wwPDB X-ray Structure Validation Summary Report (i)

Jan 31, 2016 – 08:17 PM GMT

PDB ID	:	1JK9
Title	:	Heterodimer between H48F-ySOD1 and yCCS
Authors	:	Lamb, A.L.; Torres, A.S.; O'Halloran, T.V.; Rosenzweig, A.C.
Deposited on	:	2001-07-11
Resolution	:	2.90 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

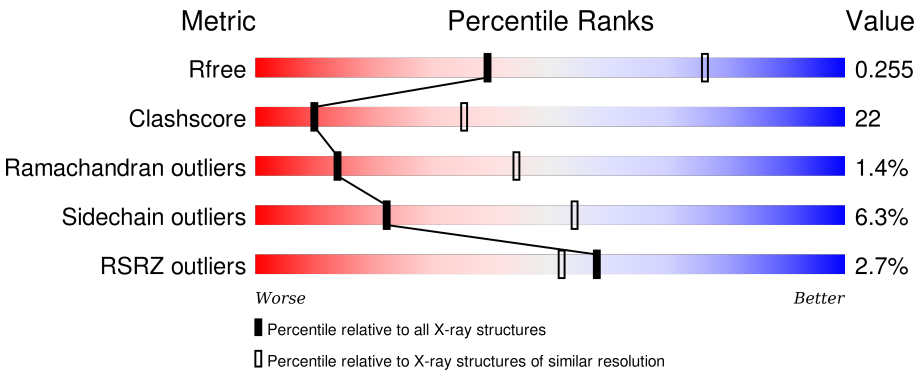
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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 <=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	153	2%, 69%, 28%, •
1	C	153	%, 70%, 27%, •
2	B	249	3%, 63%, 29%, 6%, •
2	D	249	4%, 65%, 29%, •, •

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5979 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 superoxide dismutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	153	1107	683	196	225	3	0	0	0
1	C	153	1107	683	196	225	3	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	48	PHE	HIS	ENGINEERED	UNP P00445
C	48	PHE	HIS	ENGINEERED	UNP P00445

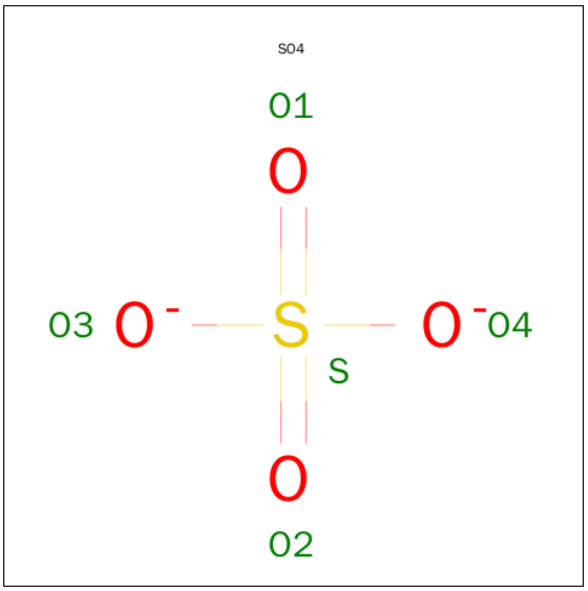
- Molecule 2 is a protein called copper chaperone for superoxide dismutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	243	1870	1172	322	367	9	0	0	0
2	D	243	1870	1172	322	367	9	0	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
3	B	1	1	1	0	0
3	A	1	1	1	0	0
3	D	1	1	1	0	0
3	C	1	1	1	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
4	A	1	5	4	1	0	0
			Total	O	S		
4	C	1	5	4	1	0	0

- Molecule 5 is water.

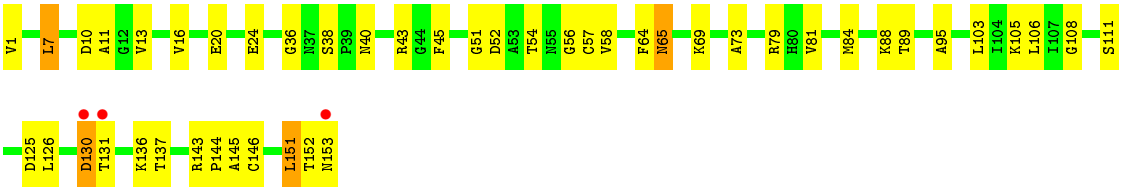
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	4	4	4	0	0
			Total	O		
5	B	2	2	2	0	0
			Total	O		
5	C	2	2	2	0	0
			Total	O		
5	D	3	3	3	0	0

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 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 (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.

