



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 24, 2024 – 01:45 PM EDT

PDB ID : 6T8X
Title : Crystal structure of MAPKAPK2 (MK2) complexed with PF-3644022 and 5-(4-bromophenyl)-N-[4-(1-piperazinyl)phenyl]-N-(2-pyridinylmethyl)-2-furan-carboxamide
Authors : Beaumont, E.J.; Barker, J.
Deposited on : 2019-10-25
Resolution : 2.81 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:

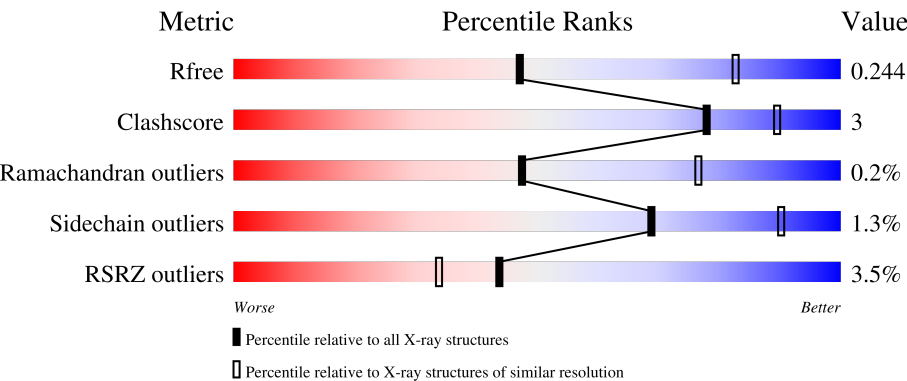
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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	3617 (2.84-2.80)
Clashscore	141614	4060 (2.84-2.80)
Ramachandran outliers	138981	3978 (2.84-2.80)
Sidechain outliers	138945	3980 (2.84-2.80)
RSRZ outliers	127900	3552 (2.84-2.80)

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	319	<div><div>2%</div><div><div></div><div>79%</div><div>9%</div><div>13%</div></div></div>
1	B	319	<div><div>3%</div><div><div></div><div>78%</div><div>8%</div><div>13%</div></div></div>
1	C	319	<div><div>3%</div><div><div></div><div>80%</div><div>7%</div><div>13%</div></div></div>
1	D	319	<div><div>4%</div><div><div></div><div>79%</div><div>9%</div><div>12%</div></div></div>

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Mol	Chain	Length	Quality of chain
1	E	319	<div><div></div><div>3%</div><div>78%</div><div>8%</div><div>13%</div></div>
1	F	319	<div><div></div><div>4%</div><div>79%</div><div>7%</div><div>13%</div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13887 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 MAP kinase-activated protein kinase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	279	Total	C	N	O	S	0	0	0
			2261	1447	387	410	17			
1	B	276	Total	C	N	O	S	0	0	0
			2233	1427	383	406	17			
1	C	276	Total	C	N	O	S	0	0	0
			2234	1429	383	405	17			
1	D	280	Total	C	N	O	S	0	0	0
			2268	1452	388	410	18			
1	E	277	Total	C	N	O	S	0	0	0
			2245	1436	384	407	18			
1	F	276	Total	C	N	O	S	0	0	0
			2239	1433	383	406	17			

There are 138 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	20	MET	-	initiating methionine	UNP P49137
A	21	GLY	-	expression tag	UNP P49137
A	22	SER	-	expression tag	UNP P49137
A	23	SER	-	expression tag	UNP P49137
A	24	HIS	-	expression tag	UNP P49137
A	25	HIS	-	expression tag	UNP P49137
A	26	HIS	-	expression tag	UNP P49137
A	27	HIS	-	expression tag	UNP P49137
A	28	HIS	-	expression tag	UNP P49137
A	29	HIS	-	expression tag	UNP P49137
A	30	SER	-	expression tag	UNP P49137
A	31	SER	-	expression tag	UNP P49137
A	32	GLY	-	expression tag	UNP P49137
A	33	GLU	-	expression tag	UNP P49137
A	34	ASN	-	expression tag	UNP P49137
A	35	LEU	-	expression tag	UNP P49137
A	36	TYR	-	expression tag	UNP P49137

