



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 07:26 am BST

PDB ID : 4V04
Title : FGFR1 in complex with ponatinib.
Authors : Tucker, J.; Klein, T.; Breed, J.; Breeze, A.; Overman, R.; Phillips, C.; Norman, R.A.
Deposited on : 2014-09-10
Resolution : 2.12 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 : 2.11
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) : 2.11

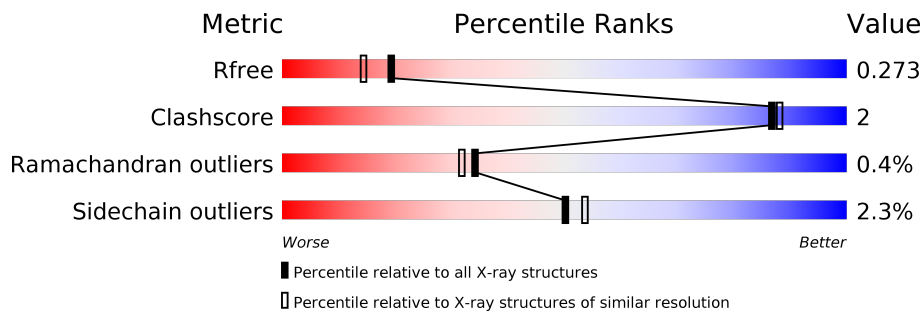
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.12 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	6241 (2.14-2.10)
Clashscore	141614	6778 (2.14-2.10)
Ramachandran outliers	138981	6705 (2.14-2.10)
Sidechain outliers	138945	6706 (2.14-2.10)

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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	309	 89% 6% 5%
1	B	309	 88% 6% 6%

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 5161 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.

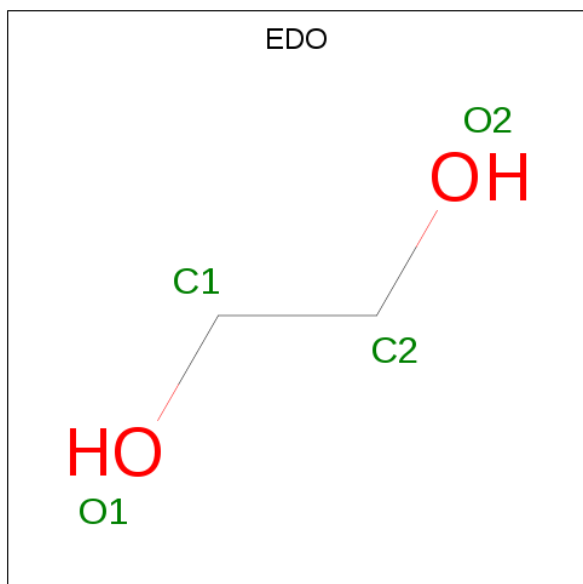
- Molecule 1 is a protein called FIBROBLAST GROWTH FACTOR RECEPTOR 1 (FMS-RELATED TYROSINE KINASE 2, PFEIFFER SYNDROME), ISOFORM CRA_B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	293	2318	1476	395	428	19	0	8	1
1	B	289	2276	1451	386	421	18	0	6	0

There are 6 discrepancies between the modelled and reference sequences:

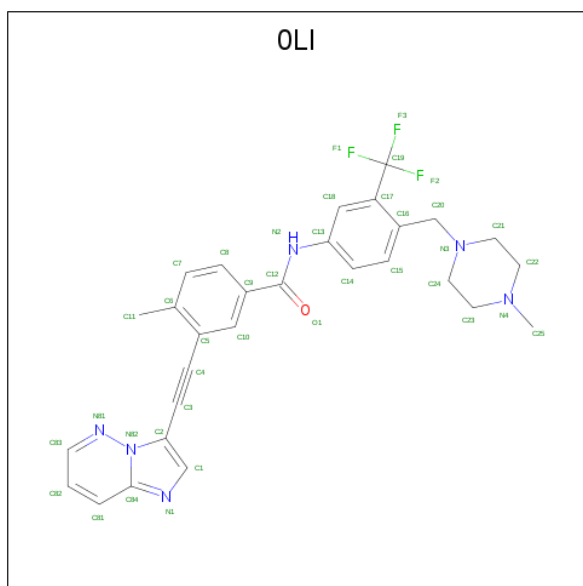
Chain	Residue	Modelled	Actual	Comment	Reference
A	457	GLY	-	expression tag	UNP D3DSX2
A	488	ALA	CYS	engineered mutation	UNP D3DSX2
A	584	SER	CYS	engineered mutation	UNP D3DSX2
B	457	GLY	-	expression tag	UNP D3DSX2
B	488	ALA	CYS	engineered mutation	UNP D3DSX2
B	584	SER	CYS	engineered mutation	UNP D3DSX2

