



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:44 AM UTC

PDB ID : 4PP7 / pdb\_00004pp7  
Title : Highly Potent and Selective 3-N-methylquinazoline-4(3H)-one Based Inhibitors of B-RafV600E Kinase  
Authors : Wenglowky, S.; Ren, L.; Grina, J.; Hansen, J.D.; Laird, E.R.; Moreno, D.; Dinkel, V.; Gloor, S.L.; Hastings, G.; Rana, S.; Rasor, K.; Sturgis, H.L.; Voegtli, W.C.; Vigers, G.P.A.; Willis, B.; Mathieu, S.; Rudolph, J.  
Deposited on : 2014-02-26  
Resolution : 3.40 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](#) ①) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtrriage (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

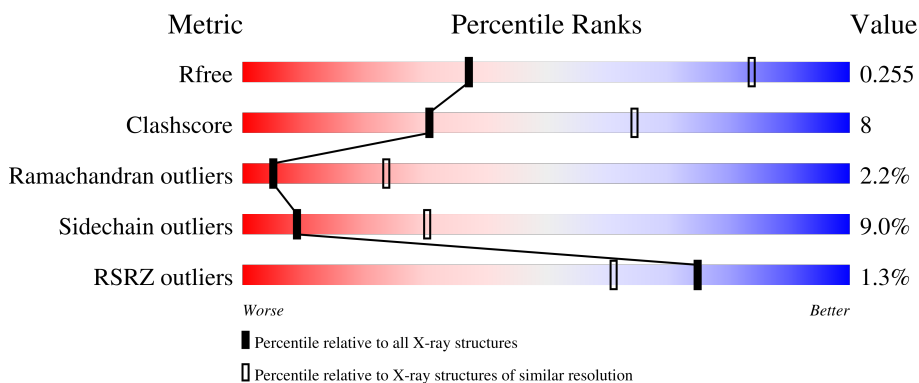
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.40 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))
$R_{free}$	180053	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	307	 % 65% 18% • 13%
1	B	307	 % 65% 21% • 11%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4390 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.

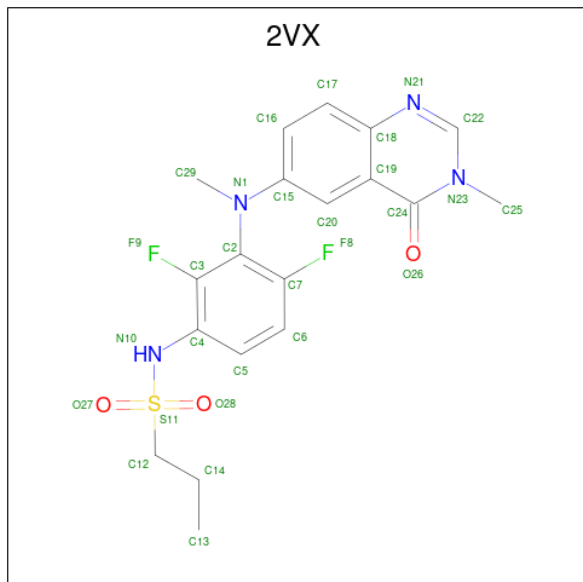
- Molecule 1 is a protein called Serine/threonine-protein kinase B-raf.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	267	Total	C	N	O	S	0	0	0
			2137	1370	373	381	13			
1	B	274	Total	C	N	O	S	0	0	0
			2195	1405	384	393	13			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	420	MET	-	expression tag	UNP P15056
A	421	ASP	-	expression tag	UNP P15056
A	422	ARG	-	expression tag	UNP P15056
A	423	GLY	-	expression tag	UNP P15056
A	424	SER	-	expression tag	UNP P15056
A	425	HIS	-	expression tag	UNP P15056
A	426	HIS	-	expression tag	UNP P15056
A	427	HIS	-	expression tag	UNP P15056
A	428	HIS	-	expression tag	UNP P15056
A	429	HIS	-	expression tag	UNP P15056
A	430	HIS	-	expression tag	UNP P15056
A	431	GLY	-	expression tag	UNP P15056
B	420	MET	-	expression tag	UNP P15056
B	421	ASP	-	expression tag	UNP P15056
B	422	ARG	-	expression tag	UNP P15056
B	423	GLY	-	expression tag	UNP P15056
B	424	SER	-	expression tag	UNP P15056
B	425	HIS	-	expression tag	UNP P15056
B	426	HIS	-	expression tag	UNP P15056
B	427	HIS	-	expression tag	UNP P15056
B	428	HIS	-	expression tag	UNP P15056
B	429	HIS	-	expression tag	UNP P15056
B	430	HIS	-	expression tag	UNP P15056
B	431	GLY	-	expression tag	UNP P15056

- Molecule 2 is N-{2,4-difluoro-3-[methyl(3-methyl-4-oxo-3,4-dihydroquinazolin-6-yl)amino]phenyl}propane-1-sulfonamide (CCD ID: 2VX) (formula: C<sub>19</sub>H<sub>20</sub>F<sub>2</sub>N<sub>4</sub>O<sub>3</sub>S).

