



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:15 PM GMT

PDB ID : 1SUL
Title : Crystal Structure of the apo-YsxC
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Deposited on : 2004-03-26
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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

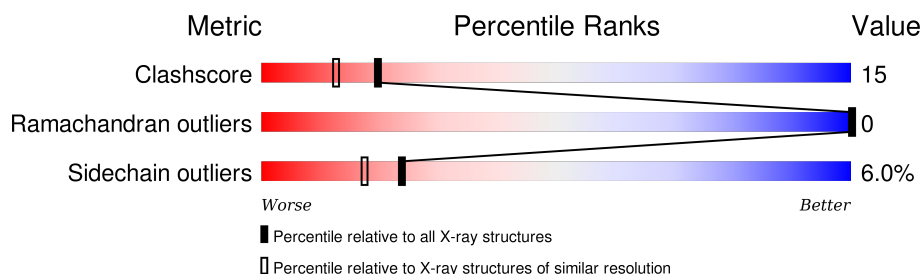
1 Overall quality at a glance

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))
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	195	 72% 20% • 5%
1	B	195	 80% 16% • •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3546 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 GTP-binding protein YsxC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	186	Total	C	N	O	S	0	0	0
			1487	953	253	277	4			
1	B	195	Total	C	N	O	S	0	0	0
			1553	992	267	290	4			

- Molecule 2 is water.

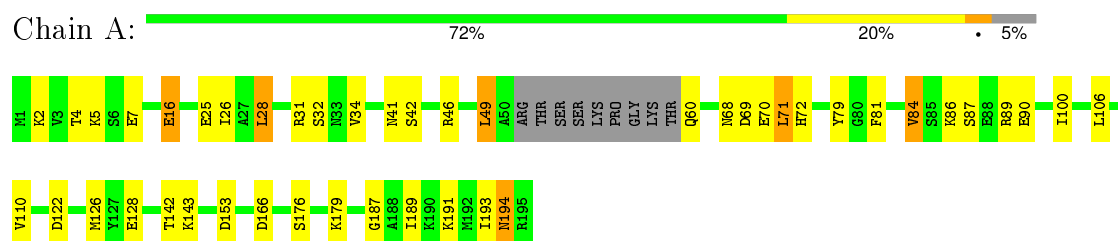
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	269	Total	O	0	0
			269	269		
2	B	237	Total	O	0	0
			237	237		

3 Residue-property plots

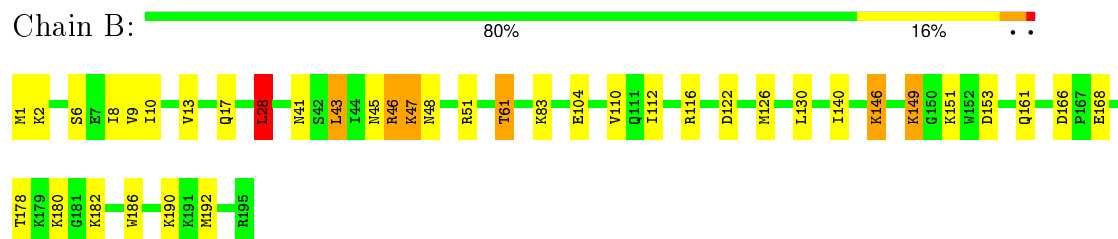
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ($\text{RSRZ} > 2$). 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.

Note EDS was not executed.

• Molecule 1: GTP-binding protein YsxC



• Molecule 1: GTP-binding protein YsxC



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	145.67Å 36.52Å 83.25Å 90.00° 122.16° 90.00°	Depositor
Resolution (Å)	10.00 – 2.00	Depositor
% Data completeness (in resolution range)	96.4 (10.00-2.00)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5	Depositor
R, R_{free}	0.173 , 0.251	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3546	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	0.63	0/1515	0.85	4/2041 (0.2%)
1	B	0.65	0/1583	0.82	4/2133 (0.2%)
All	All	0.64	0/3098	0.83	8/4174 (0.2%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	28	LEU	CA-CB-CG	7.17	131.79	115.30
1	B	46	ARG	NE-CZ-NH2	-6.93	116.83	120.30
1	A	166	ASP	CB-CG-OD2	6.39	124.05	118.30
1	B	166	ASP	CB-CG-OD2	6.25	123.93	118.30
1	A	153	ASP	CB-CG-OD2	6.02	123.72	118.30
1	A	69	ASP	CB-CG-OD2	6.00	123.70	118.30
1	B	153	ASP	CB-CG-OD2	5.81	123.53	118.30
1	A	122	ASP	CB-CG-OD1	5.15	122.94	118.30

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	1487	0	1523	54	0
1	B	1553	0	1597	38	0
2	A	269	0	0	23	4

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	237	0	0	10	4
All	All	3546	0	3120	92	4

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 15.