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0

- Molecule 3 is 3-(imidazo[1,2-b]pyridazin-3-ylethynyl)-4-methyl-N-{4-[(4-methylpiperazin-1-yl)methyl]-3-(trifluoromethyl)phenyl}benzamide (three-letter code: 0LI) (formula: C₂₉H₂₇F₃N₆O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	1
			39	29	3	6	1		
3	B	1	Total	C	F	N	O	0	1
			39	29	3	6	1		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is water.

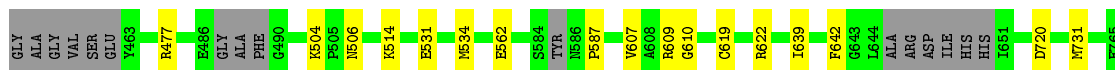
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	215	Total	O	0	0
			215	215		
5	B	205	Total	O	0	0
			205	205		

3 Residue-property plots [i](#)


These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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.

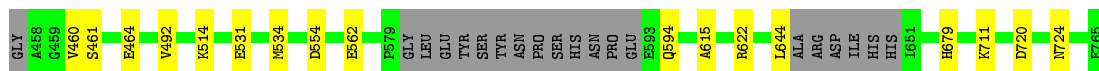
- Molecule 1: FIBROBLAST GROWTH FACTOR RECEPTOR 1 (FMS-RELATED TYROSINE KINASE 2, PFEIFFER SYNDROME), ISOFORM CRA_B

Chain A: 



- Molecule 1: FIBROBLAST GROWTH FACTOR RECEPTOR 1 (FMS-RELATED TYROSINE KINASE 2, PFEIFFER SYNDROME), ISOFORM CRA_B

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	207.25Å 58.09Å 65.23Å 90.00° 107.38° 90.00°	Depositor
Resolution (Å)	21.37 – 2.12 21.38 – 2.11	Depositor EDS
% Data completeness (in resolution range)	97.6 (21.37-2.12) 95.5 (21.38-2.11)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.32 (at 2.11Å)	Xtrriage
Refinement program	BUSTER 2.11.5	Depositor
R, R_{free}	0.186 , 0.228 0.244 , 0.273	Depositor DCC
R_{free} test set	2076 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	37.2	Xtrriage
Anisotropy	0.618	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 64.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.017 for -h-2*1,-k,l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5161	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.81% 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: SO4, OLI, EDO

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	0.52	0/2364	0.64	0/3197
1	B	0.49	0/2321	0.62	0/3139
All	All	0.50	0/4685	0.63	0/6336

There are no bond length outliers.

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	2318	0	2274	11	0
1	B	2276	0	2221	8	1
2	A	24	0	36	0	0
2	B	20	0	30	0	0
3	A	39	0	27	7	0
3	B	39	0	27	6	0
4	A	20	0	0	1	0
4	B	5	0	0	0	0
5	A	215	0	0	1	3
5	B	205	0	0	0	0
All	All	5161	0	4615	20	3

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 2.

