Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 06:32 am BST

PDB ID	:	6AD3
Title	:	Structural characterization of the condensation domain from Monacolin K polyketide synthase MokA
Authors	:	Wang, L.; Zheng, J.
Deposited on	:	2018-07-30
Resolution	:	1.79 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

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

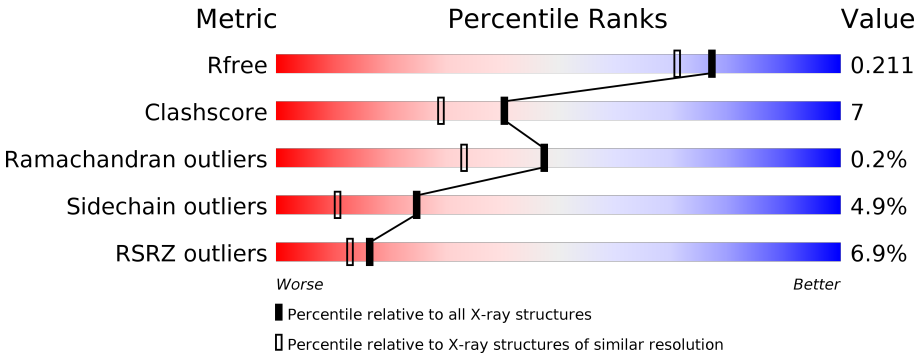
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 1.79 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}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 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% 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	519	6% 73% 13% • 13%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3788 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 Lovastatin nonaketide synthase mokA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	450	Total	C	N	O	S	0	0	0
			3528	2240	626	649	13			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP Q3S2T9
A	-18	GLY	-	expression tag	UNP Q3S2T9
A	-17	SER	-	expression tag	UNP Q3S2T9
A	-16	SER	-	expression tag	UNP Q3S2T9
A	-15	HIS	-	expression tag	UNP Q3S2T9
A	-14	HIS	-	expression tag	UNP Q3S2T9
A	-13	HIS	-	expression tag	UNP Q3S2T9
A	-12	HIS	-	expression tag	UNP Q3S2T9
A	-11	HIS	-	expression tag	UNP Q3S2T9
A	-10	HIS	-	expression tag	UNP Q3S2T9
A	-9	SER	-	expression tag	UNP Q3S2T9
A	-8	SER	-	expression tag	UNP Q3S2T9
A	-7	GLY	-	expression tag	UNP Q3S2T9
A	-6	LEU	-	expression tag	UNP Q3S2T9
A	-5	VAL	-	expression tag	UNP Q3S2T9
A	-4	PRO	-	expression tag	UNP Q3S2T9
A	-3	ARG	-	expression tag	UNP Q3S2T9
A	-2	GLY	-	expression tag	UNP Q3S2T9
A	-1	SER	-	expression tag	UNP Q3S2T9
A	0	HIS	-	expression tag	UNP Q3S2T9
A	1	MET	-	expression tag	UNP Q3S2T9
A	153	TRP	ARG	engineered mutation	UNP Q3S2T9

- Molecule 2 is water.

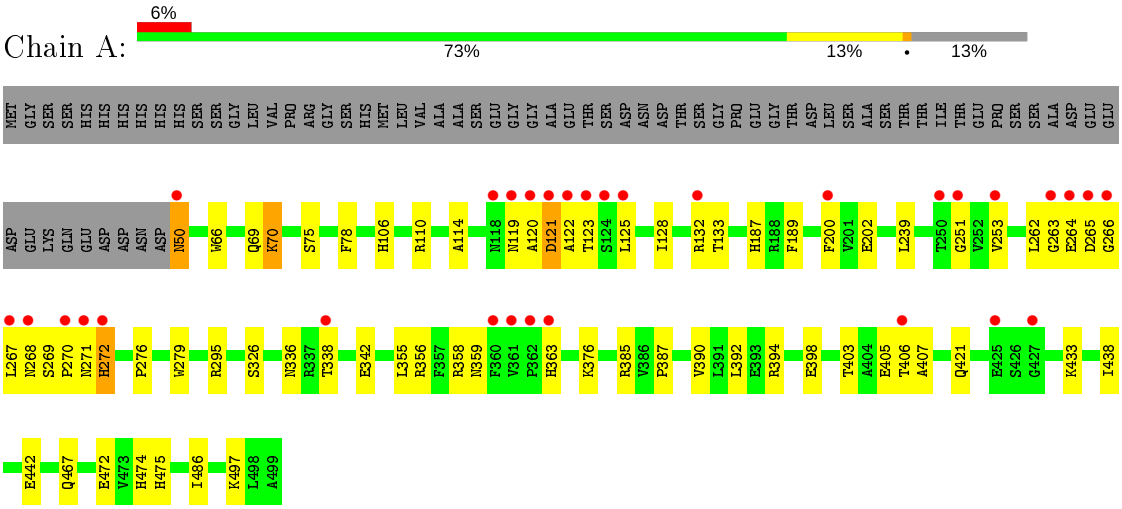
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	260	Total	O	0	0
			260	260		

