



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:44 PM UTC

PDB ID : 5FBO / pdb_00005fbo
Title : BTK-inhibitor co-structure
Authors : Fischmann, T.O.
Deposited on : 2015-12-14
Resolution : 1.89 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

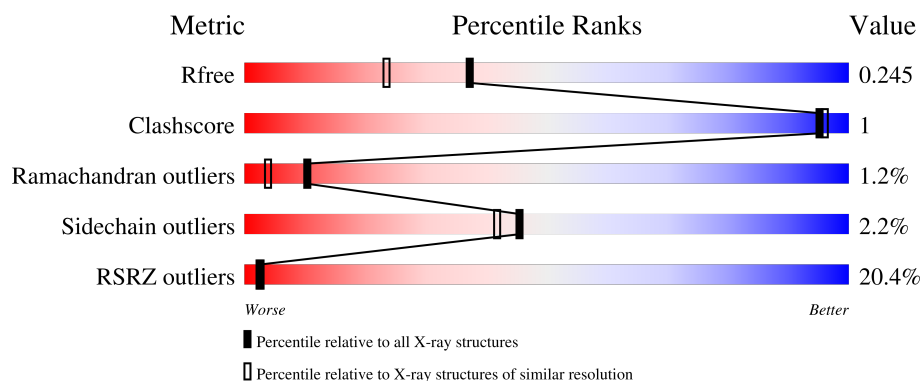
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	271	19% 89% 6%

2 Entry composition [i](#)

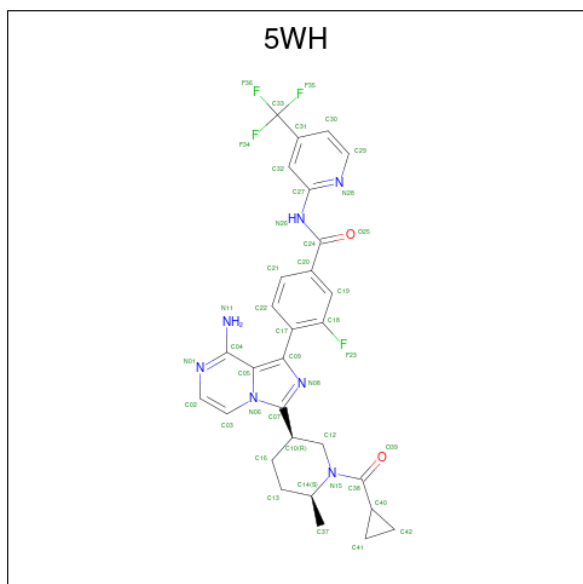
There are 4 unique types of molecules in this entry. The entry contains 2259 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 BTK.

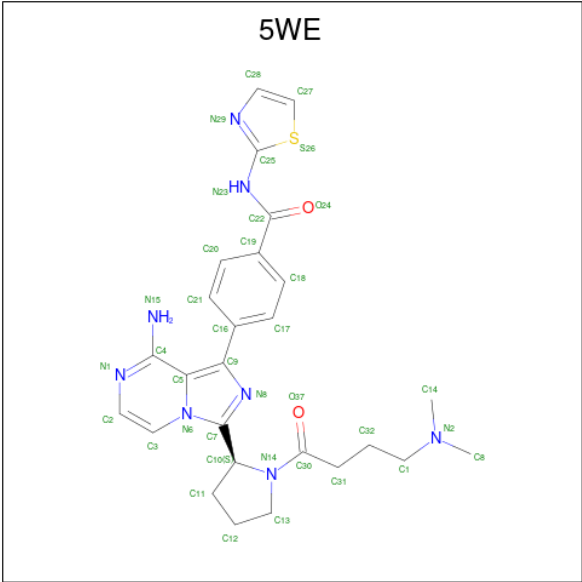
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	0	2	2
			2094	1345	346	384	19			

- Molecule 2 is 4-[8-azanyl-3-[(3 {R},6 {S})-1-cyclopropylcarbonyl-6-methyl-piperidin-3-yl]imidazo[1,5-a]pyrazin-1-yl]-3-fluoranyl- {N}-[4-(trifluoromethyl)pyridin-2-yl]benzamide (CCD ID: 5WH) (formula: C₂₉H₂₇F₄N₇O₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			42	29	4	7	2		

- Molecule 3 is 4-[8-azanyl-3-[(2 {S})-1-[4-(dimethylamino)butanoyl]pyrrolidin-2-yl]imidazo[1,5-a]pyrazin-1-yl]- {N}-(1,3-thiazol-2-yl)benzamide (CCD ID: 5WE) (formula: C₂₆H₃₀N₈O₂S).