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Chain	Residue	Modelled	Actual	Comment	Reference
A	37	PHE	-	expression tag	UNP P49137
A	38	GLN	-	expression tag	UNP P49137
A	39	GLY	-	expression tag	UNP P49137
A	40	SER	-	expression tag	UNP P49137
A	222	GLU	THR	engineered mutation	UNP P49137
A	334	GLU	THR	engineered mutation	UNP P49137
B	20	MET	-	initiating methionine	UNP P49137
B	21	GLY	-	expression tag	UNP P49137
B	22	SER	-	expression tag	UNP P49137
B	23	SER	-	expression tag	UNP P49137
B	24	HIS	-	expression tag	UNP P49137
B	25	HIS	-	expression tag	UNP P49137
B	26	HIS	-	expression tag	UNP P49137
B	27	HIS	-	expression tag	UNP P49137
B	28	HIS	-	expression tag	UNP P49137
B	29	HIS	-	expression tag	UNP P49137
B	30	SER	-	expression tag	UNP P49137
B	31	SER	-	expression tag	UNP P49137
B	32	GLY	-	expression tag	UNP P49137
B	33	GLU	-	expression tag	UNP P49137
B	34	ASN	-	expression tag	UNP P49137
B	35	LEU	-	expression tag	UNP P49137
B	36	TYR	-	expression tag	UNP P49137
B	37	PHE	-	expression tag	UNP P49137
B	38	GLN	-	expression tag	UNP P49137
B	39	GLY	-	expression tag	UNP P49137
B	40	SER	-	expression tag	UNP P49137
B	222	GLU	THR	engineered mutation	UNP P49137
B	334	GLU	THR	engineered mutation	UNP P49137
C	20	MET	-	initiating methionine	UNP P49137
C	21	GLY	-	expression tag	UNP P49137
C	22	SER	-	expression tag	UNP P49137
C	23	SER	-	expression tag	UNP P49137
C	24	HIS	-	expression tag	UNP P49137
C	25	HIS	-	expression tag	UNP P49137
C	26	HIS	-	expression tag	UNP P49137
C	27	HIS	-	expression tag	UNP P49137
C	28	HIS	-	expression tag	UNP P49137
C	29	HIS	-	expression tag	UNP P49137
C	30	SER	-	expression tag	UNP P49137
C	31	SER	-	expression tag	UNP P49137
C	32	GLY	-	expression tag	UNP P49137

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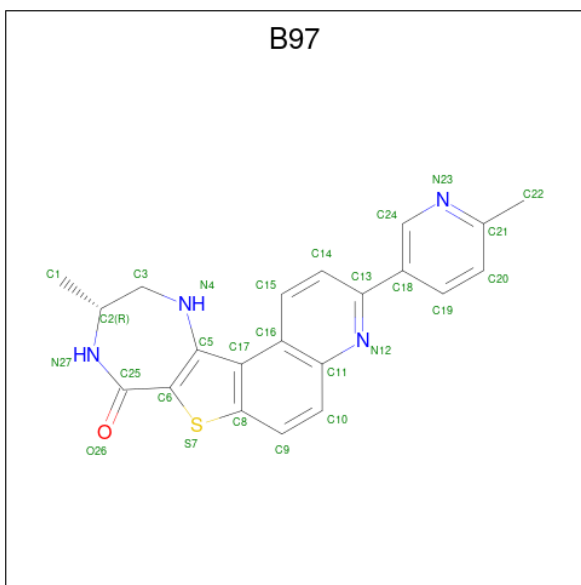
Chain	Residue	Modelled	Actual	Comment	Reference
C	33	GLU	-	expression tag	UNP P49137
C	34	ASN	-	expression tag	UNP P49137
C	35	LEU	-	expression tag	UNP P49137
C	36	TYR	-	expression tag	UNP P49137
C	37	PHE	-	expression tag	UNP P49137
C	38	GLN	-	expression tag	UNP P49137
C	39	GLY	-	expression tag	UNP P49137
C	40	SER	-	expression tag	UNP P49137
C	222	GLU	THR	engineered mutation	UNP P49137
C	334	GLU	THR	engineered mutation	UNP P49137
D	20	MET	-	initiating methionine	UNP P49137
D	21	GLY	-	expression tag	UNP P49137
D	22	SER	-	expression tag	UNP P49137
D	23	SER	-	expression tag	UNP P49137
D	24	HIS	-	expression tag	UNP P49137
D	25	HIS	-	expression tag	UNP P49137
D	26	HIS	-	expression tag	UNP P49137
D	27	HIS	-	expression tag	UNP P49137
D	28	HIS	-	expression tag	UNP P49137
D	29	HIS	-	expression tag	UNP P49137
D	30	SER	-	expression tag	UNP P49137
D	31	SER	-	expression tag	UNP P49137
D	32	GLY	-	expression tag	UNP P49137
D	33	GLU	-	expression tag	UNP P49137
D	34	ASN	-	expression tag	UNP P49137
D	35	LEU	-	expression tag	UNP P49137
D	36	TYR	-	expression tag	UNP P49137
D	37	PHE	-	expression tag	UNP P49137
D	38	GLN	-	expression tag	UNP P49137
D	39	GLY	-	expression tag	UNP P49137
D	40	SER	-	expression tag	UNP P49137
D	222	GLU	THR	engineered mutation	UNP P49137
D	334	GLU	THR	engineered mutation	UNP P49137
E	20	MET	-	initiating methionine	UNP P49137
E	21	GLY	-	expression tag	UNP P49137
E	22	SER	-	expression tag	UNP P49137
E	23	SER	-	expression tag	UNP P49137
E	24	HIS	-	expression tag	UNP P49137
E	25	HIS	-	expression tag	UNP P49137
E	26	HIS	-	expression tag	UNP P49137
E	27	HIS	-	expression tag	UNP P49137
E	28	HIS	-	expression tag	UNP P49137

