



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:11 AM UTC

PDB ID : 4C62 / pdb_00004c62
Title : Inhibitors of Jak2 Kinase domain
Authors : Read, J.; Green, I.; Pollard, H.; Howard, T.
Deposited on : 2013-09-17
Resolution : 2.75 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

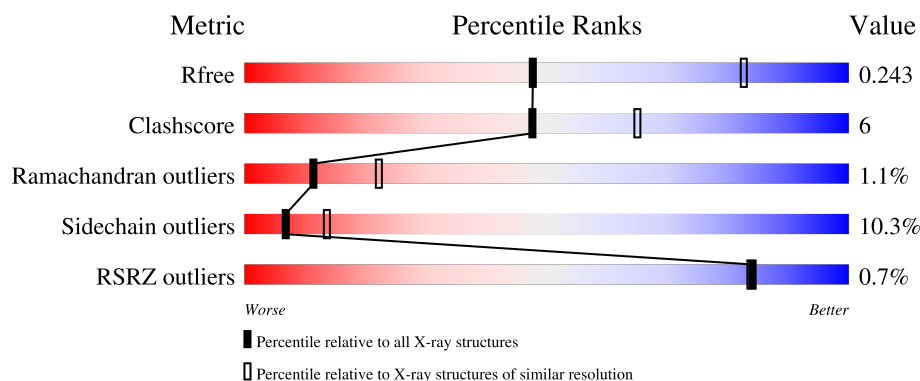
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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}	180053	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	298	 74% 13% . . 8%
1	B	298	 72% 15% . . 8%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4604 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.

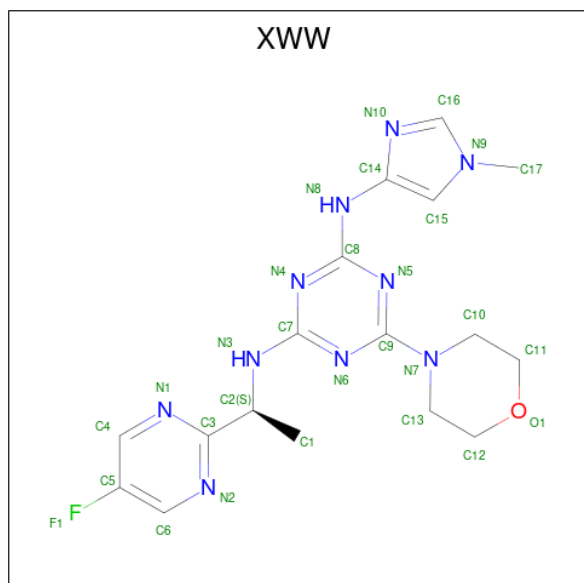
- Molecule 1 is a protein called TYROSINE-PROTEIN KINASE JAK2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	P	S	0	0	0
			2221	1411	376	418	2	14			
1	B	273	Total	C	N	O	P	S	0	0	0
			2199	1403	374	406	2	14			

There are 6 discrepancies between the modelled and reference sequences:

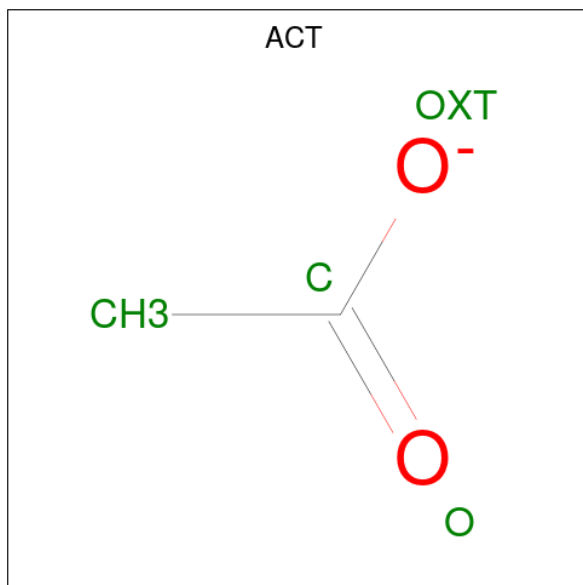
Chain	Residue	Modelled	Actual	Comment	Reference
A	943	ALA	LYS	engineered mutation	UNP O60674
A	945	ALA	LYS	engineered mutation	UNP O60674
A	1129	GLN	ASN	conflict	UNP O60674
B	943	ALA	LYS	engineered mutation	UNP O60674
B	945	ALA	LYS	engineered mutation	UNP O60674
B	1129	GLN	ASN	conflict	UNP O60674

- Molecule 2 is N2-[(1S)-1-(5-fluoropyrimidin-2-yl)ethyl]-n4-(1-methylimidazol-4-yl)-6-morpholino-1,3,5-triazine-2,4-diamine (CCD ID: XWW) (formula: C₁₇H₂₁FN₁₀O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			29	17	1	10	1		
2	B	1	Total	C	F	N	O	0	0
			29	17	1	10	1		

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



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		
3	B	1	Total	C	O	0	0
			3	2	1		

- Molecule 4 is water.

