



wwPDB X-ray Structure Validation Summary Report ⓘ

May 3, 2025 – 09:58 AM EDT

PDB ID : 3VUT / pdb_00003vut
Title : Crystal structures of non-phosphorylated MAP2K4
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Deposited on : 2012-07-05
Resolution : 3.50 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0rc1
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.43.1

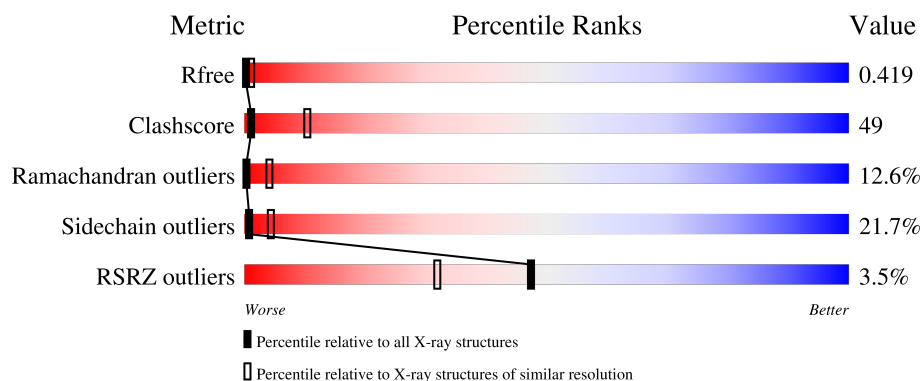
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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}	164625	1094 (3.56-3.44)
Clashscore	180529	1045 (3.54-3.46)
Ramachandran outliers	177936	1032 (3.54-3.46)
Sidechain outliers	177891	1033 (3.54-3.46)
RSRZ outliers	164620	1093 (3.56-3.44)

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 $\leq 5\%$. 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	327	<div> <div> <div></div> <div>24%</div> <div>25%</div> <div>9%</div> <div>•</div> <div>38%</div> </div> </div>
1	B	327	<div> <div>4%</div> <div>28%</div> <div>34%</div> <div>14%</div> <div>•</div> <div>20%</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3457 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 Dual specificity mitogen-activated protein kinase kinase 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	S	0	0	0
			1563	1009	251	291	12			
1	B	260	Total	C	N	O	S	0	0	0
			1894	1197	309	376	12			

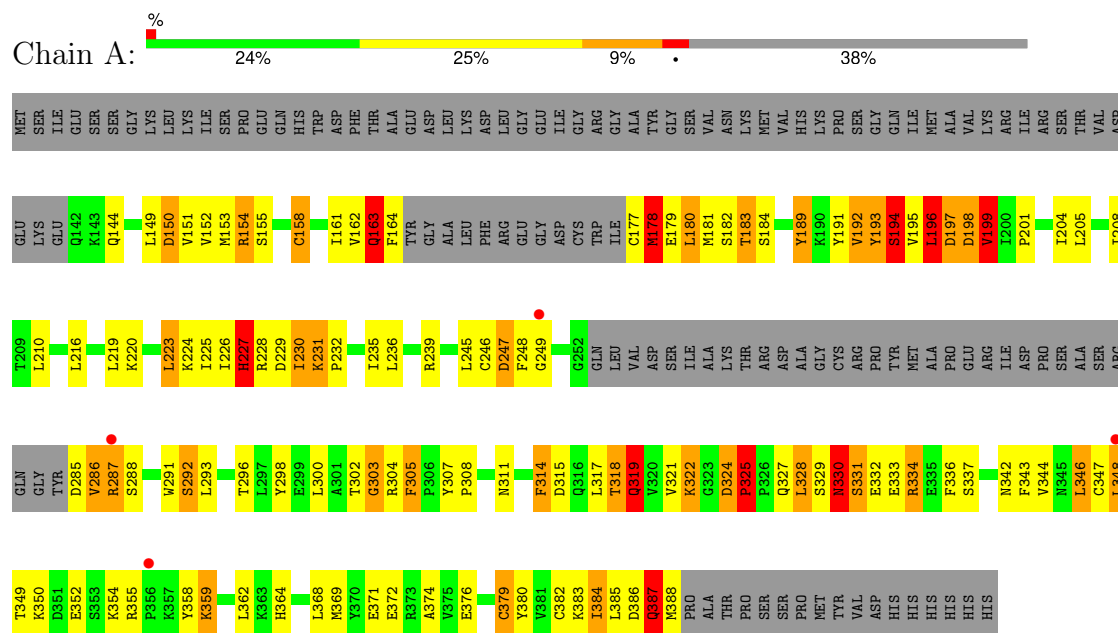
There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	79	MET	-	expression tag	UNP P45985
A	400	HIS	-	expression tag	UNP P45985
A	401	HIS	-	expression tag	UNP P45985
A	402	HIS	-	expression tag	UNP P45985
A	403	HIS	-	expression tag	UNP P45985
A	404	HIS	-	expression tag	UNP P45985
A	405	HIS	-	expression tag	UNP P45985
B	79	MET	-	expression tag	UNP P45985
B	400	HIS	-	expression tag	UNP P45985
B	401	HIS	-	expression tag	UNP P45985
B	402	HIS	-	expression tag	UNP P45985
B	403	HIS	-	expression tag	UNP P45985
B	404	HIS	-	expression tag	UNP P45985
B	405	HIS	-	expression tag	UNP P45985