All (20) 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:514:LYS:HB2	3:A:1772[A]:OLI:H11	1.62	0.80
4:A:1773:SO4:O2	5:A:2209:HOH:O	2.00	0.79
1:A:622:ARG:HH11	3:A:1772[A]:OLI:H25	1.52	0.74
1:B:514:LYS:HB2	3:B:1771[A]:OLI:H11	1.76	0.68
1:B:622:ARG:HH11	3:B:1771[A]:OLI:H25	1.62	0.64
1:A:619[A]:CYS:SG	3:A:1772[A]:OLI:H24A	2.46	0.56
1:A:562:GLU:O	3:A:1772[A]:OLI:H1	2.11	0.50
1:A:607:VAL:HG13	1:A:639[A]:ILE:HD11	1.96	0.48
1:A:531:GLU:HG3	3:A:1772[A]:OLI:H14	1.97	0.46
1:B:514:LYS:HB2	3:B:1771[A]:OLI:C11	2.45	0.46
1:B:562:GLU:O	3:B:1771[A]:OLI:H1	2.16	0.46
1:A:610:GLY:HA3	1:A:639[B]:ILE:HD12	1.97	0.46
1:A:622:ARG:NH1	3:A:1772[A]:OLI:H25	2.27	0.45
1:A:504:LYS:HG2	1:A:506:ASN:H	1.82	0.45
1:A:477:ARG:NE	1:A:477:ARG:HA	2.33	0.44
1:A:619[A]:CYS:SG	3:A:1772[A]:OLI:C24	3.07	0.43
1:B:492:VAL:HG13	3:B:1771[A]:OLI:H11A	2.02	0.42
1:B:615:ALA:HA	1:B:679:HIS:CE1	2.55	0.42
1:B:461:SER:HA	1:B:464:GLU:O	2.20	0.41
1:B:531:GLU:HG3	3:B:1771[A]:OLI:H14	2.03	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:2036:HOH:O	5:A:2036:HOH:O[2_656]	1.98	0.22
1:B:554:ASP:O	5:A:2190:HOH:O[4_556]	2.18	0.02
5:A:2034:HOH:O	5:A:2034:HOH:O[2_656]	2.18	0.02

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	291/309 (94%)	286 (98%)	4 (1%)	1 (0%)	41	40
1	B	287/309 (93%)	283 (99%)	3 (1%)	1 (0%)	41	40
All	All	578/618 (94%)	569 (98%)	7 (1%)	2 (0%)	34	40

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	460	VAL
1	A	587	PRO

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	247/269 (92%)	241 (98%)	6 (2%)	49	52
1	B	237/269 (88%)	231 (98%)	6 (2%)	47	50
All	All	484/538 (90%)	472 (98%)	12 (2%)	50	50

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	534	MET
1	A	609	ARG
1	A	642[A]	PHE
1	A	642[B]	PHE
1	A	720	ASP
1	A	731	MET
1	B	534	MET
1	B	594	GLN
1	B	644[B]	LEU
1	B	711	LYS
1	B	720	ASP

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Mol	Chain	Res	Type
1	B	724	ASN

Some 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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

18 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	EDO	A	1770	-	3,3,3	0.85	0	2,2,2	0.04	0
2	EDO	B	1770	-	3,3,3	0.69	0	2,2,2	0.30	0
2	EDO	A	1769	-	3,3,3	0.69	0	2,2,2	0.12	0
2	EDO	A	1771	-	3,3,3	0.64	0	2,2,2	0.33	0
3	OLI	B	1771[A]	-	40,43,43	0.89	1 (2%)	51,62,62	0.93	3 (5%)
2	EDO	B	1767	-	3,3,3	0.55	0	2,2,2	0.40	0
4	SO4	A	1776	-	4,4,4	0.28	0	6,6,6	0.16	0
2	EDO	A	1768	-	3,3,3	0.82	0	2,2,2	0.10	0
3	OLI	A	1772[A]	-	40,43,43	0.79	0	51,62,62	0.74	1 (1%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	A	1773	-	4,4,4	0.49	0	6,6,6	0.29	0
4	SO4	B	1772	-	4,4,4	0.32	0	6,6,6	0.35	0
4	SO4	A	1774	-	4,4,4	0.21	0	6,6,6	0.48	0
2	EDO	A	1767	-	3,3,3	0.47	0	2,2,2	0.89	0
2	EDO	A	1766	-	3,3,3	0.42	0	2,2,2	0.27	0
2	EDO	B	1768	-	3,3,3	0.54	0	2,2,2	0.47	0
4	SO4	A	1775	-	4,4,4	0.20	0	6,6,6	0.21	0
2	EDO	B	1766	-	3,3,3	0.32	0	2,2,2	0.58	0
2	EDO	B	1769	-	3,3,3	0.59	0	2,2,2	0.39	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	EDO	A	1770	-	-	0/1/1/1	-
2	EDO	B	1770	-	-	0/1/1/1	-
2	EDO	A	1769	-	-	1/1/1/1	-
2	EDO	A	1771	-	-	1/1/1/1	-
3	OLI	B	1771[A]	-	-	2/20/33/33	0/5/5/5
2	EDO	B	1767	-	-	0/1/1/1	-
2	EDO	A	1768	-	-	1/1/1/1	-
3	OLI	A	1772[A]	-	-	2/20/33/33	0/5/5/5
2	EDO	A	1767	-	-	1/1/1/1	-
2	EDO	A	1766	-	-	0/1/1/1	-
2	EDO	B	1768	-	-	0/1/1/1	-
2	EDO	B	1766	-	-	0/1/1/1	-
2	EDO	B	1769	-	-	0/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1771[A]	OLI	C82-C83	2.16	1.44	1.37

