



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 25, 2020 – 03:00 PM BST

PDB ID : 3ZMM
Title : Inhibitors of Jak2 Kinase domain
Authors : Read, J.; Green, I.; Pollard, H.; Howard, T.; Mott, R.
Deposited on : 2013-02-11
Resolution : 2.51 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 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 : 2.13
buster-report : 1.1.7 (2018)
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) : 2.13

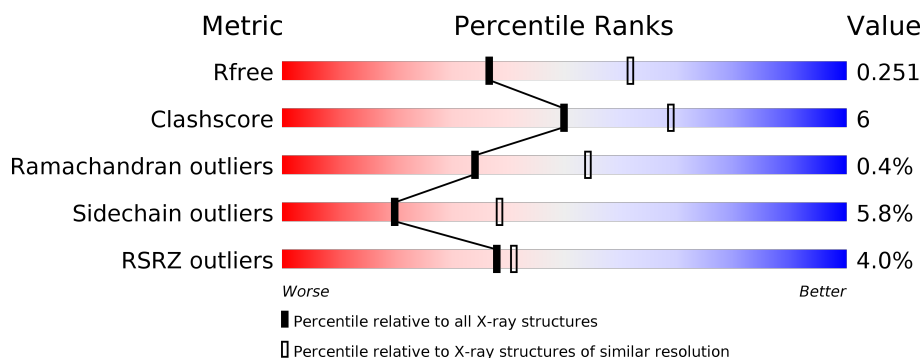
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 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	298	<div> <div>4%</div> <div> <div></div> <div>78%</div> <div>13%</div> <div>8%</div> </div> </div>
1	B	298	<div> <div>4%</div> <div> <div></div> <div>77%</div> <div>15%</div> <div>7%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4728 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.

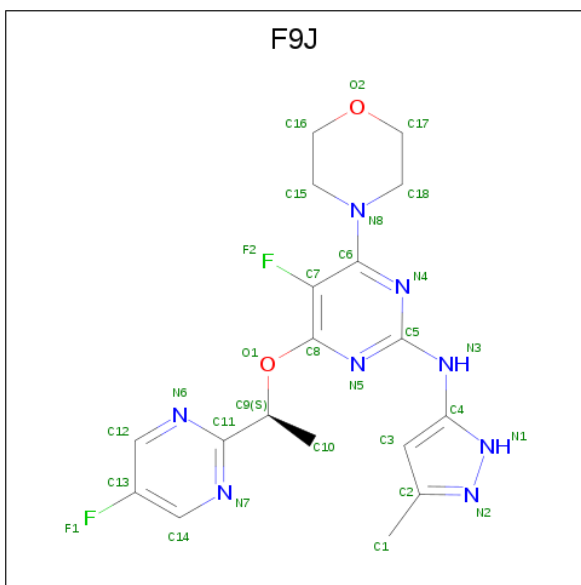
- Molecule 1 is a protein called TYROSINE-PROTEIN KINASE JAK2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	P	S	0	2	0
			2255	1439	382	418	2	14			
1	B	276	Total	C	N	O	P	S	7	1	0
			2264	1440	390	418	2	14			

There are 6 discrepancies between the modelled and reference sequences:

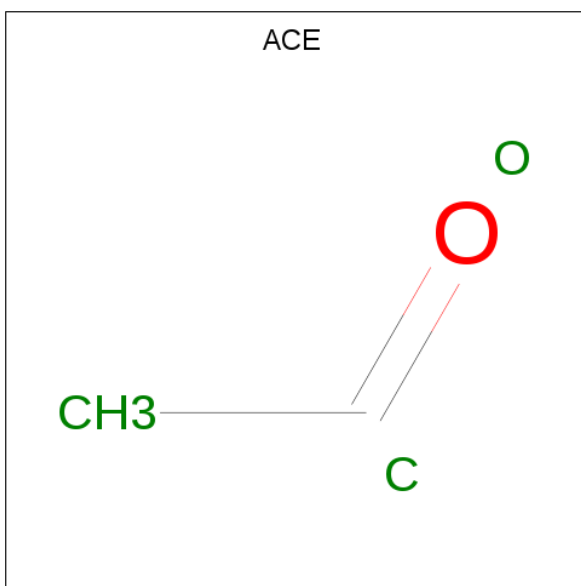
Chain	Residue	Modelled	Actual	Comment	Reference
A	943	ALA	LYS	engineered mutation	UNP O60674
A	945	ALA	LYS	engineered mutation	UNP O60674
A	1129	GLN	ASN	conflict	UNP O60674
B	943	ALA	LYS	engineered mutation	UNP O60674
B	945	ALA	LYS	engineered mutation	UNP O60674
B	1129	GLN	ASN	conflict	UNP O60674