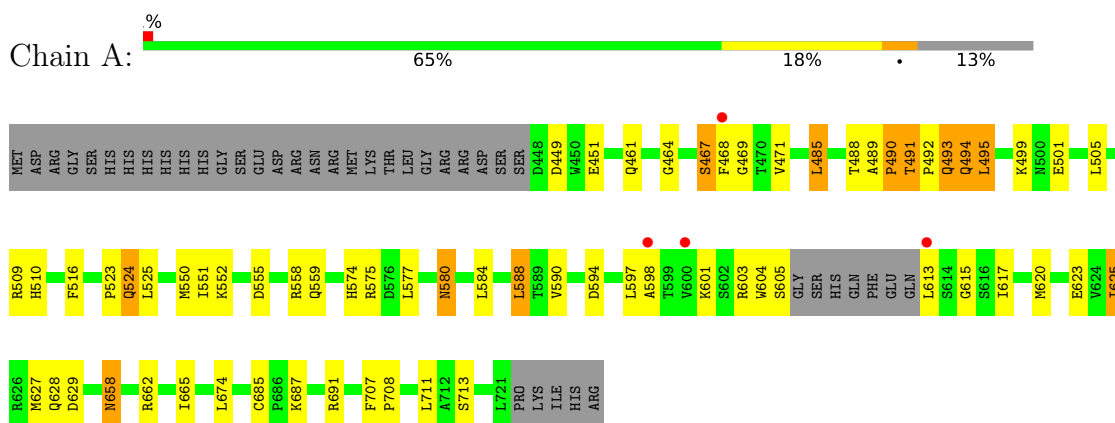


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	F	N	O			S
2	A	1	Total	19	2	4	3	1	0	0
2	B	1	Total	19	2	4	3	1	0	0

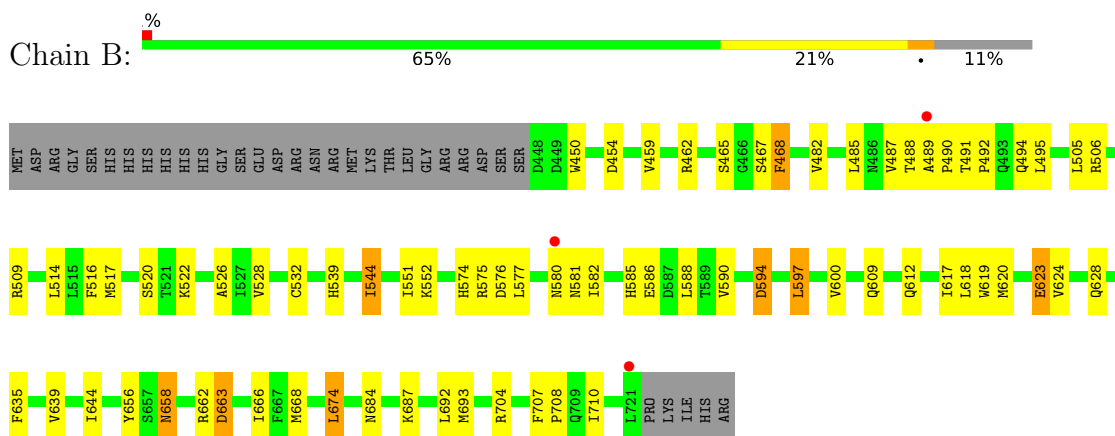
### 3 Residue-property plots [i](#)

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: Serine/threonine-protein kinase B-raf



- Molecule 1: Serine/threonine-protein kinase B-raf



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	107.97Å 107.97Å 151.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.75 – 3.40 24.75 – 3.40	Depositor EDS
% Data completeness (in resolution range)	98.5 (24.75-3.40) 98.5 (24.75-3.40)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	0.12	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.08 (at 3.38Å)	Xtrriage
Refinement program	REFMAC 5.8.0049	Depositor
R, $R_{free}$	0.195 , 0.254 0.195 , 0.255	Depositor DCC
$R_{free}$ test set	620 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	70.2	Xtrriage
Anisotropy	0.050	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 35.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4390	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	73.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% 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: 2VX

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	0.41	0/2183	0.83	4/2946 (0.1%)
1	B	0.40	0/2244	0.83	0/3029
All	All	0.41	0/4427	0.83	4/5975 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	489	ALA	CA-C-N	5.34	125.25	119.85
1	A	489	ALA	C-N-CA	5.34	125.25	119.85
1	A	490	PRO	CA-C-N	5.04	130.77	121.70
1	A	490	PRO	C-N-CA	5.04	130.77	121.70

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	2137	0	2177	35	0
1	B	2195	0	2224	39	0
2	A	29	0	20	5	0
2	B	29	0	20	5	0
All	All	4390	0	4441	73	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 8.