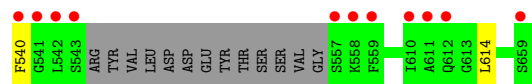


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			37	26	8	2	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	86	Total	O	0	0
			86	86		

- Molecule 1: Tyrosine-protein kinase BTK



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	72.39Å 104.47Å 38.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.08 – 1.89 38.08 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.1 (38.08-1.89) 99.8 (38.08-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 1.89Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.211 , 0.236 0.217 , 0.245	Depositor DCC
R_{free} test set	1188 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	31.6	Xtriage
Anisotropy	0.215	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2259	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.08% 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: 5WE, 5WH

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	0.85	0/2143	1.05	8/2885 (0.3%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	444	GLU	N-CA-C	-7.91	102.54	112.90
1	A	439	GLU	N-CA-C	5.99	123.56	110.80
1	A	438	SER	CA-C-N	5.75	132.53	121.54
1	A	438	SER	C-N-CA	5.75	132.53	121.54
1	A	540	PHE	CA-CB-CG	5.67	119.47	113.80
1	A	471	PHE	CA-CB-CG	5.45	119.25	113.80
1	A	439	GLU	CA-C-N	5.44	127.83	120.38
1	A	439	GLU	C-N-CA	5.44	127.83	120.38

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	2094	0	2058	3	0
2	A	42	0	0	0	0
3	A	37	0	30	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	86	0	0	0	0
All	All	2259	0	2088	5	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 1.

All (5) 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)
3:A:1002:5WE:H9	3:A:1002:5WE:H7	1.24	0.99
1:A:412:GLN:O	1:A:413:PHE:HB2	1.87	0.72
3:A:1002:5WE:H9	3:A:1002:5WE:N15	2.06	0.62
1:A:442:PHE:C	1:A:444:GLU:H	2.19	0.51
1:A:437:MET:HG2	1:A:438:SER:N	2.30	0.46

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	253/271 (93%)	240 (95%)	10 (4%)	3 (1%)	10 4

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	439	GLU
1	A	413	PHE
1	A	443	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	229/243 (94%)	224 (98%)	5 (2%)	45 42

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

Mol	Chain	Res	Type
1	A	412	GLN
1	A	445	GLU
1	A	467	GLN
1	A	518	LEU
1	A	614	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

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	5WH	A	1001	-	45,47,47	1.05	3 (6%)	55,71,71	1.15	4 (7%)
3	5WE	A	1002	-	38,41,41	1.79	8 (21%)	45,58,58	1.23	3 (6%)

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	5WH	A	1001	-	-	5/30/45/45	0/6/6/6
3	5WE	A	1002	-	-	2/26/36/36	0/5/5/5

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1002	5WE	C5-N6	4.62	1.46	1.40
2	A	1001	5WH	C17-C09	-4.30	1.41	1.48
2	A	1001	5WH	C14-N15	3.55	1.51	1.47
3	A	1002	5WE	C7-N8	3.48	1.37	1.32
3	A	1002	5WE	C30-N14	3.32	1.44	1.35
3	A	1002	5WE	C22-N23	3.19	1.43	1.37
3	A	1002	5WE	C25-N29	3.14	1.36	1.31
3	A	1002	5WE	C3-C2	3.06	1.38	1.35
2	A	1001	5WH	C03-C02	2.57	1.37	1.35
3	A	1002	5WE	C21-C16	2.23	1.42	1.39
3	A	1002	5WE	C5-C9	2.11	1.43	1.38

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	5WH	C12-N15-C14	-4.67	111.31	115.72
3	A	1002	5WE	C5-C4-N1	4.48	121.56	117.24
2	A	1001	5WH	C05-C04-N01	3.41	120.53	117.24
3	A	1002	5WE	C9-N8-C7	3.28	108.41	105.38
2	A	1001	5WH	C31-C32-C27	-3.20	118.27	122.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1002	5WE	C11-C10-N14	2.47	106.64	103.02
2	A	1001	5WH	C42-C40-C38	-2.17	114.13	116.91

There are no chirality outliers.