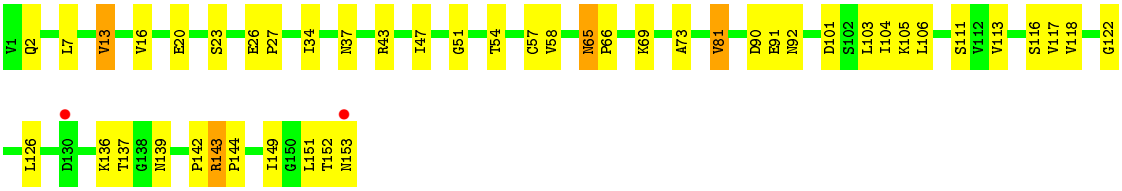
- Molecule 1: superoxide dismutase

Chain A:



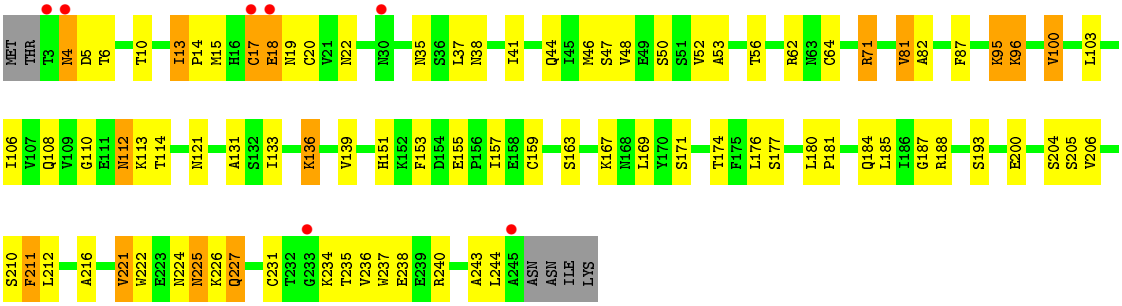
- Molecule 1: superoxide dismutase

Chain C:



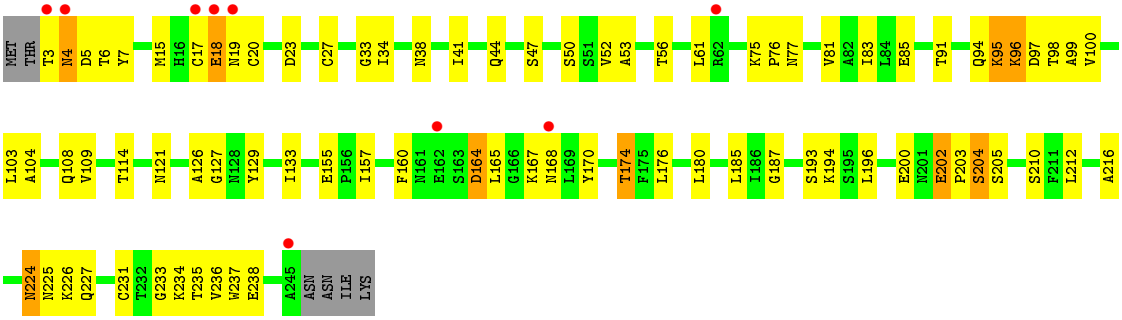
- Molecule 2: copper chaperone for superoxide dismutase

Chain B:



- Molecule 2: copper chaperone for superoxide dismutase

Chain D:


MET THR T3 N4 D5 T6 Y7 M15 H16 C17 E18 N19 C20 D23 C27 G33 I34 N38 I41 Q44 S47 S50 S51 V52 A53 T56 L61 R62 K75 P76 N77 V81 A82 I83 L84 E85 T91 Q94 K95 K96 D97 T98 A99 V100
L103 A104 Q108 V109 T114 N121 A126 G127 N128 Y129 I133 E155 P156 I157 F160 N161 E162 S163 D164 L165 G166 K167 N168 L169 Y170 T174 F175 L176 L180 L185 I186 G187 S193 K194 S195 L196 E200 N201 E202 P203 S204 S205 S210 F211 L212 A216
N224 N225 K226 Q227 C231 T232 G233 K234 T235 V236 W237 E238 A245 ASN ASN ILE LYS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	104.09Å 104.09Å 233.71Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	12.00 – 2.90 12.05 – 2.91	Depositor EDS
% Data completeness (in resolution range)	(Not available) (12.00-2.90) 95.6 (12.05-2.91)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.14 (at 2.92Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.217 , 0.260 0.215 , 0.255	Depositor DCC
R_{free} test set	3091 reflections (10.01%)	DCC
Wilson B-factor (Å^2)	32.2	Xtriage
Anisotropy	0.034	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 47.3	EDS
Estimated twinning fraction	0.035 for -h,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 32013 reflections	Xtriage
F_o,F_c correlation	0.89	EDS
Total number of atoms	5979	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, SO4

