



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2017 – 11:09 AM EDT

PDB ID : 4OWT
Title : Structural basis of SOSS1 complex assembly
Authors : Ren, W.; Sun, Q.; Tang, X.; Song, H.
Deposited on : unknown
Resolution : 2.00 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

<http://wwpdb.org/validation/2016/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) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029824

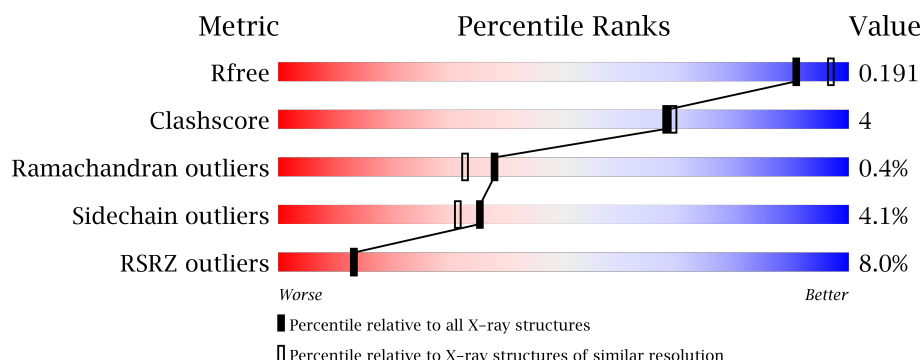
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	100719	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (2.00-2.00)

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. 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	466	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 89%, grey 10%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 1% 89% 10% </div> </div>
2	B	211	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 18%, orange 10%, yellow 5%, green 66%, grey 11%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 18% 28% 5% 66% </div> </div>
3	C	104	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 7%, green 31%, grey 60%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 1% 31% 7% 63% </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5018 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 Integrator complex subunit 3.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	466	Total	C	N	O	S	Se	0	1	0
			3743	2395	649	668	14	17			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	33	GLY	-	expression tag	UNP Q68E01
A	34	GLY	-	expression tag	UNP Q68E01

- Molecule 2 is a protein called SOSS complex subunit B1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	72	Total	C	N	O	S	0	0	0
			566	368	88	107	3			

- Molecule 3 is a protein called SOSS complex subunit C.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	39	Total	C	N	O	0	0	0
			299	192	56	51			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	347	Total	O	0	0
			347	347		
4	B	16	Total	O	0	0
			16	16		
4	C	47	Total	O	0	0
			47	47		

- Molecule 1: Integrator complex subunit 3



4 Model quality [i](#)

4.1 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	0.33	0/3804	0.51	1/5122 (0.0%)
2	B	0.26	0/571	0.41	0/769
3	C	0.33	0/307	0.49	0/417
All	All	0.32	0/4682	0.50	1/6308 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	486	ARG	NE-CZ-NH2	-5.88	117.36	120.30

There are no chirality outliers.

There are no planarity outliers.

4.2 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	3743	0	3837	30	0
2	B	566	0	576	7	0
3	C	299	0	296	4	0
4	A	347	0	0	7	3
4	B	16	0	0	2	0
4	C	47	0	0	2	1
All	All	5018	0	4709	41	3

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 4.

All (41) 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:81:LYS:NZ	4:A:501:HOH:O	2.06	0.88
1:A:222:THR:HG22	1:A:225:LEU:H	1.38	0.86
1:A:347:CYS:SG	4:A:736:HOH:O	2.42	0.77
2:B:70:LEU:O	4:B:311:HOH:O	2.05	0.74
1:A:481:LEU:HB2	1:A:486:ARG:HD2	1.76	0.67
1:A:67:SER:OG	1:A:68:GLU:N	2.29	0.66
1:A:471:HIS:O	4:A:831:HOH:O	2.13	0.65
1:A:165:MSE:HE1	1:A:203:ILE:HG23	1.81	0.62
1:A:222:THR:HG21	4:A:676:HOH:O	1.98	0.62
1:A:45:LYS:HE3	1:A:50:GLU:HG3	1.82	0.61
2:B:56:ASP:OD1	2:B:57:ASP:N	2.34	0.59
3:C:73:GLN:NE2	4:C:201:HOH:O	1.87	0.58
1:A:423:MSE:HE2	1:A:475:LEU:HD12	1.86	0.57
1:A:150:ARG:NH2	4:A:690:HOH:O	2.20	0.57
1:A:359:PRO:HB2	1:A:364:LEU:HD13	1.87	0.56
1:A:395:LEU:HD22	1:A:415:ALA:HB2	1.89	0.55
1:A:165:MSE:HE2	1:A:204:LEU:HA	1.91	0.53
1:A:351:ARG:NH1	4:A:714:HOH:O	2.34	0.49
1:A:79:VAL:HG22	1:A:86:HIS:HA	1.96	0.47
2:B:45:ASP:HB2	4:B:303:HOH:O	2.14	0.46
1:A:423:MSE:CE	1:A:475:LEU:HD12	2.46	0.46
1:A:50:GLU:HG2	1:A:54:ARG:NH1	2.30	0.45
1:A:68:GLU:HG3	1:A:106:LYS:NZ	2.32	0.45
1:A:171:ILE:HD13	1:A:184:ALA:HA	1.99	0.44
1:A:50:GLU:HG2	1:A:54:ARG:HH11	1.83	0.43
2:B:39:ARG:HA	2:B:39:ARG:HD3	1.77	0.43
1:A:416:ILE:HA	1:A:419:MSE:HE3	2.00	0.43
2:B:39:ARG:HB3	2:B:54:VAL:O	2.19	0.43
2:B:82:LEU:HD12	2:B:83:THR:H	1.83	0.43
1:A:457:PHE:HE1	1:A:495:GLU:HG3	1.83	0.43
1:A:435:ASP:OD1	1:A:488[B]:MSE:HE1	2.19	0.43
3:C:69:LYS:NZ	4:C:202:HOH:O	2.22	0.43
1:A:209:VAL:O	1:A:213:LEU:HB2	2.19	0.42
1:A:421:HIS:HD2	4:A:685:HOH:O	2.02	0.42
1:A:38:LEU:HG	1:A:138:GLN:HG2	2.00	0.42
3:C:63:HIS:N	3:C:67:GLN:HG3	2.34	0.42
3:C:63:HIS:HA	3:C:64:ILE:HA	1.89	0.41
1:A:81:LYS:HB3	1:A:81:LYS:HE2	1.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:82:LEU:HD23	2:B:108:PHE:O	2.20	0.41
1:A:377:TRP:O	1:A:381:THR:HG23	2.21	0.41
1:A:349:LEU:O	1:A:353:ILE:HG12	2.21	0.41