All (92) 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:16:GLU:N	1:A:16:GLU:OE1	1.64	1.29
1:B:149:LYS:H	1:B:149:LYS:HD2	0.97	1.11
1:B:149:LYS:N	1:B:149:LYS:HD2	1.62	1.09
2:A:342:HOH:O	1:B:151:LYS:HE2	1.55	1.03
1:B:149:LYS:CD	1:B:149:LYS:H	1.73	1.02
1:A:81:PHE:HB3	1:A:89:ARG:HD3	1.52	0.90
1:A:86:LYS:O	1:A:90:GLU:HG3	1.75	0.86
1:A:81:PHE:HA	1:A:84:VAL:CG1	2.06	0.85
1:A:194:ASN:HB3	2:A:242:HOH:O	1.78	0.82
1:A:86:LYS:NZ	2:A:457:HOH:O	2.12	0.82
1:A:4:THR:H	1:A:68:ASN:HD21	1.24	0.81
1:A:81:PHE:CD2	1:A:84:VAL:HG11	2.14	0.81
1:A:193:ILE:HG21	2:A:462:HOH:O	1.80	0.80
1:A:81:PHE:HD2	1:A:84:VAL:HG11	1.46	0.80
1:A:60:GLN:HA	2:A:425:HOH:O	1.80	0.80
1:A:81:PHE:HA	1:A:84:VAL:HG12	1.65	0.79
1:B:1:MET:HB2	1:B:190:LYS:HE2	1.66	0.78
1:A:16:GLU:H	1:A:16:GLU:CD	1.85	0.78
1:A:90:GLU:OE1	2:A:274:HOH:O	2.02	0.77
1:B:122:ASP:HB2	2:B:395:HOH:O	1.87	0.74
1:A:60:GLN:HB2	2:A:327:HOH:O	1.89	0.72
1:B:61:THR:HG21	2:B:272:HOH:O	1.89	0.71
1:B:51:ARG:HD3	2:B:225:HOH:O	1.89	0.71
1:B:9:VAL:O	1:B:10:ILE:HG23	1.94	0.68
1:A:60:GLN:CB	2:A:327:HOH:O	2.42	0.67
1:A:31:ARG:NH2	2:A:295:HOH:O	2.28	0.66
1:A:4:THR:N	1:A:68:ASN:HD21	1.93	0.65
1:A:128:GLU:OE2	2:A:233:HOH:O	2.15	0.63
1:A:70:GLU:HG3	1:A:71:LEU:HD23	1.79	0.63
1:B:48:ASN:HD22	1:B:51:ARG:HH21	1.47	0.61
1:A:81:PHE:CA	1:A:84:VAL:CG1	2.81	0.58
1:B:161:GLN:NE2	2:B:270:HOH:O	2.21	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:ARG:NH1	2:A:309:HOH:O	2.37	0.57
1:A:41:ASN:HB3	1:A:46:ARG:O	2.05	0.56
1:B:180:LYS:NZ	2:B:216:HOH:O	2.35	0.56
1:B:28:LEU:HD13	1:B:110:VAL:HB	1.88	0.56
1:A:5:LYS:NZ	2:A:449:HOH:O	2.39	0.56
1:A:84:VAL:HG22	1:A:89:ARG:HG3	1.86	0.56
1:A:100:ILE:HG21	2:A:383:HOH:O	2.07	0.55
1:A:90:GLU:OE2	2:A:354:HOH:O	2.18	0.55
1:B:13:VAL:HG22	1:B:17:GLN:HE22	1.72	0.54
1:B:168:GLU:HG2	2:B:300:HOH:O	2.06	0.54
1:B:13:VAL:H	1:B:17:GLN:NE2	2.05	0.54
1:A:100:ILE:HG13	2:A:383:HOH:O	2.08	0.54
1:B:6:SER:O	1:B:46:ARG:NH2	2.42	0.53
1:A:81:PHE:CD2	1:A:84:VAL:CG1	2.90	0.52
1:A:46:ARG:NE	2:A:309:HOH:O	2.44	0.51
1:A:70:GLU:CG	1:A:71:LEU:HD23	2.41	0.50
1:A:90:GLU:CD	2:A:354:HOH:O	2.50	0.49
1:B:45:ASN:HB3	2:B:274:HOH:O	2.12	0.49
1:A:28:LEU:HD13	1:A:110:VAL:HB	1.94	0.49
1:A:90:GLU:HG2	2:A:354:HOH:O	2.12	0.49
1:A:194:ASN:O	1:A:194:ASN:CG	2.51	0.48
1:B:41:ASN:HB3	1:B:46:ARG:O	2.13	0.48
1:B:13:VAL:H	1:B:17:GLN:HE22	1.61	0.48
1:A:60:GLN:HG2	2:A:425:HOH:O	2.13	0.48
1:B:192:MET:HE1	2:B:348:HOH:O	2.13	0.48
1:B:45:ASN:CB	2:B:274:HOH:O	2.62	0.48
1:A:60:GLN:HB3	2:A:327:HOH:O	2.12	0.48
1:B:8:ILE:O	1:B:8:ILE:HG23	2.13	0.48
1:A:71:LEU:HD13	1:A:193:ILE:HD11	1.97	0.47
1:B:192:MET:CE	1:B:192:MET:HA	2.44	0.47
1:B:178:THR:O	1:B:178:THR:HG22	2.14	0.47
1:B:45:ASN:CG	2:B:274:HOH:O	2.52	0.47
1:A:25:GLU:HG2	1:A:72:HIS:HB2	1.97	0.47
1:A:194:ASN:O	1:A:194:ASN:ND2	2.48	0.47
1:A:176:SER:O	1:A:179:LYS:HE3	2.14	0.47
1:B:146:LYS:HD2	1:B:146:LYS:N	2.29	0.47
1:B:47:LYS:HA	1:B:47:LYS:HD2	1.48	0.46
1:A:46:ARG:CZ	2:A:309:HOH:O	2.63	0.46
1:A:193:ILE:O	1:A:193:ILE:CG2	2.63	0.46
1:A:81:PHE:HD2	1:A:84:VAL:CG1	2.25	0.44
1:B:1:MET:HG2	1:B:186:TRP:CE3	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:ILE:HG13	1:A:189:ILE:HG23	2.00	0.44
1:B:112:ILE:HD13	1:B:140:ILE:HB	2.01	0.43
1:B:149:LYS:CD	1:B:149:LYS:N	2.42	0.43
1:A:86:LYS:O	1:A:90:GLU:CG	2.58	0.42
1:A:193:ILE:HD13	2:A:462:HOH:O	2.19	0.42
1:B:43:LEU:HD12	1:B:182:LYS:HG3	2.01	0.42
1:A:79:TYR:CE2	1:A:126:MET:HG2	2.55	0.42
1:A:100:ILE:HD13	1:A:106:LEU:CD2	2.50	0.41
1:B:41:ASN:ND2	1:B:48:ASN:HA	2.35	0.41
1:B:192:MET:HE2	1:B:192:MET:HA	2.00	0.41
1:A:187:GLY:O	1:A:191:LYS:HG3	2.21	0.41
1:B:116:ARG:HD3	1:B:146:LYS:CB	2.51	0.41
1:B:110:VAL:HG13	1:B:140:ILE:HG12	2.03	0.41
1:A:100:ILE:HD13	1:A:106:LEU:HD22	2.02	0.41
1:A:34:VAL:HG13	2:A:321:HOH:O	2.21	0.41
1:B:9:VAL:O	1:B:10:ILE:CG2	2.67	0.40
1:A:142:THR:OG1	1:A:143:LYS:HD2	2.20	0.40
1:B:126:MET:O	1:B:130:LEU:HG	2.21	0.40
1:A:49:LEU:O	1:A:49:LEU:HD22	2.21	0.40

