



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 17, 2024 – 11:06 PM EDT

PDB ID : 5R9G
Title : PanDDA analysis group deposition Form1 MAP kinase p38-alpha – Fragment PC587 in complex with MAP kinase p38-alpha
Authors : De Nicola, G.F.; Nichols, C.E.
Deposited on : 2020-03-04
Resolution : 1.73 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 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	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
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)	:	2.37.1

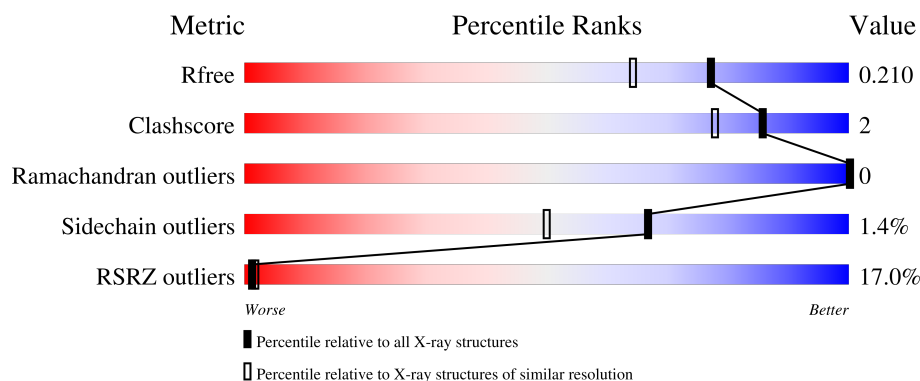
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

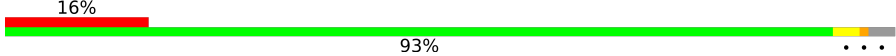
The reported resolution of this entry is 1.73 Å.

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	3764 (1.76-1.72)
Clashscore	141614	3923 (1.76-1.72)
Ramachandran outliers	138981	3878 (1.76-1.72)
Sidechain outliers	138945	3878 (1.76-1.72)
RSRZ outliers	127900	3705 (1.76-1.72)

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	360	

2 Entry composition [i](#)

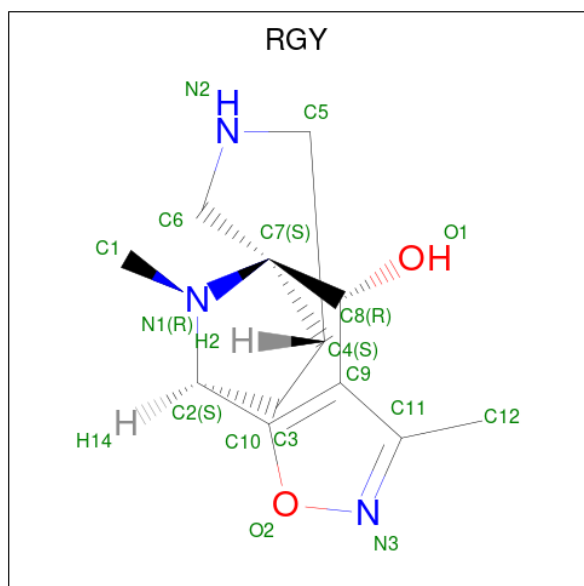
There are 7 unique types of molecules in this entry. The entry contains 3071 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	348	Total	C	N	O	S	0	3	0
			2770	1775	465	515	15			

- Molecule 2 is (4R,4aS,7aS,9S)-3,10-dimethyl-5,6,7,7a,8,9-hexahydro-4H-4a,9-epiminopyrrolo [3',4':5,6]cyclohepta[1,2-d][1,2]oxazol-4-ol (three-letter code: RGY) (formula: C₁₂H₁₇N₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			17	12	3	2		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	5	Total	Cl	0	0
			5	5		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