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Chain	Residue	Modelled	Actual	Comment	Reference
E	29	HIS	-	expression tag	UNP P49137
E	30	SER	-	expression tag	UNP P49137
E	31	SER	-	expression tag	UNP P49137
E	32	GLY	-	expression tag	UNP P49137
E	33	GLU	-	expression tag	UNP P49137
E	34	ASN	-	expression tag	UNP P49137
E	35	LEU	-	expression tag	UNP P49137
E	36	TYR	-	expression tag	UNP P49137
E	37	PHE	-	expression tag	UNP P49137
E	38	GLN	-	expression tag	UNP P49137
E	39	GLY	-	expression tag	UNP P49137
E	40	SER	-	expression tag	UNP P49137
E	222	GLU	THR	engineered mutation	UNP P49137
E	334	GLU	THR	engineered mutation	UNP P49137
F	20	MET	-	initiating methionine	UNP P49137
F	21	GLY	-	expression tag	UNP P49137
F	22	SER	-	expression tag	UNP P49137
F	23	SER	-	expression tag	UNP P49137
F	24	HIS	-	expression tag	UNP P49137
F	25	HIS	-	expression tag	UNP P49137
F	26	HIS	-	expression tag	UNP P49137
F	27	HIS	-	expression tag	UNP P49137
F	28	HIS	-	expression tag	UNP P49137
F	29	HIS	-	expression tag	UNP P49137
F	30	SER	-	expression tag	UNP P49137
F	31	SER	-	expression tag	UNP P49137
F	32	GLY	-	expression tag	UNP P49137
F	33	GLU	-	expression tag	UNP P49137
F	34	ASN	-	expression tag	UNP P49137
F	35	LEU	-	expression tag	UNP P49137
F	36	TYR	-	expression tag	UNP P49137
F	37	PHE	-	expression tag	UNP P49137
F	38	GLN	-	expression tag	UNP P49137
F	39	GLY	-	expression tag	UNP P49137
F	40	SER	-	expression tag	UNP P49137
F	222	GLU	THR	engineered mutation	UNP P49137
F	334	GLU	THR	engineered mutation	UNP P49137

- Molecule 2 is (10R)-10-methyl-3-(6-methylpyridin-3-yl)-9,10,11,12-tetrahydro-8H-[1,4]diazepino[5',6':4,5]thieno[3,2-f]quinolin-8-one (three-letter code: B97) (formula: C₂₁H₁₈N₄OS).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			27	21	4	1	1		
2	B	1	Total	C	N	O	S	0	0
			27	21	4	1	1		
2	C	1	Total	C	N	O	S	0	0
			27	21	4	1	1		
2	D	1	Total	C	N	O	S	0	0
			27	21	4	1	1		
2	E	1	Total	C	N	O	S	0	0
			27	21	4	1	1		
2	F	1	Total	C	N	O	S	0	0
			27	21	4	1	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	B	1	Total	Cl	0	0
			1	1		
3	C	1	Total	Cl	0	0
			1	1		
3	D	1	Total	Cl	0	0
			1	1		
3	E	1	Total	Cl	0	0
			1	1		
3	F	1	Total	Cl	0	0
			1	1		

-
- Chemical structure of MW8, a pyridine derivative. The structure features a central pyridine ring (N18) substituted with a 4-(4-(4-bromophenyl)-2-furyl)-2-oxo-1,2,3,4-tetrahydropyridin-5-yl group and a 4-(4-(4-bromophenyl)-2-furyl)-2-oxo-1,2,3,4-tetrahydropyridin-5-yl group. The bromine atom is labeled BR1.

- Molecule 5 is water.

