wwPDB X-ray Structure Validation Summary Report (i)

Feb 28, 2014 – 04:10 AM GMT

PDB ID	:	3MV5
Title	:	Crystal structure of Akt-1-inhibitor complexes
Authors	:	Pandit, J.
Deposited on	:	2010-05-03
Resolution	:	2.47 Å(reported)

1 Overall quality at a glance ⓘ

The reported resolution of this entry is 2.47 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	Percentile Ranks	Value
Rfree		0.281
Clashscore		8
Ramachandran outliers		0.3%
Sidechain outliers		8.8%
RSRZ outliers		0.9%

Worse ◼ Better

■ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution
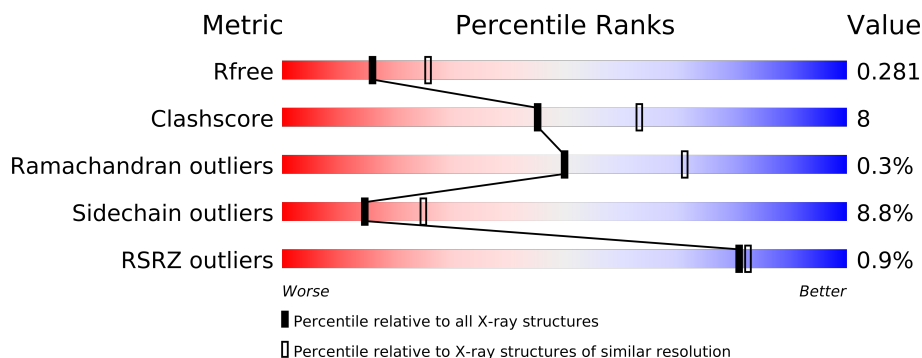
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R$_{free}$	66092	3277 (2.50-2.46)
Clashscore	79885	4136 (2.50-2.46)
Ramachandran outliers	78287	4052 (2.50-2.46)
Sidechain outliers	78261	4054 (2.50-2.46)
RSRZ outliers	66119	3279 (2.50-2.46)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	342	
2	C	10	

2 Entry composition ⓘ

There are 5 unique types of molecules in this entry. The entry contains 2749 atoms, of which 0 are hydrogen and 0 are deuterium.

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 v-akt murine thymoma viral oncogene homolog 1 (AKT1).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	316	2597	1668	437	476	1	15	0	1	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	139	GLY	-	EXPRESSION TAG	UNP B2RAM5
A	140	ALA	-	EXPRESSION TAG	UNP B2RAM5
A	141	MET	-	EXPRESSION TAG	UNP B2RAM5
A	142	ASP	-	EXPRESSION TAG	UNP B2RAM5
A	143	PRO	-	EXPRESSION TAG	UNP B2RAM5
A	473	ASP	SER	ENGINEERED	UNP B2RAM5

- Molecule 2 is a protein called GSK3-beta peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	10	79	47	16	16	0	0	0

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mn		
3	A	1	1	1	0	0

- Molecule 4 is (3R)-1-(5-METHYL-7H-PYRROLO[2,3-D]PYRIMIDIN-4-YL)PYRROLIDIN -3-AMINE (three-letter code: XFE) (formula: $C_{11}H_{15}N_5$).
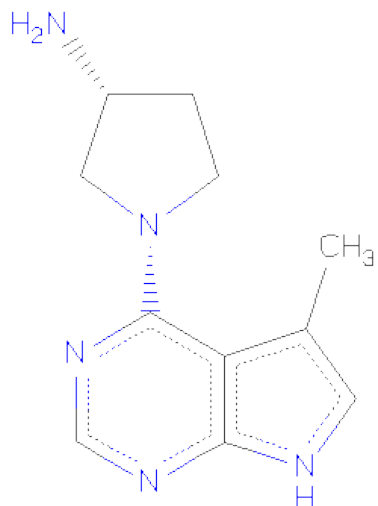
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total 16	C 11	N 5	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	55	Total 55	O 55	0	0
5	C	1	Total 1	O 1	0	0

