



wwPDB X-ray Structure Validation Summary Report ⓘ

May 3, 2025 – 10:42 AM EDT

PDB ID : 2FUM / pdb_00002fum
Title : Catalytic domain of protein kinase PknB from Mycobacterium tuberculosis in complex with mitoxantrone
Authors : Wehenkel, A.; Alzari, P.M.
Deposited on : 2006-01-27
Resolution : 2.89 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 Phenix2.0rc1
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.43.1

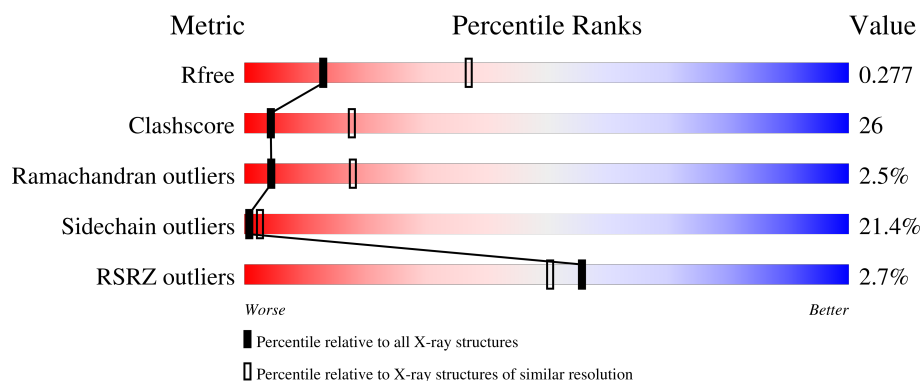
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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}	164625	2335 (2.90-2.90)
Clashscore	180529	2564 (2.90-2.90)
Ramachandran outliers	177936	2514 (2.90-2.90)
Sidechain outliers	177891	2516 (2.90-2.90)
RSRZ outliers	164620	2337 (2.90-2.90)

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	299	<div> <div>2%</div> <div> <div>43%</div> <div>34%</div> <div>9%</div> <div>12%</div> </div> </div>
1	B	299	<div> <div>4%</div> <div> <div>43%</div> <div>32%</div> <div>11%</div> <div>12%</div> </div> </div>
1	C	299	<div> <div>0%</div> <div> <div>39%</div> <div>36%</div> <div>9%</div> <div>15%</div> </div> </div>
1	D	299	<div> <div>3%</div> <div> <div>43%</div> <div>34%</div> <div>9%</div> <div>14%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7998 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 Probable serine/threonine-protein kinase pknB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			1994	1249	354	383	8			
1	B	262	Total	C	N	O	S	0	0	0
			1991	1247	357	379	8			
1	C	254	Total	C	N	O	S	0	0	0
			1931	1211	340	372	8			
1	D	258	Total	C	N	O	S	0	0	0
			1954	1225	347	374	8			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	cloning artifact	UNP P0A5S4
A	-18	GLY	-	cloning artifact	UNP P0A5S4
A	-17	SER	-	cloning artifact	UNP P0A5S4
A	-16	SER	-	cloning artifact	UNP P0A5S4
A	-15	HIS	-	cloning artifact	UNP P0A5S4
A	-14	HIS	-	cloning artifact	UNP P0A5S4
A	-13	HIS	-	cloning artifact	UNP P0A5S4
A	-12	HIS	-	cloning artifact	UNP P0A5S4
A	-11	HIS	-	cloning artifact	UNP P0A5S4
A	-10	HIS	-	cloning artifact	UNP P0A5S4
A	-9	SER	-	cloning artifact	UNP P0A5S4
A	-8	SER	-	cloning artifact	UNP P0A5S4
A	-7	GLY	-	cloning artifact	UNP P0A5S4
A	-6	LEU	-	cloning artifact	UNP P0A5S4
A	-5	VAL	-	cloning artifact	UNP P0A5S4
A	-4	PRO	-	cloning artifact	UNP P0A5S4
A	-3	ARG	-	cloning artifact	UNP P0A5S4
A	-2	GLY	-	cloning artifact	UNP P0A5S4
A	-1	SER	-	cloning artifact	UNP P0A5S4
A	0	HIS	-	cloning artifact	UNP P0A5S4
B	-19	MET	-	cloning artifact	UNP P0A5S4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	cloning artifact	UNP P0A5S4
B	-17	SER	-	cloning artifact	UNP P0A5S4
B	-16	SER	-	cloning artifact	UNP P0A5S4
B	-15	HIS	-	cloning artifact	UNP P0A5S4
B	-14	HIS	-	cloning artifact	UNP P0A5S4
B	-13	HIS	-	cloning artifact	UNP P0A5S4
B	-12	HIS	-	cloning artifact	UNP P0A5S4
B	-11	HIS	-	cloning artifact	UNP P0A5S4
B	-10	HIS	-	cloning artifact	UNP P0A5S4
B	-9	SER	-	cloning artifact	UNP P0A5S4
B	-8	SER	-	cloning artifact	UNP P0A5S4
B	-7	GLY	-	cloning artifact	UNP P0A5S4
B	-6	LEU	-	cloning artifact	UNP P0A5S4
B	-5	VAL	-	cloning artifact	UNP P0A5S4
B	-4	PRO	-	cloning artifact	UNP P0A5S4
B	-3	ARG	-	cloning artifact	UNP P0A5S4
B	-2	GLY	-	cloning artifact	UNP P0A5S4
B	-1	SER	-	cloning artifact	UNP P0A5S4
B	0	HIS	-	cloning artifact	UNP P0A5S4
C	-19	MET	-	cloning artifact	UNP P0A5S4
C	-18	GLY	-	cloning artifact	UNP P0A5S4
C	-17	SER	-	cloning artifact	UNP P0A5S4
C	-16	SER	-	cloning artifact	UNP P0A5S4
C	-15	HIS	-	cloning artifact	UNP P0A5S4
C	-14	HIS	-	cloning artifact	UNP P0A5S4
C	-13	HIS	-	cloning artifact	UNP P0A5S4
C	-12	HIS	-	cloning artifact	UNP P0A5S4
C	-11	HIS	-	cloning artifact	UNP P0A5S4
C	-10	HIS	-	cloning artifact	UNP P0A5S4
C	-9	SER	-	cloning artifact	UNP P0A5S4
C	-8	SER	-	cloning artifact	UNP P0A5S4
C	-7	GLY	-	cloning artifact	UNP P0A5S4
C	-6	LEU	-	cloning artifact	UNP P0A5S4
C	-5	VAL	-	cloning artifact	UNP P0A5S4
C	-4	PRO	-	cloning artifact	UNP P0A5S4
C	-3	ARG	-	cloning artifact	UNP P0A5S4
C	-2	GLY	-	cloning artifact	UNP P0A5S4
C	-1	SER	-	cloning artifact	UNP P0A5S4
C	0	HIS	-	cloning artifact	UNP P0A5S4
D	-19	MET	-	cloning artifact	UNP P0A5S4
D	-18	GLY	-	cloning artifact	UNP P0A5S4
D	-17	SER	-	cloning artifact	UNP P0A5S4

