry

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

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	0.41	1/684 (0.1%)	0.81	1/929 (0.1%)
2	B	0.35	0/806	0.73	2/1093 (0.2%)
3	C	0.43	0/594	0.85	2/807 (0.2%)
4	D	0.35	0/509	0.77	0/698
5	E	0.37	0/609	0.78	0/831
6	F	0.38	0/578	0.73	0/784
7	G	0.33	0/456	0.85	1/618 (0.2%)
8	H	0.33	0/2120	0.54	1/2881 (0.0%)
All	All	0.36	1/6356 (0.0%)	0.71	7/8641 (0.1%)

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
8	H	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	123	PRO	N-CD	5.08	1.54	1.47

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	75	LEU	CA-CB-CG	5.88	128.83	115.30
1	A	122	GLN	C-N-CD	5.72	140.40	128.40
7	G	61	LEU	CA-CB-CG	5.40	127.71	115.30
8	H	499	GLY	N-CA-C	-5.40	99.61	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	43	ASP	CB-CG-OD2	5.25	123.02	118.30
2	B	51	TYR	C-N-CD	-5.13	109.30	120.60
2	B	72	LEU	CA-CB-CG	5.13	127.10	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
8	H	496	LEU	Peptide

## 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	673	0	558	36	0
2	B	796	0	755	22	0
3	C	589	0	565	23	0
4	D	504	0	370	15	0
5	E	600	0	558	24	0
6	F	571	0	530	16	0
7	G	456	0	428	24	0
8	H	2089	0	1978	78	0
9	C	1	0	0	0	0
9	H	1	0	0	0	0
All	All	6280	0	5742	220	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 18.