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 (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: v-akt murine thymoma viral oncogene homolog 1 (AKT1)

Chain A:
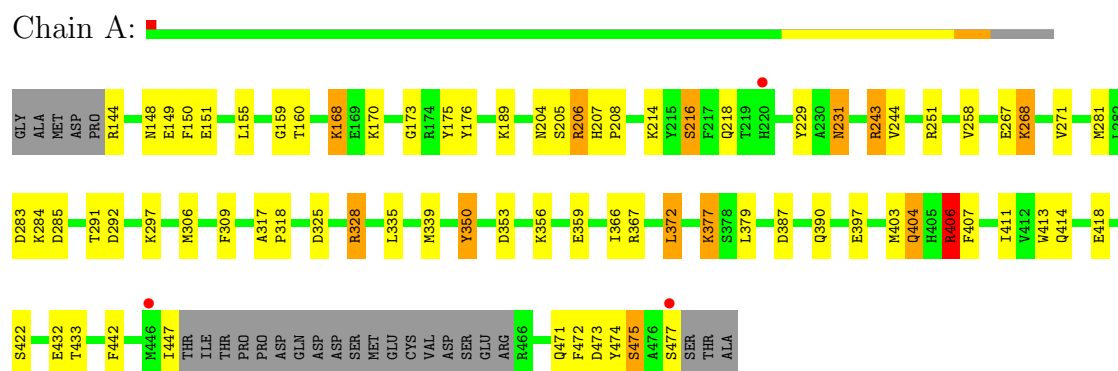
• Molecule 2: GSK3-beta peptide

Chain C:
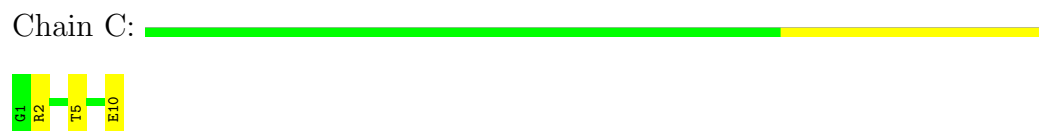
4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	42.71Å 55.42Å 93.29Å 90.00° 104.00° 90.00°	Depositor
Resolution (Å)	19.92 – 2.47 19.92 – 2.47	Depositor EDS
% Data completeness (in resolution range)	96.8 (19.92-2.47) 96.9 (19.92-2.47)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	2.10 (at 2.47Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.200 , 0.282 0.200 , 0.281	Depositor DCC
R_{free} test set	758 reflections (5.37%)	DCC
Wilson B-factor (Å^2)	34.2	Xtriage
Anisotropy	0.138	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 20.7	EDS
Estimated twinning fraction	0.013 for h,-k,-h-l	Xtriage
L-test for twinning	$<|L|> = 0.49, <L^2> = 0.32$	Xtriage
Outliers	0 of 14880 reflections	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	2749	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.08% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: TPO, MN, XFE

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	1.15	7/2647 (0.3%)	1.07	10/3558 (0.3%)
2	C	1.12	0/80	0.90	0/105
All	All	1.15	7/2727 (0.3%)	1.06	10/3663 (0.3%)

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	350	TYR	CD2-CE2	8.96	1.52	1.39
1	A	350	TYR	CE2-CZ	6.59	1.47	1.38
1	A	243	ARG	CG-CD	6.08	1.67	1.51
1	A	229	TYR	CD2-CE2	5.57	1.47	1.39
1	A	397	GLU	CG-CD	5.55	1.60	1.51