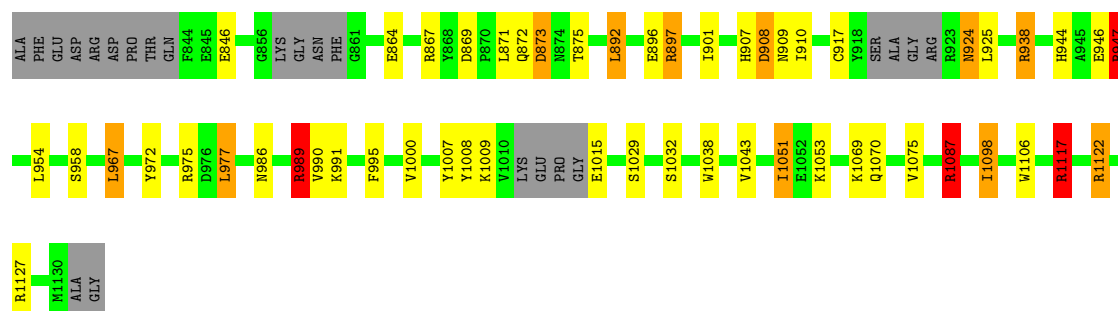
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	44	Total	O	0	0
			44	44		
4	B	73	Total	O	0	0
			73	73		

3 Residue-property plots

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.

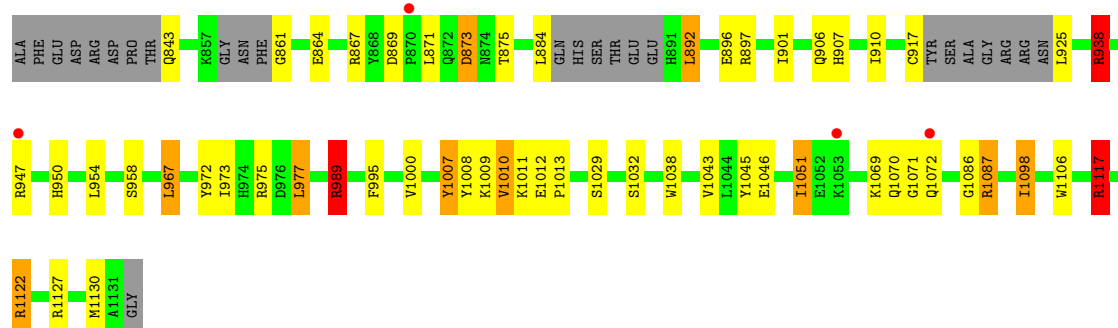
• Molecule 1: TYROSINE-PROTEIN KINASE JAK2

Chain A: 



• Molecule 1: TYROSINE-PROTEIN KINASE JAK2

Chain B: 



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 126.89Å 135.88Å 90.00° 97.53° 90.00°	Depositor
Resolution (Å)	21.00 – 2.75 21.00 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.4 (21.00-2.75) 99.4 (21.00-2.75)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.75Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.196 , 0.246 0.200 , 0.243	Depositor DCC
R_{free} test set	991 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	46.5	Xtriage
Anisotropy	0.031	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 47.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.026 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4604	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACT, XWW, PTR

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	0.65	0/2231	1.09	24/3012 (0.8%)
1	B	0.70	0/2210	1.11	27/2983 (0.9%)
All	All	0.68	0/4441	1.10	51/5995 (0.9%)

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	A	0	6
1	B	0	4
All	All	0	10

There are no bond length outliers.

