



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 17, 2024 – 10:24 AM EST

PDB ID : 4JI9  
Title : JAK2 kinase (JH1 domain) in complex with TG101209  
Authors : Eigenbrot, C.; Ultsch, M.  
Deposited on : 2013-03-05  
Resolution : 2.40 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

<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>.

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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 : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 1.20.1  
EDS : 3.0  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.003 (Gargrove)  
Density-Fitness : 1.0.11  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.39

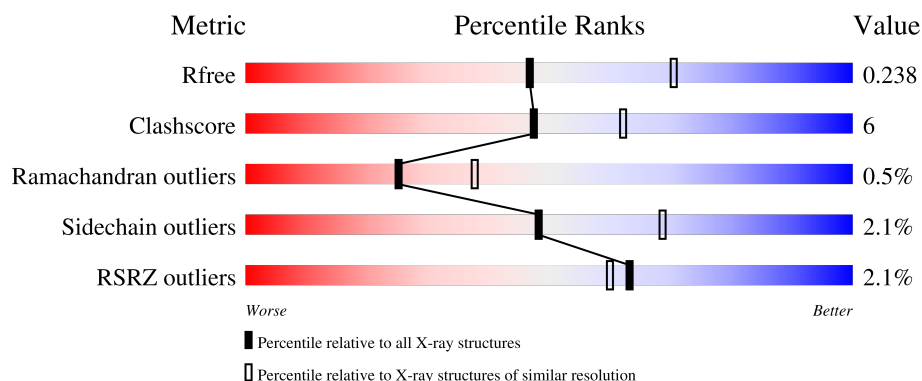
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	4642 (2.40-2.40)
Clashscore	180529	5218 (2.40-2.40)
Ramachandran outliers	177936	5158 (2.40-2.40)
Sidechain outliers	177891	5159 (2.40-2.40)
RSRZ outliers	164620	4642 (2.40-2.40)

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  $\geq 3$ , 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	300	
1	B	300	

## 2 Entry composition [i](#)

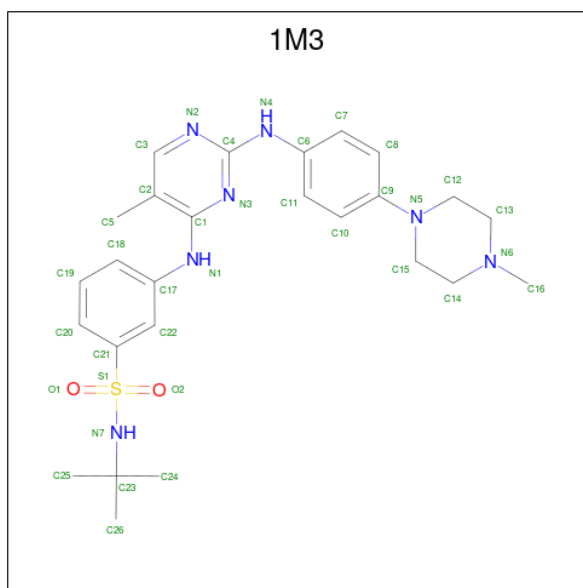
There are 3 unique types of molecules in this entry. The entry contains 4961 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	286	Total	C	N	O	P	S	0	0	0
			2373	1506	408	443	2	14			
1	B	290	Total	C	N	O	P	S	0	0	0
			2406	1525	415	450	2	14			

- Molecule 2 is N-tert-butyl-3-[(5-methyl-2-{[4-(4-methylpiperazin-1-yl)phenyl]amino}pyrimidin-4-yl)amino]benzenesulfonamide (three-letter code: 1M3) (formula: C<sub>26</sub>H<sub>35</sub>N<sub>7</sub>O<sub>2</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			36	26	7	2	1		
2	B	1	Total	C	N	O	S	0	0
			36	26	7	2	1		