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1771[A]	OLI	C19-C17-C16	2.58	124.11	121.22
3	B	1771[A]	OLI	C82-C81-C84	-2.17	117.97	120.22
3	A	1772[A]	OLI	C82-C81-C84	-2.12	118.02	120.22
3	B	1771[A]	OLI	C8-C9-C12	2.01	127.14	120.62

There are no chirality outliers.

All (8) torsion outliers are listed below:

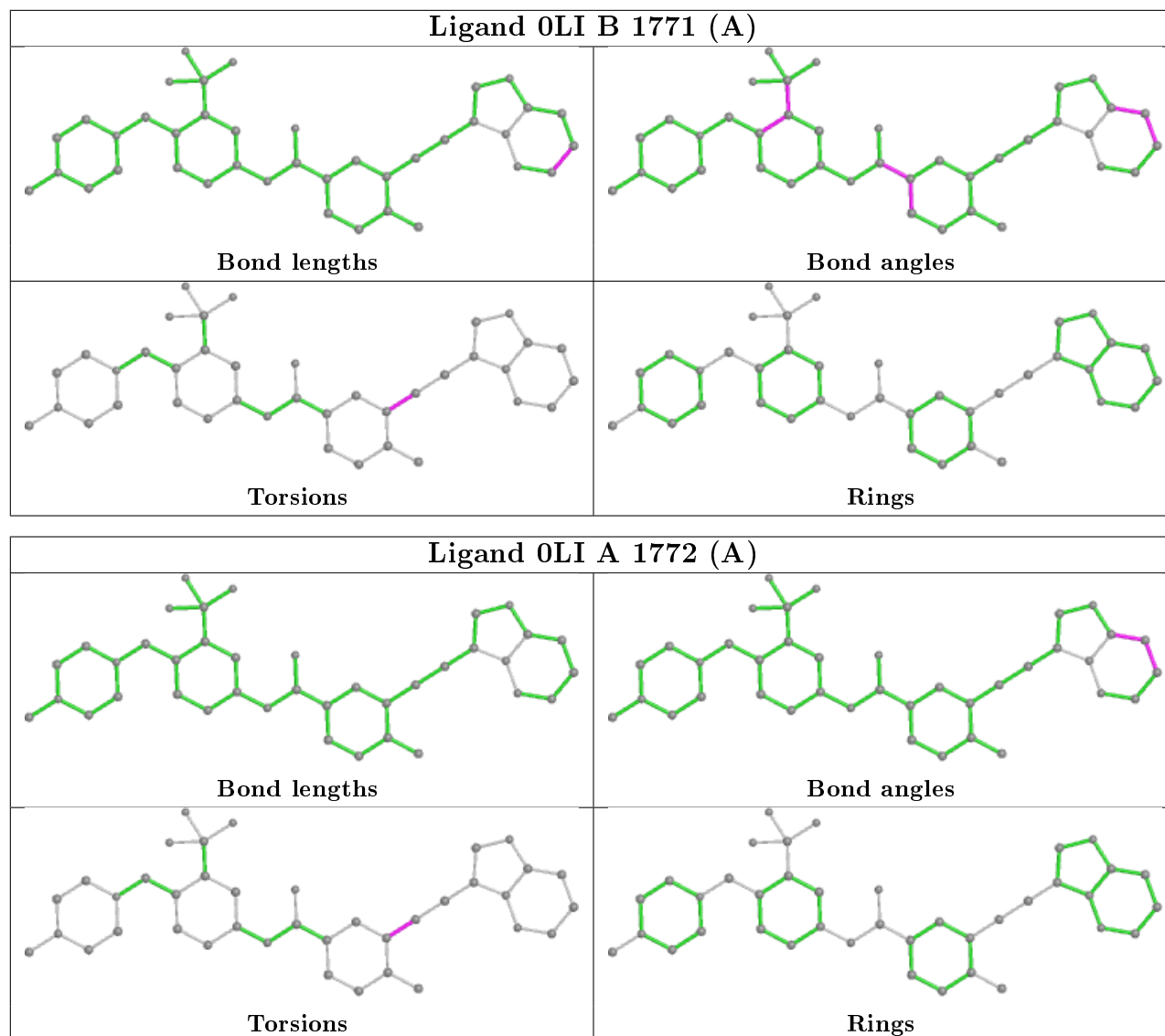
Mol	Chain	Res	Type	Atoms
3	A	1772[A]	0LI	C3-C4-C5-C6
3	A	1772[A]	0LI	C3-C4-C5-C10
3	B	1771[A]	0LI	C3-C4-C5-C10
2	A	1769	EDO	O1-C1-C2-O2
2	A	1768	EDO	O1-C1-C2-O2
3	B	1771[A]	0LI	C3-C4-C5-C6
2	A	1771	EDO	O1-C1-C2-O2
2	A	1767	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1771[A]	0LI	6	0
3	A	1772[A]	0LI	7	0
4	A	1773	SO4	1	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](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