All (220) 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:72:ASP:HA	1:A:75:ALA:HA	1.29	1.07
5:E:73:LEU:HD13	5:E:78:ILE:HD11	1.27	1.05
1:A:72:ASP:HA	1:A:75:ALA:CA	1.92	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:30:LEU:HD12	3:C:31:GLN:N	1.80	0.96
3:C:4:PRO:HG3	6:F:40:SER:CB	1.96	0.95
5:E:73:LEU:O	5:E:73:LEU:HD12	1.73	0.89
3:C:42:SER:O	3:C:43:ASP:OD1	1.95	0.84
5:E:43:VAL:HB	5:E:73:LEU:HD11	1.61	0.82
6:F:47:VAL:HG23	6:F:73:LEU:HB2	1.62	0.80
2:B:16:VAL:HG13	2:B:71:TYR:HB2	1.63	0.79
5:E:45:LEU:HD13	5:E:73:LEU:HD23	1.65	0.79
7:G:39:ARG:HG2	7:G:39:ARG:HH11	1.49	0.78
8:H:495:ASN:O	8:H:496:LEU:HD12	1.84	0.77
4:D:17:ILE:HG23	4:D:79:ILE:HG22	1.66	0.77
7:G:39:ARG:HG2	7:G:39:ARG:NH1	2.01	0.75
1:A:72:ASP:HA	1:A:75:ALA:N	2.02	0.74
1:A:74:TYR:O	1:A:75:ALA:HB3	1.88	0.73
8:H:539:ILE:O	8:H:543:LEU:HD23	1.88	0.73
3:C:73:VAL:O	6:F:74:ARG:HG3	1.87	0.73
1:A:121:ASP:O	1:A:124:LEU:HB2	1.88	0.73
6:F:25:ASN:HB3	6:F:33:LEU:HD11	1.69	0.73
8:H:632:ASN:HD21	8:H:703:TYR:HB2	1.55	0.71
5:E:21:LEU:HD13	5:E:86:LYS:CB	2.20	0.71
8:H:621:VAL:O	8:H:622:SER:HB3	1.92	0.70
3:C:45:VAL:HG12	3:C:62:ARG:HB3	1.72	0.70
8:H:605:PHE:HD2	8:H:644:ILE:HD12	1.55	0.69
7:G:38:ILE:O	7:G:49:ILE:HA	1.92	0.69
4:D:11:LYS:HE3	4:D:31:ASN:HA	1.74	0.69
7:G:38:ILE:HG22	7:G:40:VAL:HG23	1.75	0.69
1:A:68:LEU:HD21	1:A:71:PHE:CB	2.23	0.68
3:C:30:LEU:HD12	3:C:31:GLN:H	1.54	0.68
8:H:659:ILE:O	8:H:663:ASN:ND2	2.27	0.68
8:H:489:ILE:HG21	8:H:535:ILE:HD11	1.76	0.66
7:G:38:ILE:HB	7:G:50:GLY:O	1.95	0.66
5:E:21:LEU:HD12	5:E:31:GLU:HB3	1.78	0.65
1:A:121:ASP:HA	1:A:124:LEU:HD12	1.78	0.65
5:E:73:LEU:HD13	5:E:78:ILE:CD1	2.16	0.65
2:B:73:ASN:N	2:B:73:ASN:OD1	2.30	0.64
8:H:515:ASP:OD1	8:H:517:SER:HB3	1.97	0.63
8:H:533:ILE:HG21	8:H:587:PHE:HA	1.81	0.62
1:A:120:GLU:O	1:A:123:PRO:HG2	2.00	0.62
5:E:73:LEU:CD1	5:E:78:ILE:HD11	2.19	0.60
4:D:25:ILE:O	4:D:25:ILE:HD12	2.01	0.60
7:G:39:ARG:CG	7:G:39:ARG:HH11	2.15	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:ASP:HA	1:A:75:ALA:H	1.65	0.59
8:H:492:LEU:HB3	8:H:542:PHE:CE2	2.38	0.59
8:H:568:SER:HB2	8:H:627:SER:HB2	1.83	0.59
1:A:72:ASP:CA	1:A:75:ALA:HA	2.19	0.59
3:C:24:ARG:HG2	3:C:48:ILE:HG12	1.86	0.58
8:H:511:LEU:HD11	8:H:539:ILE:HD11	1.86	0.58
8:H:622:SER:HA	8:H:625:THR:HG23	1.84	0.58
8:H:731:ILE:HG23	8:H:776:LEU:HD11	1.86	0.58
7:G:39:ARG:HA	7:G:49:ILE:HG13	1.85	0.58
7:G:66:THR:H	7:G:90:GLY:HA2	1.68	0.58
4:D:17:ILE:HG12	4:D:79:ILE:HB	1.86	0.58
4:D:73:GLY:O	4:D:76:ILE:HG12	2.03	0.58