- Molecule 2 is 5-FLUORO-4-[(1S)-1-(5-FLUOROPYRIMIDIN-2-YL)ETHOXY]-N-(5-METHYL-1H-PYRAZOL-3-YL)-6-MORPHOLINO-PYRIMIDIN-2-AMINE (three-letter code: F9J) (formula: C₁₈H₂₀F₂N₈O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0
			30	18	2	8	2	
2	B	1	Total	C	F	N	O	0
			30	18	2	8	2	

- Molecule 3 is ACETYL GROUP (three-letter code: ACE) (formula: C₂H₄O).



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

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

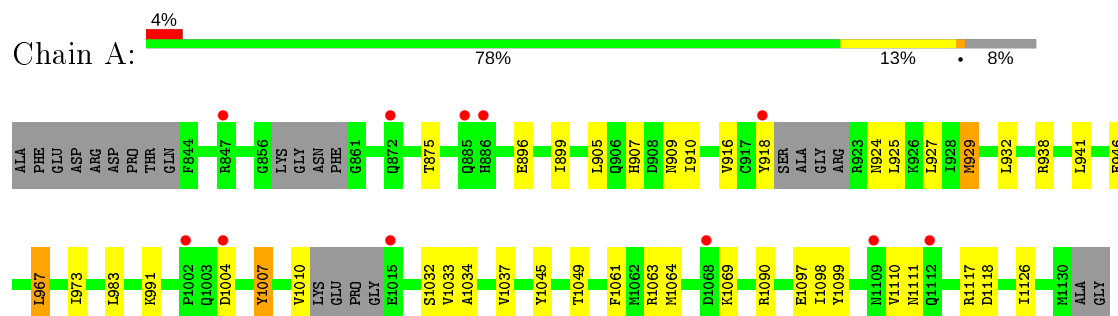
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	53	Total	O	0	0
			53	53		
4	B	87	Total	O	0	0
			87	87		

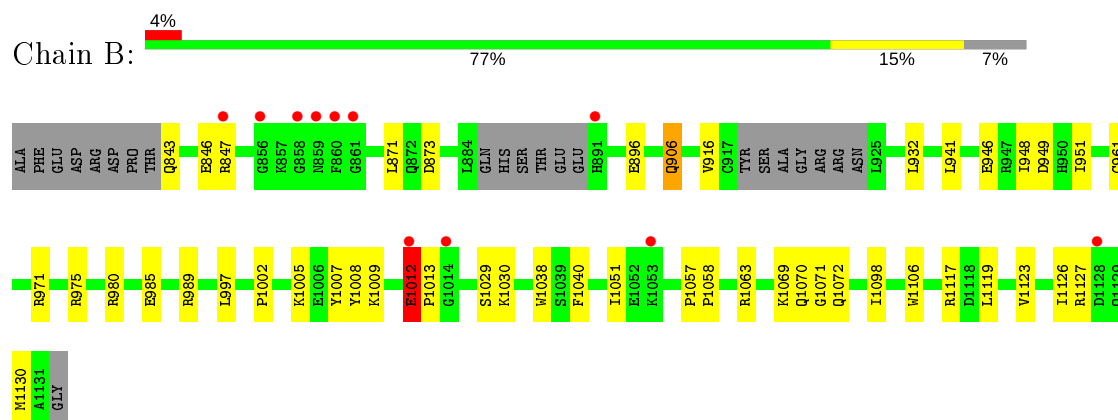
3 Residue-property plots [i](#)

