



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:49 AM GMT

PDB ID : 2IVS  
Title : CRYSTAL STRUCTURE OF NON-PHOSPHORYLATED RET TYROSINE KINASE DOMAIN  
Authors : Knowles, P.P.; Murray-Rust, J.; Mcdonald, N.Q.  
Deposited on : 2006-06-16  
Resolution : 2.00 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](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) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

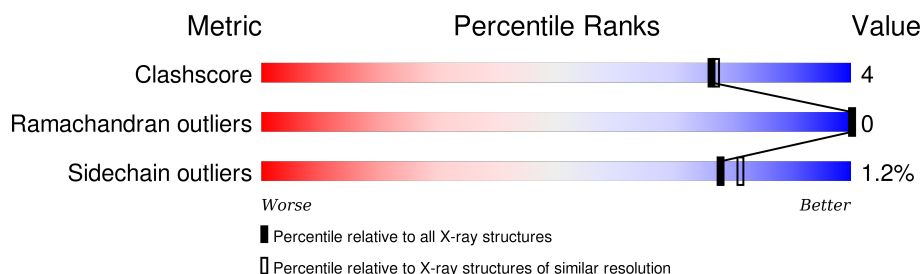
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	314	
1	B	314	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	FMT	A	3015	-	-	X	-

## 2 Entry composition [i](#)

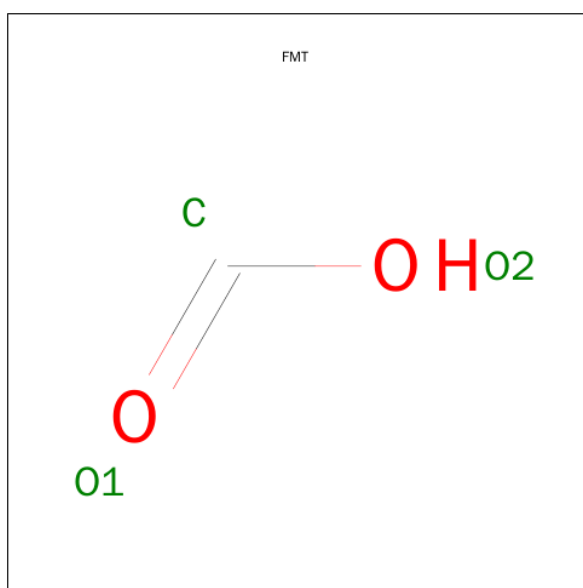
There are 4 unique types of molecules in this entry. The entry contains 4647 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROTO-ONCOGENE TYROSINE-PROTEIN KINASE RECEPTOR RET.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	284	Total	C	N	O	S	1	2	0
			2204	1422	376	392	14			
1	B	277	Total	C	N	O	S	0	0	0
			2119	1373	359	373	14			

- Molecule 2 is FORMIC ACID (three-letter code: FMT) (formula: CH<sub>2</sub>O<sub>2</sub>).



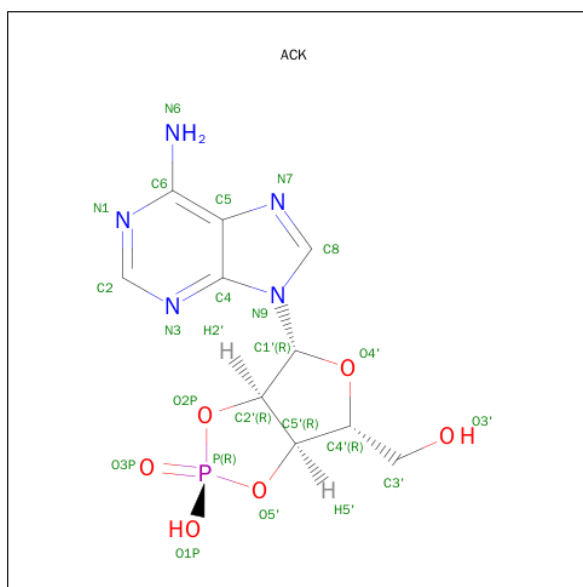
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			3	1	2		
2	A	1	Total	C	O	0	0
			3	1	2		
2	A	1	Total	C	O	0	0
			3	1	2		
2	B	1	Total	C	O	0	0
			3	1	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			3	1	2		
2	B	1	Total	C	O	0	0
			3	1	2		

- Molecule 3 is (3AR,4R,6R,6AR)-4-(6-AMINO-9H-PURIN-9-YL)-6-(HYDROXYMETHYL) TETRAHYDROFURO[3,4-D][1,3,2]DIOXAPHOSPHOL-2-OL 2-OXIDE (three-letter code: ACK) (formula: C<sub>10</sub>H<sub>12</sub>N<sub>5</sub>O<sub>6</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
3	B	1	Total	C	N	O	P	0	0
			22	10	5	6	1		

- Molecule 4 is water.