1:A:108:ASN:HD21	4:D:77:LYS:HA	1.69	0.58
3:C:75:LEU:HD11	6:F:67:PHE:HE2	1.69	0.58
8:H:520:VAL:HG12	8:H:561:HIS:ND1	2.18	0.58
3:C:63:CYS:SG	3:C:66:VAL:HG11	2.45	0.57
3:C:76:ILE:HB	6:F:72:PHE:HB3	1.86	0.57
1:A:68:LEU:HD11	1:A:71:PHE:CB	2.35	0.56
8:H:495:ASN:C	8:H:496:LEU:HD12	2.25	0.56
1:A:72:ASP:C	1:A:75:ALA:H	2.09	0.56
2:B:10:LEU:HD11	2:B:72:LEU:HD21	1.86	0.56
1:A:74:TYR:O	1:A:75:ALA:CB	2.52	0.55
1:A:103:MET:HB2	4:D:79:ILE:HG12	1.89	0.55
6:F:35:SER:O	6:F:52:ALA:HA	2.05	0.55
2:B:54:LEU:O	2:B:57:VAL:HG12	2.06	0.54
7:G:89:LEU:HD11	7:G:92:THR:HG22	1.88	0.54
3:C:30:LEU:C	3:C:30:LEU:HD12	2.28	0.54
5:E:28:ARG:NH2	6:F:57:GLU:OE2	2.40	0.54
7:G:89:LEU:HD11	7:G:92:THR:CG2	2.38	0.54
3:C:16:VAL:N	3:C:28:GLY:O	2.32	0.53
5:E:46:GLU:HG3	5:E:70:ARG:HG2	1.89	0.53
8:H:543:LEU:HD12	8:H:547:GLN:HB3	1.90	0.53
8:H:552:LEU:HD22	8:H:598:PHE:CD2	2.44	0.53
8:H:673:LEU:O	8:H:676:LYS:N	2.41	0.53
2:B:19:LEU:HD22	2:B:66:THR:HG23	1.89	0.52
3:C:30:LEU:HA	3:C:41:LEU:HD23	1.92	0.52
8:H:536:MET:HA	8:H:539:ILE:HG22	1.89	0.52
8:H:729:ASP:O	8:H:733:ALA:N	2.42	0.52
4:D:25:ILE:HG12	4:D:71:ILE:CD1	2.39	0.52
6:F:75:GLY:O	6:F:78:VAL:HG22	2.10	0.52
8:H:475:LYS:O	8:H:479:LEU:HD23	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:ASP:CA	1:A:75:ALA:H	2.22	0.52
8:H:482:ILE:HG21	8:H:531:LYS:HD3	1.91	0.52
1:A:120:GLU:O	1:A:123:PRO:CG	2.58	0.52
7:G:64:ASP:OD1	7:G:65:ASP:N	2.43	0.52
8:H:474:GLY:O	8:H:475:LYS:HB3	2.11	0.51
5:E:41:VAL:HG11	7:G:95:ARG:HD2	1.93	0.51
7:G:30:LEU:HD11	7:G:55:TYR:CD1	2.45	0.51
7:G:38:ILE:HG22	7:G:40:VAL:CG2	2.39	0.51
7:G:38:ILE:CB	7:G:50:GLY:O	2.58	0.51
2:B:63:ARG:O	2:B:66:THR:HG22	2.11	0.51
2:B:21:ASN:O	2:B:22:ASP:HB2	2.11	0.51
8:H:682:PRO:HG2	8:H:687:ASN:HD21	1.77	0.50
2:B:-5:TYR:HE1	2:B:9:THR:HG21	1.76	0.50
3:C:68:ILE:HG23	3:C:68:ILE:O	2.12	0.50
2:B:79:THR:HB	3:C:31:GLN:OE1	2.12	0.50
8:H:588:GLN:HG3	8:H:592:LEU:HD12	1.94	0.50
1:A:121:ASP:O	1:A:124:LEU:N	2.45	0.50
8:H:568:SER:HB3	8:H:628:LYS:H	1.77	0.50
1:A:82:CYS:SG	1:A:102:PHE:HB2	2.51	0.50
6:F:56:TYR:C	6:F:56:TYR:CD1	2.85	0.50
8:H:479:LEU:HD21	8:H:529:PHE:CE2	2.46	0.50
8:H:486:TYR:CD2	8:H:535:ILE:HD12	2.47	0.49
8:H:488:GLU:HG2	8:H:507:MET:HA	1.94	0.49
8:H:623:PHE:C	8:H:623:PHE:CD1	2.85	0.49
3:C:10:LEU:HD22	8:H:479:LEU:HB3	1.95	0.49
5:E:45:LEU:HD13	5:E:73:LEU:CD2	2.38	0.49
1:A:82:CYS:N	1:A:99:ARG:O	2.46	0.49
4:D:75:PHE:CD1	7:G:100:VAL:HG22	2.47	0.49
8:H:568:SER:HB3	8:H:628:LYS:N	2.27	0.49
1:A:113:GLY:HA3	2:B:60:ILE:HD13	1.94	0.49