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: Lovastatin nonaketide synthase mokA

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 2 21	Depositor
Cell constants a, b, c, α, β, γ	51.10Å 64.04Å 173.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.79 42.88 – 1.79	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-1.79) 99.8 (42.88-1.79)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.63 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.176 , 0.206 0.189 , 0.211	Depositor DCC
R_{free} test set	2681 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å^2)	20.9	Xtriage
Anisotropy	0.091	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 43.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	3788	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.78% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.68	0/3610	0.66	2/4897 (0.0%)

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
1	A	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	269	SER	C-N-CD	5.83	140.63	128.40
1	A	263	GLY	N-CA-C	-5.21	100.09	113.10

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	200	PHE	Sidechain
1	A	251	GLY	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	3528	0	3469	50	1
2	A	260	0	0	3	0
All	All	3788	0	3469	50	1

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 7.

All (50) 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:50:ASN:HB3	2:A:656:HOH:O	1.67	0.94
1:A:265:ASP:CB	1:A:266:GLY:HA2	2.08	0.83
1:A:120:ALA:HB3	1:A:121:ASP:C	2.01	0.80
1:A:356:ARG:O	1:A:376:LYS:NZ	2.16	0.77
1:A:120:ALA:HB1	1:A:123:THR:H	1.50	0.75
1:A:120:ALA:HB1	1:A:123:THR:OG1	1.86	0.75
1:A:264:GLU:HA	1:A:474:HIS:CE1	2.23	0.74
1:A:70:LYS:HD3	1:A:122:ALA:HB1	1.69	0.73
1:A:270:PRO:HB3	1:A:272:HIS:CE1	2.25	0.71
1:A:120:ALA:N	1:A:121:ASP:HA	2.06	0.70
1:A:114:ALA:CB	1:A:128:ILE:HD11	2.22	0.70
1:A:279:TRP:HA	1:A:467:GLN:HE22	1.56	0.69
1:A:114:ALA:HB2	1:A:128:ILE:HD11	1.79	0.65
1:A:270:PRO:CB	1:A:272:HIS:CE1	2.80	0.64
1:A:187:HIS:HD2	1:A:189:PHE:H	1.45	0.63
1:A:120:ALA:CB	1:A:123:THR:H	2.12	0.61
1:A:264:GLU:HA	1:A:474:HIS:HE1	1.65	0.61
1:A:120:ALA:O	1:A:123:THR:HG23	2.01	0.59
1:A:69:GLN:HE22	1:A:187:HIS:CE1	2.23	0.57
1:A:276:PRO:HA	1:A:407:ALA:HB2	1.85	0.57
1:A:106:HIS:HE1	1:A:202:GLU:OE2	1.88	0.56
1:A:326:SER:O	1:A:358:ARG:O	2.25	0.55
1:A:120:ALA:HB3	1:A:121:ASP:O	2.06	0.54
1:A:472:GLU:H	1:A:475:HIS:HD1	1.57	0.51
1:A:110:ARG:HB3	1:A:133:THR:HA	1.92	0.51
1:A:114:ALA:HB2	1:A:128:ILE:CD1	2.39	0.51
1:A:114:ALA:HB3	1:A:128:ILE:HD11	1.93	0.51
1:A:69:GLN:HG3	1:A:75:SER:CB	2.40	0.51
1:A:66:TRP:CD2	1:A:125:LEU:HD22	2.47	0.50
1:A:187:HIS:CD2	1:A:189:PHE:H	2.25	0.50
1:A:265:ASP:CB	1:A:266:GLY:CA	2.85	0.50
1:A:106:HIS:HD2	2:A:695:HOH:O	1.95	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:120:ALA:N	1:A:121:ASP:CA	2.74	0.49
1:A:120:ALA:HB1	1:A:123:THR:N	2.22	0.49
1:A:239:LEU:HD22	1:A:394:ARG:HD3	1.96	0.47
1:A:403:THR:OG1	1:A:406:THR:HG22	2.14	0.47
1:A:69:GLN:HE22	1:A:187:HIS:HE1	1.62	0.47
1:A:270:PRO:HB2	1:A:272:HIS:CE1	2.49	0.46
1:A:69:GLN:HG3	1:A:75:SER:HB2	1.99	0.45
1:A:70:LYS:CD	1:A:122:ALA:CB	2.94	0.45
1:A:338:THR:N	1:A:342:GLU:OE2	2.40	0.45
1:A:70:LYS:HD3	1:A:122:ALA:CB	2.43	0.45
1:A:295:ARG:HD2	2:A:510:HOH:O	2.17	0.44
1:A:78:PHE:HA	1:A:442:GLU:HB2	2.00	0.44
1:A:385:ARG:O	1:A:387:PRO:HD3	2.17	0.44
1:A:70:LYS:CD	1:A:122:ALA:HB1	2.43	0.43
1:A:358:ARG:HG3	1:A:359:ASN:N	2.33	0.43
1:A:66:TRP:CZ2	1:A:70:LYS:HD3	2.55	0.42
1:A:270:PRO:O	1:A:271:ASN:C	2.59	0.42
1:A:363:HIS:O	1:A:497:LYS:HD2	2.21	0.41