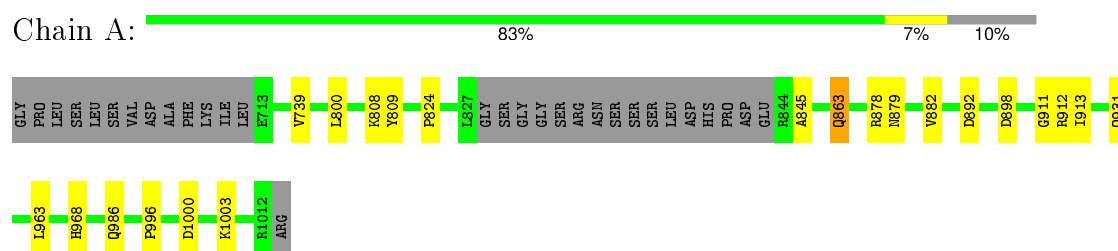
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	147	Total	O	0	0
			147	147		
4	B	115	Total	O	0	0
			115	115		

### 3 Residue-property plots [i](#)

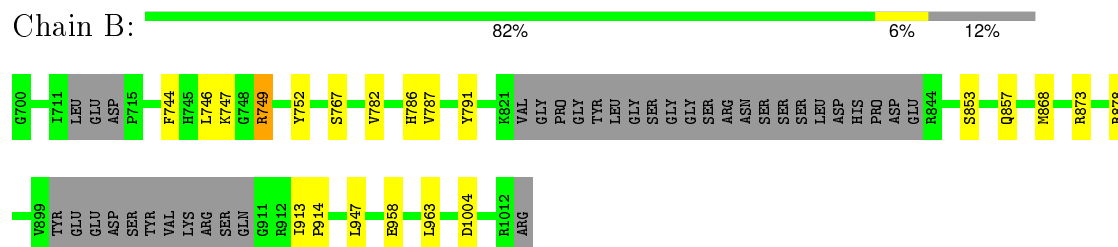
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

#### • Molecule 1: PROTO-ONCOGENE TYROSINE-PROTEIN KINASE RECEPTOR RET



#### • Molecule 1: PROTO-ONCOGENE TYROSINE-PROTEIN KINASE RECEPTOR RET



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	50.40 Å   80.22 Å   79.68 Å 90.00°   100.09°   90.00°	Depositor
Resolution (Å)	50.00 – 2.00	Depositor
% Data completeness (in resolution range)	97.7 (50.00-2.00)	Depositor
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.180 , 0.229	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4647	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

## 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: ACK, FMT

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	1.58	2/2261 (0.1%)	1.06	6/3062 (0.2%)
1	B	0.69	0/2167	0.74	3/2933 (0.1%)
All	All	1.23	2/4428 (0.0%)	0.92	9/5995 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	863[A]	GLN	CD-NE2	47.44	2.51	1.32
1	A	863[B]	GLN	CD-NE2	47.44	2.51	1.32

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	863[A]	GLN	CG-CD-NE2	-26.59	52.89	116.70
1	A	863[B]	GLN	CG-CD-NE2	-26.59	52.89	116.70
1	A	863[A]	GLN	OE1-CD-NE2	14.38	154.97	121.90
1	A	863[B]	GLN	OE1-CD-NE2	14.38	154.97	121.90
1	A	912	ARG	NE-CZ-NH1	-9.94	115.33	120.30
1	A	912	ARG	NE-CZ-NH2	6.84	123.72	120.30
1	B	749	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	B	947	LEU	CA-CB-CG	5.30	127.49	115.30
1	B	749	ARG	NE-CZ-NH2	-5.25	117.67	120.30

There are no chirality outliers.

There are no planarity outliers.

## 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	2204	0	2146	22	0
1	B	2119	0	2066	12	0
2	A	9	0	3	7	0
2	B	9	0	3	0	0
3	A	22	0	11	0	0
3	B	22	0	11	0	0
4	A	147	0	0	5	0
4	B	115	0	0	1	0
All	All	4647	0	4240	33	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 4.