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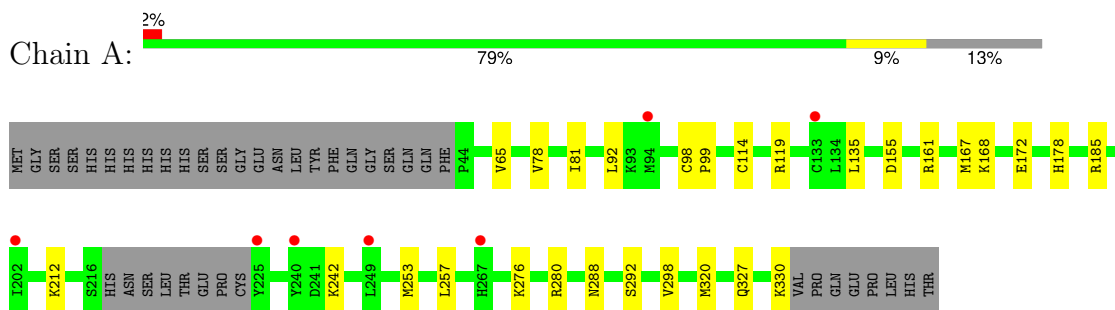
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	E	2	Total	O	0	0
			2	2		
5	F	5	Total	O	0	0
			5	5		

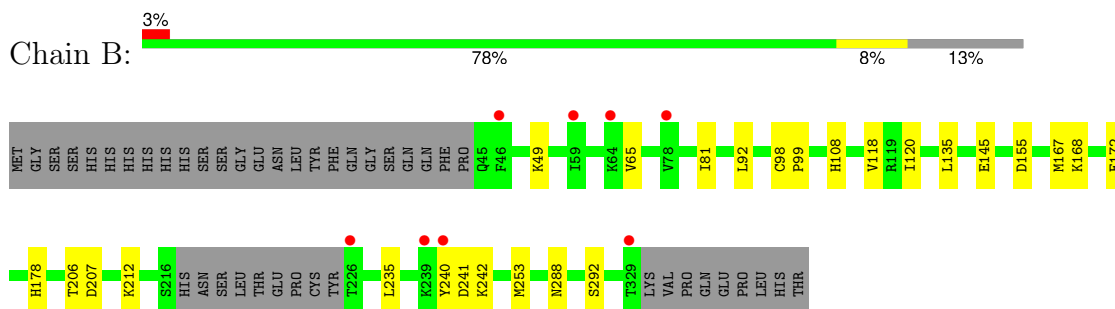
3 Residue-property plots

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.

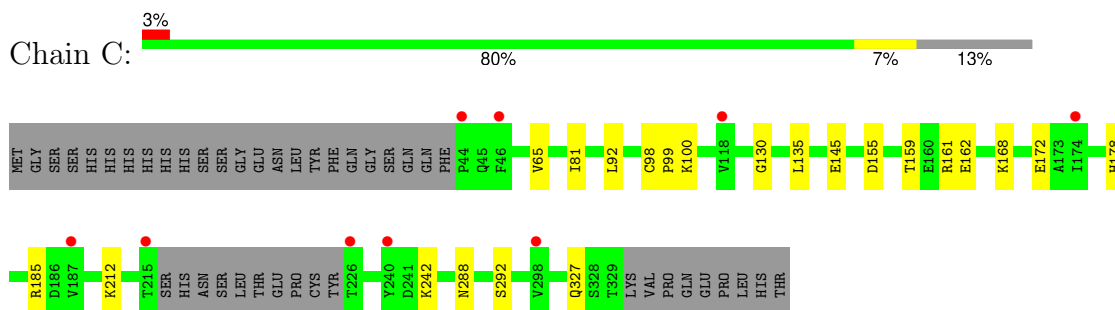
- Molecule 1: MAP kinase-activated protein kinase 2



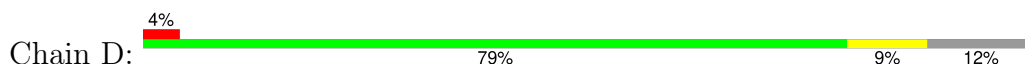
- Molecule 1: MAP kinase-activated protein kinase 2