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	0.39	0/1127	0.66	0/1524
1	C	0.38	0/1127	0.65	0/1524
2	B	0.39	0/1906	0.66	0/2584
2	D	0.40	0/1906	0.65	0/2584
All	All	0.39	0/6066	0.65	0/8216

There are no bond length outliers.

There are no bond angle outliers.

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	1107	0	1074	36	0
1	C	1107	0	1074	38	0
2	B	1870	0	1843	111	0
2	D	1870	0	1843	105	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	5	0	0	0	0
5	A	4	0	0	0	0
5	B	2	0	0	0	0
5	C	2	0	0	0	0
5	D	3	0	0	0	0
All	All	5979	0	5834	259	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 22.

The worst 5 of 259 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)
2:B:96:LYS:H	2:B:96:LYS:HE2	1.05	1.12
2:D:96:LYS:HE2	2:D:96:LYS:H	1.18	1.09
2:D:108:GLN:HG3	2:D:114:THR:HG22	1.43	0.96
2:B:95:LYS:HD2	2:B:96:LYS:HZ1	1.36	0.90
2:B:95:LYS:HD2	2:B:96:LYS:NZ	1.87	0.89

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	151/153 (99%)	134 (89%)	15 (10%)	2 (1%)	15 46
1	C	151/153 (99%)	144 (95%)	7 (5%)	0	100 100
2	B	241/249 (97%)	218 (90%)	21 (9%)	2 (1%)	24 60
2	D	241/249 (97%)	209 (87%)	25 (10%)	7 (3%)	6 23
All	All	784/804 (98%)	705 (90%)	68 (9%)	11 (1%)	14 44

5 of 11 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	131	THR
2	B	5	ASP
2	D	165	LEU
2	D	204	SER
2	B	205	SER

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	118/118 (100%)	111 (94%)	7 (6%)	24	58
1	C	118/118 (100%)	111 (94%)	7 (6%)	24	58
2	B	208/214 (97%)	192 (92%)	16 (8%)	16	42
2	D	208/214 (97%)	197 (95%)	11 (5%)	28	63
All	All	652/664 (98%)	611 (94%)	41 (6%)	22	54

5 of 41 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	200	GLU
2	B	227	GLN
2	D	200	GLU
2	B	211	PHE
2	B	221	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 26 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	224	ASN
1	C	65	ASN
2	D	224	ASN
2	B	225	ASN
1	C	37	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)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	SO4	A	401	-	4,4,4	0.30	0	6,6,6	0.15	0
4	SO4	C	402	-	4,4,4	0.24	0	6,6,6	0.10	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SO4	A	401	-	-	0/0/0/0	0/0/0/0
4	SO4	C	402	-	-	0/0/0/0	0/0/0/0

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 (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, 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	153/153 (100%)	-0.46	3 (1%)	68	64	10, 24, 46, 56	0
1	C	153/153 (100%)	-0.46	2 (1%)	79	78	8, 22, 43, 51	0
2	B	243/249 (97%)	-0.20	7 (2%)	55	49	9, 28, 49, 71	0
2	D	243/249 (97%)	-0.22	9 (3%)	45	38	6, 25, 51, 67	0
All	All	792/804 (98%)	-0.31	21 (2%)	58	52	6, 25, 48, 71	0

The worst 5 of 21 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	18	GLU	4.5
2	B	3	THR	4.4
2	B	4	ASN	3.5
2	D	19	ASN	3.1
1	A	153	ASN	3.1

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th 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	LLDF	B-factors(Å^2)	Q<0.9
3	ZN	C	302	1/1	0.99	0.07	-1.58	19,19,19,19	0
3	ZN	A	301	1/1	0.99	0.07	-1.64	18,18,18,18	0
4	SO4	C	402	5/5	0.98	0.10	-1.95	25,26,28,29	0
4	SO4	A	401	5/5	0.98	0.09	-2.00	23,25,25,26	0
3	ZN	B	303	1/1	0.99	0.07	-2.97	22,22,22,22	0
3	ZN	D	304	1/1	0.97	0.05	-5.21	25,25,25,25	0

6.5 Other polymers (i)

There are no such residues in this entry.