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	cloning artifact	UNP P0A5S4
D	-15	HIS	-	cloning artifact	UNP P0A5S4
D	-14	HIS	-	cloning artifact	UNP P0A5S4
D	-13	HIS	-	cloning artifact	UNP P0A5S4
D	-12	HIS	-	cloning artifact	UNP P0A5S4
D	-11	HIS	-	cloning artifact	UNP P0A5S4
D	-10	HIS	-	cloning artifact	UNP P0A5S4
D	-9	SER	-	cloning artifact	UNP P0A5S4
D	-8	SER	-	cloning artifact	UNP P0A5S4
D	-7	GLY	-	cloning artifact	UNP P0A5S4
D	-6	LEU	-	cloning artifact	UNP P0A5S4
D	-5	VAL	-	cloning artifact	UNP P0A5S4
D	-4	PRO	-	cloning artifact	UNP P0A5S4
D	-3	ARG	-	cloning artifact	UNP P0A5S4
D	-2	GLY	-	cloning artifact	UNP P0A5S4
D	-1	SER	-	cloning artifact	UNP P0A5S4
D	0	HIS	-	cloning artifact	UNP P0A5S4

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- The chemical structure is a complex molecule featuring a central naphthalene ring system. The naphthalene core is substituted with several side chains and functional groups. On the left side, there is a side chain starting with a blue nitrogen atom (NH) bonded to a green carbon atom (CAN), which is further bonded to a green carbon atom (GAL) and a red hydroxyl group (OH). This side chain is attached to the naphthalene ring at position 1. On the right side, there is a side chain starting with a blue nitrogen atom (NH) bonded to a green carbon atom (CAN), which is further bonded to a green carbon atom (GAL) and a red hydroxyl group (OH). This side chain is attached to the naphthalene ring at position 8. The naphthalene ring itself has several other substituents: a red hydroxyl group (OH) at position 2, a red hydroxyl group (OH) at position 3, a red hydroxyl group (OH) at position 4, a red hydroxyl group (OH) at position 5, a red hydroxyl group (OH) at position 6, a red hydroxyl group (OH) at position 7, a red hydroxyl group (OH) at position 9, and a red hydroxyl group (OH) at position 10. The naphthalene ring is also substituted with several other groups: a green carbon atom (CAN) at position 1, a green carbon atom (CAN) at position 8, a green carbon atom (CAN) at position 2, a green carbon atom (CAN) at position 3, a green carbon atom (CAN) at position 4, a green carbon atom (CAN) at position 5, a green carbon atom (CAN) at position 6, a green carbon atom (CAN) at position 7, a green carbon atom (CAN) at position 9, and a green carbon atom (CAN) at position 10. The naphthalene ring is also substituted with several other groups: a green carbon atom (CAN) at position 1, a green carbon atom (CAN) at position 8, a green carbon atom (CAN) at position 2, a green carbon atom (CAN) at position 3, a green carbon atom (CAN) at position 4, a green carbon atom (CAN) at position 5, a green carbon atom (CAN) at position 6, a green carbon atom (CAN) at position 7, a green carbon atom (CAN) at position 9, and a green carbon atom (CAN) at position 10.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 32	C 22	N 4	O 6	0	0
2	B	1	Total 32	C 22	N 4	O 6	0	0