All (73) 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:490:PRO:HA	1:B:491:THR:OG1	1.57	1.05
1:A:603:ARG:HH11	1:A:603:ARG:HG3	1.45	0.81
1:A:490:PRO:HA	1:A:491:THR:OG1	1.82	0.79
1:B:658:ASN:HD22	1:B:658:ASN:H	1.33	0.72
1:A:490:PRO:HB3	1:A:494:GLN:HB2	1.76	0.66
1:A:594:ASP:H	2:A:801:2VX:H3	1.42	0.64
1:A:603:ARG:HG3	1:A:603:ARG:NH1	2.13	0.63
1:A:467:SER:O	1:A:469:GLY:N	2.29	0.63
1:B:454:ASP:H	1:B:522:LYS:HD3	1.65	0.61
1:A:464:GLY:HA3	1:A:471:VAL:HB	1.85	0.58
1:A:505:LEU:HD22	2:A:801:2VX:H9	1.85	0.58
1:A:492:PRO:HA	1:A:493:GLN:C	2.28	0.58
1:B:594:ASP:H	2:B:801:2VX:H3	1.48	0.58
1:A:523:PRO:O	1:A:524:GLN:HB2	2.05	0.57
1:A:598:ALA:HA	1:A:601:LYS:HB2	1.87	0.56
1:B:551:ILE:HD12	1:B:551:ILE:H	1.71	0.56
1:B:467:SER:O	1:B:468:PHE:HB3	2.07	0.53
1:B:582:ILE:HG23	1:B:590:VAL:HG13	1.92	0.52
1:B:620:MET:HE3	1:B:624:VAL:HG12	1.90	0.52
2:A:801:2VX:O27	2:A:801:2VX:H1	2.10	0.52
1:A:658:ASN:H	1:A:658:ASN:HD22	1.57	0.52
1:A:495:LEU:HD22	1:A:499:LYS:HE2	1.92	0.51
1:A:551:ILE:H	1:A:551:ILE:HD12	1.76	0.51
1:B:520:SER:HB3	1:B:526:ALA:HB3	1.91	0.51
1:B:586:GLU:O	1:B:588:LEU:HD12	2.11	0.50
1:B:454:ASP:N	1:B:522:LYS:HD3	2.26	0.50
1:A:524:GLN:HG2	1:A:525:LEU:H	1.77	0.49
1:A:707:PHE:O	1:A:711:LEU:HB2	2.13	0.49
1:B:450:TRP:HE1	1:B:506:ARG:HD2	1.77	0.49
1:A:620:MET:HE3	1:A:625:ILE:HD13	1.94	0.49
1:A:524:GLN:HG2	1:A:525:LEU:N	2.27	0.49
1:A:601:LYS:O	1:A:605:SER:HB3	2.12	0.48
1:B:532:CYS:O	2:B:801:2VX:H14	2.13	0.48
1:B:576:ASP:O	1:B:581:ASN:ND2	2.46	0.48
1:A:617:ILE:HA	1:A:620:MET:HE2	1.95	0.48
1:A:501:GLU:HG3	1:A:597:LEU:HA	1.95	0.48
1:A:509:ARG:HD3	1:B:516:PHE:O	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:510:HIS:HA	1:B:517:MET:HE2	1.96	0.48
1:A:516:PHE:O	1:B:509:ARG:HD3	2.14	0.48
1:B:692:LEU:HD11	1:B:710:ILE:HG23	1.96	0.47
1:B:514:LEU:HD23	2:B:801:2VX:H9	1.97	0.47
1:A:471:VAL:HG21	2:A:801:2VX:H19	1.97	0.46
2:B:801:2VX:H13	2:B:801:2VX:H18	1.70	0.46
1:B:468:PHE:HZ	1:B:597:LEU:HD23	1.81	0.46
1:A:485:LEU:H	1:A:485:LEU:HD13	1.81	0.46
1:B:482:VAL:HG22	1:B:528:VAL:HG22	1.98	0.46
1:B:532:CYS:HA	1:B:585:HIS:HD2	1.81	0.45
1:B:707:PHE:O	1:B:708:PRO:C	2.60	0.45
1:A:707:PHE:N	1:A:708:PRO:HD2	2.32	0.44
1:B:492:PRO:HA	1:B:494:GLN:H	1.82	0.44
1:B:544:ILE:H	1:B:544:ILE:HG13	1.60	0.44
1:A:580:ASN:ND2	1:A:580:ASN:H	2.16	0.44
1:B:574:HIS:O	1:B:575:ARG:HB2	2.18	0.43
1:B:619:TRP:HA	1:B:644:ILE:HG13	2.00	0.43
1:B:635:PHE:O	1:B:639:VAL:HG23	2.18	0.43
1:B:617:ILE:HG13	1:B:618:LEU:HD12	2.00	0.43
1:B:623:GLU:OE1	1:B:704:ARG:NH2	2.45	0.43
1:A:559:GLN:HB3	1:A:590:VAL:HB	1.99	0.42
1:B:505:LEU:HD22	2:B:801:2VX:H4	2.00	0.42
1:A:588:LEU:HD13	1:B:586:GLU:HG3	2.01	0.42
1:B:693:MET:HE3	1:B:693:MET:HB2	1.95	0.42
1:B:468:PHE:CZ	1:B:597:LEU:HD23	2.55	0.42
1:A:555:ASP:OD1	1:A:558:ARG:NH1	2.52	0.42
2:A:801:2VX:H18	2:A:801:2VX:H13	1.68	0.42
1:B:663:ASP:HA	1:B:666:ILE:HD12	2.02	0.42
1:A:580:ASN:H	1:A:580:ASN:HD22	1.68	0.41
1:A:574:HIS:O	1:A:575:ARG:HB2	2.21	0.41
1:B:462:ARG:HH11	1:B:465:SER:HB3	1.86	0.41
1:A:628:GLN:NE2	1:A:629:ASP:OD1	2.54	0.41
1:B:585:HIS:CE1	1:B:586:GLU:HG2	2.56	0.41
1:B:668:MET:HE2	1:B:674:LEU:HB2	2.02	0.41
1:B:539:HIS:ND1	1:B:544:ILE:HD11	2.36	0.41
1:A:550:MET:CE	1:A:685:CYS:HA	2.51	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	263/307 (86%)	237 (90%)	18 (7%)	8 (3%)	3	18
1	B	272/307 (89%)	242 (89%)	26 (10%)	4 (2%)	8	30
All	All	535/614 (87%)	479 (90%)	44 (8%)	12 (2%)	5	24