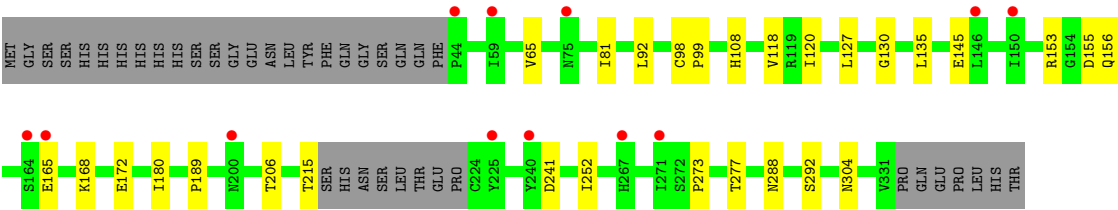


- Molecule 1: MAP kinase-activated protein kinase 2

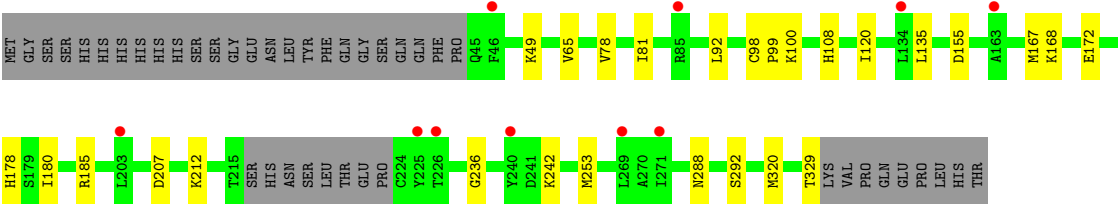
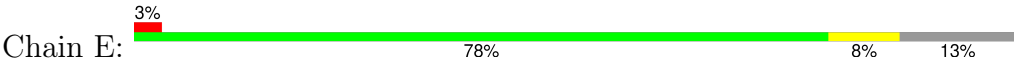


- Molecule 1: MAP kinase-activated protein kinase 2

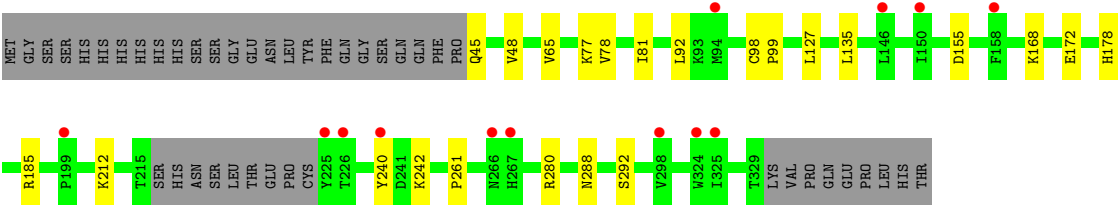
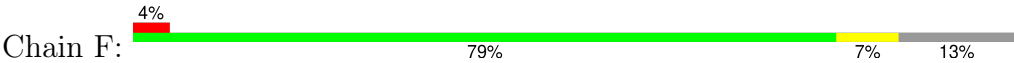




● Molecule 1: MAP kinase-activated protein kinase 2



● Molecule 1: MAP kinase-activated protein kinase 2



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	131.53Å 287.34Å 148.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	120.00 – 2.81 119.60 – 2.81	Depositor EDS
% Data completeness (in resolution range)	99.8 (120.00-2.81) 99.8 (119.60-2.81)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.54 (at 2.82Å)	Xtriage
Refinement program	BUSTER 2.11.7 (3-OCT-2019)	Depositor
R, R_{free}	0.228 , 0.247 0.230 , 0.244	Depositor DCC
R_{free} test set	3441 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	83.0	Xtriage
Anisotropy	0.594	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 70.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13887	wwPDB-VP
Average B, all atoms (Å ²)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% 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: CL, MW8, B97

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	0.44	0/2312	0.58	0/3119
1	B	0.43	0/2282	0.57	0/3079
1	C	0.43	0/2284	0.57	0/3082
1	D	0.45	0/2319	0.59	0/3129
1	E	0.41	0/2295	0.55	0/3097
1	F	0.39	0/2289	0.56	0/3089
All	All	0.43	0/13781	0.57	0/18595

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	2261	0	2280	16	0
1	B	2233	0	2250	12	0
1	C	2234	0	2253	11	0
1	D	2268	0	2289	14	0
1	E	2245	0	2259	14	0
1	F	2239	0	2254	11	0
2	A	27	0	18	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	27	0	18	1	0
2	C	27	0	18	1	0
2	D	27	0	18	1	0
2	E	27	0	18	2	0
2	F	27	0	18	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	A	34	0	0	1	0
4	B	34	0	0	0	0
4	C	34	0	0	1	0
4	D	34	0	0	0	0
4	E	34	0	0	1	0
4	F	34	0	0	1	0
5	A	4	0	0	0	0
5	B	9	0	0	0	0
5	C	9	0	0	0	0
5	D	6	0	0	0	0
5	E	2	0	0	0	0
5	F	5	0	0	0	0
All	All	13887	0	13693	81	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 3.