All (3) 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)
4:A:558:HOH:O	4:A:567:HOH:O[3_564]	1.94	0.26
4:A:519:HOH:O	4:A:519:HOH:O[5_675]	2.05	0.15
4:A:583:HOH:O	4:C:205:HOH:O[5_675]	2.10	0.10

4.3 Torsion angles ⓘ

4.3.1 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	465/466 (100%)	452 (97%)	12 (3%)	1 (0%)	51	48
2	B	62/211 (29%)	56 (90%)	5 (8%)	1 (2%)	11	5
3	C	37/104 (36%)	36 (97%)	1 (3%)	0	100	100
All	All	564/781 (72%)	544 (96%)	18 (3%)	2 (0%)	38	33

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	67	SER
2	B	39	ARG

4.3.2 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	418/401 (104%)	405 (97%)	13 (3%)	45	44
2	B	66/179 (37%)	60 (91%)	6 (9%)	11	6
3	C	30/85 (35%)	28 (93%)	2 (7%)	19	13
All	All	514/665 (77%)	493 (96%)	21 (4%)	35	31

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	LEU
1	A	41	SER
1	A	111	LEU
1	A	158	LEU
1	A	167	PHE
1	A	199	LEU
1	A	213	LEU
1	A	222	THR
1	A	289	LEU
1	A	345	LEU
1	A	364	LEU
1	A	395	LEU
1	A	397	TYR
2	B	5	THR
2	B	19	LEU
2	B	38	VAL
2	B	39	ARG
2	B	57	ASP
2	B	61	LEU
3	C	84	ILE
3	C	98	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

4.3.3 RNA [i](#)

There are no RNA molecules in this entry.

4.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

4.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

4.6 Ligand geometry [i](#)

There are no ligands in this entry.

4.7 Other polymers [i](#)

There are no such residues in this entry.

4.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

5 Fit of model and data ⓘ

5.1 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, 95th 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 ²)	Q<0.9
1	A	450/466 (96%)	-0.38	6 (1%) 77 77	15, 30, 62, 106	0
2	B	72/211 (34%)	2.13	38 (52%) 0 1	35, 77, 100, 121	0
3	C	39/104 (37%)	-0.50	1 (2%) 56 56	15, 26, 65, 103	0
All	All	561/781 (71%)	-0.07	45 (8%) 13 13	15, 31, 86, 121	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	56	ASP	5.9
2	B	38	VAL	5.6
2	B	40	THR	5.5
2	B	55	TRP	5.3
1	A	33	GLY	5.2
2	B	108	PHE	5.1
1	A	69	ARG	5.0
2	B	10	ILE	5.0
2	B	85	TYR	4.8
2	B	6	PHE	4.8
2	B	41	CYS	4.8
2	B	84	LEU	4.7
2	B	5	THR	4.2
2	B	39	ARG	4.1
2	B	70	LEU	4.0
2	B	106	PRO	3.8
2	B	82	LEU	3.7
2	B	83	THR	3.7
3	C	63	HIS	3.6
2	B	52	ILE	3.5
2	B	21	PHE	3.3
2	B	54	VAL	3.3
2	B	9	ASP	3.2

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Mol	Chain	Res	Type	RSRZ
2	B	57	ASP	3.2
2	B	7	VAL	3.2
1	A	65	GLY	3.2
2	B	8	LYS	3.1
2	B	105	VAL	3.0
2	B	19	LEU	3.0
2	B	92	LEU	2.9
2	B	53	SER	2.8
1	A	68	GLU	2.7
2	B	61	LEU	2.7
2	B	58	VAL	2.7
2	B	20	ILE	2.6
2	B	86	THR	2.6
2	B	27	GLY	2.6
2	B	59	GLY	2.6
2	B	50	ILE	2.4
2	B	62	ILE	2.4
2	B	60	ASN	2.3
1	A	365	SER	2.3
2	B	91	ASP	2.3
2	B	25	GLU	2.3
1	A	70	GLU	2.2

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.4 Ligands ⓘ

There are no ligands in this entry.

5.5 Other polymers ⓘ

There are no such residues in this entry.