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	897	ARG	NE-CZ-NH2	-10.57	109.68	119.20
1	A	897	ARG	CD-NE-CZ	10.19	138.66	124.40
1	B	975	ARG	NE-CZ-NH2	-9.78	110.39	119.20
1	B	1117	ARG	NE-CZ-NH2	-9.16	110.95	119.20
1	A	947	ARG	NE-CZ-NH2	-9.13	110.98	119.20
1	B	989	ARG	NE-CZ-NH2	-9.07	111.03	119.20
1	A	947	ARG	CD-NE-CZ	9.04	137.05	124.40
1	A	1087	ARG	NE-CZ-NH2	-8.57	111.48	119.20
1	A	1087	ARG	CD-NE-CZ	8.57	136.40	124.40
1	B	938	ARG	NE-CZ-NH2	-8.33	111.71	119.20
1	B	1010	VAL	N-CA-C	8.25	119.12	111.45
1	A	1122	ARG	NE-CZ-NH2	-8.22	111.80	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1117	ARG	NE-CZ-NH2	-8.14	111.87	119.20
1	B	989	ARG	CD-NE-CZ	7.99	135.59	124.40
1	A	1117	ARG	CD-NE-CZ	7.86	135.40	124.40
1	B	897	ARG	NE-CZ-NH2	-7.72	112.25	119.20
1	B	975	ARG	CD-NE-CZ	7.69	135.17	124.40
1	B	938	ARG	CD-NE-CZ	7.68	135.15	124.40
1	A	975	ARG	NE-CZ-NH2	-7.66	112.30	119.20
1	A	1122	ARG	CD-NE-CZ	7.60	135.03	124.40
1	A	989	ARG	NE-CZ-NH2	-7.31	112.62	119.20
1	B	1122	ARG	CD-NE-CZ	7.02	134.23	124.40
1	B	1122	ARG	NE-CZ-NH2	-6.95	112.94	119.20
1	B	1117	ARG	CD-NE-CZ	6.91	134.07	124.40
1	A	938	ARG	NE-CZ-NH2	-6.88	113.01	119.20
1	B	947	ARG	NE-CZ-NH2	-6.87	113.02	119.20
1	A	1087	ARG	NE-CZ-NH1	6.77	128.27	121.50
1	B	1087	ARG	NE-CZ-NH2	-6.40	113.44	119.20
1	A	938	ARG	CD-NE-CZ	6.38	133.34	124.40
1	A	989	ARG	CD-NE-CZ	6.35	133.29	124.40
1	B	897	ARG	CD-NE-CZ	6.25	133.16	124.40
1	B	989	ARG	NE-CZ-NH1	6.24	127.74	121.50
1	B	1087	ARG	CD-NE-CZ	6.21	133.09	124.40
1	A	989	ARG	CG-CD-NE	-6.13	98.52	112.00
1	B	1117	ARG	NE-CZ-NH1	6.10	127.60	121.50
1	B	1122	ARG	NE-CZ-NH1	6.10	127.60	121.50
1	B	975	ARG	NE-CZ-NH1	6.05	127.55	121.50
1	B	938	ARG	NE-CZ-NH1	6.02	127.52	121.50
1	A	975	ARG	CD-NE-CZ	5.98	132.77	124.40
1	A	1117	ARG	NE-CZ-NH1	5.88	127.38	121.50
1	A	975	ARG	NE-CZ-NH1	5.86	127.36	121.50
1	B	989	ARG	CG-CD-NE	5.84	124.85	112.00
1	B	947	ARG	NE-CZ-NH1	5.74	127.24	121.50
1	B	1087	ARG	NE-CZ-NH1	5.71	127.21	121.50
1	A	1122	ARG	NE-CZ-NH1	5.63	127.13	121.50
1	B	897	ARG	NE-CZ-NH1	5.62	127.12	121.50
1	B	1117	ARG	CG-CD-NE	-5.40	100.13	112.00
1	A	938	ARG	NE-CZ-NH1	5.39	126.89	121.50
1	A	897	ARG	CG-CD-NE	5.33	123.73	112.00
1	B	897	ARG	CG-CD-NE	-5.33	100.28	112.00
1	A	908	ASP	N-CA-C	5.01	117.47	111.71

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1087	ARG	Sidechain
1	A	1117	ARG	Sidechain
1	A	1122	ARG	Sidechain
1	A	897	ARG	Sidechain
1	A	947	ARG	Sidechain
1	A	989	ARG	Sidechain
1	B	1117	ARG	Sidechain
1	B	1122	ARG	Sidechain
1	B	938	ARG	Sidechain
1	B	989	ARG	Sidechain