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	468	PHE
1	A	488	THR
1	A	524	GLN
1	B	468	PHE
1	A	494	GLN
1	B	489	ALA
1	A	467	SER
1	A	491	THR
1	A	615	GLY
1	A	627	MET
1	B	487	VAL
1	B	488	THR

### 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	235/271 (87%)	213 (91%)	22 (9%)	8	29

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	241/271 (89%)	220 (91%)	21 (9%)	9	32
All	All	476/542 (88%)	433 (91%)	43 (9%)	9	30

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	449	ASP
1	A	451	GLU
1	A	461	GLN
1	A	485	LEU
1	A	493	GLN
1	A	495	LEU
1	A	552	LYS
1	A	577	LEU
1	A	580	ASN
1	A	584	LEU
1	A	588	LEU
1	A	604	TRP
1	A	613	LEU
1	A	623	GLU
1	A	625	ILE
1	A	658	ASN
1	A	662	ARG
1	A	665	ILE
1	A	674	LEU
1	A	687	LYS
1	A	691	ARG
1	A	713	SER
1	B	459	VAL
1	B	485	LEU
1	B	495	LEU
1	B	544	ILE
1	B	552	LYS
1	B	577	LEU
1	B	580	ASN
1	B	594	ASP
1	B	597	LEU
1	B	600	VAL
1	B	609	GLN
1	B	612	GLN
1	B	623	GLU
1	B	628	GLN

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Mol	Chain	Res	Type
1	B	656	TYR
1	B	658	ASN
1	B	662	ARG
1	B	663	ASP
1	B	674	LEU
1	B	684	ASN
1	B	687	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	494	GLN
1	A	524	GLN
1	A	539	HIS
1	A	574	HIS
1	A	580	ASN
1	A	585	HIS
1	A	631	ASN
1	A	658	ASN
1	B	493	GLN
1	B	524	GLN
1	B	580	ASN
1	B	581	ASN
1	B	585	HIS
1	B	609	GLN
1	B	612	GLN
1	B	628	GLN
1	B	631	ASN
1	B	653	GLN
1	B	658	ASN