All (7) torsion outliers are listed below:

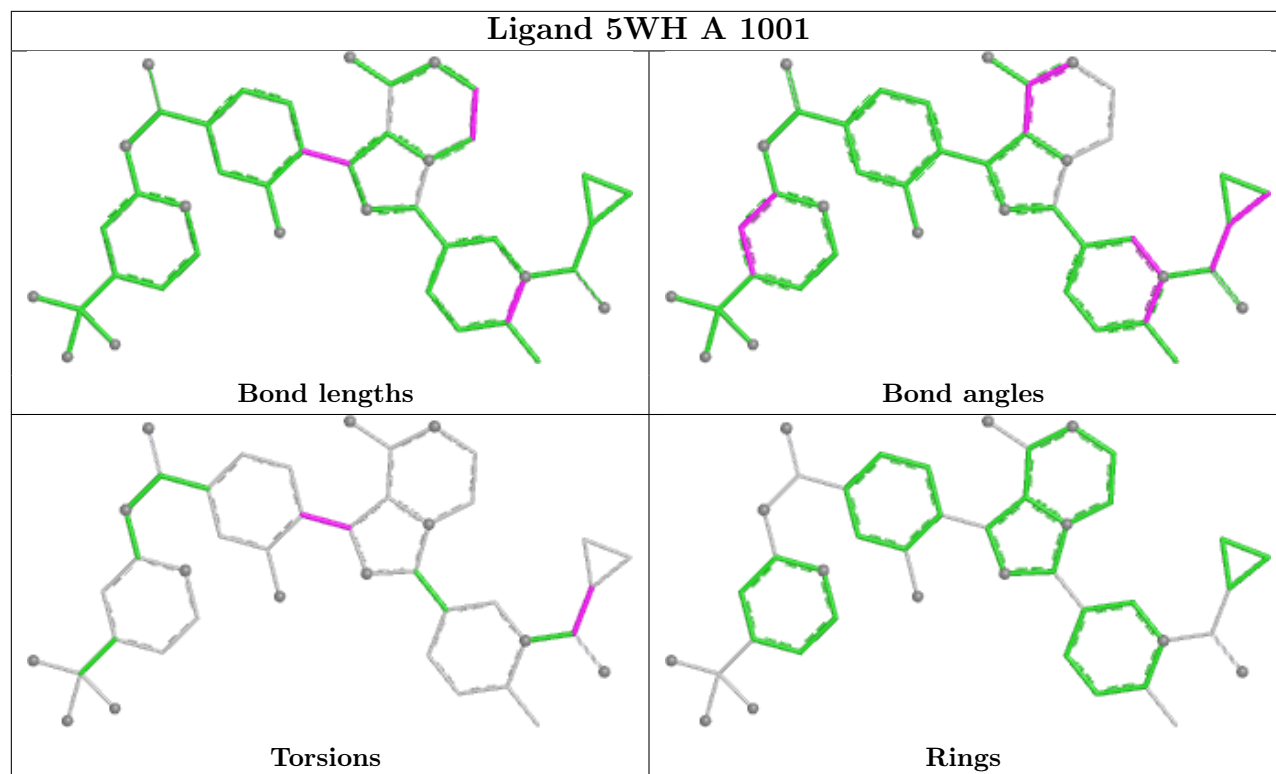
Mol	Chain	Res	Type	Atoms
2	A	1001	5WH	N15-C38-C40-C41
2	A	1001	5WH	N15-C38-C40-C42
3	A	1002	5WE	C32-C1-N2-C8
3	A	1002	5WE	C32-C1-N2-C14
2	A	1001	5WH	O39-C38-C40-C41
2	A	1001	5WH	O39-C38-C40-C42
2	A	1001	5WH	N08-C09-C17-C18