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	2221	0	2127	25	0
1	B	2199	0	2130	23	0
2	A	29	0	21	1	0
2	B	29	0	21	2	0
3	A	6	0	6	0	0
3	B	3	0	3	0	0
4	A	44	0	0	6	0
4	B	73	0	0	2	0
All	All	4604	0	4308	49	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 (49) 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:947:ARG:NH2	4:A:2016:HOH:O	2.11	0.84
2:B:2135:XWW:H15	2:B:2135:XWW:N4	2.02	0.74
1:A:967:LEU:HD23	1:A:972:TYR:HB2	1.69	0.74
1:B:967:LEU:HD23	1:B:972:TYR:HB2	1.70	0.74
1:A:892:LEU:HD12	1:A:892:LEU:O	2.03	0.59
1:A:871:LEU:O	1:A:873:ASP:HB3	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:2135:XWW:N4	2:B:2135:XWW:C15	2.61	0.57
1:B:892:LEU:HD12	1:B:892:LEU:O	2.06	0.54
1:A:1032:SER:HB2	4:A:2031:HOH:O	2.09	0.51
1:B:871:LEU:O	1:B:873:ASP:HB3	2.10	0.51
1:B:977:LEU:HD22	1:B:1043:VAL:HG21	1.92	0.51
1:B:954:LEU:HD21	1:B:1098:ILE:HD13	1.93	0.51
1:A:989:ARG:HH12	1:B:1117:ARG:CZ	2.25	0.49
1:A:947:ARG:CZ	4:A:2016:HOH:O	2.58	0.47
1:A:977:LEU:HB2	4:A:2025:HOH:O	2.14	0.47
1:A:954:LEU:HD21	1:A:1098:ILE:HD13	1.95	0.47
1:A:967:LEU:HD23	1:A:972:TYR:CB	2.43	0.47
1:B:967:LEU:HD23	1:B:972:TYR:CB	2.44	0.46
1:A:1051:ILE:HG22	1:A:1051:ILE:O	2.14	0.46
1:B:1007:PTR:CD1	1:B:1007:PTR:C	2.92	0.46
1:A:907:HIS:HB3	1:A:910:ILE:HD12	1.97	0.46
1:B:1010:VAL:O	1:B:1012:GLU:N	2.49	0.46
1:B:907:HIS:HB3	1:B:910:ILE:HD12	1.98	0.45
1:A:944:HIS:CG	4:A:2016:HOH:O	2.70	0.44
1:A:977:LEU:HD22	1:A:1043:VAL:HG21	1.98	0.44
1:A:1007:PTR:CD1	1:A:1007:PTR:C	2.96	0.44
1:A:924:ASN:HD22	1:A:924:ASN:HA	1.59	0.44
1:A:989:ARG:NH1	1:B:1117:ARG:NH2	2.67	0.43
1:A:967:LEU:HD11	1:A:995:PHE:HE1	1.82	0.43
1:B:973:ILE:HD12	1:B:1032:SER:C	2.44	0.43
1:B:1086:GLY:C	1:B:1087:ARG:HG2	2.44	0.43
1:B:867:ARG:HG2	1:B:869:ASP:HB2	2.01	0.42
1:B:1117:ARG:HD3	4:B:2070:HOH:O	2.18	0.42
1:A:954:LEU:HB2	1:A:1127:ARG:HD3	2.02	0.42
1:A:989:ARG:HG3	1:A:990:VAL:N	2.34	0.42
1:A:909:ASN:HA	1:A:991:LYS:HG2	2.02	0.41
1:A:986:ASN:OD1	1:A:986:ASN:C	2.63	0.41
1:B:967:LEU:HD11	1:B:995:PHE:HE1	1.86	0.41
1:A:867:ARG:HG2	1:A:869:ASP:HB2	2.02	0.41
1:A:1053:LYS:N	4:A:2036:HOH:O	2.54	0.41
1:B:861:GLY:HA2	1:B:884:LEU:HA	2.02	0.41
1:B:843:GLN:N	4:B:2001:HOH:O	2.52	0.41
1:B:1051:ILE:HG22	1:B:1051:ILE:O	2.19	0.41
1:B:954:LEU:HB2	1:B:1127:ARG:HD3	2.03	0.41
1:B:1038:TRP:CE3	1:B:1106:TRP:HA	2.56	0.41
1:B:950:HIS:HB2	1:B:1130:MET:HE1	2.03	0.40
1:B:1045:TYR:O	1:B:1046:GLU:C	2.63	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1038:TRP:CE3	1:A:1106:TRP:HA	2.57	0.40
2:A:2133:XWW:N4	2:A:2133:XWW:H15	2.36	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	265/298 (89%)	255 (96%)	8 (3%)	2 (1%)	16	30
1	B	263/298 (88%)	255 (97%)	4 (2%)	4 (2%)	8	15
All	All	528/596 (89%)	510 (97%)	12 (2%)	6 (1%)	11	22

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1011	LYS
1	B	1013	PRO
1	B	1071	GLY
1	A	872	GLN
1	A	1051	ILE
1	B	1051	ILE

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	234/267 (88%)	207 (88%)	27 (12%)	5	10
1	B	232/267 (87%)	211 (91%)	21 (9%)	9	17
All	All	466/534 (87%)	418 (90%)	48 (10%)	7	13