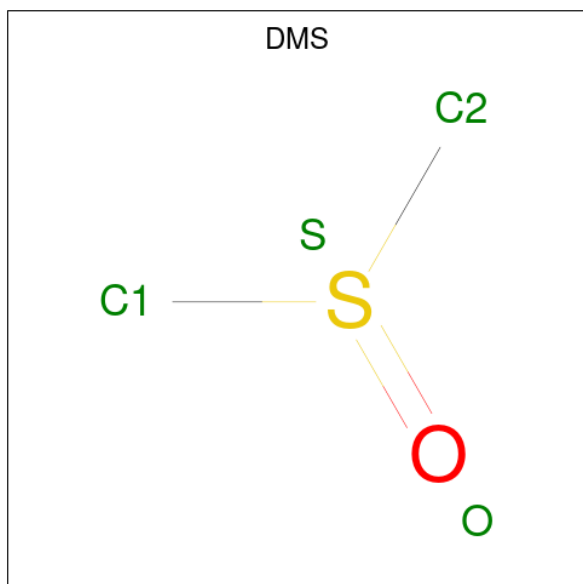
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C₂H₆OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	O	S	0	0
			4	2	1	1		

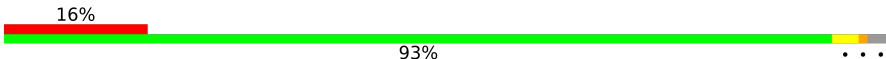
- Molecule 7 is water.

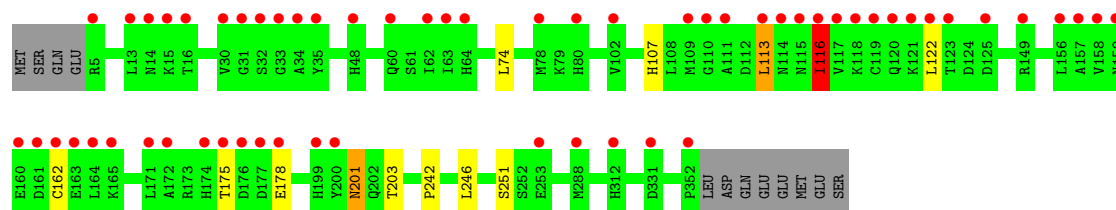
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	269	Total	O	0	0
			269	269		

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: Mitogen-activated protein kinase 14

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.91Å 85.94Å 127.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.34 – 1.73 29.32 – 1.73	Depositor EDS
% Data completeness (in resolution range)	99.2 (29.34-1.73) 99.2 (29.32-1.73)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 1.73Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.176 , 0.195 0.206 , 0.210	Depositor DCC
R_{free} test set	2645 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.021	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 54.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3071	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.08% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: DMS, CL, SO4, MG, RGY

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	0.92	1/2836 (0.0%)	0.83	0/3859

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	116	ILE	C-N	13.40	1.64	1.34

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2770	0	2690	11	0
2	A	17	0	0	1	0
3	A	5	0	0	0	0
4	A	1	0	0	0	0
5	A	5	0	0	1	0
6	A	4	0	6	0	0
7	A	269	0	0	2	0
All	All	3071	0	2696	12	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 2.

All (12) 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)
5:A:408:SO4:O3	7:A:501:HOH:O	2.01	0.78
1:A:162:CYS:SG	2:A:401:RGY:N2	2.82	0.53
1:A:116:ILE:HD12	1:A:122:LEU:HD21	1.90	0.51
1:A:201:ASN:HD22	1:A:201:ASN:C	2.14	0.50
1:A:175:THR:HG22	1:A:178:GLU:CG	2.43	0.48
1:A:242:PRO:HB3	1:A:246:LEU:HD23	1.94	0.48
1:A:107:HIS:HD2	7:A:534:HOH:O	1.95	0.48
1:A:74:LEU:HD13	1:A:74:LEU:C	2.36	0.45
1:A:113:LEU:HD23	1:A:113:LEU:HA	1.81	0.42
1:A:201:ASN:HD22	1:A:203:THR:H	1.67	0.41
1:A:201:ASN:ND2	1:A:203:THR:OG1	2.54	0.41
1:A:175:THR:HG23	1:A:178:GLU:H	1.85	0.40

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	349/360 (97%)	339 (97%)	10 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	294/319 (92%)	290 (99%)	4 (1%)	67	50