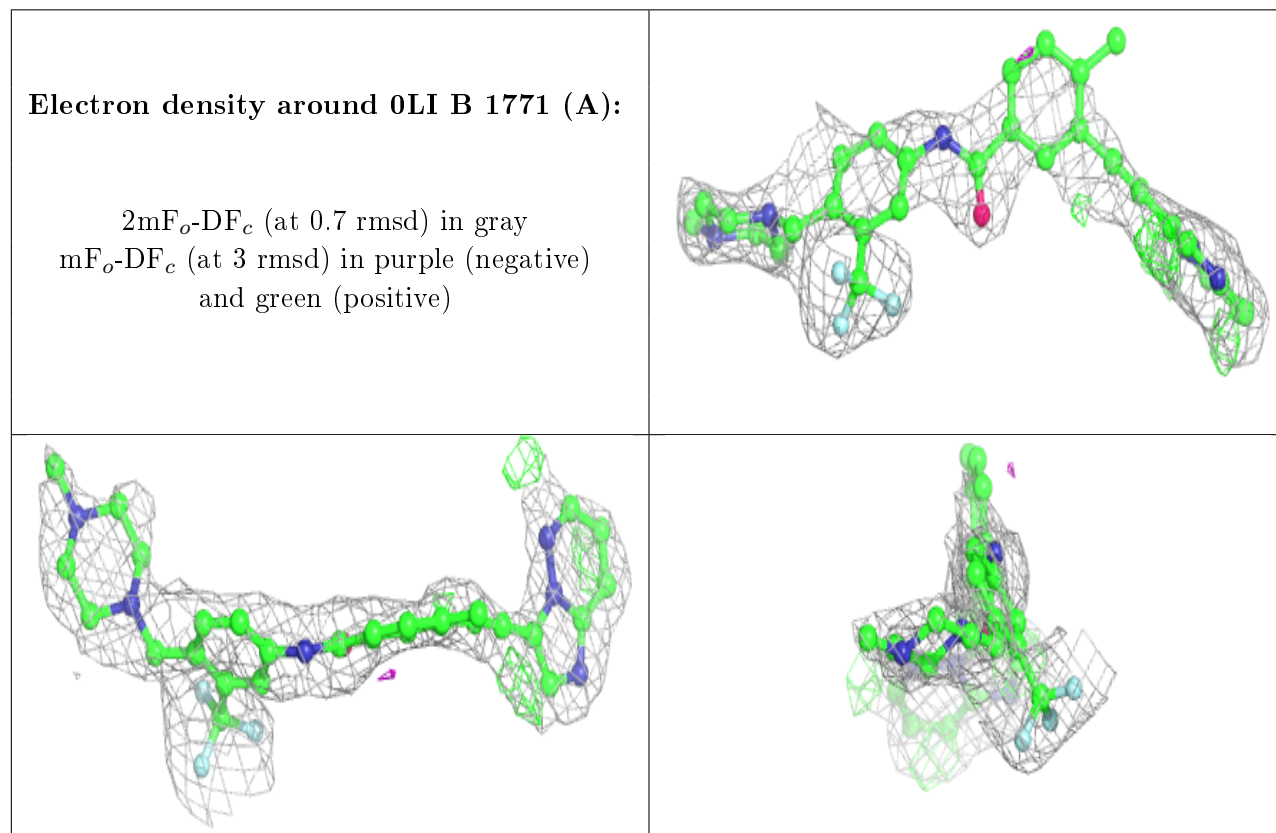
6.3 Carbohydrates [i](#)