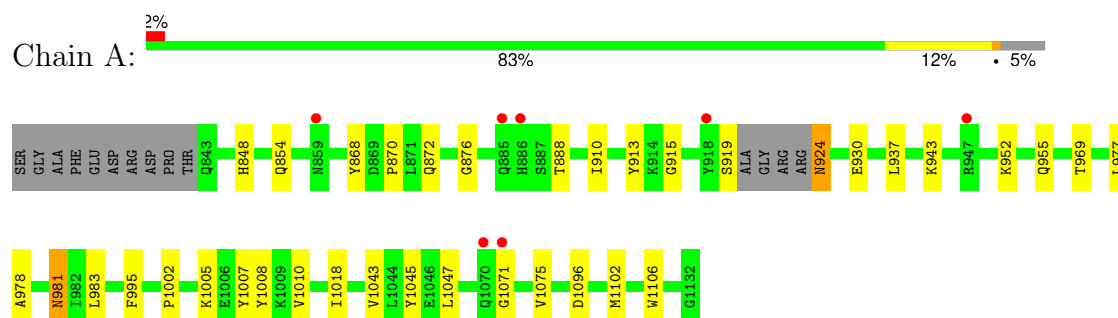
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	53	Total 53	O 53	0	0
3	B	57	Total 57	O 57	0	0

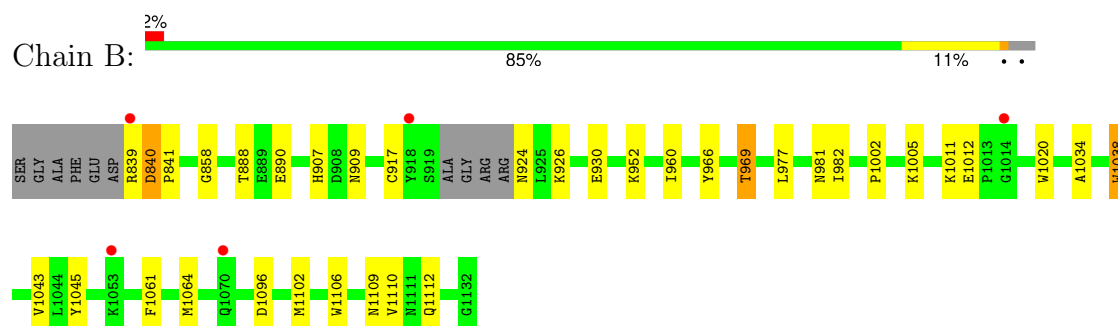
### 3 Residue-property plots [i](#)

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	P 41	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	111.99Å 111.99Å 70.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 50.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.9 (50.00-2.40) 98.9 (50.00-2.40)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.92 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, $R_{free}$	0.202 , 0.245 0.196 , 0.238	Depositor DCC
$R_{free}$ test set	852 reflections (2.49%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.9	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 32.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.030 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4961	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.13% 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, 0.333 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: PTR, 1M3

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	0/2389	0.55	0/3211
1	B	0.46	2/2423 (0.1%)	0.56	0/3258
All	All	0.46	2/4812 (0.0%)	0.55	0/6469

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1038	TRP	CD2-CE2	5.22	1.47	1.41
1	B	1020	TRP	CD2-CE2	5.08	1.47	1.41

