wwPDB X-ray Structure Validation Summary Report (i)

Feb 3, 2024 – 08:15 PM EST

PDB ID	:	1MR1
Title	:	Crystal Structure of a Smad4-Ski Complex
Authors	:	Wu, J.-W.; Krawitz, A.R.; Chai, J.; Li, W.; Zhang, F.; Luo, K.; Shi, Y.
Deposited on	:	2002-09-17
Resolution	:	2.85 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

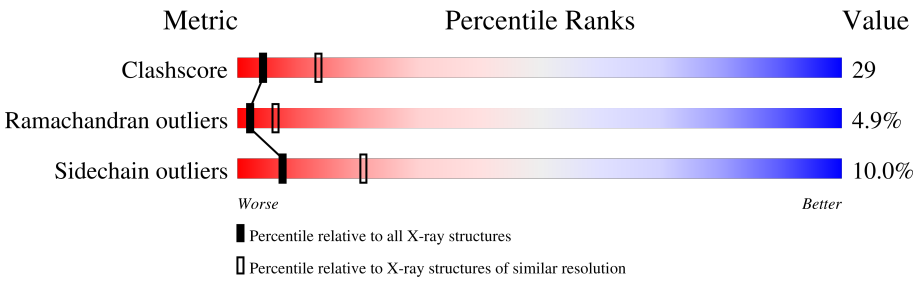
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.85 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	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)

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 of the lower bar indicate the fraction of residues that contain outliers for >=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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	235	46% 33% • 17%
1	B	235	39% 37% 5% • 18%
2	C	99	51% 41% 6% •
2	D	99	41% 42% 10% • •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4684 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 Mothers against decapentaplegic homolog 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	195	1549	978	278	281	12	0	0	0
1	B	193	1545	979	275	279	12	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	318	MET	-	initiating methionine	UNP Q13485
B	318	MET	-	initiating methionine	UNP Q13485

- Molecule 2 is a protein called Ski oncogene.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	97	789	499	141	140	9	0	0	0
2	D	96	783	496	140	138	9	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	215	GLY	-	cloning artifact	UNP P12755
C	216	SER	-	cloning artifact	UNP P12755
C	217	HIS	-	cloning artifact	UNP P12755
C	218	MET	-	cloning artifact	UNP P12755
D	215	GLY	-	cloning artifact	UNP P12755
D	216	SER	-	cloning artifact	UNP P12755
D	217	HIS	-	cloning artifact	UNP P12755
D	218	MET	-	cloning artifact	UNP P12755

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

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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	7	7	7	0	0
			Total	O		
4	B	2	2	2	0	0
			Total	O		
4	C	6	6	6	0	0
			Total	O		
4	D	1	1	1	0	0

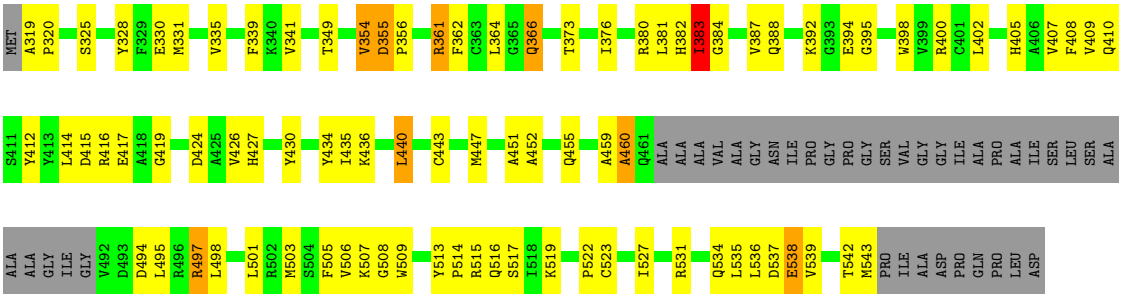
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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. 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. 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.

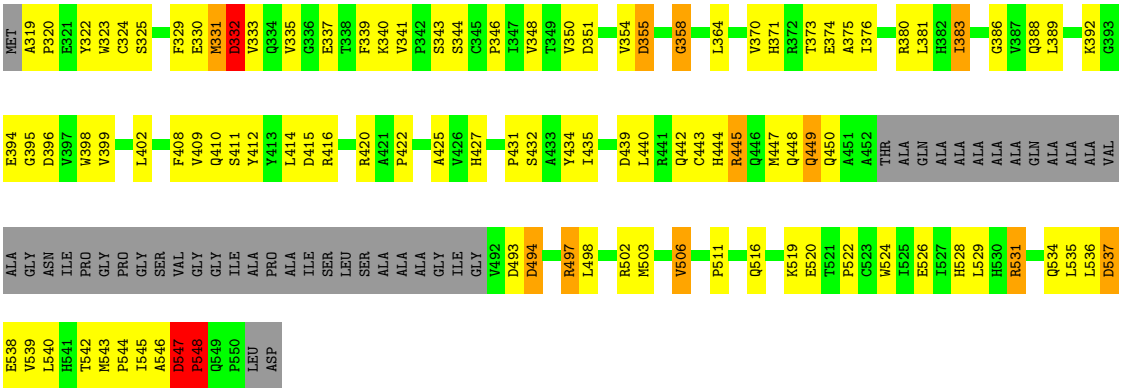
• Molecule 1: Mothers against decapentaplegic homolog 4