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: TYROSINE-PROTEIN KINASE JAK2



• Molecule 1: TYROSINE-PROTEIN KINASE JAK2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	44.51Å 126.69Å 135.59Å 90.00° 97.22° 90.00°	Depositor
Resolution (Å)	17.57 – 2.51 17.57 – 2.52	Depositor EDS
% Data completeness (in resolution range)	95.5 (17.57-2.51) 96.1 (17.57-2.52)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 2.52Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.193 , 0.246 0.199 , 0.251	Depositor DCC
R_{free} test set	1239 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	41.8	Xtriage
Anisotropy	0.328	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 41.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.022 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4728	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.84% 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: F9J, PTR, ACE

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	0.45	0/2273	0.65	0/3064
1	B	0.48	0/2279	0.71	1/3069 (0.0%)
All	All	0.47	0/4552	0.68	1/6133 (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	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	949	ASP	CB-CG-OD1	5.07	122.86	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	1012	GLU	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	2255	0	2192	23	0
1	B	2264	0	2226	23	0
2	A	30	0	20	3	0
2	B	30	0	20	6	0
3	A	6	0	6	0	0
3	B	3	0	3	0	0
4	A	53	0	0	1	0
4	B	87	0	0	2	0
All	All	4728	0	4467	52	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 6.

All (52) 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:932:LEU:H	2:A:2133:F9J:H1	1.19	0.91
1:B:1098:ILE:HD11	1:B:1126:ILE:HG21	1.55	0.87
1:B:932:LEU:H	2:B:2134:F9J:H1	1.34	0.75
1:A:1010:VAL:O	1:A:1010:VAL:HG12	1.86	0.74
1:A:1010:VAL:O	1:A:1010:VAL:CG1	2.39	0.70
1:A:1098:ILE:HD11	1:A:1126:ILE:HG21	1.79	0.63
2:A:2133:F9J:F2	2:A:2133:F9J:H181	1.88	0.63
1:B:1051:ILE:O	1:B:1051:ILE:HG22	2.05	0.57
2:B:2134:F9J:H3	2:B:2134:F9J:N5	2.21	0.56
2:B:2134:F9J:C18	2:B:2134:F9J:F2	2.44	0.55
1:B:1038:TRP:CE3	1:B:1106:TRP:HA	2.42	0.54
1:B:871:LEU:HB2	1:B:873:ASP:OD1	2.08	0.53
1:A:973:ILE:HD12	1:A:1032:SER:C	2.31	0.51
1:A:909:ASN:HA	1:A:991:LYS:HG2	1.93	0.50
1:B:843:GLN:N	4:B:2001:HOH:O	2.43	0.50
1:A:1033:VAL:O	1:A:1037:VAL:HG23	2.12	0.49
1:A:907:HIS:HB3	1:A:910:ILE:HD13	1.95	0.49
1:A:929:MET:HE1	4:A:2014:HOH:O	2.13	0.49
1:B:1007:PTR:CD1	1:B:1007:PTR:C	2.91	0.48
1:A:932:LEU:HD12	1:A:983:LEU:CB	2.44	0.47
2:B:2134:F9J:N5	2:B:2134:F9J:C3	2.77	0.47
1:A:1034:ALA:CB	1:A:1110:VAL:HG13	2.44	0.47
1:B:975:ARG:HD3	1:B:997:LEU:O	2.14	0.47
1:A:899:ILE:HG12	1:A:927:LEU:HD13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1063[B]:ARG:CZ	1:A:1063[B]:ARG:HB3	2.47	0.45
1:B:961:CYS:SG	1:B:1040:PHE:CZ	3.10	0.45
1:A:1004:ASP:OD1	1:A:1004:ASP:N	2.43	0.44
1:A:1061:PHE:HA	1:A:1064:MET:HE3	1.99	0.44
2:B:2134:F9J:H182	2:B:2134:F9J:F2	2.07	0.44
1:A:1045:TYR:O	1:A:1049:THR:HG23	2.17	0.44
2:B:2134:F9J:N7	2:B:2134:F9J:H3	2.33	0.44
1:A:932:LEU:HD12	1:A:983:LEU:HB2	1.99	0.43
1:B:1051:ILE:N	1:B:1051:ILE:HD13	2.33	0.43
1:A:905:LEU:HD11	1:A:967:LEU:HD21	2.00	0.43
1:B:951:ILE:HG23	1:B:1127:ARG:HD2	1.99	0.43
1:B:971:ARG:NH2	1:B:1005:LYS:O	2.52	0.43
1:A:875:THR:HB	1:B:906:GLN:NE2	2.34	0.43
1:B:941:LEU:HD23	1:B:948:ILE:HD13	2.00	0.43
1:B:932:LEU:HD22	1:B:985:GLU:HG3	2.00	0.42
1:A:1117:ARG:O	1:A:1118:ASP:C	2.57	0.42
1:B:1012:GLU:N	1:B:1013:PRO:CD	2.82	0.42
1:B:1098:ILE:N	1:B:1098:ILE:HD13	2.34	0.42
1:A:1090:ARG:HG3	1:A:1099:TYR:HB2	2.02	0.41
1:B:1002:PRO:HG2	1:B:1005:LYS:HB2	2.01	0.41
1:B:1130:MET:HB3	1:B:1130:MET:HE2	1.93	0.41
1:A:1007:PTR:C	1:A:1007:PTR:CD1	2.97	0.41
2:A:2133:F9J:H3	2:A:2133:F9J:N5	2.36	0.41
1:B:1030:LYS:HE2	4:B:2051:HOH:O	2.20	0.41
1:B:1098:ILE:CD1	1:B:1126:ILE:HD13	2.51	0.41
1:B:1057:PRO:HB2	1:B:1058:PRO:HD3	2.03	0.41
1:B:1119:LEU:O	1:B:1123:VAL:HG23	2.21	0.41
1:A:905:LEU:CD1	1:A:967:LEU:HD21	2.52	0.40