### 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	2VX	A	801	-	30,31,31	1.98	7 (23%)	43,46,46	1.64	10 (23%)
2	2VX	B	801	-	30,31,31	1.95	6 (20%)	43,46,46	1.53	7 (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	2VX	A	801	-	-	7/17/17/17	0/3/3/3
2	2VX	B	801	-	-	2/17/17/17	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	2VX	C4-N10	-5.51	1.34	1.42
2	B	801	2VX	S11-N10	5.26	1.73	1.62
2	A	801	2VX	S11-N10	4.95	1.72	1.62
2	B	801	2VX	C4-N10	-4.92	1.35	1.42
2	B	801	2VX	O27-S11	3.06	1.47	1.43
2	A	801	2VX	O27-S11	2.86	1.47	1.43
2	B	801	2VX	O28-S11	2.82	1.47	1.43
2	B	801	2VX	C15-N1	-2.71	1.35	1.42
2	A	801	2VX	O28-S11	2.58	1.47	1.43
2	A	801	2VX	C15-N1	-2.29	1.36	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	2VX	C12-S11	2.14	1.85	1.78
2	B	801	2VX	C12-S11	2.14	1.84	1.78
2	A	801	2VX	C6-C7	2.04	1.42	1.37

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	2VX	C15-N1-C2	4.79	125.94	119.98
2	B	801	2VX	C15-N1-C2	4.30	125.34	119.98
2	B	801	2VX	C7-C2-C3	3.62	119.14	115.25
2	A	801	2VX	C4-N10-S11	3.29	131.58	123.58
2	B	801	2VX	O26-C24-N23	3.25	124.08	119.70
2	B	801	2VX	C2-C3-C4	-3.13	120.43	124.75
2	A	801	2VX	O26-C24-N23	3.07	123.84	119.70
2	A	801	2VX	C2-C3-C4	-2.96	120.66	124.75
2	A	801	2VX	C18-N21-C22	2.95	119.69	116.69
2	A	801	2VX	C7-C2-C3	2.88	118.34	115.25
2	A	801	2VX	C3-C4-N10	-2.86	117.43	120.73
2	B	801	2VX	C4-N10-S11	2.84	130.48	123.58
2	B	801	2VX	C12-S11-N10	-2.71	102.88	106.72
2	A	801	2VX	C12-S11-N10	-2.52	103.14	106.72
2	B	801	2VX	C18-N21-C22	2.37	119.11	116.69
2	A	801	2VX	C5-C4-C3	2.25	120.85	117.72
2	A	801	2VX	O27-S11-C12	2.02	111.07	107.86

There are no chirality outliers.

All (9) torsion outliers are listed below:

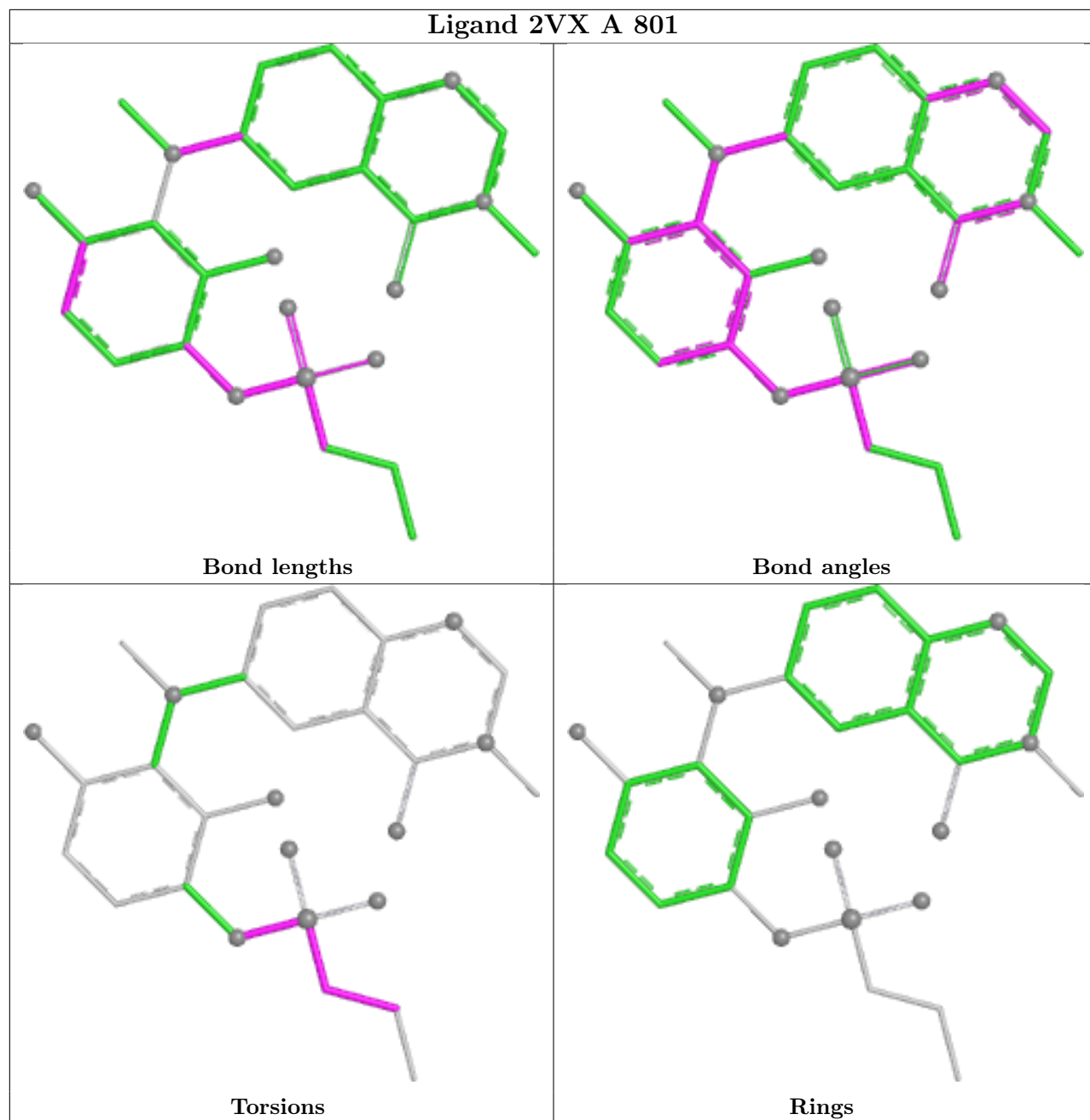
Mol	Chain	Res	Type	Atoms
2	A	801	2VX	C14-C12-S11-N10
2	A	801	2VX	C14-C12-S11-O27
2	A	801	2VX	C14-C12-S11-O28
2	A	801	2VX	S11-C12-C14-C13
2	B	801	2VX	S11-C12-C14-C13
2	A	801	2VX	C4-N10-S11-O27
2	A	801	2VX	C4-N10-S11-O28
2	B	801	2VX	C7-C2-N1-C29
2	A	801	2VX	C4-N10-S11-C12