3:C:39:ILE:O	3:C:68:ILE:HG22	2.13	0.49
8:H:543:LEU:HD12	8:H:547:GLN:CB	2.42	0.48
8:H:683:PRO:O	8:H:687:ASN:ND2	2.45	0.48
1:A:122:GLN:N	1:A:123:PRO:HD2	2.27	0.48
3:C:7:LEU:O	3:C:10:LEU:HB2	2.13	0.48
4:D:45:GLU:O	4:D:63:ALA:N	2.47	0.48
5:E:16:ILE:HG12	5:E:34:LEU:HD22	1.96	0.48
5:E:11:VAL:O	5:E:15:THR:HG23	2.14	0.48
1:A:108:ASN:ND2	4:D:77:LYS:HA	2.28	0.48
8:H:556:PHE:CZ	8:H:637:LEU:HD21	2.48	0.48
8:H:559:LEU:HD21	8:H:633:LEU:HD23	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:43:GLY:C	6:F:45:MET:H	2.15	0.48
7:G:40:VAL:N	7:G:48:VAL:O	2.42	0.48
8:H:605:PHE:CD2	8:H:644:ILE:HD12	2.43	0.47
8:H:716:LYS:O	8:H:720:GLN:HG3	2.15	0.47
8:H:709:ALA:HB2	8:H:779:PHE:HA	1.95	0.47
8:H:733:ALA:O	8:H:737:GLU:N	2.48	0.47
1:A:121:ASP:OD1	1:A:124:LEU:HD12	2.15	0.47
1:A:65:PHE:O	1:A:82:CYS:HA	2.14	0.47
7:G:38:ILE:O	7:G:49:ILE:HG23	2.15	0.47
8:H:495:ASN:C	8:H:497:ARG:O	2.53	0.47
4:D:4:LEU:O	4:D:8:THR:OG1	2.31	0.47
5:E:73:LEU:HD12	5:E:73:LEU:C	2.33	0.47
8:H:536:MET:HG3	8:H:591:ILE:HD11	1.96	0.47
5:E:22:ILE:CG2	5:E:78:ILE:HD12	2.46	0.46
7:G:39:ARG:HB2	7:G:105:ALA:HB2	1.97	0.46
8:H:612:LEU:O	8:H:616:ILE:HG12	2.15	0.46
5:E:40:PHE:O	5:E:41:VAL:HG12	2.15	0.46
8:H:474:GLY:C	8:H:476:LYS:H	2.17	0.46
2:B:51:TYR:HA	2:B:52:PRO:HD2	1.73	0.46
8:H:632:ASN:O	8:H:636:ILE:HG13	2.16	0.46
1:A:103:MET:SD	4:D:79:ILE:HD11	2.55	0.46
8:H:674:GLU:HG3	8:H:722:ILE:HG21	1.98	0.46
2:B:73:ASN:HB2	2:B:75:ASN:OD1	2.15	0.46
1:A:121:ASP:O	1:A:124:LEU:CB	2.63	0.46
2:B:44:SER:C	2:B:45:CYS:SG	2.94	0.46
5:E:43:VAL:HG12	5:E:45:LEU:HD12	1.98	0.46
5:E:71:MET:CB	6:F:82:SER:HB2	2.45	0.46
8:H:481:LEU:HB3	8:H:526:MET:HE2	1.98	0.45
8:H:630:GLY:O	8:H:634:ILE:HG12	2.15	0.45
1:A:120:GLU:O	1:A:124:LEU:HG	2.16	0.45
2:B:75:ASN:N	2:B:75:ASN:OD1	2.47	0.45
5:E:43:VAL:O	5:E:73:LEU:HG	2.17	0.45
8:H:546:GLN:HA	8:H:549:LEU:HB3	1.98	0.45
3:C:17:TYR:HB3	3:C:77:SER:O	2.17	0.45
8:H:719:HIS:O	8:H:723:ILE:HG12	2.17	0.45
6:F:47:VAL:CG2	6:F:73:LEU:HB2	2.42	0.45
3:C:63:CYS:SG	3:C:66:VAL:CG1	3.05	0.44
8:H:622:SER:HA	8:H:625:THR:CG2	2.47	0.44
5:E:43:VAL:HB	5:E:73:LEU:CD1	2.40	0.44
8:H:546:GLN:O	8:H:550:LYS:HB2	2.17	0.44
8:H:482:ILE:HD11	8:H:532:GLY:CA	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:545:LYS:HD2	8:H:602:ASN:OD1	2.17	0.44
8:H:620:ASN:O	8:H:623:PHE:HB3	2.18	0.44
1:A:120:GLU:C	1:A:123:PRO:HD2	2.37	0.44
1:A:73:GLN:O	1:A:74:TYR:HD1	2.00	0.44
5:E:26:SER:OG	5:E:27:ASN:N	2.50	0.44
7:G:66:THR:N	7:G:90:GLY:HA2	2.32	0.44