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	267/298 (90%)	255 (96%)	12 (4%)	0	100	100
1	B	269/298 (90%)	256 (95%)	11 (4%)	2 (1%)	22	39
All	All	536/596 (90%)	511 (95%)	23 (4%)	2 (0%)	34	54

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1071	GLY
1	B	1012	GLU

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	241/267 (90%)	228 (95%)	13 (5%)	22	42
1	B	243/267 (91%)	227 (93%)	16 (7%)	16	32
All	All	484/534 (91%)	455 (94%)	29 (6%)	20	37

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	896	GLU
1	A	916	VAL
1	A	918	TYR
1	A	924	ASN
1	A	925	LEU
1	A	929	MET
1	A	938	ARG
1	A	941	LEU
1	A	946	GLU
1	A	967	LEU
1	A	1069	LYS
1	A	1097	GLU
1	A	1111	ASN
1	B	846	GLU

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Mol	Chain	Res	Type
1	B	847	ARG
1	B	896	GLU
1	B	906	GLN
1	B	916	VAL
1	B	946	GLU
1	B	980	ARG
1	B	989	ARG
1	B	1009	LYS
1	B	1029	SER
1	B	1063[A]	ARG
1	B	1063[B]	ARG
1	B	1069	LYS
1	B	1070	GLN
1	B	1072	GLN
1	B	1117	ARG

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

Mol	Chain	Res	Type
1	A	924	ASN
1	A	942	GLN
1	B	859	ASN
1	B	891	HIS
1	B	906	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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
1	PTR	A	1007	1	15,16,17	0.72	0	19,22,24	1.04	1 (5%)
1	PTR	A	1008	1	15,16,17	0.67	0	19,22,24	0.87	0
1	PTR	B	1008	1	15,16,17	0.68	0	19,22,24	1.30	2 (10%)
1	PTR	B	1007	1	15,16,17	0.75	0	19,22,24	1.12	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
1	PTR	A	1007	1	-	0/10/11/13	0/1/1/1
1	PTR	A	1008	1	-	0/10/11/13	0/1/1/1
1	PTR	B	1008	1	-	1/10/11/13	0/1/1/1
1	PTR	B	1007	1	-	0/10/11/13	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1008	PTR	P-OH-CZ	3.00	133.36	123.75
1	B	1008	PTR	CG-CB-CA	-2.93	108.16	114.10
1	A	1007	PTR	O3P-P-O2P	2.05	115.47	107.64