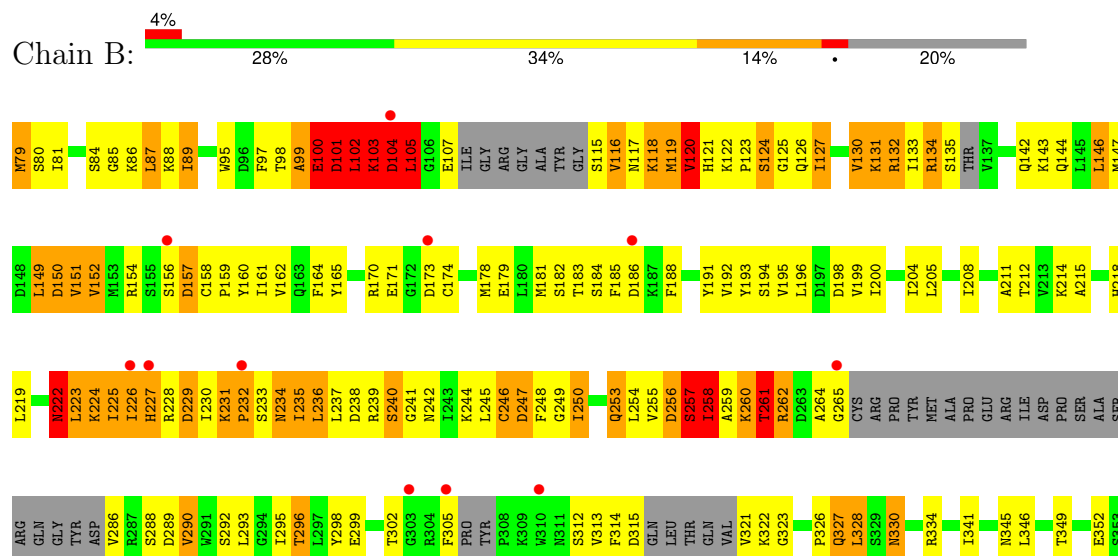
3 Residue-property plots [i](#)

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 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: Dual specificity mitogen-activated protein kinase kinase 4



- Molecule 1: Dual specificity mitogen-activated protein kinase kinase 4



K354	K355	P356	K357	Y358	K359	E360	L361	L362	K363	H364		I367	L368		E371	E372	R373	ALA	VAL	GLU	VAL	VAL	ALA	CYS	TYR	VAL	CYS	LYS	ILE	LEU	ASP	GLN	MET	PRO	ALA	THR	PRO	SER	SER	PRO	PRO	MET	TYR	VAL	ASP	HIS	HIS	HIS	HIS	HIS	HIS
------	------	------	------	------	------	------	------	------	------	------	--	------	------	--	------	------	------	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.77Å 87.36Å 118.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.50 20.00 – 3.50	Depositor EDS
% Data completeness (in resolution range)	98.2 (20.00-3.50) 97.7 (20.00-3.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.34 (at 3.48Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.331 , 0.407 0.331 , 0.419	Depositor DCC
R_{free} test set	593 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	120.1	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 77.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	3457	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.14% 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

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	0.74	0/1590	1.13	8/2153 (0.4%)
1	B	0.83	2/1919 (0.1%)	1.28	18/2597 (0.7%)
All	All	0.79	2/3509 (0.1%)	1.22	26/4750 (0.5%)

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	B	0	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	257	SER	N-CA	-5.56	1.39	1.46
1	B	261	THR	CA-C	5.04	1.58	1.52

The worst 5 of 26 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	158	CYS	CA-C-N	12.33	132.01	119.56
1	A	158	CYS	C-N-CA	12.33	132.01	119.56
1	B	118	LYS	N-CA-C	-10.32	95.07	110.24
1	B	257	SER	N-CA-C	-8.86	91.93	110.80
1	B	260	LYS	N-CA-C	7.54	120.30	110.43

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	100	GLU	Peptide

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Mol	Chain	Res	Type	Group
1	B	103	LYS	Peptide
1	B	105	LEU	Peptide
1	B	244	LYS	Peptide

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	1563	0	1540	99	0
1	B	1894	0	1765	231	0
All	All	3457	0	3305	330	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 49.

The worst 5 of 330 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:B:219:LEU:O	1:B:222:ASN:O	1.54	1.25
1:B:115:SER:O	1:B:131:LYS:O	1.53	1.24
1:B:133:ILE:CB	1:B:174:CYS:O	1.87	1.22
1:B:223:LEU:HG	1:B:224:LYS:CB	1.69	1.20
1:B:256:ASP:OD2	1:B:261:THR:HG23	1.47	1.13

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	197/327 (60%)	142 (72%)	26 (13%)	29 (15%)	0	3
1	B	248/327 (76%)	169 (68%)	52 (21%)	27 (11%)	0	5
All	All	445/654 (68%)	311 (70%)	78 (18%)	56 (13%)	0	4

5 of 56 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	178	MET
1	A	247	ASP
1	A	287	ARG
1	A	330	ASN
1	A	334	ARG

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	172/296 (58%)	130 (76%)	42 (24%)	0	3
1	B	196/296 (66%)	158 (81%)	38 (19%)	1	6
All	All	368/592 (62%)	288 (78%)	80 (22%)	1	4

5 of 80 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	204	ILE
1	B	261	THR
1	B	222	ASN
1	B	246	CYS
1	B	328	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	253	GLN

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Mol	Chain	Res	Type
1	B	345	ASN
1	B	126	GLN
1	B	218	HIS
1	B	222	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	203/327 (62%)	0.29	4 (1%) 64 45	64, 98, 143, 166	0
1	B	260/327 (79%)	0.44	12 (4%) 38 27	58, 99, 134, 171	0
All	All	463/654 (70%)	0.37	16 (3%) 47 32	58, 99, 137, 171	0

The worst 5 of 16 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	104	ASP	5.2
1	B	173	ASP	4.1
1	B	227	HIS	3.4
1	B	303	GLY	3.0
1	A	249	GLY	2.9

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 monosaccharides 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.