1:A:82:CYS:HB2	1:A:99:ARG:CB	2.48	0.44
5:E:57:ASP:O	5:E:60:ARG:N	2.48	0.44
3:C:4:PRO:CG	6:F:40:SER:CB	2.84	0.44
8:H:527:LEU:HD12	8:H:527:LEU:HA	1.84	0.44
5:E:75:GLY:HA2	5:E:78:ILE:HG12	1.99	0.43
8:H:515:ASP:OD1	8:H:517:SER:N	2.50	0.43
8:H:533:ILE:O	8:H:590:ILE:HG21	2.18	0.43
8:H:612:LEU:HG	8:H:637:LEU:HB3	2.01	0.43
8:H:662:TRP:HA	8:H:665:ILE:HD12	2.00	0.43
8:H:673:LEU:O	8:H:677:ILE:HG13	2.18	0.43
8:H:623:PHE:O	8:H:623:PHE:HD1	2.01	0.43
3:C:42:SER:O	3:C:64:GLU:O	2.37	0.43
7:G:53:LYS:HG3	7:G:64:ASP:HB2	1.98	0.43
8:H:564:ILE:HD11	8:H:584:VAL:HA	2.01	0.43
8:H:578:LEU:HA	8:H:581:LEU:HD12	2.00	0.43
8:H:564:ILE:HG21	8:H:587:PHE:CD2	2.54	0.43
8:H:620:ASN:OD1	8:H:623:PHE:HB2	2.19	0.43
2:B:11:VAL:HA	2:B:29:LEU:HD23	2.01	0.42
1:A:56:VAL:HG22	1:A:112:LEU:HB3	2.00	0.42
8:H:661:THR:HG22	8:H:665:ILE:HD11	2.01	0.42
1:A:73:GLN:C	1:A:74:TYR:HD1	2.22	0.42
6:F:42:ASP:CB	6:F:46:ASN:HB2	2.49	0.42
1:A:74:TYR:CD1	1:A:74:TYR:N	2.86	0.42
3:C:75:LEU:HD11	6:F:67:PHE:CE2	2.53	0.42
8:H:660:SER:HA	8:H:663:ASN:HD22	1.84	0.42
2:B:19:LEU:HD12	2:B:23:ILE:HB	2.01	0.42
8:H:691:MET:HE2	8:H:699:MET:HG2	2.00	0.42
1:A:68:LEU:HD12	1:A:69:ARG:H	1.85	0.42
2:B:57:VAL:HG21	2:B:60:ILE:HD11	2.02	0.42
7:G:63:LEU:HD13	7:G:94:ILE:HG12	2.00	0.42
8:H:475:LYS:HG3	8:H:529:PHE:HZ	1.84	0.42
8:H:570:TYR:OH	8:H:679:LEU:O	2.23	0.41
8:H:629:ILE:HA	8:H:629:ILE:HD13	1.78	0.41
4:D:45:GLU:OE2	7:G:39:ARG:NH2	2.52	0.41
2:B:4:PHE:O	2:B:8:LYS:HG3	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:495:ASN:O	8:H:497:ARG:O	2.39	0.41
8:H:632:ASN:HD22	8:H:632:ASN:HA	1.76	0.41
2:B:-3:GLN:HG2	2:B:-2:GLY:N	2.36	0.41
8:H:604:ASN:OD1	8:H:605:PHE:N	2.54	0.41
8:H:482:ILE:HD12	8:H:526:MET:HB2	2.03	0.40
2:B:8:LYS:HE2	2:B:8:LYS:HB3	1.63	0.40
4:D:17:ILE:HB	4:D:25:ILE:CD1	2.50	0.40
2:B:83:GLN:O	2:B:87:ARG:HG3	2.22	0.40
7:G:55:TYR:C	7:G:55:TYR:CD2	2.93	0.40
1:A:79:LEU:HB2	1:A:82:CYS:SG	2.61	0.40
8:H:511:LEU:HD11	8:H:539:ILE:CD1	2.51	0.40
2:B:-7:ASN:CB	2:B:81:LEU:HD21	2.51	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	92/101 (91%)	88 (96%)	3 (3%)	1 (1%)	16	58
2	B	102/105 (97%)	93 (91%)	8 (8%)	1 (1%)	17	59
3	C	77/89 (86%)	75 (97%)	2 (3%)	0	100	100
4	D	77/114 (68%)	73 (95%)	4 (5%)	0	100	100
5	E	81/93 (87%)	77 (95%)	4 (5%)	0	100	100
6	F	75/86 (87%)	71 (95%)	4 (5%)	0	100	100
7	G	65/115 (56%)	62 (95%)	3 (5%)	0	100	100
8	H	269/328 (82%)	253 (94%)	16 (6%)	0	100	100
All	All	838/1031 (81%)	792 (94%)	44 (5%)	2 (0%)	49	83