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	846	GLU
1	A	864	GLU
1	A	873	ASP
1	A	875	THR
1	A	892	LEU
1	A	896	GLU
1	A	901	ILE
1	A	908	ASP
1	A	917	CYS
1	A	924	ASN
1	A	925	LEU
1	A	938	ARG
1	A	946	GLU
1	A	947	ARG
1	A	958	SER
1	A	967	LEU
1	A	977	LEU
1	A	1000	VAL
1	A	1009	LYS
1	A	1015	GLU
1	A	1029	SER
1	A	1069	LYS
1	A	1070	GLN
1	A	1075	VAL
1	A	1087	ARG
1	A	1098	ILE
1	A	1117	ARG
1	B	864	GLU
1	B	873	ASP
1	B	875	THR
1	B	892	LEU
1	B	896	GLU
1	B	901	ILE
1	B	906	GLN
1	B	917	CYS

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Mol	Chain	Res	Type
1	B	925	LEU
1	B	938	ARG
1	B	958	SER
1	B	967	LEU
1	B	977	LEU
1	B	989	ARG
1	B	1000	VAL
1	B	1009	LYS
1	B	1029	SER
1	B	1069	LYS
1	B	1070	GLN
1	B	1072	GLN
1	B	1098	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	848	HIS
1	A	854	GLN
1	A	909	ASN
1	A	924	ASN
1	A	1084	ASN
1	B	848	HIS
1	B	854	GLN
1	B	944	HIS

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	0.72	0	17,22,24	1.17	2 (11%)
1	PTR	A	1007	1	15,16,17	0.73	0	17,22,24	1.10	0
1	PTR	B	1008	1	15,16,17	0.67	0	17,22,24	1.34	3 (17%)
1	PTR	B	1007	1	15,16,17	0.70	0	17,22,24	1.09	1 (5%)

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	A	1007	1	-	0/10/11/13	0/1/1/1
1	PTR	B	1008	1	-	0/10/11/13	0/1/1/1
1	PTR	B	1007	1	-	1/10/11/13	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1008	PTR	O3P-P-O2P	2.47	117.07	107.80
1	B	1008	PTR	P-OH-CZ	2.35	132.25	123.88
1	B	1008	PTR	O3P-P-OH	-2.25	98.66	105.32
1	A	1008	PTR	O3P-P-OH	-2.07	99.20	105.32
1	B	1007	PTR	O3P-P-OH	-2.05	99.27	105.32
1	A	1008	PTR	O3P-P-O2P	2.00	115.32	107.80

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	1007	PTR	CZ-OH-P-O2P

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 oligosaccharides 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	XWW	A	2133	-	30,32,32	1.34	2 (6%)	37,44,44	2.79	12 (32%)
3	ACT	A	2135	-	1,2,3	1.32	0	0,1,3	-	-
3	ACT	B	2134	-	1,2,3	1.42	0	0,1,3	-	-
3	ACT	A	2134	-	1,2,3	1.25	0	0,1,3	-	-
2	XWW	B	2135	-	30,32,32	1.42	2 (6%)	37,44,44	2.83	14 (37%)

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	XWW	A	2133	-	-	3/16/24/24	0/4/4/4
2	XWW	B	2135	-	-	3/16/24/24	0/4/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2135	XWW	C7-N3	5.85	1.41	1.34
2	A	2133	XWW	C7-N3	5.04	1.40	1.34
2	B	2135	XWW	C9-N7	4.02	1.44	1.35
2	A	2133	XWW	C9-N7	3.98	1.44	1.35

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2133	XWW	C15-N9-C16	6.76	109.38	106.71
2	B	2135	XWW	C15-N9-C16	6.61	109.32	106.71
2	A	2133	XWW	C13-N7-C9	-6.13	109.56	121.67
2	A	2133	XWW	N9-C16-N10	-6.12	110.32	112.66
2	A	2133	XWW	C10-N7-C9	-6.09	109.63	121.67
2	B	2135	XWW	C9-N5-C8	5.61	118.58	113.84
2	A	2133	XWW	C9-N6-C7	5.48	118.46	113.84
2	B	2135	XWW	C13-N7-C9	-5.43	110.94	121.67
2	B	2135	XWW	C1-C2-C3	-5.42	104.40	110.92
2	B	2135	XWW	C10-N7-C9	-5.34	111.11	121.67
2	B	2135	XWW	C9-N6-C7	4.99	118.05	113.84
2	A	2133	XWW	C9-N5-C8	4.97	118.04	113.84
2	B	2135	XWW	N9-C16-N10	-4.42	110.97	112.66
2	B	2135	XWW	N3-C7-N4	3.59	122.73	117.09
2	B	2135	XWW	C7-N3-C2	-3.37	117.03	122.67
2	B	2135	XWW	N6-C9-N5	-3.27	120.46	126.27
2	A	2133	XWW	C7-N3-C2	-3.12	117.44	122.67
2	B	2135	XWW	N2-C3-N1	2.99	129.87	126.04
2	A	2133	XWW	N6-C9-N5	-2.98	120.98	126.27
2	A	2133	XWW	C11-C10-N7	2.86	115.33	109.93
2	B	2135	XWW	N3-C7-N6	-2.77	112.73	117.09
2	A	2133	XWW	N3-C7-N4	2.57	121.12	117.09
2	B	2135	XWW	C12-C13-N7	2.49	114.65	109.93
2	A	2133	XWW	N6-C7-N4	-2.19	122.67	126.26
2	B	2135	XWW	N6-C9-N7	2.13	120.16	117.12
2	A	2133	XWW	C3-C2-N3	-2.02	108.84	111.70

There are no chirality outliers.

