



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 02:49 PM GMT

PDB ID : 4ANB  
Title : Crystal structures of human MEK1 with carboxamide-based allosteric inhibitor XL518 (GDC-0973), or related analogs.  
Authors : Rice, K.D.; Aay, N.; Anand, N.K.; Blazey, C.M.; Bowles, O.J.; Bussenius, J.; Costanzo, S.; Curtis, J.K.; Defina, S.C.; Dubenko, L.; Engst, S.; Joshi, A.A.; Kennedy, A.R.; Kim, A.I.; Koltun, E.S.; Loughheed, J.C.; Manalo, J.C.L.; Martini, J.F.; Nuss, J.M.; Peto, C.J.; Tsang, T.H.; Yu, P.; Johnston, S.  
Deposited on : 2012-03-16  
Resolution : 2.20 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](mailto: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.

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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

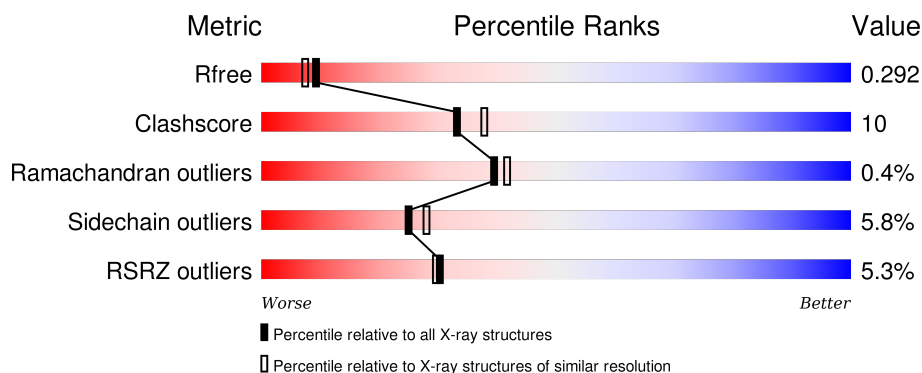
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


The reported resolution of this entry is 2.20 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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  $\geq 3$ , 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	301	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 2324 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.

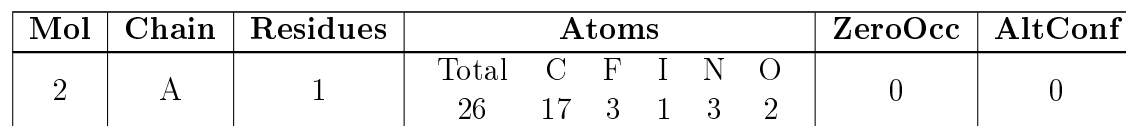
- Molecule 1 is a protein called DUAL SPECIFICITY MITOGEN-ACTIVATED PROTEIN KINASE KINASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	282	Total	C	N	O	S	0	0	0
			2209	1412	376	405	16			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	58	MET	-	EXPRESSION TAG	UNP Q02750
A	59	ALA	-	EXPRESSION TAG	UNP Q02750
A	60	LEU	-	EXPRESSION TAG	UNP Q02750
A	393	GLU	-	EXPRESSION TAG	UNP Q02750
A	394	GLY	-	EXPRESSION TAG	UNP Q02750
A	395	HIS	-	EXPRESSION TAG	UNP Q02750
A	396	HIS	-	EXPRESSION TAG	UNP Q02750
A	397	HIS	-	EXPRESSION TAG	UNP Q02750
A	398	HIS	-	EXPRESSION TAG	UNP Q02750
A	399	HIS	-	EXPRESSION TAG	UNP Q02750
A	400	HIS	-	EXPRESSION TAG	UNP Q02750
A	218	GLU	SER	ENGINEERED MUTATION	UNP Q02750
A	222	GLU	SER	ENGINEERED MUTATION	UNP Q02750
A	328	ALA	GLY	ENGINEERED MUTATION	UNP Q02750

- Molecule 2 is [3-(AMINOMETHYL)-3-OXIDANYL-AZETIDIN-1-YL]-[3,4-BIS(FLUORANYL)-2-[(2-FLUORANYL-4-IODANYL-PHENYL)AMINO]PHENYL]METHANONE (three-letter code: YQY) (formula: C<sub>17</sub>H<sub>15</sub>F<sub>3</sub>IN<sub>3</sub>O<sub>2</sub>).



- 
- ACP
- The image displays a chemical structure of a nucleotide derivative, specifically a dinucleotide. The structure is composed of a purine base, a ribose sugar, and a pyrophosphate group. The purine base is shown in blue, with atoms labeled N1, N3, N7, N9, C2, C4, C6, and C8. The ribose sugar is shown in red, with atoms labeled C1', C2', C3', C4', and C5'. The pyrophosphate group is shown in pink, with atoms labeled P1, P2, and P3. The structure is labeled with atom names and numbers, and is titled 'ACP'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 31	C 11	N 5	O 12	P 3	0	0