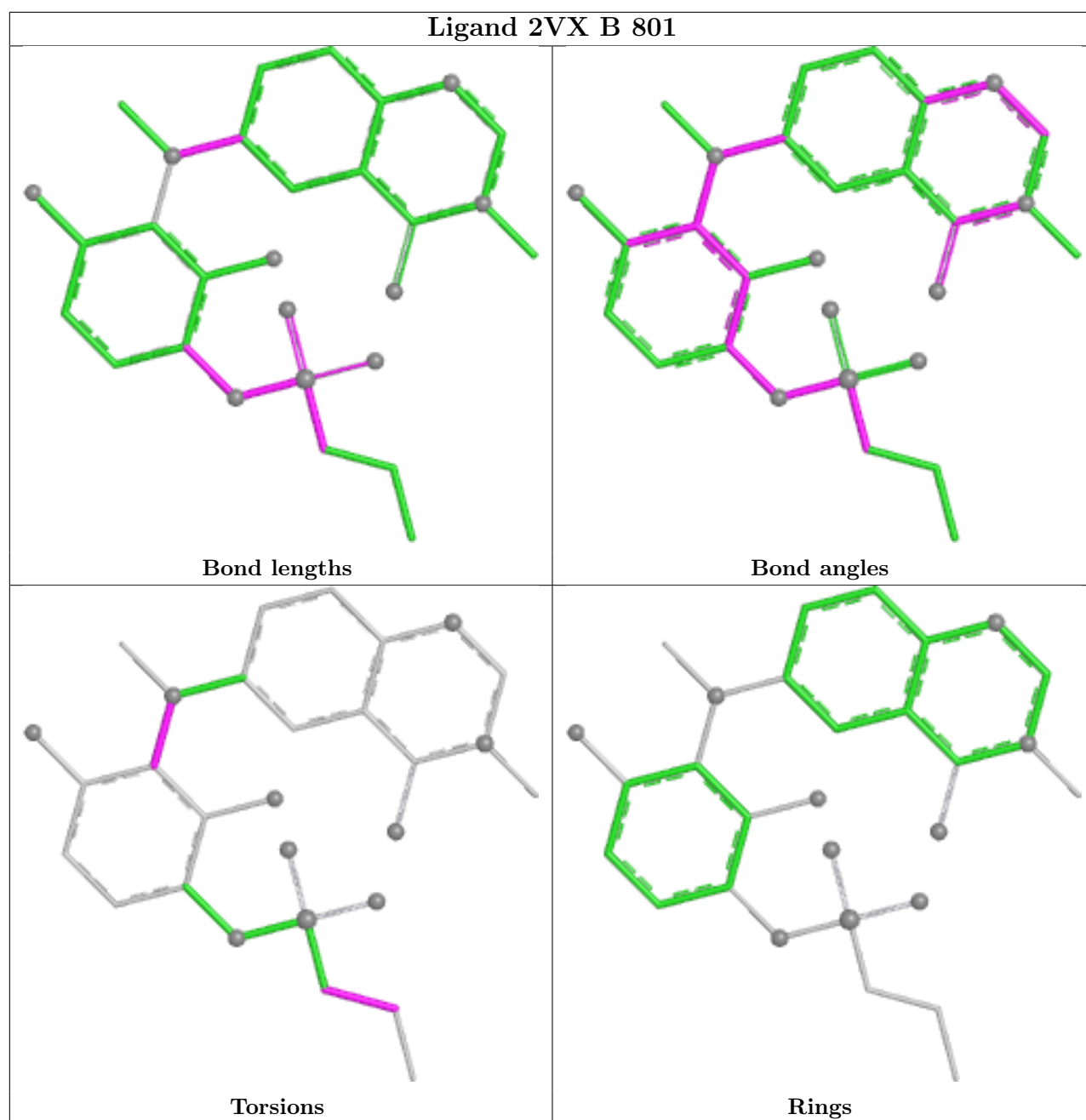
There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	2VX	5	0
2	B	801	2VX	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	267/307 (86%)	-0.12	4 (1%) 72 57	44, 67, 130, 152	0
1	B	274/307 (89%)	-0.21	3 (1%) 78 65	44, 67, 111, 129	0
All	All	541/614 (88%)	-0.16	7 (1%) 75 61	44, 67, 121, 152	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	721	LEU	5.4
1	B	580	ASN	2.7
1	A	600	VAL	2.6
1	A	613	LEU	2.5
1	A	468	PHE	2.5
1	A	598	ALA	2.2
1	B	489	ALA	2.0