All (2) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	A	100	GLY
2	B	52	PRO

### 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	53/91 (58%)	47 (89%)	6 (11%)	6	33
2	B	84/99 (85%)	80 (95%)	4 (5%)	28	64
3	C	59/81 (73%)	56 (95%)	3 (5%)	26	63
4	D	29/104 (28%)	26 (90%)	3 (10%)	8	37
5	E	58/84 (69%)	56 (97%)	2 (3%)	40	72
6	F	58/75 (77%)	55 (95%)	3 (5%)	25	62
7	G	39/103 (38%)	39 (100%)	0	100	100
8	H	204/307 (66%)	191 (94%)	13 (6%)	19	56
All	All	584/944 (62%)	550 (94%)	34 (6%)	22	59

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	ARG
1	A	102	PHE
1	A	109	VAL
1	A	115	VAL
1	A	121	ASP
1	A	122	GLN
2	B	-3	GLN
2	B	9	THR
2	B	16	VAL
2	B	45	CYS
3	C	62	ARG
3	C	63	CYS
3	C	75	LEU
4	D	25	ILE

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Mol	Chain	Res	Type
4	D	45	GLU
4	D	68	GLU
5	E	21	LEU
5	E	28	ARG
6	F	17	SER
6	F	47	VAL
6	F	60	ASN
8	H	486	TYR
8	H	497	ARG
8	H	530	ASP
8	H	531	LYS
8	H	535	ILE
8	H	540	PHE
8	H	565	ILE
8	H	585	ASP
8	H	612	LEU
8	H	623	PHE
8	H	637	LEU
8	H	672	SER
8	H	773	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	ASN
1	A	108	ASN
8	H	588	GLN
8	H	632	ASN
8	H	663	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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 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

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	94/101 (93%)	0.46	8 (8%) 11 8	102, 179, 212, 225	0
2	B	104/105 (99%)	0.10	1 (0%) 82 72	92, 125, 187, 198	0
3	C	79/89 (88%)	-0.11	0 100 100	50, 100, 192, 209	0
4	D	81/114 (71%)	0.25	8 (9%) 7 6	157, 181, 208, 213	0
5	E	83/93 (89%)	-0.22	0 100 100	71, 133, 177, 183	0
6	F	77/86 (89%)	-0.14	0 100 100	79, 106, 145, 153	0
7	G	69/115 (60%)	-0.24	2 (2%) 51 38	84, 157, 190, 219	0
8	H	279/328 (85%)	-0.32	1 (0%) 92 88	50, 95, 163, 231	0
All	All	866/1031 (83%)	-0.08	20 (2%) 60 48	50, 126, 197, 231	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	79	LEU	4.5
4	D	32	VAL	4.1
1	A	102	PHE	3.8
7	G	26	ALA	3.6
4	D	82	GLN	3.4
7	G	27	ILE	3.1
4	D	79	ILE	3.0
4	D	8	THR	3.0
1	A	64	LEU	2.9
1	A	90	GLU	2.7
4	D	25	ILE	2.7
4	D	15	MET	2.4
1	A	68	LEU	2.4
1	A	103	MET	2.3
2	B	70	VAL	2.2
1	A	113	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	114	GLU	2.1
4	D	4	LEU	2.1
4	D	78	PHE	2.1
8	H	472	SER	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](#)

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. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	B-factors( $\text{\AA}^2$ )	Q<0.9
9	CO	H	1080	1/1	0.67	0.42	189,189,189,189	0
9	CO	C	1080	1/1	0.86	0.41	144,144,144,144	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.