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	2373	0	2343	24	0
1	B	2406	0	2374	23	0
2	A	36	0	35	5	0
2	B	36	0	35	5	0
3	A	53	0	0	1	0
3	B	57	0	0	1	0
All	All	4961	0	4787	53	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 (53) 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:1045:TYR:HB2	1:B:1102:MET:HE3	1.71	0.72
1:A:1002:PRO:HG2	1:A:1005:LYS:HB2	1.78	0.65
1:B:1002:PRO:HG2	1:B:1005:LYS:HB2	1.80	0.64
1:A:1102:MET:HE3	1:A:1106:TRP:CH2	2.33	0.63
1:A:1008:PTR:HE1	1:A:1010:VAL:HB	1.81	0.62
2:A:1201:1M3:H21	2:A:1201:1M3:N3	2.13	0.62
1:B:952:LYS:HG3	3:B:1341:HOH:O	2.00	0.60
1:B:839:ARG:HG3	1:B:840:ASP:N	2.17	0.60
1:B:1034:ALA:CB	1:B:1110:VAL:HG13	2.33	0.58
1:A:1102:MET:HE3	1:A:1106:TRP:HH2	1.68	0.58
2:A:1201:1M3:H21	2:A:1201:1M3:H13	1.85	0.58
2:B:1201:1M3:N3	2:B:1201:1M3:H21	2.18	0.57
1:B:930:GLU:O	2:B:1201:1M3:H16	2.06	0.56
1:A:1018:ILE:HD13	1:A:1075:VAL:HG22	1.88	0.56
1:A:952:LYS:CE	1:A:955:GLN:NE2	2.69	0.55
1:A:952:LYS:CE	1:A:955:GLN:HE22	2.20	0.54
1:A:952:LYS:HE2	1:A:955:GLN:HE22	1.73	0.53
1:B:888:THR:HG22	1:B:890:GLU:H	1.73	0.53
1:A:952:LYS:HE2	1:A:955:GLN:NE2	2.24	0.52
1:B:1109:ASN:HD22	1:B:1112:GLN:HE21	1.55	0.52
1:A:977:LEU:HG	1:A:1043:VAL:HG21	1.92	0.51
1:A:848:HIS:HD2	1:A:872:GLN:HE21	1.59	0.50
1:B:977:LEU:HG	1:B:1043:VAL:HG21	1.94	0.48
2:A:1201:1M3:N3	2:A:1201:1M3:H13	2.27	0.48
1:B:888:THR:HG22	1:B:890:GLU:N	2.28	0.48
1:A:930:GLU:O	2:A:1201:1M3:H16	2.13	0.48
1:A:1102:MET:CE	1:A:1106:TRP:CH2	2.96	0.47
1:A:943:LYS:HE3	3:A:1344:HOH:O	2.14	0.47
1:B:1061:PHE:HA	1:B:1064:MET:HE3	1.97	0.47
2:B:1201:1M3:N3	2:B:1201:1M3:H13	2.29	0.46
1:B:858:GLY:HA3	2:B:1201:1M3:H2	1.97	0.46
1:B:966:TYR:O	1:B:969:THR:HB	2.16	0.45
1:B:977:LEU:HD11	1:B:982:ILE:HD11	1.99	0.45
1:B:1034:ALA:HB3	1:B:1110:VAL:HG13	2.00	0.44
1:A:937:LEU:HD21	1:A:1047:LEU:HD21	2.00	0.44
1:A:1045:TYR:HB2	1:A:1102:MET:HE3	2.00	0.44
1:A:919:SER:HA	1:A:924:ASN:ND2	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:868:TYR:O	1:A:876:GLY:HA3	2.18	0.43
1:A:848:HIS:CG	1:A:870:PRO:HA	2.54	0.43
1:B:1102:MET:HE3	1:B:1106:TRP:CH2	2.54	0.42
1:B:1045:TYR:HD1	1:B:1102:MET:HE1	1.85	0.42
1:B:960:ILE:HG12	1:B:982:ILE:HD13	2.01	0.42
1:A:978:ALA:H	1:A:981:ASN:HD21	1.68	0.42
1:A:983:LEU:HD11	2:A:1201:1M3:C1	2.50	0.42
1:B:1038:TRP:CE3	1:B:1106:TRP:HA	2.54	0.42
1:B:907:HIS:HE1	1:B:909:ASN:HD22	1.69	0.41
1:A:952:LYS:HE3	1:A:955:GLN:NE2	2.36	0.41
1:B:1102:MET:HE3	1:B:1106:TRP:HH2	1.85	0.41
1:A:913:TYR:CZ	1:A:915:GLY:HA2	2.55	0.40
1:A:910:ILE:HG23	1:A:995:PHE:CZ	2.57	0.40
1:B:840:ASP:HA	1:B:841:PRO:HD3	1.93	0.40
1:B:917:CYS:HB3	1:B:926:LYS:HB2	2.04	0.40
2:B:1201:1M3:N3	2:B:1201:1M3:C18	2.84	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	280/300 (93%)	268 (96%)	11 (4%)	1 (0%)	30	44
1	B	284/300 (95%)	275 (97%)	7 (2%)	2 (1%)	19	29
All	All	564/600 (94%)	543 (96%)	18 (3%)	3 (0%)	25	38