All (81) 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:E:236:GLY:HA3	1:F:280:ARG:HE	1.67	0.60
1:D:168:LYS:O	1:D:172:GLU:HG3	2.02	0.59
1:A:92:LEU:HD11	1:A:135:LEU:HB3	1.86	0.58
1:F:92:LEU:HD11	1:F:135:LEU:HB3	1.86	0.57
1:D:288:ASN:O	1:D:292:SER:HB2	2.05	0.57
1:A:280:ARG:HH12	1:D:215:THR:CG2	2.17	0.57
1:A:288:ASN:O	1:A:292:SER:HB2	2.05	0.56
1:B:92:LEU:HD11	1:B:135:LEU:HB3	1.87	0.56
1:D:92:LEU:HD11	1:D:135:LEU:HB3	1.87	0.56
1:E:92:LEU:HD11	1:E:135:LEU:HB3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:401:B97:HN4	2:B:401:B97:H15	1.71	0.55
2:E:401:B97:HN4	2:E:401:B97:H15	1.70	0.55
1:A:178:HIS:CE1	1:A:242:LYS:HB3	2.42	0.55
1:E:178:HIS:CE1	1:E:242:LYS:HB3	2.41	0.54
1:B:288:ASN:O	1:B:292:SER:HB2	2.07	0.54
1:E:288:ASN:O	1:E:292:SER:HB2	2.07	0.54
2:A:401:B97:HN4	2:A:401:B97:H15	1.72	0.54
1:B:212:LYS:HZ3	1:B:240:TYR:HD1	1.52	0.54
1:C:178:HIS:CE1	1:C:242:LYS:HB3	2.43	0.54
1:D:127:LEU:HB2	1:E:49:LYS:HB2	1.90	0.54
1:D:189:PRO:HD3	1:D:252:ILE:HG12	1.90	0.53
1:A:276:LYS:O	1:A:280:ARG:HG2	2.09	0.53
1:F:288:ASN:O	1:F:292:SER:HB2	2.08	0.53
1:B:178:HIS:CE1	1:B:242:LYS:HB3	2.43	0.53
1:F:168:LYS:O	1:F:172:GLU:HG3	2.09	0.53
1:A:168:LYS:O	1:A:172:GLU:HG3	2.08	0.52
1:C:288:ASN:O	1:C:292:SER:HB2	2.09	0.52
1:D:153:ARG:O	1:D:156:GLN:HB3	2.10	0.52
1:B:168:LYS:O	1:B:172:GLU:HG3	2.10	0.51
1:C:168:LYS:O	1:C:172:GLU:HG3	2.11	0.51
1:E:168:LYS:O	1:E:172:GLU:HG3	2.11	0.50
2:D:401:B97:HN4	2:D:401:B97:H15	1.76	0.50
2:C:401:B97:HN4	2:C:401:B97:H15	1.77	0.50
1:D:130:GLY:HA3	1:E:180:ILE:HB	1.92	0.49
1:B:49:LYS:HB2	1:F:127:LEU:HB2	1.95	0.48
4:E:403:MW8:C10	4:E:403:MW8:C23	2.92	0.48
1:A:185:ARG:HH11	1:A:212:LYS:HB2	1.79	0.48
1:A:172:GLU:HG2	1:A:320:MET:CE	2.44	0.47
1:E:65:VAL:HA	1:E:81:ILE:HG22	1.95	0.47
1:D:65:VAL:HA	1:D:81:ILE:HG22	1.96	0.47
1:E:185:ARG:HH11	1:E:212:LYS:HB2	1.79	0.47
1:C:98:CYS:HB2	1:C:99:PRO:HD2	1.95	0.47
1:A:167:MET:HG3	1:A:253:MET:HE2	1.97	0.47
1:B:65:VAL:HA	1:B:81:ILE:HG22	1.97	0.46
4:C:403:MW8:C10	4:C:403:MW8:C23	2.94	0.46
1:F:78:VAL:HG21	2:F:401:B97:C5	2.46	0.46
1:B:240:TYR:CD2	1:B:241:ASP:N	2.83	0.46
1:C:65:VAL:HA	1:C:81:ILE:HG22	1.96	0.46
1:F:178:HIS:CE1	1:F:242:LYS:HB3	2.51	0.45
2:F:401:B97:HN4	2:F:401:B97:H15	1.81	0.45
1:B:167:MET:HG3	1:B:253:MET:HE2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:98:CYS:HB2	1:D:99:PRO:HD2	1.98	0.45
1:F:185:ARG:HH21	1:F:212:LYS:HB2	1.81	0.45
1:F:98:CYS:HB2	1:F:99:PRO:HD2	1.98	0.45
1:A:65:VAL:HA	1:A:81:ILE:HG22	1.98	0.45
1:B:98:CYS:HB2	1:B:99:PRO:HD2	1.99	0.45
1:C:185:ARG:HH21	1:C:212:LYS:HB2	1.81	0.44
1:D:108:HIS:CG	1:D:120:ILE:HD11	2.52	0.44
1:A:98:CYS:HB2	1:A:99:PRO:HD2	1.99	0.44
1:E:167:MET:HG3	1:E:253:MET:HE2	2.00	0.43
1:D:273:PRO:O	1:D:277:THR:HG23	2.16	0.43
1:A:257:LEU:HD11	1:A:298:VAL:HG21	1.99	0.43
1:F:65:VAL:HA	1:F:81:ILE:HG22	1.99	0.43
1:F:261:PRO:HD3	4:F:403:MW8:C8	2.49	0.43
1:A:161:ARG:HD2	1:A:327:GLN:HE21	1.84	0.43
1:E:98:CYS:HB2	1:E:99:PRO:HD2	2.00	0.43
1:E:78:VAL:HG21	2:E:401:B97:C5	2.49	0.42
1:C:130:GLY:HA3	1:D:180:ILE:HB	2.00	0.42
1:C:92:LEU:HD11	1:C:135:LEU:HB3	2.01	0.42
1:A:119:ARG:HB2	1:A:119:ARG:HH11	1.83	0.42
1:C:98:CYS:SG	1:C:100:LYS:HG2	2.58	0.42
1:B:118:VAL:HB	1:B:206:THR:HG22	2.01	0.42
1:D:118:VAL:HB	1:D:206:THR:HG22	2.02	0.42
1:C:161:ARG:HD2	1:C:327:GLN:HE21	1.86	0.41
4:A:403:MW8:C10	4:A:403:MW8:C23	2.98	0.41
1:E:172:GLU:HG2	1:E:320:MET:CE	2.51	0.41
1:A:78:VAL:HG21	2:A:401:B97:C6	2.51	0.41
1:B:108:HIS:CG	1:B:120:ILE:HD11	2.56	0.40
1:E:108:HIS:CG	1:E:120:ILE:HD11	2.56	0.40
1:A:78:VAL:HG21	2:A:401:B97:C5	2.51	0.40
1:C:159:THR:OG1	1:C:162:GLU:HG2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	275/319 (86%)	265 (96%)	10 (4%)	0	100	100
1	B	272/319 (85%)	263 (97%)	7 (3%)	2 (1%)	22	51
1	C	272/319 (85%)	259 (95%)	13 (5%)	0	100	100
1	D	276/319 (86%)	268 (97%)	8 (3%)	0	100	100
1	E	273/319 (86%)	261 (96%)	11 (4%)	1 (0%)	34	64
1	F	272/319 (85%)	259 (95%)	12 (4%)	1 (0%)	34	64
All	All	1640/1914 (86%)	1575 (96%)	61 (4%)	4 (0%)	47	76