All (33) 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:863[A]:GLN:NE2	1:A:863[A]:GLN:HG2	1.15	1.44
1:A:863[A]:GLN:CG	1:A:863[A]:GLN:NE2	2.00	1.24
1:A:931[B]:GLN:NE2	4:A:2100:HOH:O	1.97	0.98
1:A:1000:ASP:OD2	4:A:2137:HOH:O	1.93	0.86
1:B:749:ARG:NH2	4:B:2020:HOH:O	2.19	0.74
1:A:879:ASN:ND2	2:A:3015:FMT:C	2.56	0.67
1:A:863[A]:GLN:NE2	1:A:863[A]:GLN:CD	2.51	0.64
1:A:892:ASP:OD2	2:A:3015:FMT:H	2.00	0.61
1:A:800:LEU:HG	1:B:958:GLU:HG3	1.83	0.60
1:A:879:ASN:ND2	2:A:3015:FMT:H	2.17	0.59
1:B:782:VAL:HG11	1:B:787:VAL:HG11	1.84	0.58
1:A:878:ARG:NH2	4:A:2073:HOH:O	2.41	0.53
1:A:879:ASN:HD21	2:A:3015:FMT:C	2.21	0.53
1:A:863[A]:GLN:NE2	4:A:2067:HOH:O	2.41	0.52
1:B:744:PHE:HB2	1:B:752:TYR:CE1	2.46	0.51
1:A:911:GLY:HA3	4:A:2099:HOH:O	2.10	0.51
1:A:879:ASN:HD22	2:A:3015:FMT:H	1.75	0.50
1:A:913:ILE:N	1:A:913:ILE:HD12	2.28	0.49
1:A:892:ASP:OD2	2:A:3015:FMT:C	2.60	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:747:LYS:HE2	1:B:791:TYR:O	2.13	0.48
1:B:873:ARG:HB3	1:B:913:ILE:HD11	1.96	0.48
1:A:892:ASP:CG	2:A:3015:FMT:H	2.33	0.47
1:B:853:SER:O	1:B:857:GLN:HG3	2.13	0.47
1:A:808:LYS:HE2	1:A:809:TYR:CZ	2.50	0.46
1:A:809:TYR:HB2	1:A:882:VAL:HB	1.98	0.45
1:A:986:GLN:HB3	1:A:996:PRO:HD3	1.97	0.44
1:B:868:MET:HB3	1:B:868:MET:HE2	1.96	0.43
1:A:963:LEU:HG	1:A:968:HIS:HB2	2.01	0.42
1:B:746:LEU:HD11	1:B:791:TYR:HB3	2.01	0.42
1:B:786:HIS:CE1	1:B:857:GLN:OE1	2.72	0.42
1:B:963:LEU:HD23	1:B:963:LEU:HA	1.88	0.41
1:B:878:ARG:NH2	1:B:914:PRO:HG2	2.35	0.41
1:A:824:PRO:HG2	1:A:845:ALA:HB2	2.02	0.41

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	282/314 (90%)	277 (98%)	5 (2%)	0	100	100
1	B	269/314 (86%)	263 (98%)	6 (2%)	0	100	100
All	All	551/628 (88%)	540 (98%)	11 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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	222/272 (82%)	219 (99%)	3 (1%)	74	77
1	B	212/272 (78%)	210 (99%)	2 (1%)	84	88
All	All	434/544 (80%)	429 (99%)	5 (1%)	78	81

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	739	VAL
1	A	898	ASP
1	A	1003	LYS
1	B	767	SER
1	B	1004	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	860	GLN
1	A	879	ASN
1	B	796	GLN
1	B	860	GLN
1	B	879	ASN
1	B	926	HIS

### 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

8 ligands are 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	$\# Z  > 2$
2	FMT	A	3013	-	0,2,2	0.00	-	0,1,1	0.00	-
2	FMT	A	3014	-	0,2,2	0.00	-	0,1,1	0.00	-
2	FMT	A	3015	-	0,2,2	0.00	-	0,1,1	0.00	-
3	ACK	A	3016	-	18,25,25	1.14	1 (5%)	18,39,39	2.24	5 (27%)
2	FMT	B	3013	-	0,2,2	0.00	-	0,1,1	0.00	-
2	FMT	B	3014	-	0,2,2	0.00	-	0,1,1	0.00	-
2	FMT	B	3015	-	0,2,2	0.00	-	0,1,1	0.00	-
3	ACK	B	3016	-	18,25,25	0.99	1 (5%)	18,39,39	2.39	6 (33%)

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	FMT	A	3013	-	-	0/0/0/0	0/0/0/0
2	FMT	A	3014	-	-	0/0/0/0	0/0/0/0
2	FMT	A	3015	-	-	0/0/0/0	0/0/0/0
3	ACK	A	3016	-	-	0/2/32/32	0/4/4/4
2	FMT	B	3013	-	-	0/0/0/0	0/0/0/0
2	FMT	B	3014	-	-	0/0/0/0	0/0/0/0
2	FMT	B	3015	-	-	0/0/0/0	0/0/0/0
3	ACK	B	3016	-	-	0/2/32/32	0/4/4/4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	3016	ACK	C5-C4	2.61	1.46	1.40
3	A	3016	ACK	C5-C4	2.95	1.47	1.40

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	3016	ACK	N3-C2-N1	-8.12	122.68	128.89
3	A	3016	ACK	N3-C2-N1	-7.21	123.37	128.89
3	A	3016	ACK	C1'-N9-C4	-2.90	122.57	126.94
3	A	3016	ACK	O1P-P-O2P	-2.70	103.63	108.46
3	B	3016	ACK	C4-C5-N7	-2.17	107.48	109.48
3	B	3016	ACK	C2-N1-C6	2.15	122.61	118.77
3	B	3016	ACK	C4'-O4'-C1'	2.24	112.18	109.72
3	B	3016	ACK	O4'-C1'-N9	2.29	112.90	108.10
3	B	3016	ACK	O1P-P-O5'	2.56	113.04	108.46
3	A	3016	ACK	N6-C6-N1	2.60	124.77	119.20
3	A	3016	ACK	C2-N1-C6	2.64	123.48	118.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	3015	FMT	7	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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.