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1071	GLY
1	B	840	ASP

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Mol	Chain	Res	Type
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	260/270 (96%)	254 (98%)	6 (2%)	45	66
1	B	264/270 (98%)	259 (98%)	5 (2%)	52	72
All	All	524/540 (97%)	513 (98%)	11 (2%)	48	69

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	854	GLN
1	A	888	THR
1	A	924	ASN
1	A	969	THR
1	A	981	ASN
1	A	1096	ASP
1	B	924	ASN
1	B	969	THR
1	B	981	ASN
1	B	1011	LYS
1	B	1096	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	848	HIS
1	A	909	ASN
1	A	924	ASN
1	A	955	GLN
1	A	974	HIS
1	A	981	ASN

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Mol	Chain	Res	Type
1	B	909	ASN
1	B	955	GLN
1	B	974	HIS
1	B	981	ASN
1	B	1112	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	1008	1	15,16,17	0.77	0	17,22,24	1.06	0
1	PTR	B	1007	1	15,16,17	0.76	0	17,22,24	0.86	0
1	PTR	A	1007	1	15,16,17	0.79	0	17,22,24	0.88	1 (5%)
1	PTR	B	1008	1	15,16,17	0.76	0	17,22,24	0.98	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	1008	1	-	0/10/11/13	0/1/1/1
1	PTR	B	1007	1	-	0/10/11/13	0/1/1/1
1	PTR	A	1007	1	-	0/10/11/13	0/1/1/1
1	PTR	B	1008	1	-	1/10/11/13	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1007	PTR	OH-P-O1P	-2.30	101.81	109.48

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	1008	PTR	C-CA-CB-CG

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1008	PTR	1	0

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

2 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	1M3	A	1201	-	39,39,39	1.75	2 (5%)	56,57,57	1.58	12 (21%)
2	1M3	B	1201	-	39,39,39	1.78	3 (7%)	56,57,57	1.38	9 (16%)

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	1M3	A	1201	-	-	0/24/34/34	0/4/4/4
2	1M3	B	1201	-	-	0/24/34/34	0/4/4/4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1201	1M3	S1-N7	9.86	1.76	1.61
2	A	1201	1M3	S1-N7	9.72	1.76	1.61
2	A	1201	1M3	C23-N7	-2.41	1.45	1.49
2	B	1201	1M3	C23-N7	-2.15	1.45	1.49
2	B	1201	1M3	C6-N4	-2.13	1.36	1.40

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1201	1M3	N2-C4-N3	-4.92	121.67	126.42
2	A	1201	1M3	O1-S1-N7	3.92	116.32	107.36
2	B	1201	1M3	N2-C4-N3	-3.64	122.90	126.42
2	A	1201	1M3	C2-C3-N2	-3.47	121.61	125.09
2	B	1201	1M3	C2-C3-N2	-3.42	121.65	125.09
2	B	1201	1M3	C2-C1-N3	-3.11	119.83	123.46
2	B	1201	1M3	C15-C14-N6	-2.94	106.13	110.86
2	A	1201	1M3	C12-N5-C9	2.91	126.05	118.11
2	A	1201	1M3	C10-C9-N5	-2.75	117.56	121.39
2	A	1201	1M3	O1-S1-O2	-2.73	116.20	119.52
2	B	1201	1M3	C12-N5-C9	2.73	125.55	118.11
2	A	1201	1M3	C3-N2-C4	2.58	119.35	115.81
2	A	1201	1M3	C2-C1-N3	-2.49	120.55	123.46
2	A	1201	1M3	C22-C21-S1	-2.49	116.52	119.06
2	B	1201	1M3	O2-S1-C21	-2.31	105.06	107.98
2	A	1201	1M3	C21-S1-N7	-2.27	104.06	108.11
2	A	1201	1M3	C20-C21-S1	2.19	122.17	119.76
2	B	1201	1M3	O1-S1-N7	2.16	112.30	107.36
2	A	1201	1M3	C14-C15-N5	2.08	115.16	110.78
2	B	1201	1M3	C23-N7-S1	-2.07	124.67	128.22
2	B	1201	1M3	C21-S1-N7	-2.04	104.48	108.11