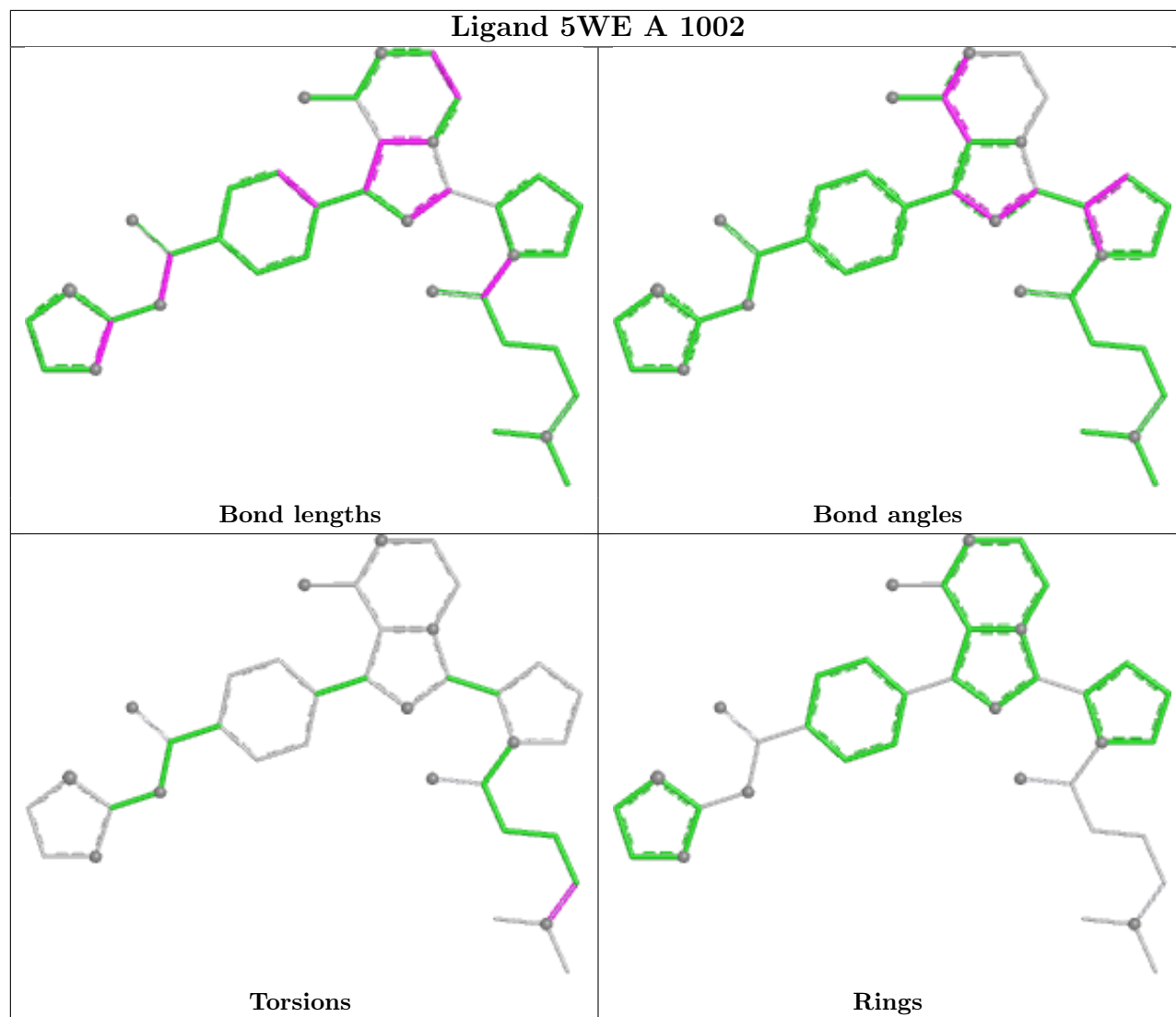
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1002	5WE	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 ⓘ

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	255/271 (94%)	0.91	52 (20%) 3 2	15, 36, 92, 125	2 (0%)

