



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2017 – 11:09 pm GMT

PDB ID : 4RC3
Title : Crystal structure of Ser/Thr kinase Pim1 in complex with Mitoxantrone derivatives
Authors : Zhang, W.; Wan, X.; Huang, N.
Deposited on : 2014-09-14
Resolution : 2.34 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
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)	:	recalc28949

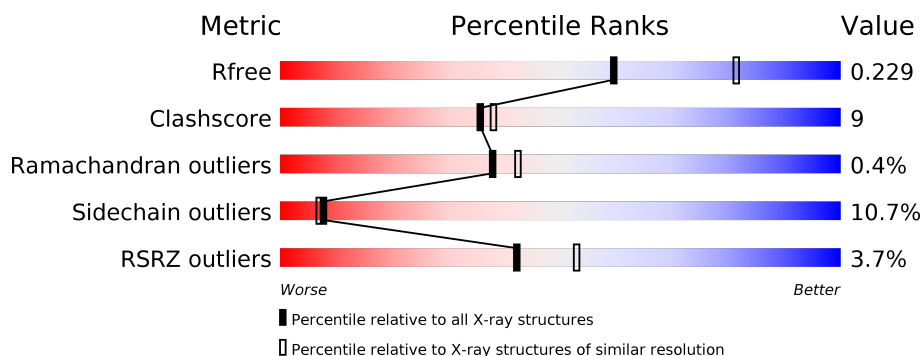
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.34 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	1570 (2.36-2.32)
Clashscore	112137	1673 (2.36-2.32)
Ramachandran outliers	110173	1654 (2.36-2.32)
Sidechain outliers	110143	1655 (2.36-2.32)
RSRZ outliers	101464	1576 (2.36-2.32)

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	289	<div> <div>3%</div> <div>71%</div> <div>21%</div> <div>6%</div> </div>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 2256 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.

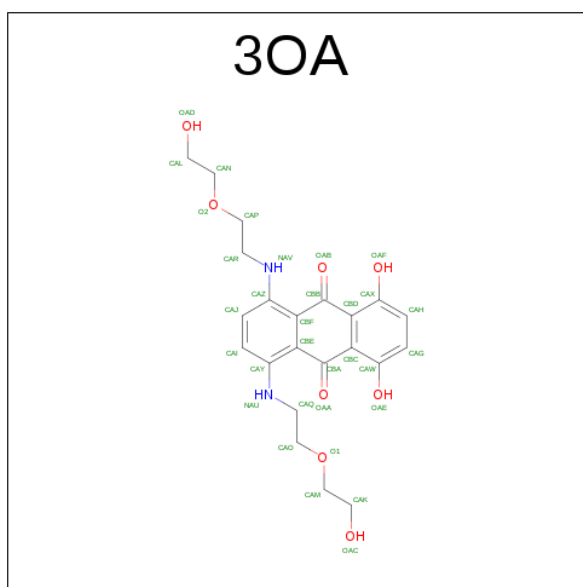
- Molecule 1 is a protein called Serine/threonine-protein kinase pim-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	272	2224	1421	388	406	1	8	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25	GLY	-	EXPRESSION TAG	UNP P11309
A	26	PRO	-	EXPRESSION TAG	UNP P11309
A	27	HIS	-	EXPRESSION TAG	UNP P11309
A	28	MET	-	EXPRESSION TAG	UNP P11309

- Molecule 2 is 1,4-DIHYDROXY-5,8-BIS{[2-(2-HYDROXYETHOXY)ETHYL]AMINO}AN THRACENE-9,10-DIONE (three-letter code: 3OA) (formula: C₂₂H₂₆N₂O₈).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			32	22	2	8		

- Molecule 1: Serine/threonine-protein kinase pim-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	97.87Å 97.87Å 80.83Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	36.48 – 2.34 36.48 – 2.34	Depositor EDS
% Data completeness (in resolution range)	99.6 (36.48-2.34) 99.6 (36.48-2.34)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.53 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.191 , 0.222 0.198 , 0.229	Depositor DCC
R_{free} test set	951 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	45.8	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.049 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2256	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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