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	328	ARG	NE-CZ-NH2	-7.32	116.64	120.30
1	A	350	TYR	CB-CG-CD1	-7.29	116.62	121.00
1	A	350	TYR	CB-CG-CD2	6.71	125.03	121.00
1	A	325	ASP	CB-CG-OD1	6.66	124.29	118.30
1	A	367	ARG	NE-CZ-NH1	5.75	123.17	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens

added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2597	0	2560	44	0
2	C	79	0	77	5	0
3	A	1	0	0	0	0
4	A	16	0	15	0	0
5	A	55	0	0	5	0
5	C	1	0	0	0	0
All	All	2749	0	2652	45	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

The worst 5 of 45 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:309:PHE:HD2	2:C:10:GLU:HG2	1.13	1.05
1:A:390:GLN:HG3	5:A:40:HOH:O	1.65	0.95
1:A:309:PHE:CD2	2:C:10:GLU:HG2	2.06	0.88
1:A:404:GLN:HB2	5:A:7:HOH:O	1.72	0.87
1:A:309:PHE:HD2	2:C:10:GLU:CG	1.90	0.84

There are no symmetry-related clashes.

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	312/342 (91%)	298 (96%)	13 (4%)	1 (0%)	50	71
2	C	8/10 (80%)	8 (100%)	0	0	100	100
All	All	320/352 (91%)	306 (96%)	13 (4%)	1 (0%)	50	71

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	292	ASP

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	277/299 (93%)	252 (91%)	25 (9%)	14	24
2	C	8/8 (100%)	8 (100%)	0	100	100
All	All	285/307 (93%)	260 (91%)	25 (9%)	14	25

5 of 25 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	267	GLU
1	A	284	LYS
1	A	475	SER
1	A	268	LYS
1	A	285	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	204	ASN
1	A	231	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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	TPO	A	308	1	10,10,11	7.19	1 (10%)	12,14,16	1.96	5 (41%)

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	TPO	A	308	1	-	0/9/11/13	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	308	TPO	O-C	22.56	1.27	1.11

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	308	TPO	CB-CA-N	3.63	114.67	109.60
1	A	308	TPO	O3P-P-O1P	3.07	120.47	110.44
1	A	308	TPO	OG1-CB-CG2	-2.46	106.01	110.13
1	A	308	TPO	O3P-P-OG1	-2.44	100.06	107.09
1	A	308	TPO	O3P-P-O2P	2.41	116.99	107.61

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5   Carbohydrates (i)

There are no carbohydrates in this entry.

5.6   Ligand geometry (i)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	XFE	A	999	-	18,18,18	1.39	2 (11%)	24,26,26	3.35	11 (45%)

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	XFE	A	999	-	-	0/4/13/13	0/1/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	999	XFE	C14-C10	4.22	1.50	1.42
4	A	999	XFE	C8-N9	3.00	1.38	1.32

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	999	XFE	N9-C8-N7	-8.87	121.29	128.71
4	A	999	XFE	C10-C11-N12	6.84	113.50	106.54
4	A	999	XFE	C14-C10-C11	-6.43	98.93	106.08
4	A	999	XFE	C8-N7-C6	4.47	121.22	111.53
4	A	999	XFE	C15-C14-C13	-3.85	119.49	127.91

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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, 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	316/342 (92%)	-0.19	3 (0%)	81	82	11, 30, 53, 65	0
2	C	10/10 (100%)	0.50	0	100	100	34, 38, 45, 52	0
All	All	326/352 (92%)	-0.17	3 (0%)	81	82	11, 30, 52, 65	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	477	SER	4.3
1	A	220	HIS	3.1
1	A	446	MET	2.3

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

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	RSR	LLDF	B-factors(Å^2)	Q<0.9
1	TPO	A	308	11/12	0.11	-0.73	33,35,44,46	0

6.3   Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å^2)	Q<0.9
3	MN	A	1	1/1	0.13	-0.53	63,63,63,63	0
4	XFE	A	999	16/16	0.11	-0.90	21,23,25,27	0

6.5 Other polymers ⓘ

There are no such residues in this entry.