All (6) torsion outliers are listed below:

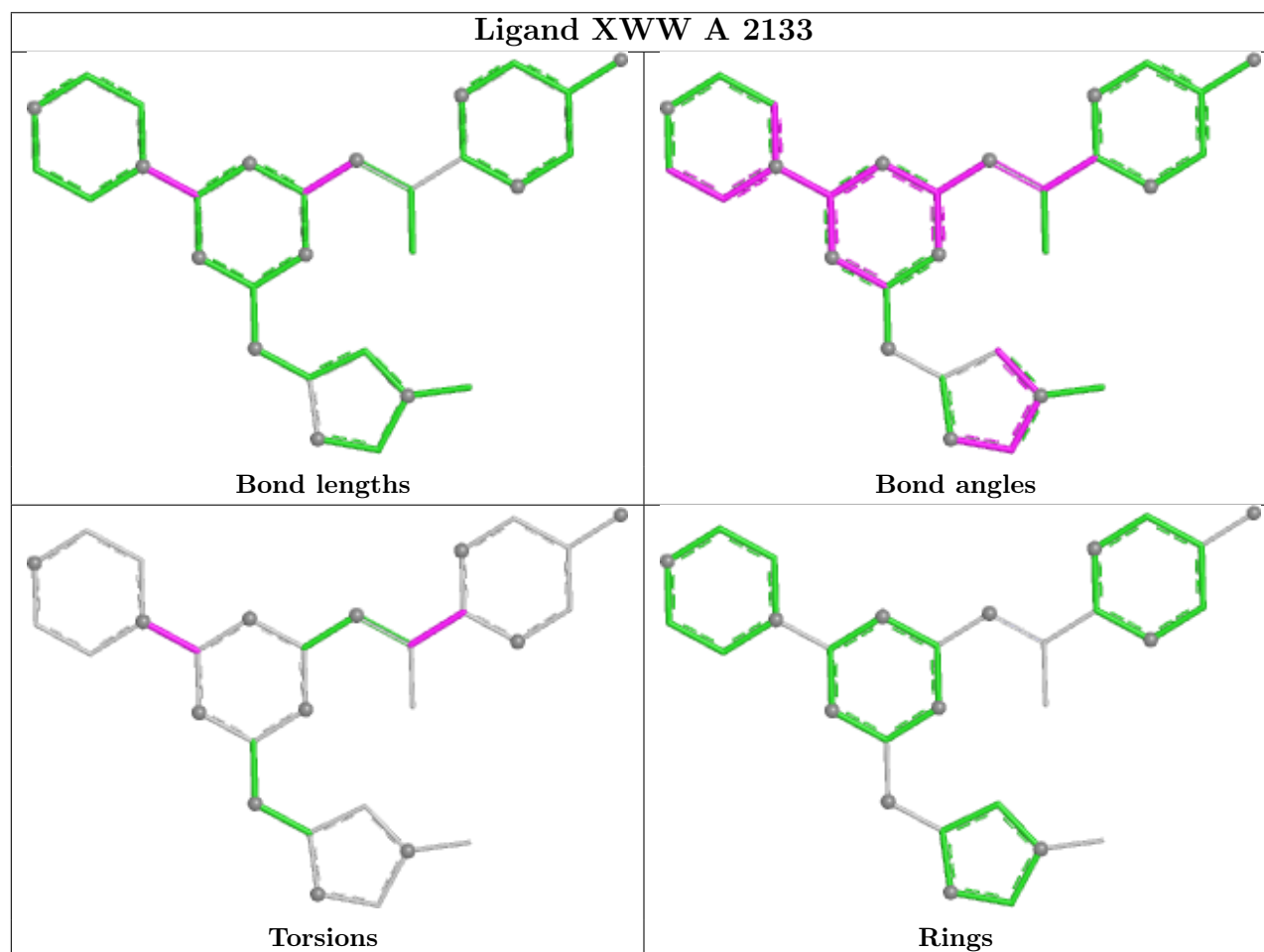
Mol	Chain	Res	Type	Atoms
2	B	2135	XWW	N6-C9-N7-C10
2	B	2135	XWW	N5-C9-N7-C10
2	A	2133	XWW	N3-C2-C3-N2
2	B	2135	XWW	N3-C2-C3-N2
2	A	2133	XWW	N6-C9-N7-C10
2	A	2133	XWW	N5-C9-N7-C10