All (1) 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)
1:A:267:LEU:O	1:A:385:ARG:NH2[1_655]	2.06	0.14

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	448/519 (86%)	434 (97%)	13 (3%)	1 (0%)	47 33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	268	ASN

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	368/430 (86%)	350 (95%)	18 (5%)	25 11

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	ASN
1	A	70	LYS
1	A	119	ASN
1	A	121	ASP
1	A	132	ARG
1	A	253	VAL
1	A	262	LEU
1	A	272	HIS
1	A	336	ASN
1	A	355	LEU
1	A	390	VAL
1	A	392	LEU
1	A	398	GLU
1	A	405	GLU
1	A	421	GLN
1	A	433	LYS
1	A	438	ILE
1	A	486	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	GLN
1	A	106	HIS
1	A	154	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	187	HIS
1	A	198	ASN
1	A	272	HIS
1	A	277	ASN
1	A	336	ASN
1	A	467	GLN
1	A	474	HIS

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

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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^{th} 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	450/519 (86%)	0.31	31 (6%) 16 13	12, 22, 48, 115	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	120	ALA	14.0
1	A	121	ASP	14.0
1	A	267	LEU	13.3
1	A	123	THR	7.1
1	A	122	ALA	6.9
1	A	119	ASN	5.2
1	A	124	SER	4.9
1	A	266	GLY	4.6
1	A	270	PRO	4.5
1	A	361	VAL	4.5
1	A	263	GLY	4.5
1	A	118	ASN	3.8
1	A	251	GLY	3.6
1	A	264	GLU	3.5
1	A	360	PHE	3.4
1	A	425	GLU	3.4
1	A	338	THR	3.4
1	A	265	ASP	3.1
1	A	271	ASN	3.1
1	A	253	VAL	3.0
1	A	50	ASN	2.8
1	A	268	ASN	2.8
1	A	363	HIS	2.7
1	A	132	ARG	2.6
1	A	362	PRO	2.6
1	A	250	THR	2.4
1	A	427	GLY	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	406	THR	2.1
1	A	200	PHE	2.1
1	A	272	HIS	2.1
1	A	125	LEU	2.1

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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.