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

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

### 6.3 Carbohydrates [i](#)

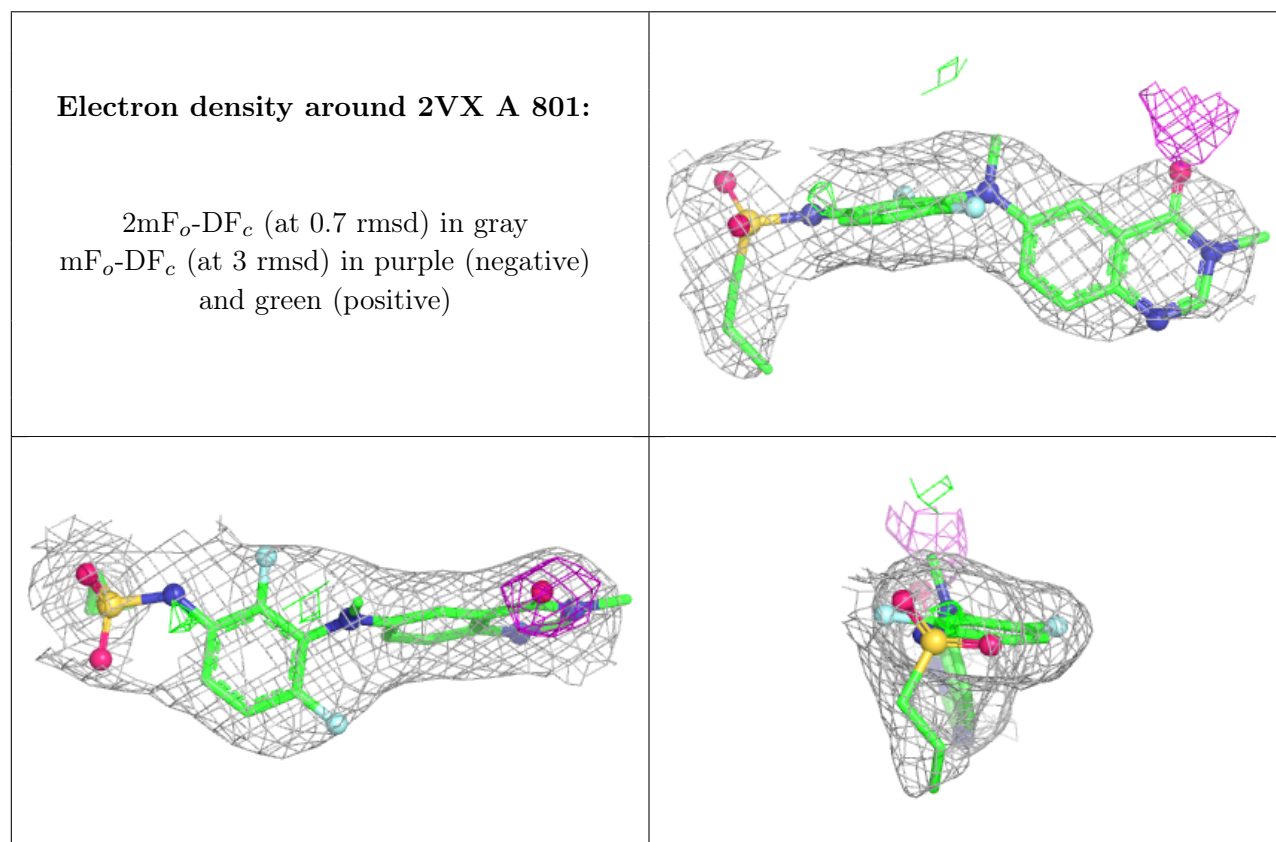
There are no oligosaccharides in this entry.