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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			32	22	4	6		
2	D	1	Total	C	N	O	0	0
			32	22	4	6		

- Molecule 1: Probable serine/threonine-protein kinase pknB



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	116.86Å 116.86Å 260.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	69.84 – 2.89 69.84 – 2.89	Depositor EDS
% Data completeness (in resolution range)	68.1 (69.84-2.89) 68.1 (69.84-2.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.51 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.218 , 0.278 0.214 , 0.277	Depositor DCC
R_{free} test set	1426 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	97.6	Xtriage
Anisotropy	0.212	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 119.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7998	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.21% 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: MIX

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	0.93	2/2035 (0.1%)	1.25	13/2775 (0.5%)
1	B	1.01	2/2032 (0.1%)	1.28	18/2770 (0.6%)
1	C	0.87	1/1972 (0.1%)	1.21	13/2691 (0.5%)
1	D	0.79	1/1994 (0.1%)	1.11	8/2719 (0.3%)
All	All	0.91	6/8033 (0.1%)	1.22	52/10955 (0.5%)

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	2

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	179	THR	CA-C	10.39	1.63	1.53
1	B	22	MET	CA-C	6.77	1.61	1.52
1	D	116	ILE	CG1-CD1	6.40	1.76	1.51
1	A	179	THR	CA-CB	5.95	1.63	1.53
1	C	4	PRO	N-CA	5.66	1.52	1.46

The worst 5 of 52 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	179	THR	N-CA-C	13.96	128.94	109.11
1	B	180	ALA	N-CA-C	10.02	132.14	110.80
1	A	179	THR	N-CA-C	9.78	131.64	110.80
1	B	181	GLN	N-CA-C	8.56	121.56	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	157	PHE	N-CA-C	8.50	121.36	110.65

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	178	GLY	Peptide
1	A	180	ALA	Peptide

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	1994	0	1940	112	0
1	B	1991	0	1944	109	0
1	C	1931	0	1868	92	0
1	D	1954	0	1908	104	0
2	A	32	0	26	3	0
2	B	32	0	27	4	0
2	C	32	0	27	3	0
2	D	32	0	28	8	0
All	All	7998	0	7768	415	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 26.

The worst 5 of 415 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:D:116:ILE:CD1	1:D:116:ILE:CG1	1.76	1.63
1:B:99:THR:HG21	1:B:142:ALA:HA	1.16	1.16
1:C:189:ARG:HB2	1:C:189:ARG:NH1	1.64	1.12
1:B:180:ALA:HB2	1:B:221:PRO:HA	1.39	1.04
1:C:181:GLN:HG3	1:C:224:VAL:HG11	1.39	1.02

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	259/299 (87%)	222 (86%)	32 (12%)	5 (2%)	6	24
1	B	258/299 (86%)	223 (86%)	29 (11%)	6 (2%)	5	20
1	C	250/299 (84%)	216 (86%)	26 (10%)	8 (3%)	3	13
1	D	254/299 (85%)	217 (85%)	30 (12%)	7 (3%)	4	16
All	All	1021/1196 (85%)	878 (86%)	117 (12%)	26 (2%)	4	18

5 of 26 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	22	MET
1	A	44	ALA
1	A	180	ALA
1	B	22	MET
1	B	44	ALA

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	208/242 (86%)	159 (76%)	49 (24%)	0	1
1	B	208/242 (86%)	168 (81%)	40 (19%)	1	3
1	C	203/242 (84%)	155 (76%)	48 (24%)	0	1
1	D	205/242 (85%)	166 (81%)	39 (19%)	1	4
All	All	824/968 (85%)	648 (79%)	176 (21%)	1	2

5 of 176 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	135	ILE
1	D	27	LEU
1	C	184	SER
1	C	243	LEU
1	D	62	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 30 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	277	HIS
1	D	136	HIS
1	C	68	HIS
1	D	181	GLN
1	D	127	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](#)

4 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	MIX	D	3539	-	34,34,34	1.10	4 (11%)	46,46,46	0.79	1 (2%)
2	MIX	B	1539	-	34,34,34	1.28	4 (11%)	46,46,46	1.46	7 (15%)
2	MIX	A	539	-	34,34,34	1.49	4 (11%)	46,46,46	1.01	4 (8%)
2	MIX	C	2539	-	34,34,34	1.09	4 (11%)	46,46,46	1.02	2 (4%)