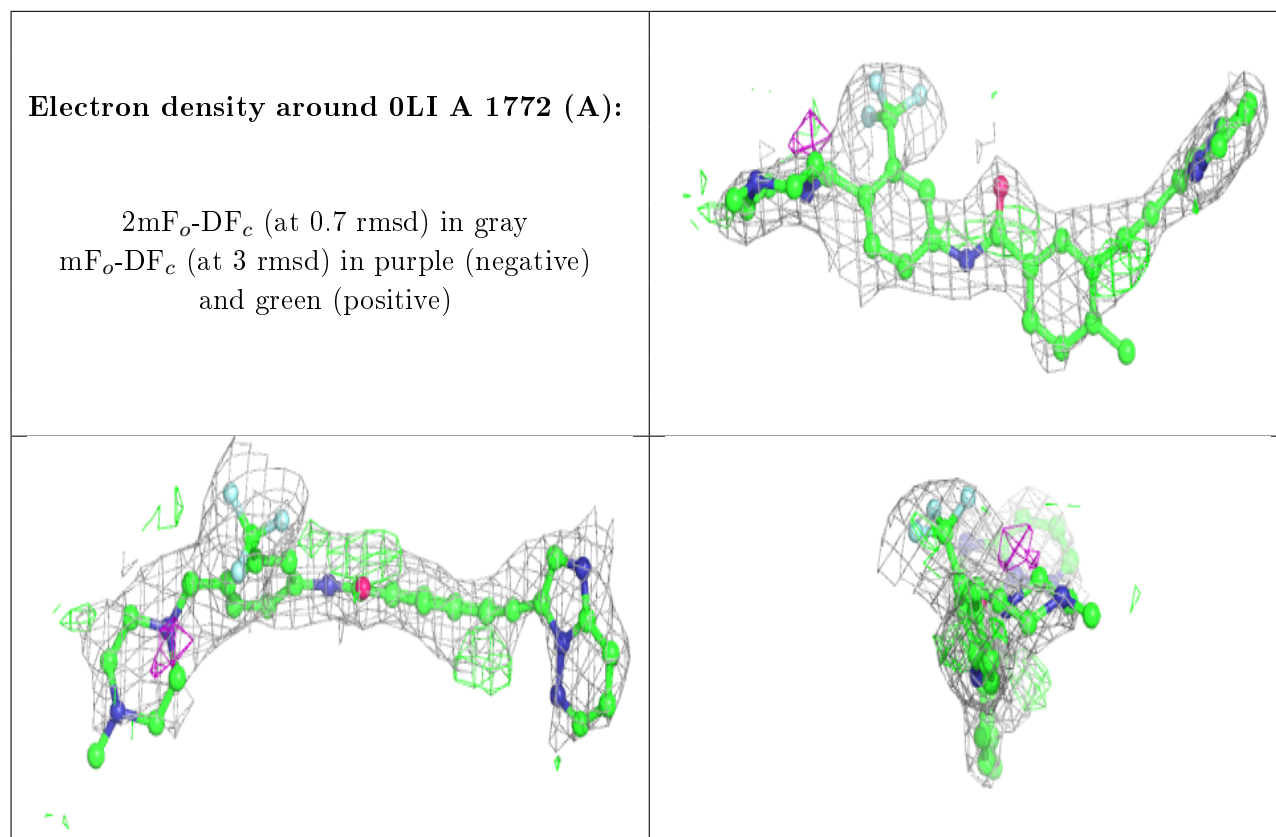
Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

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.





6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.