



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:38 PM GMT

PDB ID : 4IZ7
Title : Structure of Non-Phosphorylated ERK2 bound to the PEA-15 Death Effector Domain
Authors : Mace, P.D.; Robinson, H.; Riedl, S.J.
Deposited on : 2013-01-29
Resolution : 1.80 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 (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

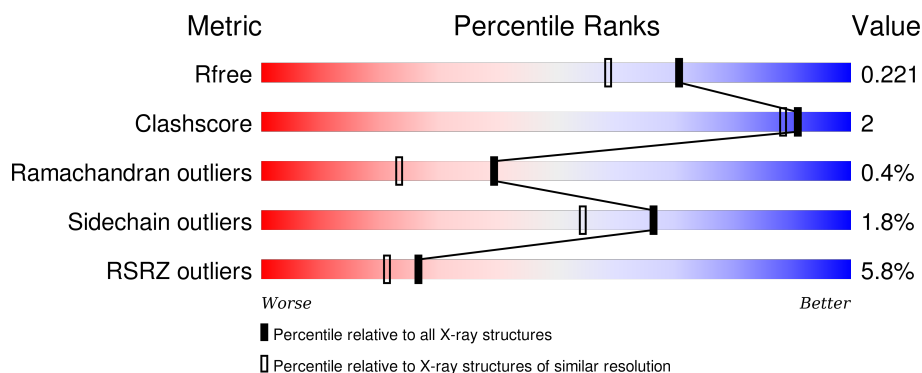
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (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. 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	356	<div> <div>6%</div> <div>91%</div> <div>5% . .</div> </div>
1	C	356	<div> <div>5%</div> <div>87%</div> <div>8% . .</div> </div>
2	B	97	<div> <div>7%</div> <div>81%</div> <div>11% 7%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6926 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 Mitogen-activated protein kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	340	Total	C	N	O	S	0	6	0
			2768	1782	478	493	15			
1	A	344	Total	C	N	O	S	0	5	0
			2804	1798	483	508	15			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	5	GLY	-	EXPRESSION TAG	UNP P28482
C	6	PRO	-	EXPRESSION TAG	UNP P28482
C	7	GLY	-	EXPRESSION TAG	UNP P28482
A	5	GLY	-	EXPRESSION TAG	UNP P28482
A	6	PRO	-	EXPRESSION TAG	UNP P28482
A	7	GLY	-	EXPRESSION TAG	UNP P28482

- Molecule 2 is a protein called Astrocytic phosphoprotein PEA-15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	90	Total	C	N	O	S	0	4	0
			704	440	112	148	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	GLY	-	EXPRESSION TAG	UNP Q9Z297

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

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

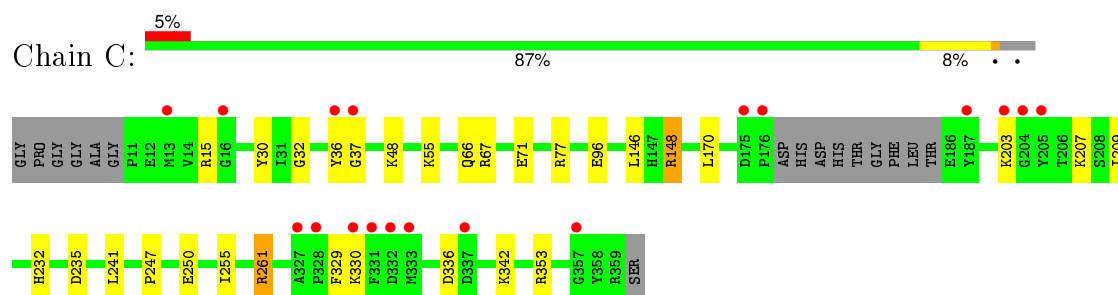
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	297	Total 297	O 297	0	0
4	A	284	Total 284	O 284	0	0
4	B	68	Total 68	O 68	0	0