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	1008	PTR	CZ-OH-P-O3P

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1007	PTR	1	0
1	B	1007	PTR	1	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 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	F9J	B	2134	-	31,33,33	1.10	4 (12%)	34,46,46	1.37	6 (17%)
3	ACE	A	2135	-	1,2,2	1.41	0	1,1,1	0.47	0
3	ACE	A	2134	-	1,2,2	1.34	0	1,1,1	0.37	0
2	F9J	A	2133	-	31,33,33	0.99	1 (3%)	34,46,46	1.20	3 (8%)
3	ACE	B	2135	-	1,2,2	1.39	0	1,1,1	0.02	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	F9J	B	2134	-	-	3/14/24/24	0/4/4/4
2	F9J	A	2133	-	-	2/14/24/24	0/4/4/4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2134	F9J	C6-N8	2.55	1.44	1.37
2	B	2134	F9J	C2-N2	2.50	1.37	1.34
2	A	2133	F9J	C6-N8	2.49	1.43	1.37
2	B	2134	F9J	C5-N3	2.15	1.40	1.36
2	B	2134	F9J	C4-N3	-2.13	1.34	1.38

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2133	F9J	C18-N8-C15	3.14	118.44	111.52
2	B	2134	F9J	O1-C8-N5	2.58	123.01	118.65
2	B	2134	F9J	C4-C3-C2	-2.55	104.60	107.23
2	B	2134	F9J	N4-C6-N8	2.54	120.93	116.80
2	A	2133	F9J	C3-C2-N2	-2.53	106.22	110.81
2	B	2134	F9J	N7-C11-N6	2.35	129.05	125.95
2	B	2134	F9J	C3-C2-N2	-2.32	106.61	110.81
2	B	2134	F9J	C8-O1-C9	-2.13	113.73	116.79
2	A	2133	F9J	O1-C8-N5	2.09	122.18	118.65

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	2134	F9J	C7-C6-N8-C15
2	A	2133	F9J	N6-C11-C9-O1
2	B	2134	F9J	N4-C6-N8-C15
2	B	2134	F9J	N4-C6-N8-C18
2	A	2133	F9J	N4-C6-N8-C18

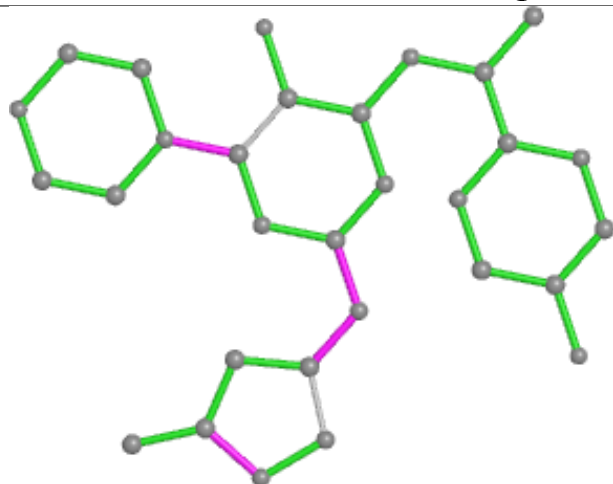
There are no ring outliers.