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	MIX	D	3539	-	-	9/14/30/30	0/3/3/3
2	MIX	B	1539	-	-	11/14/30/30	0/3/3/3
2	MIX	A	539	-	-	7/14/30/30	0/3/3/3
2	MIX	C	2539	-	-	6/14/30/30	0/3/3/3

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1539	MIX	CBC-CBA	-3.98	1.38	1.47
2	A	539	MIX	CBD-CBB	-3.87	1.38	1.47
2	A	539	MIX	CBE-CBA	-3.69	1.39	1.47
2	A	539	MIX	CBC-CBA	-3.64	1.39	1.47
2	A	539	MIX	CBF-CBB	-3.19	1.40	1.47

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1539	MIX	CAJ-CAZ-NAV	-4.73	113.94	121.75
2	B	1539	MIX	CBF-CAZ-NAV	4.35	126.53	121.34
2	B	1539	MIX	OAA-CBA-CBC	-2.82	116.74	121.44
2	A	539	MIX	CAJ-CAZ-NAV	-2.69	117.31	121.75
2	A	539	MIX	CAP-CAR-NAV	-2.67	106.04	111.54

There are no chirality outliers.

5 of 33 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1539	MIX	OAD-CAL-CAN-NAT
2	C	2539	MIX	OAC-CAK-CAM-NAS

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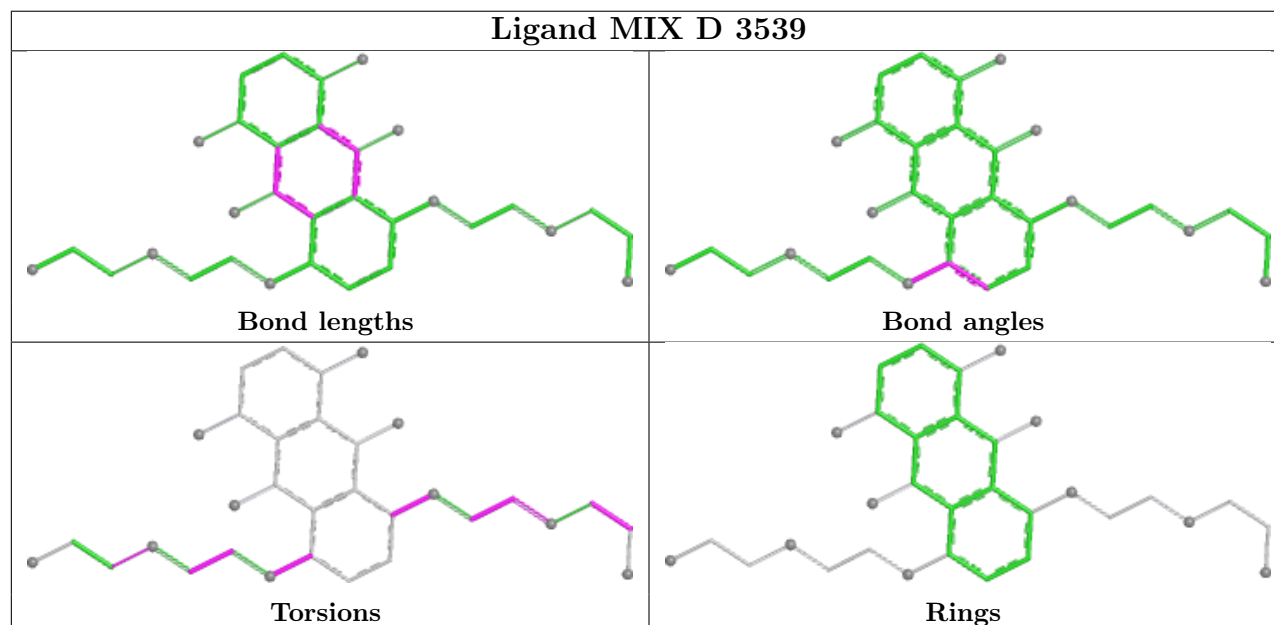
Mol	Chain	Res	Type	Atoms
2	C	2539	MIX	CAK-CAM-NAS-CAO
2	B	1539	MIX	CAJ-CAZ-NAV-CAR
2	A	539	MIX	CAJ-CAZ-NAV-CAR