There are no chirality outliers.

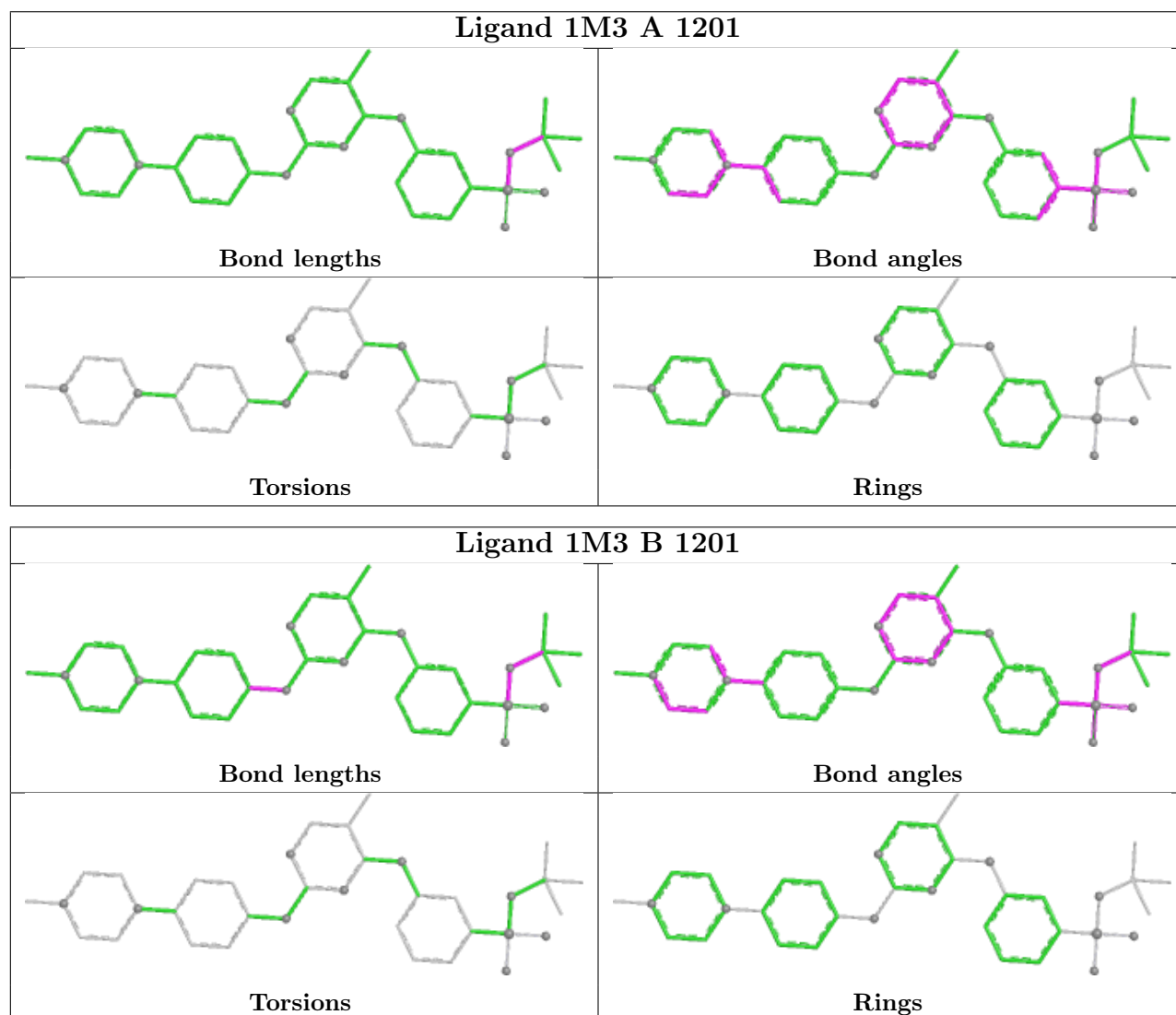
There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1201	1M3	5	0
2	B	1201	1M3	5	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.