2 monomers are involved in 9 short contacts:

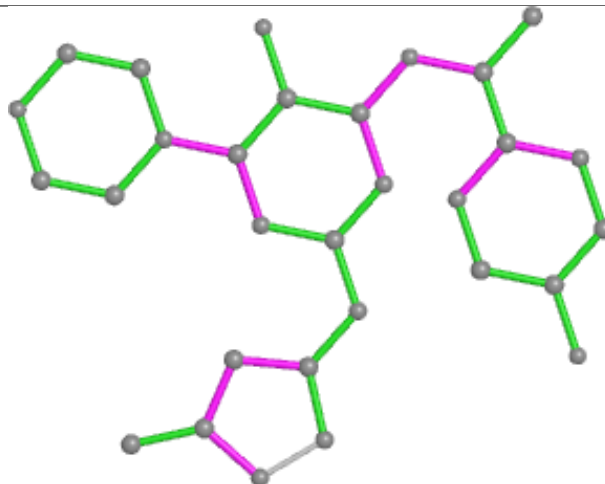
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2134	F9J	6	0
2	A	2133	F9J	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

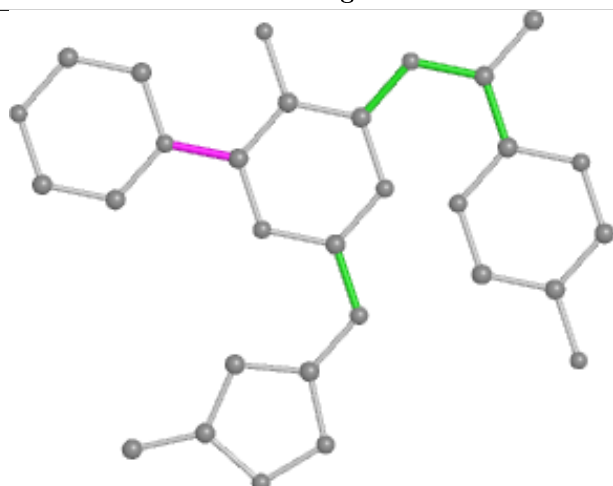
Ligand F9J B 2134



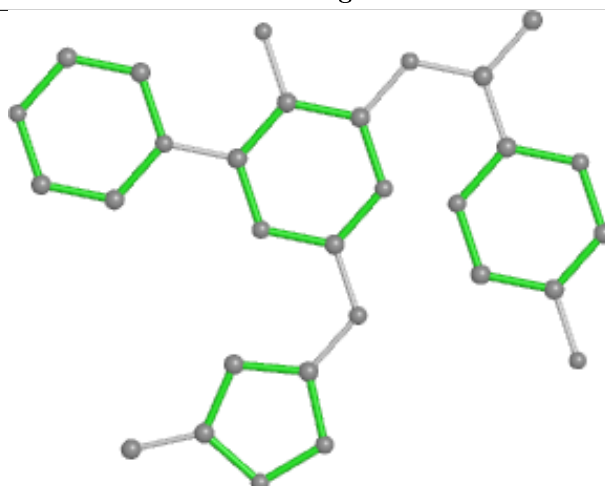
Bond lengths



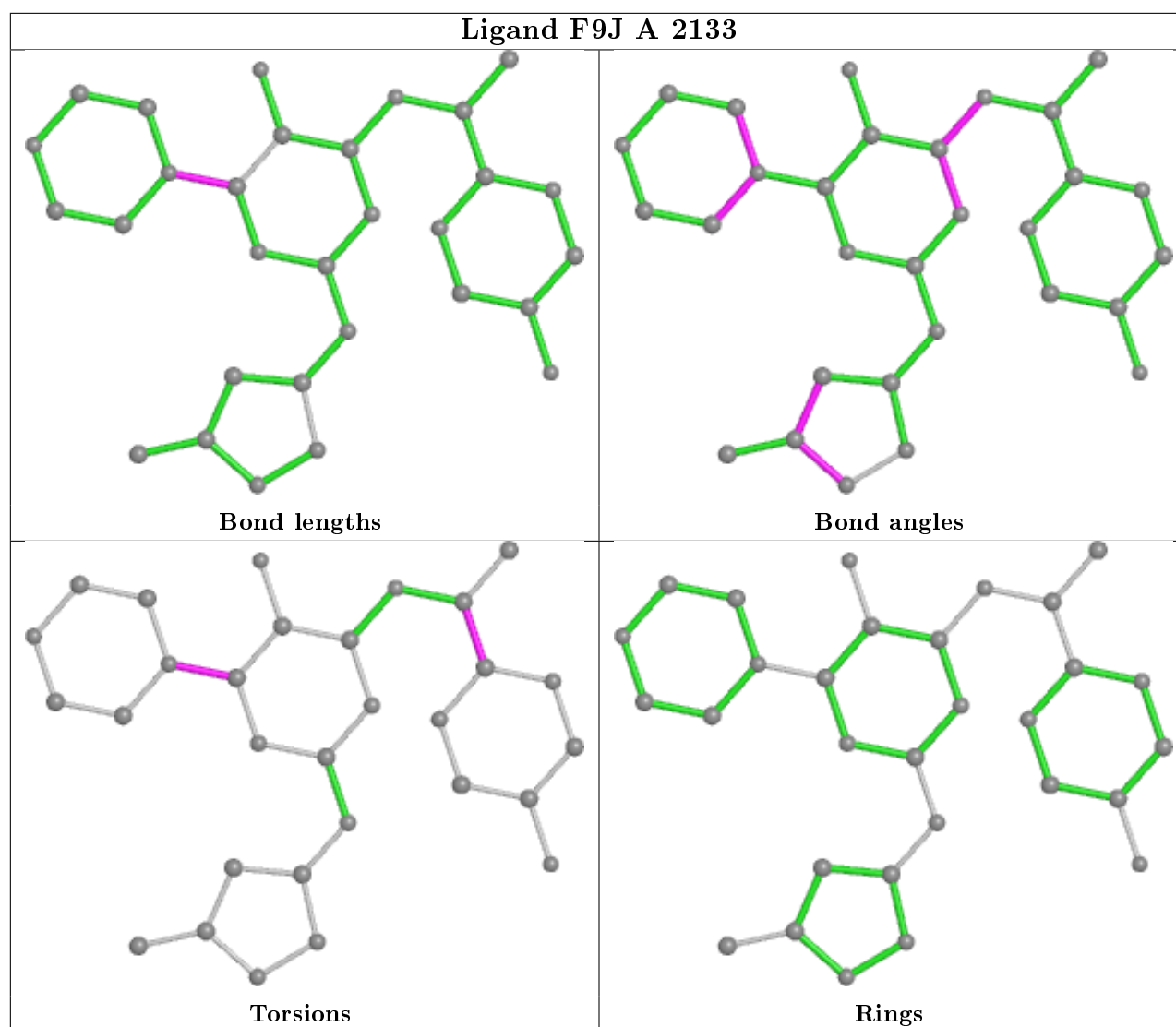
Bond angles



Torsions



Rings



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	273/298 (91%)	0.02	11 (4%) 38 41	41, 63, 96, 111	0
1	B	274/298 (91%)	-0.07	11 (4%) 38 41	35, 56, 95, 122	1 (0%)
All	All	547/596 (91%)	-0.02	22 (4%) 38 41	35, 59, 95, 122	1 (0%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	861	GLY	4.7
1	B	859	ASN	4.4
1	B	860	PHE	3.9
1	A	872	GLN	3.4
1	B	1012	GLU	3.2
1	B	1014	GLY	2.9
1	A	1068	ASP	2.8
1	B	858	GLY	2.7
1	A	1004	ASP	2.7
1	A	886	HIS	2.7
1	B	1128	ASP	2.6
1	B	847	ARG	2.6
1	A	1015	GLU	2.5
1	A	885	GLN	2.4
1	A	918	TYR	2.4
1	A	847	ARG	2.3
1	A	1112	GLN	2.3
1	B	891	HIS	2.1
1	A	1002	PRO	2.1
1	B	1053	LYS	2.1
1	A	1109	ASN	2.0
1	B	856	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	PTR	A	1007	16/17	0.85	0.23	80,88,102,106	0
1	PTR	A	1008	16/17	0.85	0.20	75,87,108,113	0
1	PTR	B	1008	16/17	0.94	0.12	51,60,74,76	0
1	PTR	B	1007	16/17	0.96	0.13	55,58,69,71	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