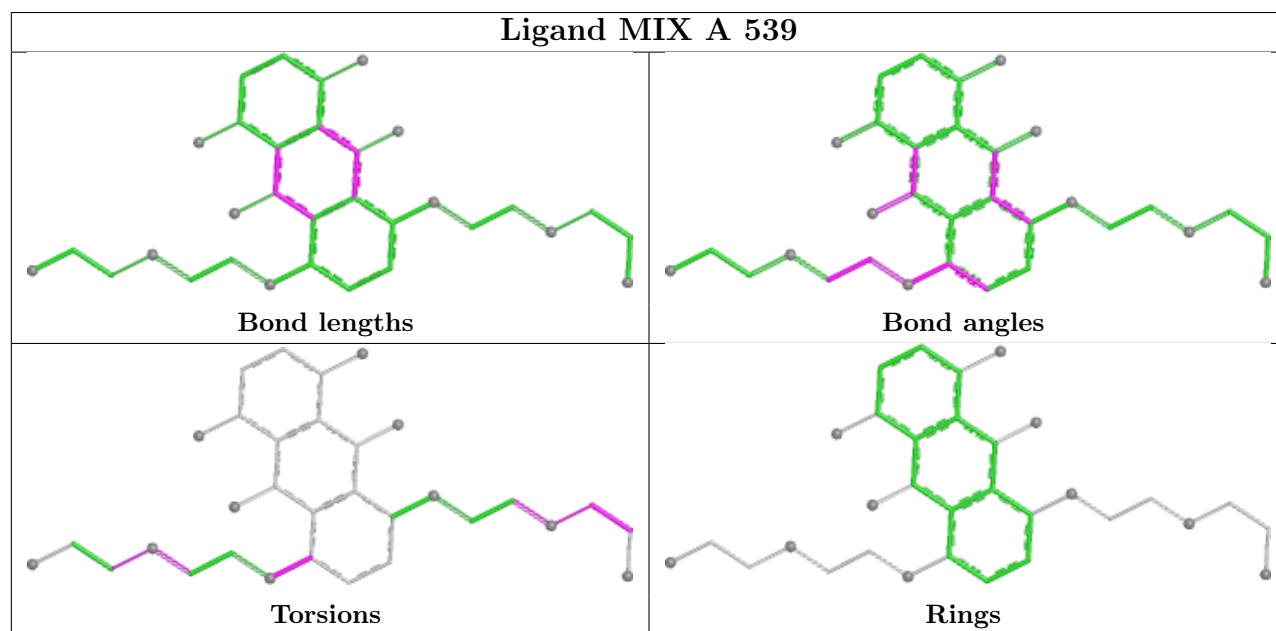
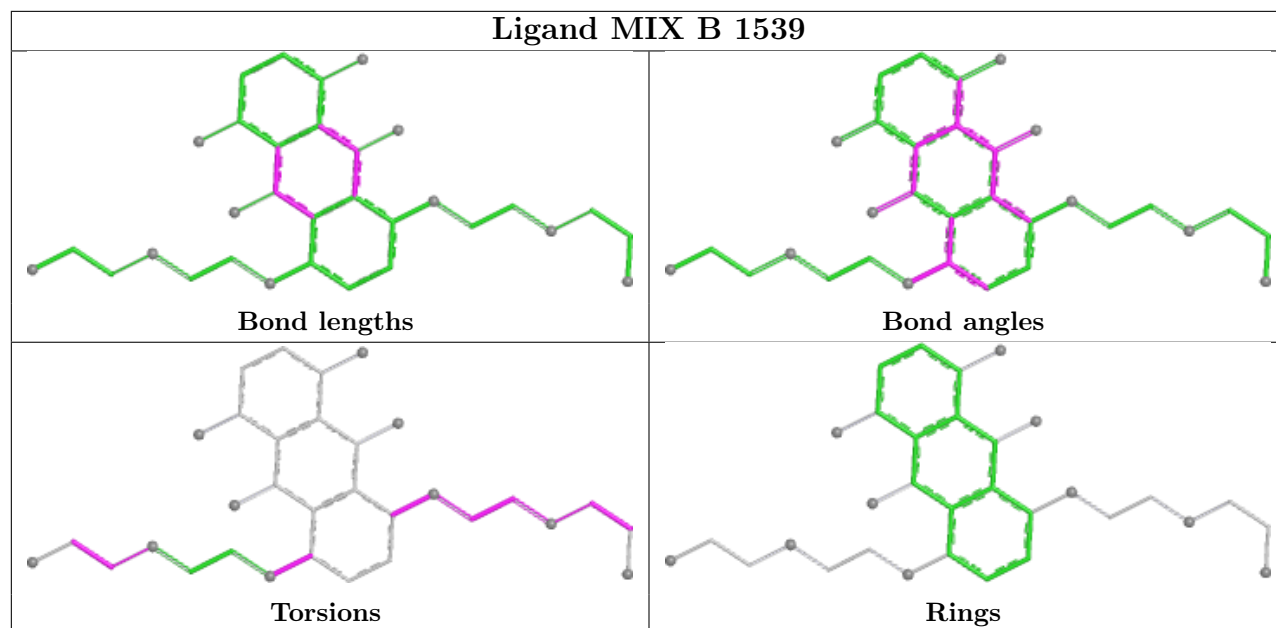
There are no ring outliers.