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	240	TYR
1	E	207	ASP
1	B	207	ASP
1	B	235	LEU

5.3.2 Protein sidechains ⓘ

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	250/287 (87%)	247 (99%)	3 (1%)	71	91
1	B	247/287 (86%)	245 (99%)	2 (1%)	81	94
1	C	247/287 (86%)	245 (99%)	2 (1%)	81	94
1	D	251/287 (88%)	246 (98%)	5 (2%)	55	83
1	E	248/287 (86%)	245 (99%)	3 (1%)	71	91
1	F	247/287 (86%)	243 (98%)	4 (2%)	62	87
All	All	1490/1722 (86%)	1471 (99%)	19 (1%)	69	90

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114	CYS
1	A	155	ASP
1	A	330	LYS
1	B	145	GLU
1	B	155	ASP
1	C	145	GLU
1	C	155	ASP
1	D	145	GLU
1	D	155	ASP
1	D	165	GLU
1	D	241	ASP
1	D	304	ASN
1	E	100	LYS
1	E	155	ASP
1	E	329	THR
1	F	45	GLN
1	F	48	VAL
1	F	77	LYS
1	F	155	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	96	GLN
1	A	156	GLN
1	A	181	ASN
1	A	327	GLN
1	C	327	GLN
1	D	156	GLN
1	E	156	GLN
1	F	156	GLN
1	F	327	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 6 are monoatomic - leaving 12 for Mogul analysis.