Bond lengths and bond angles in the following residue types are not validated in this section: 3OA, SEP

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	0.97	0/2272	1.03	8/3081 (0.3%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	278	ARG	NE-CZ-NH1	7.92	124.26	120.30
1	A	186	ASP	CB-CG-OD1	6.97	124.57	118.30
1	A	156	ARG	NE-CZ-NH2	-6.71	116.95	120.30
1	A	156	ARG	NE-CZ-NH1	6.45	123.53	120.30
1	A	290	MET	CG-SD-CE	-6.40	89.95	100.20
1	A	185	ILE	N-CA-C	5.98	127.14	111.00
1	A	170	ASP	CB-CG-OD1	5.50	123.25	118.30
1	A	159	HIS	CB-CA-C	5.05	120.50	110.40

There are no chirality outliers.

There are no planarity outliers.

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	2224	0	2165	38	0
2	A	32	0	25	6	0
All	All	2256	0	2190	41	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 9.

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:80:LEU:HD23	1:A:94:LYS:HE2	1.54	0.88
1:A:150:GLN:HE22	1:A:181:GLU:HA	1.50	0.74
1:A:127:GLN:HG3	2:A:401:3OA:CAK	2.18	0.73
1:A:259:VAL:HG12	1:A:263:CYS:HB3	1.74	0.68
1:A:79:GLU:OE2	1:A:83:GLY:HA2	1.94	0.67
1:A:80:LEU:HD11	1:A:86:VAL:CG1	2.27	0.65
1:A:127:GLN:HG3	2:A:401:3OA:H2	1.77	0.65
1:A:80:LEU:HD12	1:A:80:LEU:H	1.63	0.63
2:A:401:3OA:OAF	2:A:401:3OA:OAB	2.16	0.63
1:A:126:VAL:HA	1:A:175:ILE:O	2.03	0.58
1:A:259:VAL:CG1	1:A:263:CYS:HB3	2.34	0.57
1:A:80:LEU:HD11	1:A:86:VAL:HG13	1.87	0.56
1:A:112:ARG:HB3	1:A:113:PRO:CD	2.37	0.55
2:A:401:3OA:NAV	2:A:401:3OA:OAB	2.41	0.54
1:A:80:LEU:HD11	1:A:86:VAL:HG11	1.92	0.52
1:A:100:PHE:O	1:A:101:SER:HB3	2.10	0.50
1:A:37:GLN:O	1:A:58:VAL:HG22	2.12	0.50
2:A:401:3OA:NAU	2:A:401:3OA:OAA	2.41	0.49
1:A:106:LEU:HA	1:A:120:LEU:HD23	1.93	0.49
1:A:269:TRP:CD1	1:A:279:PRO:HD3	2.48	0.49
1:A:38:TYR:CE1	1:A:57:ARG:HG3	2.48	0.48
1:A:150:GLN:NE2	1:A:182:LEU:H	2.11	0.47
1:A:152:LEU:CD1	1:A:290:MET:HE2	2.45	0.46
1:A:72:ASP:OD1	1:A:72:ASP:N	2.44	0.45
1:A:86:VAL:HB	1:A:87:PRO:CD	2.46	0.45
1:A:127:GLN:HA	2:A:401:3OA:H2	1.98	0.45
1:A:58:VAL:O	1:A:59:SER:C	2.54	0.44
1:A:112:ARG:HB3	1:A:113:PRO:HD2	1.99	0.44
1:A:80:LEU:O	1:A:81:PRO:C	2.54	0.44
1:A:152:LEU:CD1	1:A:290:MET:CE	2.96	0.43
1:A:214:ARG:HB3	1:A:214:ARG:CZ	2.48	0.43
1:A:215:TYR:O	1:A:217:ARG:HG2	2.19	0.42
1:A:293:VAL:HG22	1:A:294:LEU:N	2.34	0.42
1:A:58:VAL:O	1:A:60:ASP:N	2.52	0.42
1:A:80:LEU:HD23	1:A:94:LYS:CE	2.39	0.41
1:A:86:VAL:HB	1:A:87:PRO:HD2	2.03	0.41
1:A:176:ASP:C	1:A:176:ASP:OD1	2.59	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:VAL:HG12	1:A:41:GLY:N	2.36	0.40
1:A:148:PHE:CE2	1:A:152:LEU:HD11	2.56	0.40
1:A:293:VAL:HG22	1:A:294:LEU:O	2.20	0.40
1:A:79:GLU:OE2	1:A:83:GLY:CA	2.68	0.40