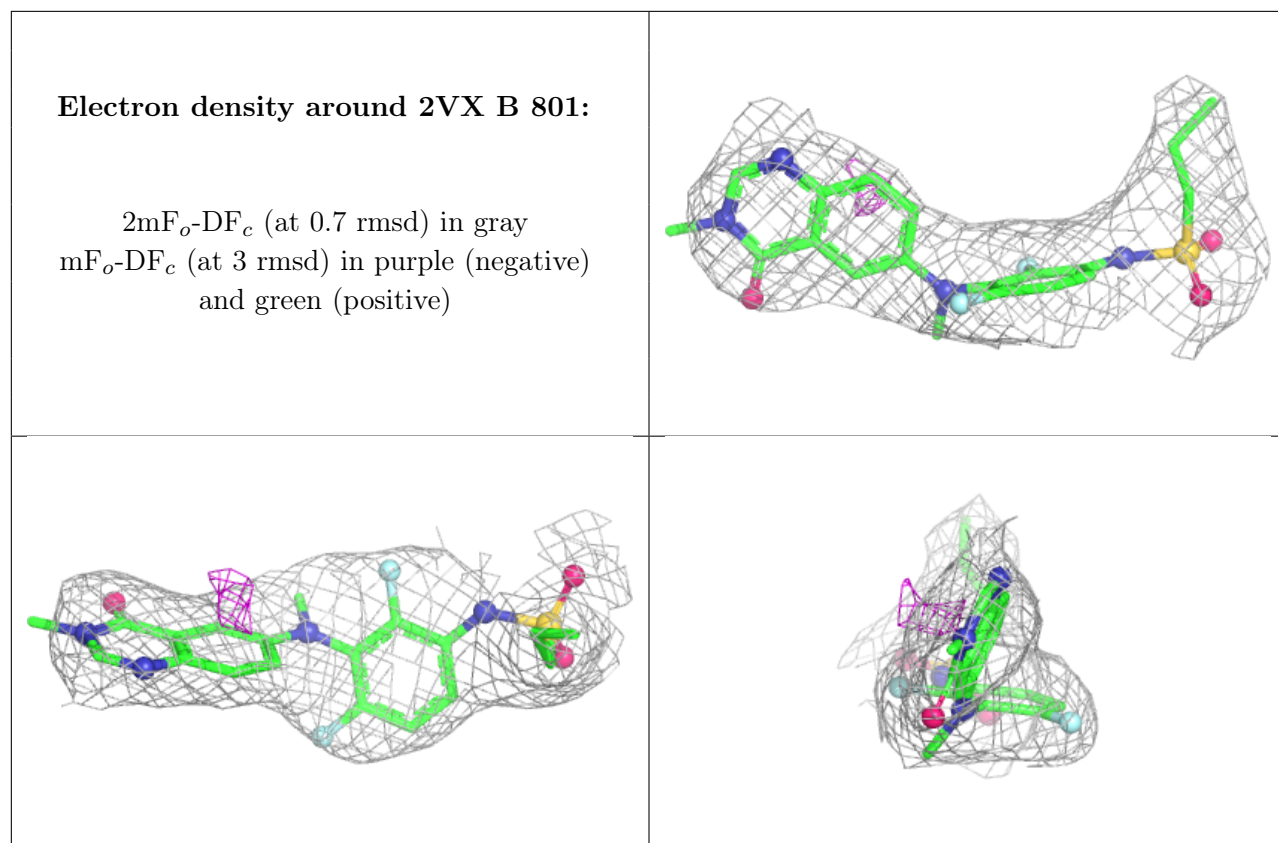
### 6.4 Ligands [i](#)

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	2VX	A	801	29/29	0.96	0.09	51,57,64,66	0
2	2VX	B	801	29/29	0.97	0.09	47,56,67,70	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.





## 6.5 Other polymers [i](#)

There are no such residues in this entry.