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	LEU
1	A	116	ILE
1	A	201	ASN
1	A	251	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	HIS
1	A	115	ASN
1	A	128	GLN
1	A	201	ASN
1	A	272	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 6 are monoatomic - leaving 3 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
6	DMS	A	409	-	3,3,3	0.30	0	3,3,3	0.09	0
2	RGY	A	401	-	11,20,20	1.62	3 (27%)	8,33,33	1.79	2 (25%)
5	SO4	A	408	-	4,4,4	0.35	0	6,6,6	0.22	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
2	RGY	A	401	-	-	-	0/5/4/4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	RGY	C9-C8	-3.10	1.49	1.51
2	A	401	RGY	C2-N1	3.01	1.52	1.48
2	A	401	RGY	C1-N1	2.18	1.49	1.46

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	RGY	C5-N2-C6	3.46	113.58	105.42
2	A	401	RGY	O1-C8-C7	-2.52	107.92	111.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	RGY	1	0
5	A	408	SO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	116:ILE	C	117:VAL	N	1.64

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	348/360 (96%)	1.10	59 (16%) ⓘ ⓘ	13, 25, 50, 83	33 (9%)

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	164	LEU	21.6
1	A	156	LEU	20.9
1	A	117	VAL	20.0
1	A	113	LEU	19.2
1	A	78[A]	MET	15.5
1	A	157	ALA	13.7
1	A	80	HIS	13.3
1	A	122	LEU	12.9
1	A	158	VAL	11.7
1	A	162	CYS	11.7
1	A	114	ASN	11.5
1	A	171	LEU	11.2
1	A	116	ILE	10.5
1	A	165	LYS	10.4
1	A	118	LYS	10.3
1	A	288[A]	MET	9.9
1	A	119	CYS	9.9
1	A	123	THR	9.0
1	A	312	HIS	8.9
1	A	34	ALA	8.0
1	A	149	ARG	7.8
1	A	33	GLY	7.4
1	A	163	GLU	6.2
1	A	125[A]	ASP	6.1
1	A	115	ASN	6.0
1	A	161	ASP	5.6
1	A	352	PRO	5.6

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Mol	Chain	Res	Type	RSRZ
1	A	35	TYR	5.6
1	A	175	THR	5.3
1	A	159	ASN	5.0
1	A	110	GLY	4.8
1	A	32	SER	4.6
1	A	15	LYS	4.6
1	A	178	GLU	4.2
1	A	120	GLN	4.1
1	A	63	ILE	4.0
1	A	174	HIS	3.8
1	A	160	GLU	3.6
1	A	177	ASP	3.4
1	A	14	ASN	3.4
1	A	13	LEU	3.3
1	A	30	VAL	3.1
1	A	172	ALA	3.1
1	A	31	GLY	3.0
1	A	109	MET	3.0
1	A	121	LYS	3.0
1	A	199	HIS	2.9
1	A	111	ALA	2.9
1	A	16	THR	2.7
1	A	176	ASP	2.5
1	A	5	ARG	2.5
1	A	331	ASP	2.5
1	A	253	GLU	2.4
1	A	200	TYR	2.4
1	A	62	ILE	2.3
1	A	102	VAL	2.3
1	A	48	HIS	2.2
1	A	60	GLN	2.2
1	A	64	HIS	2.2

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no monosaccharides 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. 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	B-factors(Å ²)	Q<0.9
2	RGY	A	401	17/17	0.37	0.34	38,39,40,41	17
6	DMS	A	409	4/4	0.83	0.14	73,77,79,87	0
5	SO4	A	408	5/5	0.90	0.21	52,56,61,82	0
3	CL	A	404	1/1	0.91	0.06	50,50,50,50	0
3	CL	A	405	1/1	0.97	0.06	38,38,38,38	0
4	MG	A	407	1/1	0.97	0.30	31,31,31,31	0
3	CL	A	402	1/1	0.99	0.11	43,43,43,43	0
3	CL	A	403	1/1	0.99	0.03	42,42,42,42	0
3	CL	A	406	1/1	0.99	0.05	33,33,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.