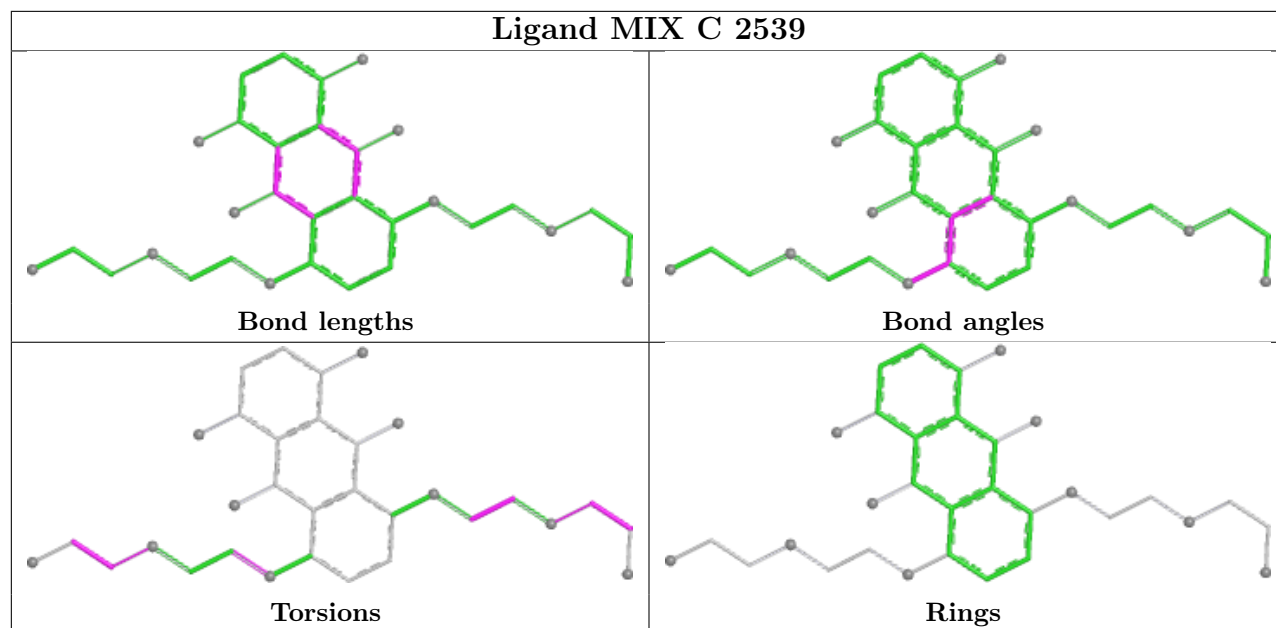
4 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	3539	MIX	8	0
2	B	1539	MIX	4	0
2	A	539	MIX	3	0
2	C	2539	MIX	3	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	263/299 (87%)	-0.03	5 (1%) 66 60	76, 79, 83, 89	0
1	B	262/299 (87%)	0.16	11 (4%) 41 35	75, 79, 83, 89	0
1	C	254/299 (84%)	0.03	4 (1%) 70 64	77, 79, 81, 86	0
1	D	258/299 (86%)	0.26	8 (3%) 51 46	76, 79, 82, 86	0
All	All	1037/1196 (86%)	0.11	28 (2%) 56 50	75, 79, 83, 89	0

The worst 5 of 28 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	189	ARG	4.5
1	D	177	ILE	4.4
1	D	7	LEU	4.4
1	D	159	ILE	4.3
1	B	84	ALA	3.7

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](#)