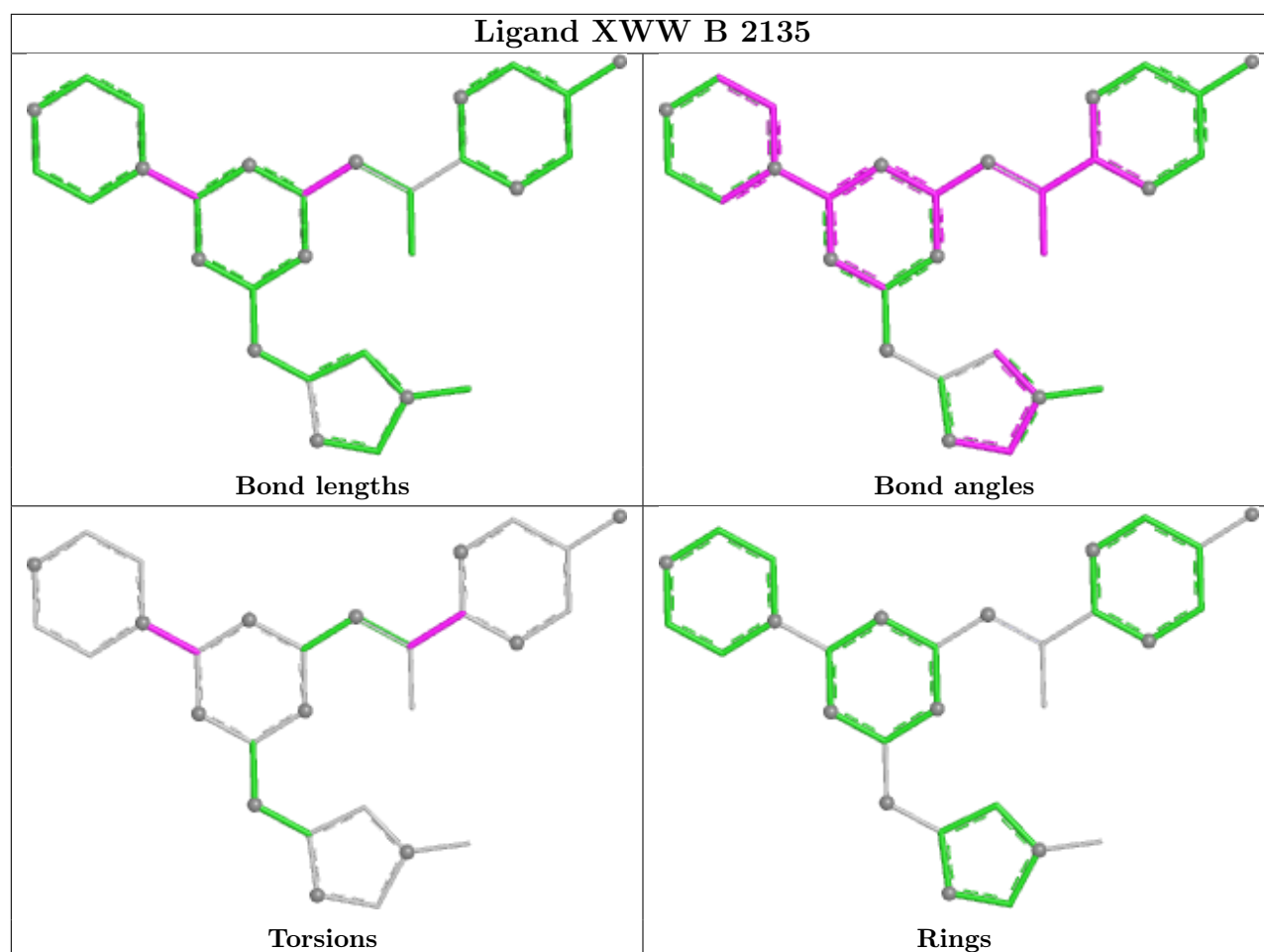
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2133	XWW	1	0
2	B	2135	XWW	2	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, 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%)	-0.10	0 100 100	26, 52, 90, 102	1 (0%)
1	B	271/298 (90%)	-0.19	4 (1%) 72 71	22, 45, 84, 104	0
All	All	544/596 (91%)	-0.14	4 (0%) 84 84	22, 48, 88, 104	1 (0%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1053	LYS	2.5
1	B	870	PRO	2.2
1	B	1072	GLN	2.2
1	B	947	ARG	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, 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	0.88	0.12	70,86,100,102	0
1	PTR	A	1008	16/17	0.92	0.10	71,78,93,95	0
1	PTR	B	1008	16/17	0.95	0.07	41,45,50,50	0
1	PTR	B	1007	16/17	0.96	0.06	41,49,57,60	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