## 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	284/300 (94%)	-0.52	7 (2%) 58 55	20, 34, 89, 114	0
1	B	288/300 (96%)	-0.54	5 (1%) 69 65	22, 37, 69, 112	0
All	All	572/600 (95%)	-0.53	12 (2%) 63 60	20, 36, 80, 114	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	918	TYR	3.9
1	B	1070	GLN	3.7
1	A	918	TYR	3.3
1	A	859	ASN	3.2
1	B	839	ARG	2.6
1	A	886	HIS	2.6
1	B	1014	GLY	2.3
1	A	947	ARG	2.3
1	A	1070	GLN	2.2
1	A	1071	GLY	2.2
1	B	1053	LYS	2.2
1	A	885	GLN	2.1

### 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, 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	B-factors(Å <sup>2</sup> )	Q<0.9
1	PTR	B	1008	16/17	0.90	0.09	38,49,81,84	0
1	PTR	A	1007	16/17	0.91	0.09	39,45,97,118	0

*Continued on next page...*

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
1	PTR	B	1007	16/17	0.92	0.08	38,50,79,87	0
1	PTR	A	1008	16/17	0.93	0.08	34,48,69,71	0

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [i](#)

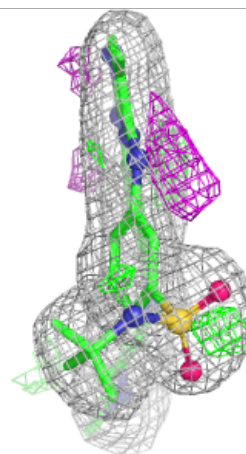
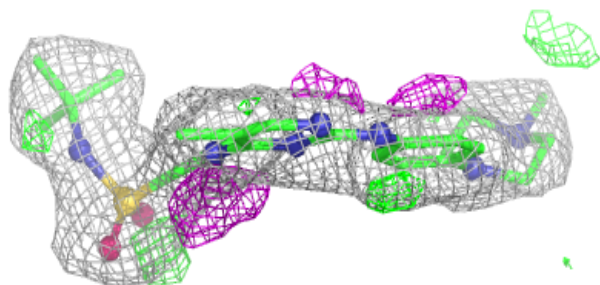
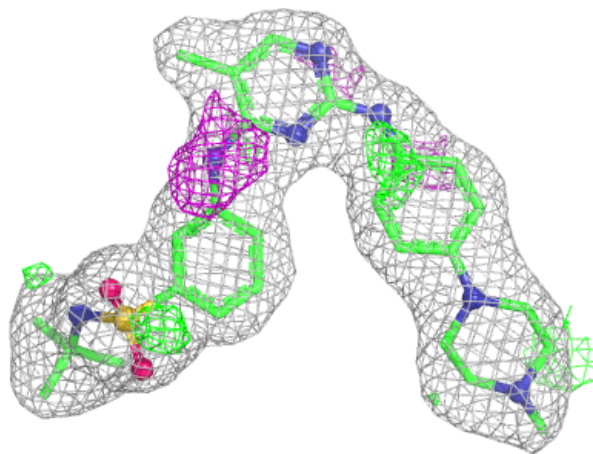
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