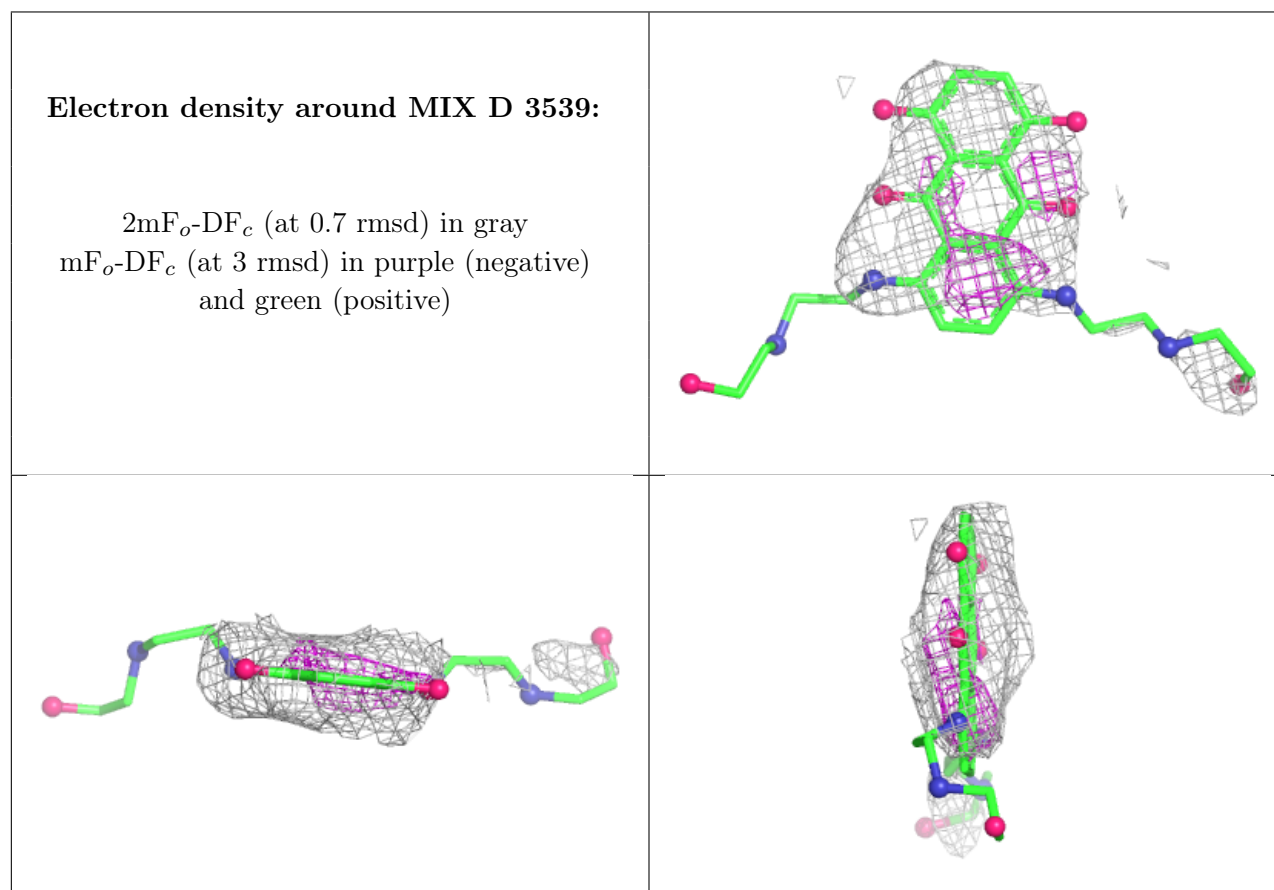
There are no monosaccharides in this entry.

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, 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.

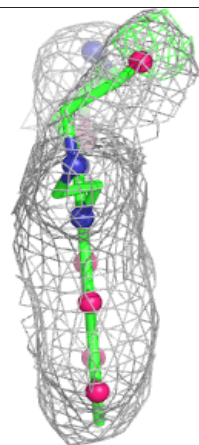
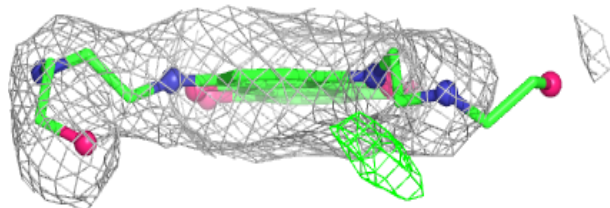
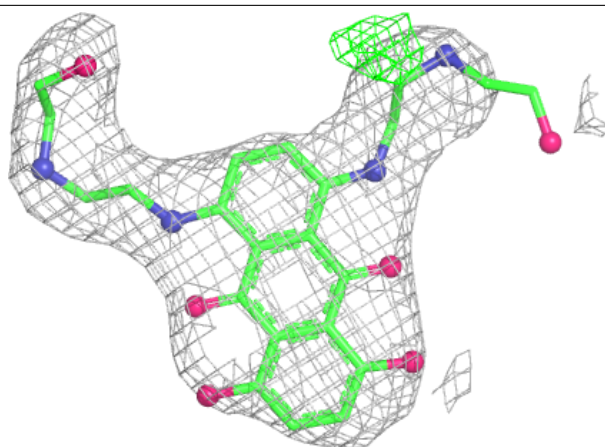
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	MIX	D	3539	32/32	0.79	0.17	158,163,164,164	0
2	MIX	C	2539	32/32	0.87	0.17	122,127,130,132	0
2	MIX	B	1539	32/32	0.90	0.11	80,81,87,87	0
2	MIX	A	539	32/32	0.91	0.12	91,91,103,105	0

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.