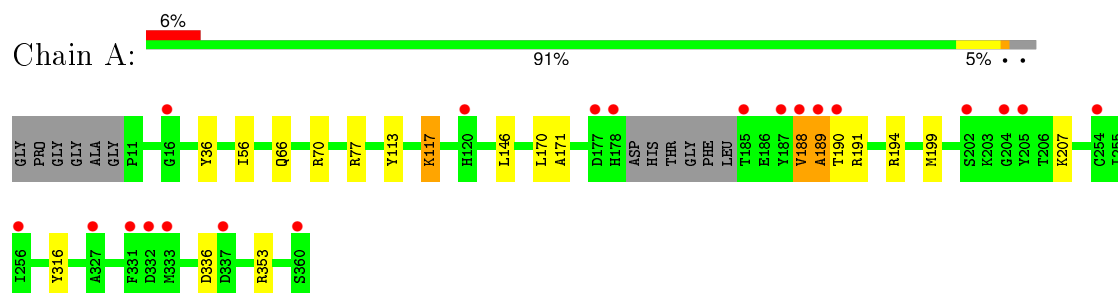
3 Residue-property plots [i](#)

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.

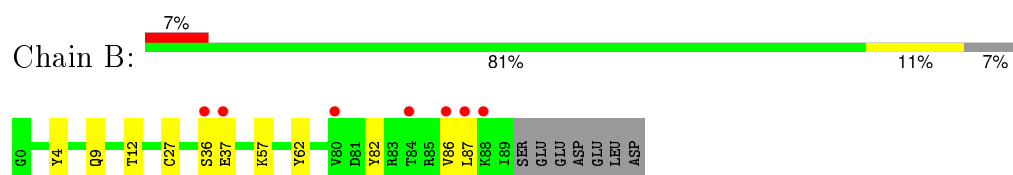
- Molecule 1: Mitogen-activated protein kinase 1



- Molecule 1: Mitogen-activated protein kinase 1



- Molecule 2: Astrocytic phosphoprotein PEA-15



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	73.48 Å 205.33 Å 57.75 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.70 – 1.80 29.67 – 1.80	Depositor EDS
% Data completeness (in resolution range)	98.8 (29.70-1.80) 98.9 (29.67-1.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 1.80 Å)	Xtriage
Refinement program	REFMAC 5.6.0107	Depositor
R, R_{free}	0.185 , 0.222 0.184 , 0.221	Depositor DCC
R_{free} test set	4107 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	22.4	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 81104 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6926	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.25% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

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

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	0.63	0/2899	0.66	0/3929
1	C	0.65	0/2864	0.75	4/3881 (0.1%)
2	B	0.64	0/735	0.68	0/998
All	All	0.64	0/6498	0.70	4/8808 (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 (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	261	ARG	NE-CZ-NH2	-10.98	114.81	120.30
1	C	261	ARG	NE-CZ-NH1	10.87	125.73	120.30
1	C	15	ARG	NE-CZ-NH1	5.60	123.10	120.30
1	C	148	ARG	NE-CZ-NH2	-5.34	117.63	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	188	VAL	Peptide
1	A	189	ALA	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	2804	0	2773	10	1
1	C	2768	0	2770	13	1
2	B	704	0	660	4	0
3	A	1	0	0	0	0
4	A	284	0	0	2	0
4	B	68	0	0	0	0
4	C	297	0	0	0	0
All	All	6926	0	6203	27	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 2.

All (27) 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:70[B]:ARG:HD3	1:A:171:ALA:O	1.66	0.95
1:A:70[A]:ARG:HD3	1:A:171:ALA:O	1.67	0.95
1:A:194:ARG:HD2	1:A:199:MET:HG3	1.73	0.70
1:C:148:ARG:HD3	1:C:170:LEU:O	1.92	0.68
1:C:146:LEU:HD22	1:C:207:LYS:HA	1.77	0.65
2:B:27[B]:CYS:HB3	2:B:62:TYR:OH	1.98	0.64
1:C:37:GLY:HA3	1:C:55:LYS:O	2.00	0.61
1:C:77:ARG:HD2	1:C:329:PHE:CD2	2.40	0.57
1:A:170:LEU:HD13	1:A:188:VAL:HB	1.89	0.54
1:C:36:TYR:OH	1:C:67:ARG:HD2	2.08	0.53
1:A:191:ARG:HD3	4:A:715:HOH:O	2.09	0.51
1:A:36:TYR:HB2	1:A:56:ILE:HG23	1.93	0.49
1:C:232:HIS:CD2	1:C:235:ASP:H	2.30	0.49
1:A:113:TYR:CZ	1:A:117:LYS:HE3	2.47	0.49
1:C:241:LEU:HD22	1:C:247:PRO:HD3	1.96	0.47
1:A:66:GLN:CD	1:A:336:ASP:HA	2.34	0.47
1:C:66:GLN:OE1	1:C:336:ASP:HA	2.15	0.46
1:A:146:LEU:HD22	1:A:207:LYS:HA	1.97	0.46
1:C:255:ILE:O	1:C:261:ARG:HD2	2.16	0.45
2:B:82:TYR:O	2:B:86:VAL:HB	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:232:HIS:HD2	1:C:235:ASP:H	1.64	0.44
2:B:9:GLN:O	2:B:12[B]:THR:HG22	2.19	0.43
1:A:117:LYS:HE2	4:A:512:HOH:O	2.19	0.43
1:C:30:TYR:CZ	1:C:32:GLY:HA2	2.53	0.43
1:C:148:ARG:HG2	1:C:209:ILE:HD11	2.01	0.43
2:B:4:TYR:CD1	2:B:57:LYS:HA	2.56	0.41
1:C:96:GLU:H	1:C:96:GLU:CD	2.24	0.40