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

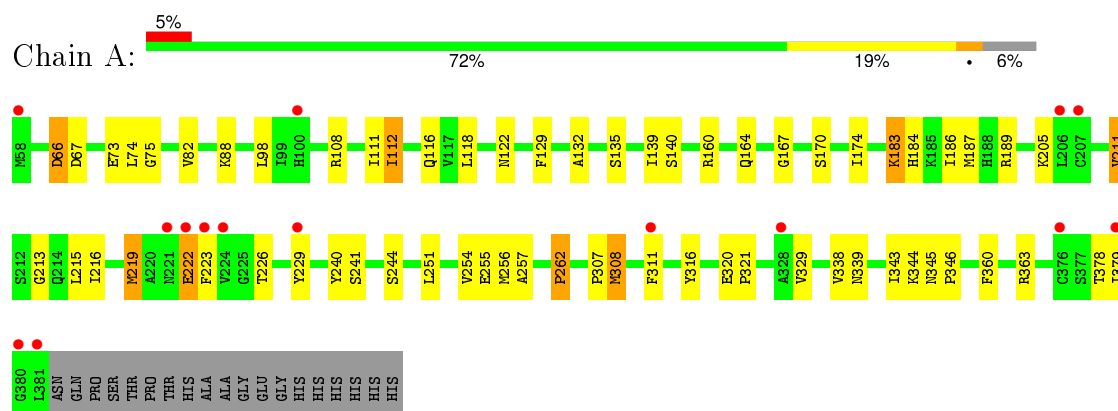
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	57	Total	O	0	0
			57	57		

### 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: DUAL SPECIFICITY MITOGEN-ACTIVATED PROTEIN KINASE KINASE 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	109.00 Å   109.00 Å   48.08 Å 90.00°   90.00°   120.00°	Depositor
Resolution (Å)	94.49 – 2.20 35.68 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.4 (94.49-2.20) 99.5 (35.68-2.20)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.04 (at 2.20 Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.226 , 0.296 0.223 , 0.292	Depositor DCC
$R_{free}$ test set	858 reflections (5.43%)	DCC
Wilson B-factor (Å <sup>2</sup> )	56.8	Xtriage
Anisotropy	0.093	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 49.0	EDS
Estimated twinning fraction	0.039 for h,-h-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 16672 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	2324	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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: MG, ACP, YQY

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	0.93	1/2254 (0.0%)	0.91	3/3038 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	262	PRO	C-N	13.59	1.65	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	98	LEU	CA-CB-CG	6.85	131.04	115.30
1	A	256	MET	CG-SD-CE	6.40	110.44	100.20
1	A	262	PRO	O-C-N	-6.32	112.59	122.70

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	2209	0	2237	43	0
2	A	26	0	15	3	0
3	A	31	0	14	0	0
4	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	57	0	0	5	2
All	All	2324	0	2266	46	2

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 10.

The worst 5 of 46 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:67:ASP:O	1:A:88:LYS:HE2	1.70	0.89
1:A:170:SER:O	1:A:174:ILE:HG12	1.79	0.81
1:A:311:PHE:CD2	5:A:2048:HOH:O	2.35	0.80
1:A:183:LYS:HB3	1:A:184:HIS:CD2	2.23	0.72
1:A:229:TYR:OH	1:A:255:GLU:OE1	2.09	0.70

All (2) 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)
5:A:2025:HOH:O	5:A:2047:HOH:O[5_554]	1.88	0.32
5:A:2026:HOH:O	5:A:2047:HOH:O[5_554]	2.16	0.04

## 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	280/301 (93%)	266 (95%)	13 (5%)	1 (0%)	39 42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	308	MET

### 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	243/258 (94%)	229 (94%)	14 (6%)	25 28

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	211	VAL
1	A	219	MET
1	A	329	VAL
1	A	183	LYS
1	A	244	SER

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	122	ASN
1	A	164	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	$\# Z  > 2$
2	YQY	A	1382	-	24,28,28	1.26	3 (12%)	30,42,42	1.94	11 (36%)
3	ACP	A	1383	4	25,33,33	1.45	4 (16%)	31,52,52	2.19	6 (19%)

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	YQY	A	1382	-	-	0/12/25/25	0/3/3/3
3	ACP	A	1383	4	-	0/15/38/38	0/3/3/3

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1382	YQY	F26-C9	-2.30	1.31	1.35
3	A	1383	ACP	PG-O3G	-2.14	1.49	1.54
3	A	1383	ACP	C5-C4	2.05	1.45	1.40
3	A	1383	ACP	C2-N3	2.28	1.36	1.32
2	A	1382	YQY	C3-C4	2.43	1.43	1.38

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1383	ACP	N3-C2-N1	-7.34	123.27	128.89
3	A	1383	ACP	C2'-C1'-N9	-5.28	106.23	114.29
2	A	1382	YQY	C3-C2-C1	-3.59	117.05	121.22
2	A	1382	YQY	F22-C10-C9	-3.44	113.65	118.33
3	A	1383	ACP	C4-C5-N7	-3.21	106.53	109.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1382	YQY	3	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	282/301 (93%)	0.18	15 (5%) 30 29	40, 58, 80, 96	8 (2%)

The worst 5 of 15 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	224	VAL	4.8
1	A	379	ILE	4.5
1	A	381	LEU	4.4
1	A	380	GLY	4.2
1	A	221	ASN	3.6

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	YQY	A	1382	26/26	0.99	0.15	0.00	40,45,51,55	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ACP	A	1383	31/31	0.97	0.13	-0.30	45,52,62,63	0
4	MG	A	1384	1/1	0.69	0.15	-	58,58,58,58	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.