All (4) 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)
2:A:284:HOH:O	2:B:281:HOH:O[3_454]	2.09	0.11
2:A:211:HOH:O	2:B:373:HOH:O[3_454]	2.11	0.09
2:A:281:HOH:O	2:B:307:HOH:O[4_557]	2.17	0.03
2:A:344:HOH:O	2:B:340:HOH:O[2_667]	2.17	0.03

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	182/195 (93%)	178 (98%)	4 (2%)	0	100	100
1	B	193/195 (99%)	190 (98%)	3 (2%)	0	100	100
All	All	375/390 (96%)	368 (98%)	7 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	162/170 (95%)	151 (93%)	11 (7%)	20	13
1	B	170/170 (100%)	161 (95%)	9 (5%)	28	22
All	All	332/340 (98%)	312 (94%)	20 (6%)	24	17

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	7	GLU
1	A	16	GLU
1	A	28	LEU
1	A	32	SER
1	A	42	SER
1	A	49	LEU
1	A	71	LEU
1	A	84	VAL
1	A	87	SER
1	A	194	ASN
1	B	2	LYS
1	B	28	LEU
1	B	43	LEU
1	B	47	LYS
1	B	61	THR
1	B	83	LYS
1	B	104	GLU

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Mol	Chain	Res	Type
1	B	146	LYS
1	B	149	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	A	60	GLN
1	A	63	ASN
1	A	68	ASN
1	A	111	GLN
1	A	125	GLN
1	A	164	ASN
1	B	17	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 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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.