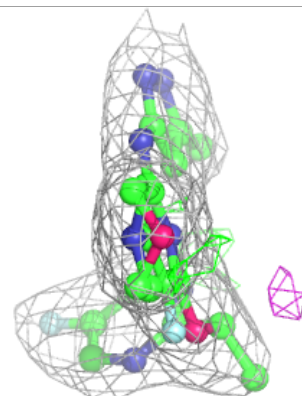
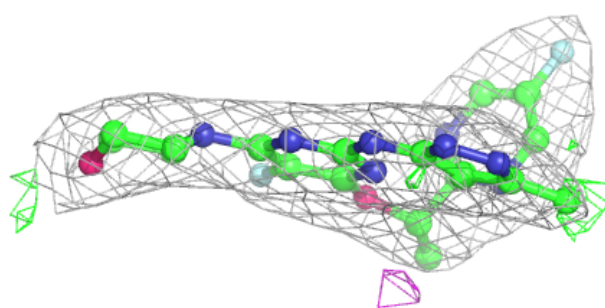
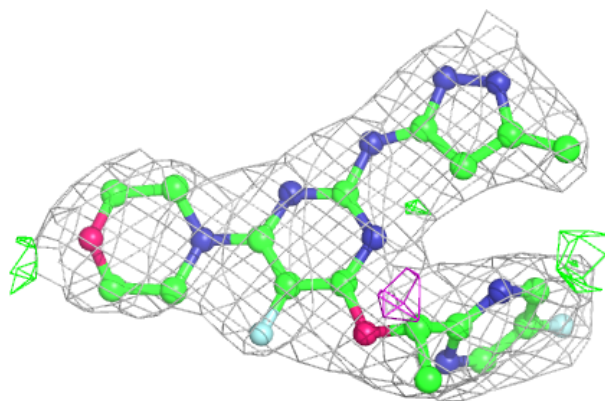
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	ACE	A	2134	3/3	0.78	0.21	67,67,74,75	0
3	ACE	A	2135	3/3	0.81	0.16	74,74,76,78	0
2	F9J	B	2134	30/30	0.93	0.13	45,54,65,67	0
3	ACE	B	2135	3/3	0.93	0.19	48,48,54,58	0
2	F9J	A	2133	30/30	0.96	0.12	39,45,50,56	0

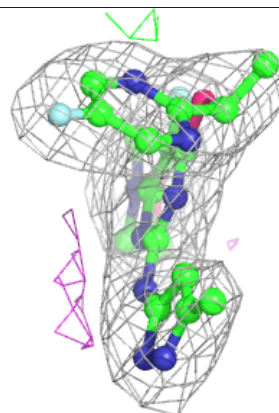
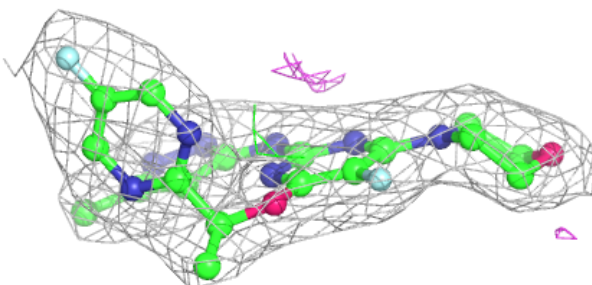
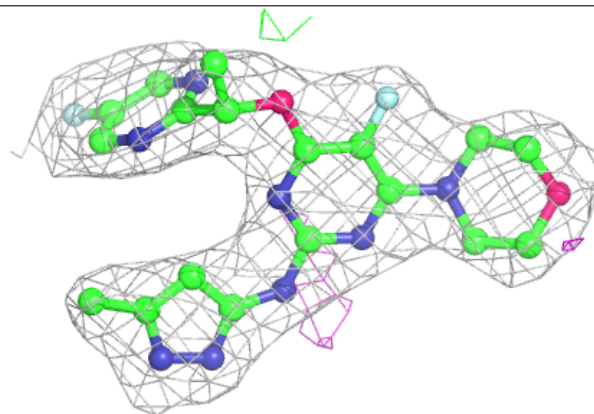
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around F9J B 2134:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around F9J A 2133:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers

There are no such residues in this entry.