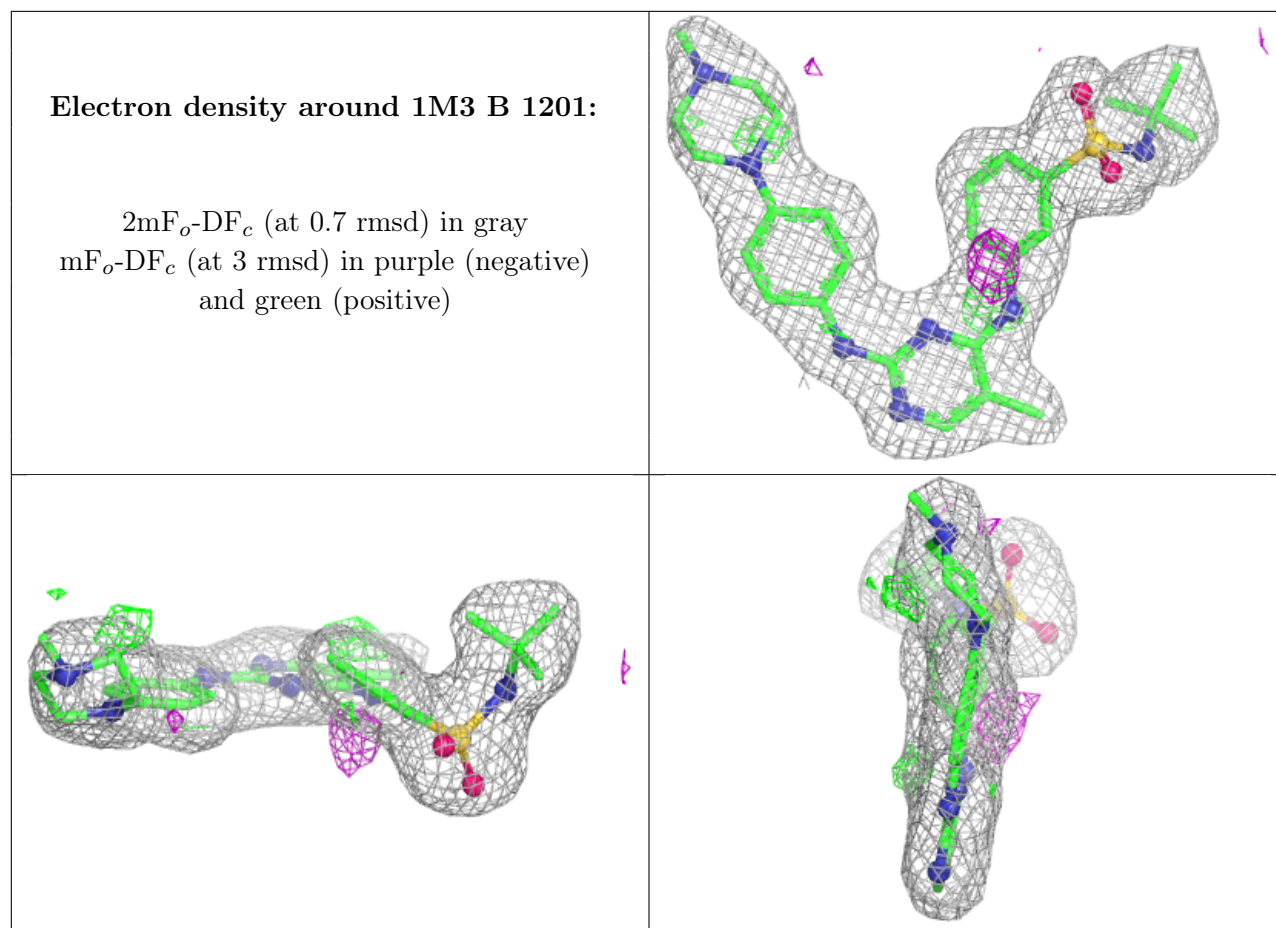
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	1M3	A	1201	36/36	0.92	0.10	20,28,53,64	0
2	1M3	B	1201	36/36	0.93	0.09	25,31,57,61	0

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 1M3 A 1201:**

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





## 6.5 Other polymers [i](#)

There are no such residues in this entry.