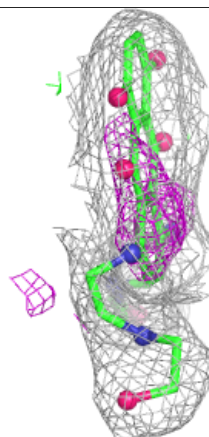
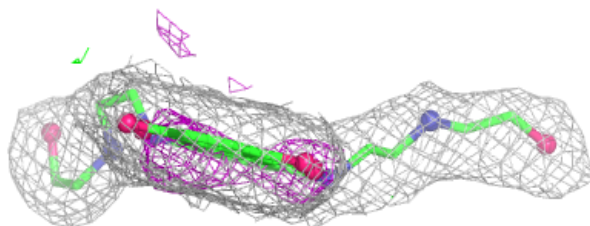
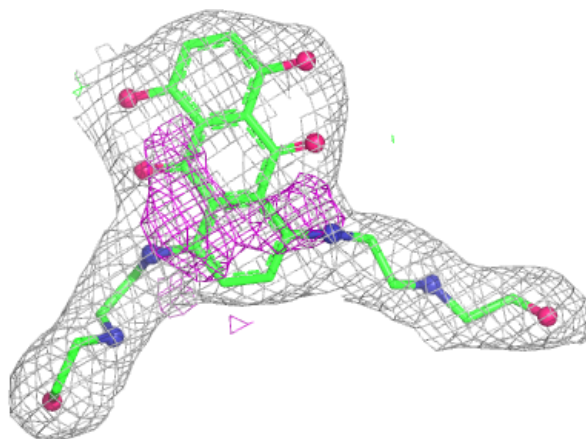
Electron density around MIX C 2539:

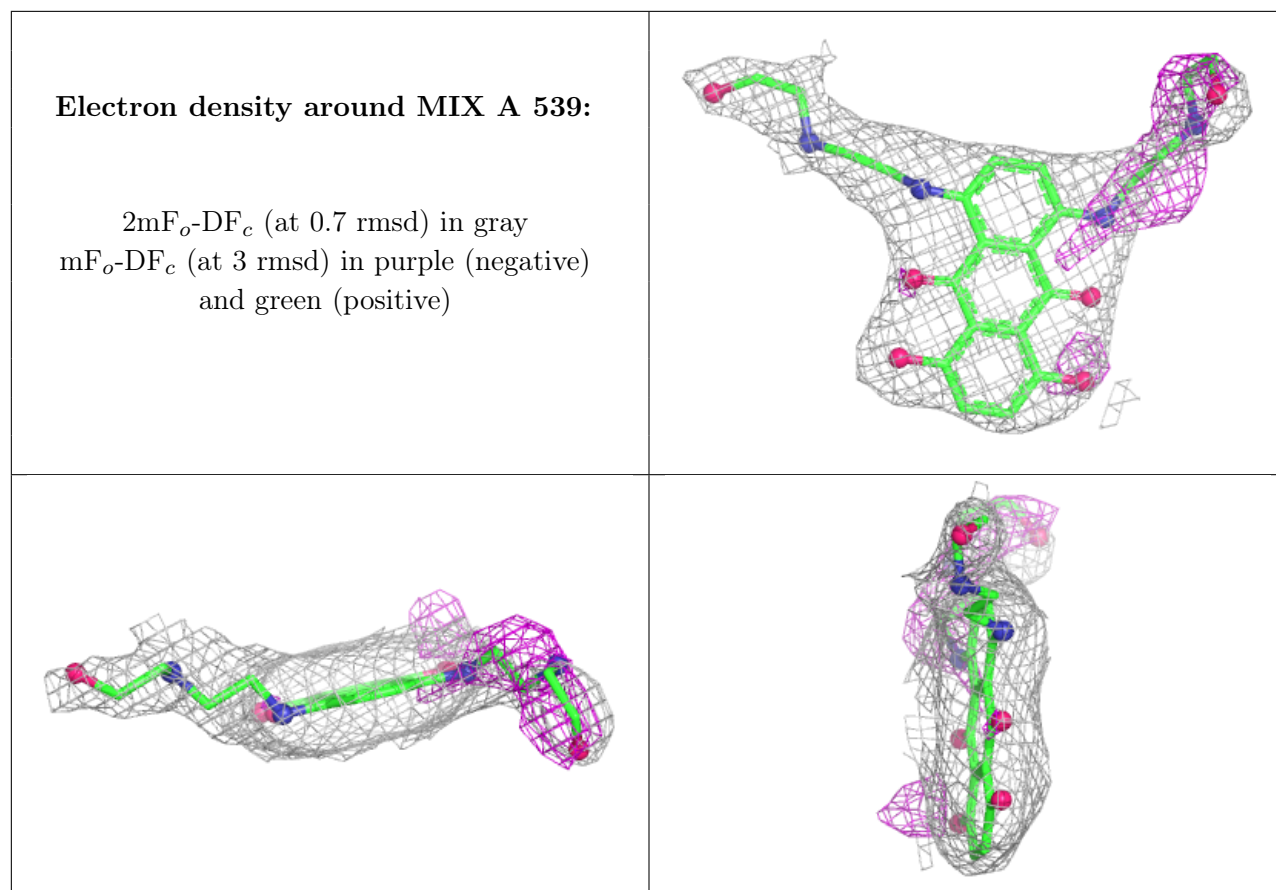
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around MIX B 1539:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.