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:C:261:ARG:NH2	1:A:316:TYR:O[2_655]	2.10	0.10

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	345/356 (97%)	334 (97%)	10 (3%)	1 (0%)	46	29
1	C	342/356 (96%)	328 (96%)	14 (4%)	0	100	100
2	B	92/97 (95%)	87 (95%)	3 (3%)	2 (2%)	8	1
All	All	779/809 (96%)	749 (96%)	27 (4%)	3 (0%)	39	23

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	189	ALA
2	B	36	SER
2	B	87	LEU

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	308/318 (97%)	304 (99%)	4 (1%)	76	68
1	C	303/318 (95%)	296 (98%)	7 (2%)	58	42
2	B	80/91 (88%)	79 (99%)	1 (1%)	76	68
All	All	691/727 (95%)	679 (98%)	12 (2%)	66	57

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	48	LYS
1	C	71	GLU
1	C	203	LYS
1	C	250	GLU
1	C	330	LYS
1	C	342	LYS
1	C	353	ARG
1	A	77	ARG
1	A	117	LYS
1	A	190	THR
1	A	353	ARG
2	B	37	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	47	ASN
1	C	61	HIS
1	C	123	ASN
1	C	232	HIS
1	C	238	ASN
1	A	47	ASN
1	A	123	ASN
1	A	276	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 1 ligands modelled in this entry, 1 is 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 ⓘ

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	344/356 (96%)	0.22	20 (5%)	26 21	11, 22, 41, 54	1 (0%)
1	C	340/356 (95%)	0.05	18 (5%)	30 25	9, 20, 39, 55	0
2	B	90/97 (92%)	0.11	7 (7%)	16 12	14, 22, 42, 51	0
All	All	774/809 (95%)	0.14	45 (5%)	26 21	9, 21, 41, 55	1 (0%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	189	ALA	9.3
1	C	333	MET	5.8
2	B	37	GLU	5.2
1	C	332	ASP	4.8
1	C	176	PRO	4.8
1	A	332	ASP	4.7
1	C	16	GLY	4.0
2	B	36	SER	3.8
1	C	331	PHE	3.8
1	A	178	HIS	3.7
1	A	333	MET	3.7
1	C	175	ASP	3.6
1	C	357	GLY	3.6
1	A	360	SER	3.5
1	A	254	CYS	3.5
1	A	177	ASP	3.4
1	C	203	LYS	3.3
1	C	204	GLY	3.1
1	C	36	TYR	3.1
1	A	187	TYR	3.0
1	A	16	GLY	2.9
1	A	185	THR	2.8
1	A	337	ASP	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	187	TYR	2.7
1	A	190	THR	2.6
1	A	204	GLY	2.6
2	B	86	VAL	2.5
1	A	205	TYR	2.5
1	C	205	TYR	2.4
1	C	327	ALA	2.4
1	C	337	ASP	2.3
1	A	331	PHE	2.3
1	A	256	ILE	2.3
2	B	80	VAL	2.3
2	B	84	THR	2.3
2	B	88	LYS	2.3
1	A	188	VAL	2.2
1	A	202	SER	2.2
1	A	120	HIS	2.2
1	C	37	GLY	2.1
2	B	87	LEU	2.1
1	C	330	LYS	2.1
1	C	328	PRO	2.1
1	C	13	MET	2.1
1	A	327	ALA	2.0

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(\AA^2)	Q<0.9
3	NA	A	401	1/1	0.96	0.22	1.42	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.