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	267/289 (92%)	261 (98%)	5 (2%)	1 (0%)	38	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	59	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	243/257 (95%)	217 (89%)	26 (11%)	8	7

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	59	SER
1	A	62	LEU
1	A	72	ASP
1	A	76	ASP
1	A	80	LEU
1	A	82	ASN
1	A	84	THR
1	A	104	ILE
1	A	105	ARG
1	A	124	GLU
1	A	128	ASP
1	A	139	LEU
1	A	159	HIS
1	A	160	ASN
1	A	177	LEU
1	A	189	SER
1	A	196	THR
1	A	200	ASP
1	A	217	ARG
1	A	239	ASP
1	A	243	GLU
1	A	252	GLN
1	A	256	ARG
1	A	298	GLU
1	A	305	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	150	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](#)

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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	$\# Z > 2$
1	SEP	A	261	1	9,9,10	0.68	0	9,12,14	1.23	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
1	SEP	A	261	1	-	0/5/8/10	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.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

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
2	3OA	A	401	-	34,34,34	2.89	8 (23%)	46,46,46	1.83	13 (28%)

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
2	3OA	A	401	-	-	0/14/30/30	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	3OA	CBF-CBB	-3.50	1.39	1.47
2	A	401	3OA	CBC-CBA	-3.24	1.39	1.47
2	A	401	3OA	CBD-CBB	-3.04	1.40	1.47
2	A	401	3OA	CBD-CAX	-2.62	1.37	1.41
2	A	401	3OA	CBD-CBC	-2.23	1.36	1.41
2	A	401	3OA	CAR-CAP	2.36	1.59	1.50
2	A	401	3OA	OAA-CBA	9.43	1.39	1.22
2	A	401	3OA	OAB-CBB	10.84	1.42	1.22

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	3OA	CBE-CAY-NAU	-4.69	115.84	121.35
2	A	401	3OA	CBF-CAZ-NAV	-3.88	116.79	121.35
2	A	401	3OA	OAF-CAX-CBD	-3.57	114.32	121.14
2	A	401	3OA	OAA-CBA-CBC	-2.44	117.47	121.49
2	A	401	3OA	CAO-CAQ-NAU	-2.34	105.62	111.18
2	A	401	3OA	CBD-CBC-CBA	-2.25	116.85	120.01
2	A	401	3OA	CBE-CBF-CBB	-2.20	116.92	120.01
2	A	401	3OA	OAB-CBB-CBD	-2.07	118.09	121.49
2	A	401	3OA	CAJ-CAZ-NAV	2.02	125.22	121.78
2	A	401	3OA	CBE-CBF-CAZ	2.03	121.26	119.78
2	A	401	3OA	CBD-CBB-CBF	2.42	122.28	118.07
2	A	401	3OA	CBC-CBA-CBE	2.64	122.65	118.07
2	A	401	3OA	CAP-CAR-NAV	4.57	122.02	111.18

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	3OA	6	0

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	271/289 (93%)	0.03	10 (3%)	42 53	29, 48, 81, 105	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	81	PRO	4.6
1	A	80	LEU	2.7
1	A	83	GLY	2.7
1	A	85	ARG	2.6
1	A	82	ASN	2.6
1	A	173	ILE	2.5
1	A	58	VAL	2.3
1	A	59	SER	2.1
1	A	73	ARG	2.0
1	A	84	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	SEP	A	261	10/11	0.88	0.20	-	15,22,45,58	0

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(\AA^2)	Q<0.9
2	3OA	A	401	32/32	0.90	0.21	1.96	46,71,118,127	0

6.5 Other polymers [i](#)

There are no such residues in this entry.