Chain A: 46% 33% • 17%



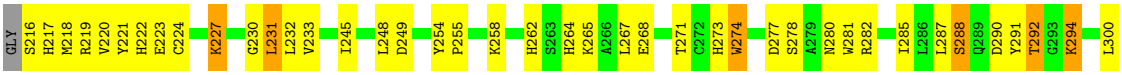
• Molecule 1: Mothers against decapentaplegic homolog 4

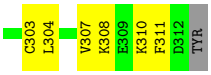
Chain B: 39% 37% 5% • 18%



• Molecule 2: Ski oncogene

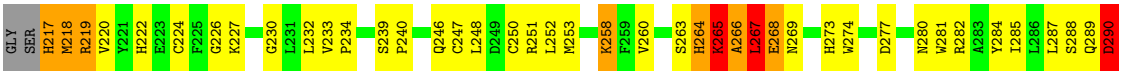
Chain C: 51% 41% 6% •

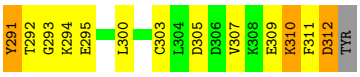




- Molecule 2: Ski oncogene

Chain D:





4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	109.80Å 109.80Å 141.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.85	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.85)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.231 , 0.280	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4684	wwPDB-VP
Average B, all atoms (Å^2)	49.0	wwPDB-VP

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

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	0.44	0/1586	0.68	0/2149
1	B	0.38	0/1585	0.63	0/2150
2	C	0.48	0/811	0.78	1/1093 (0.1%)
2	D	0.54	0/805	0.90	3/1085 (0.3%)
All	All	0.45	0/4787	0.72	4/6477 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	266	ALA	N-CA-C	-5.94	94.97	111.00
2	D	226	GLY	N-CA-C	-5.62	99.05	113.10
2	C	294	LYS	N-CA-C	5.54	125.95	111.00
2	D	267	LEU	N-CA-C	-5.04	97.40	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	1549	0	1509	83	0
1	B	1545	0	1505	93	0
2	C	789	0	757	44	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	783	0	752	63	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	7	0	0	0	0
4	B	2	0	0	0	0
4	C	6	0	0	1	0
4	D	1	0	0	0	0
All	All	4684	0	4523	270	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 29.

The worst 5 of 270 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:366:GLN:HE21	1:A:366:GLN:N	1.29	1.28
1:A:366:GLN:H	1:A:366:GLN:NE2	1.45	1.14
2:D:218:MET:HB2	2:D:232:LEU:HB3	1.42	1.02
1:B:420:ARG:HH11	1:B:425:ALA:HB2	1.21	1.00
2:D:264:HIS:O	2:D:266:ALA:N	2.01	0.92

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	191/235 (81%)	164 (86%)	21 (11%)	6 (3%)	4 14
1	B	189/235 (80%)	159 (84%)	19 (10%)	11 (6%)	1 4
2	C	95/99 (96%)	80 (84%)	12 (13%)	3 (3%)	4 13
2	D	94/99 (95%)	74 (79%)	12 (13%)	8 (8%)	1 1

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	569/668 (85%)	477 (84%)	64 (11%)	28 (5%)	2 6

5 of 28 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	384	GLY
1	B	332	ASP
1	B	355	ASP
1	B	547	ASP
2	C	294	LYS

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	164/187 (88%)	151 (92%)	13 (8%)	12 31
1	B	167/187 (89%)	150 (90%)	17 (10%)	7 20
2	C	86/87 (99%)	79 (92%)	7 (8%)	11 30
2	D	85/87 (98%)	72 (85%)	13 (15%)	2 7
All	All	502/548 (92%)	452 (90%)	50 (10%)	7 21

5 of 50 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	547	ASP
2	C	288	SER
2	D	312	ASP
1	B	548	PRO
2	C	271	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 22 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	528	HIS
2	C	280	ASN
2	C	222	HIS
2	D	217	HIS
1	B	382	HIS

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 monosaccharides 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 (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.