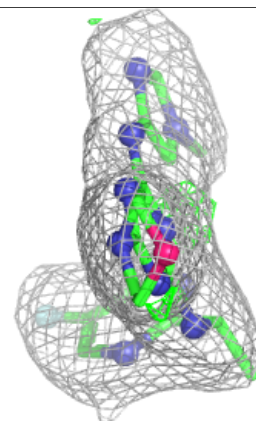
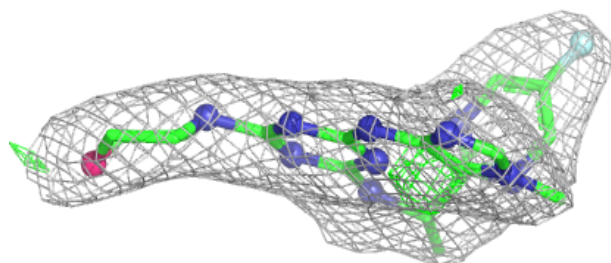
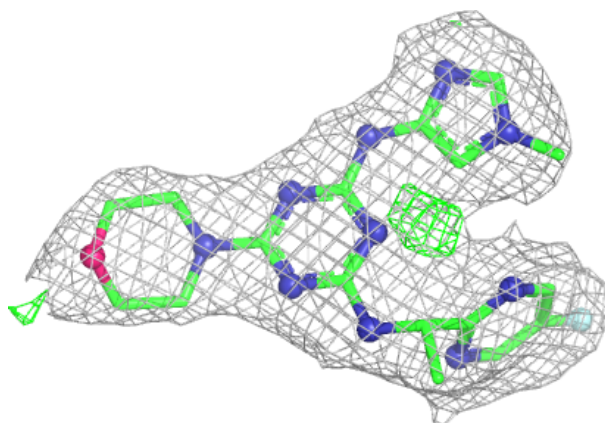
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists 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	ACT	A	2135	3/4	0.81	0.16	57,57,62,63	0
3	ACT	A	2134	3/4	0.89	0.14	58,58,61,64	0
2	XWW	B	2135	29/29	0.92	0.08	33,40,44,48	0
3	ACT	B	2134	3/4	0.96	0.08	40,40,41,41	0
2	XWW	A	2133	29/29	0.97	0.05	24,27,34,38	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.

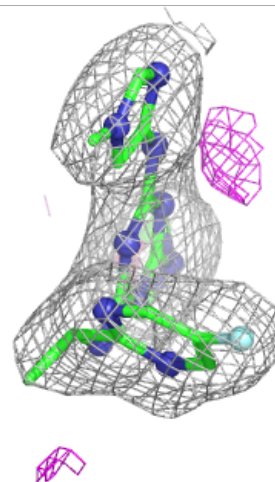
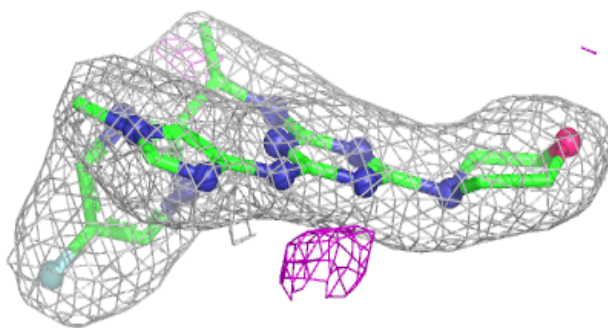
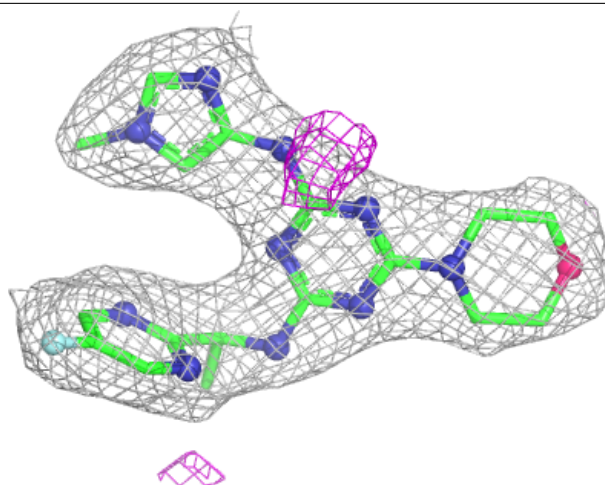
Electron density around XWW B 2135:

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 XWW 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 [i](#)

There are no such residues in this entry.