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	543	SER	10.5
1	A	659	SER	7.3
1	A	557	SER	7.2
1	A	443	ILE	7.2
1	A	542	LEU	6.5
1	A	442	PHE	6.2
1	A	413	PHE	6.0
1	A	541	GLY	5.2
1	A	392	TYR	4.6
1	A	432	ILE	4.4
1	A	438	SER	4.2
1	A	489	MET	4.0
1	A	437	MET	3.9
1	A	466	LYS	3.9
1	A	540	PHE	3.9
1	A	404	PHE	3.7
1	A	440	ASP	3.7
1	A	412	GLN	3.5
1	A	439	GLU	3.4
1	A	423	GLY	3.4
1	A	467	GLN	3.3
1	A	435	GLY	3.3
1	A	469	PRO	3.1
1	A	405	LEU	2.9
1	A	431	MET	2.9
1	A	434	GLU	2.9
1	A	448	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	408	LEU	2.8
1	A	411	GLY	2.7
1	A	446	ALA	2.6
1	A	611	ALA	2.6
1	A	410	THR	2.5
1	A	395	TRP	2.5
1	A	518	LEU	2.4
1	A	559	PHE	2.4
1	A	445	GLU	2.4
1	A	468	ARG	2.4
1	A	436	SER	2.4
1	A	539	ASP	2.4
1	A	496	GLN	2.4
1	A	441	GLU	2.3
1	A	393	GLY	2.3
1	A	415	VAL	2.3
1	A	444	GLU	2.3
1	A	610	ILE	2.3
1	A	447	LYS	2.2
1	A	455	GLU	2.2
1	A	470	ILE	2.2
1	A	425	TYR	2.1
1	A	517	PHE	2.1
1	A	612	GLN	2.1
1	A	558	LYS	2.1

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

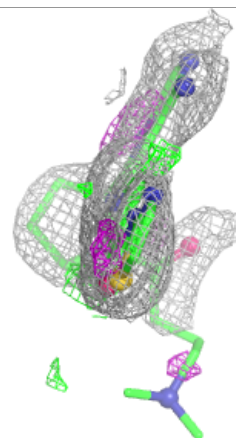
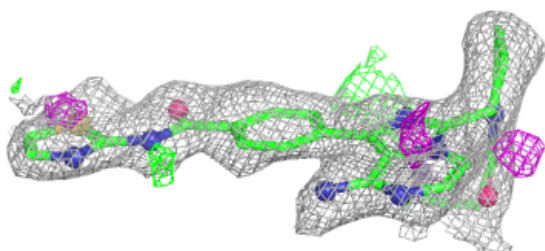
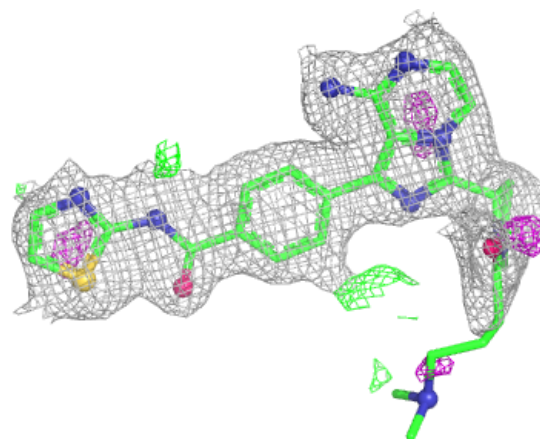
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	5WE	A	1002	37/37	0.77	0.16	58,61,78,78	0
2	5WH	A	1001	42/42	0.89	0.13	27,45,90,93	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.

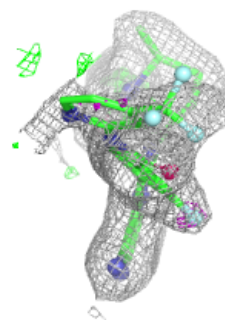
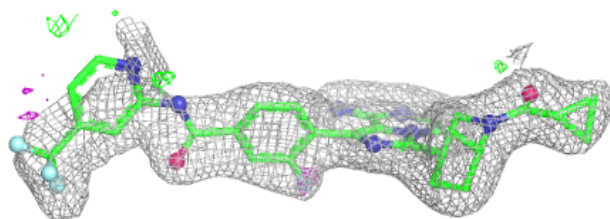
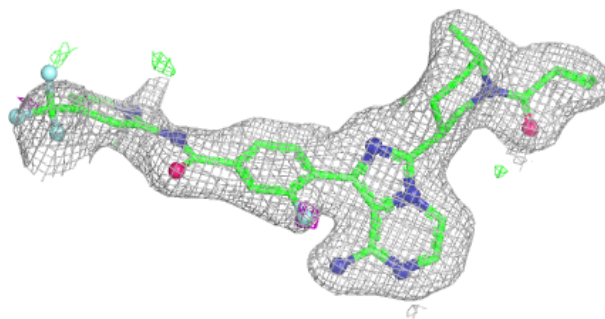
Electron density around 5WE A 1002:

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 5WH A 1001:

$2mF_o-DF_c$ (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.