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
4	MW8	B	403	-	36,38,38	1.03	1 (2%)	44,52,52	1.37	4 (9%)
2	B97	B	401	-	25,31,31	1.92	4 (16%)	26,46,46	1.18	2 (7%)
2	B97	A	401	-	25,31,31	1.96	4 (16%)	26,46,46	1.19	2 (7%)
2	B97	E	401	-	25,31,31	1.78	4 (16%)	26,46,46	1.14	1 (3%)
2	B97	C	401	-	25,31,31	1.84	5 (20%)	26,46,46	1.13	1 (3%)
4	MW8	E	403	-	36,38,38	1.11	3 (8%)	44,52,52	1.21	4 (9%)
4	MW8	A	403	-	36,38,38	1.08	1 (2%)	44,52,52	1.31	5 (11%)
2	B97	F	401	-	25,31,31	2.03	4 (16%)	26,46,46	1.02	2 (7%)
4	MW8	C	403	-	36,38,38	1.08	2 (5%)	44,52,52	1.32	7 (15%)
4	MW8	D	403	-	36,38,38	1.12	2 (5%)	44,52,52	1.31	5 (11%)
4	MW8	F	403	-	36,38,38	1.01	2 (5%)	44,52,52	1.03	4 (9%)
2	B97	D	401	-	25,31,31	2.04	4 (16%)	26,46,46	1.13	1 (3%)

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
4	MW8	B	403	-	-	2/18/32/32	0/5/5/5
2	B97	B	401	-	-	0/0/17/17	0/5/5/5
2	B97	A	401	-	-	0/0/17/17	0/5/5/5
2	B97	E	401	-	-	0/0/17/17	0/5/5/5
2	B97	C	401	-	-	0/0/17/17	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MW8	E	403	-	-	0/18/32/32	0/5/5/5
4	MW8	A	403	-	-	0/18/32/32	0/5/5/5
2	B97	F	401	-	-	0/0/17/17	0/5/5/5
4	MW8	C	403	-	-	0/18/32/32	0/5/5/5
4	MW8	D	403	-	-	1/18/32/32	0/5/5/5
4	MW8	F	403	-	-	2/18/32/32	0/5/5/5
2	B97	D	401	-	-	0/0/17/17	0/5/5/5

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	401	B97	C25-N27	7.29	1.39	1.34
2	D	401	B97	C25-N27	7.01	1.39	1.34
2	A	401	B97	C25-N27	6.66	1.39	1.34
2	B	401	B97	C25-N27	6.60	1.39	1.34
2	E	401	B97	C25-N27	5.72	1.38	1.34
2	C	401	B97	C25-N27	5.32	1.38	1.34
2	C	401	B97	C5-C17	4.83	1.48	1.41
2	D	401	B97	C5-C17	4.75	1.48	1.41
4	A	403	MW8	C11-C13	-4.68	1.42	1.49
4	D	403	MW8	C11-C13	-4.54	1.43	1.49
2	F	401	B97	C5-C17	4.47	1.47	1.41
2	A	401	B97	C5-C17	4.34	1.47	1.41
2	B	401	B97	C5-C17	4.24	1.47	1.41
2	E	401	B97	C5-C17	4.12	1.47	1.41
4	C	403	MW8	C11-C13	-4.01	1.43	1.49
4	F	403	MW8	C11-C13	-3.65	1.44	1.49
4	B	403	MW8	C11-C13	-3.60	1.44	1.49
4	E	403	MW8	C11-C13	-3.36	1.44	1.49
2	A	401	B97	C14-C13	-2.62	1.37	1.41
4	C	403	MW8	C5-C8	-2.61	1.42	1.46
2	C	401	B97	C2-N27	-2.52	1.44	1.47
2	D	401	B97	C10-C11	-2.48	1.37	1.41
2	C	401	B97	C10-C11	-2.48	1.37	1.41
4	F	403	MW8	C13-N15	2.46	1.40	1.36
2	A	401	B97	C10-C11	-2.44	1.37	1.41
2	D	401	B97	C14-C13	-2.40	1.37	1.41
4	E	403	MW8	C13-N15	2.38	1.40	1.36
2	E	401	B97	C14-C13	-2.33	1.38	1.41
2	B	401	B97	C10-C11	-2.33	1.38	1.41
2	E	401	B97	C10-C11	-2.32	1.38	1.41
2	C	401	B97	C14-C13	-2.20	1.38	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	B97	C14-C13	-2.19	1.38	1.41
4	D	403	MW8	C5-C8	-2.14	1.42	1.46
2	F	401	B97	C10-C11	-2.11	1.38	1.41
2	F	401	B97	C14-C13	-2.07	1.38	1.41
4	E	403	MW8	C5-C8	-2.02	1.42	1.46

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	403	MW8	C9-C8-C5	5.10	133.54	128.77
4	D	403	MW8	C9-C8-C5	4.52	133.00	128.77
2	A	401	B97	C17-C16-C11	3.96	121.27	117.96
4	C	403	MW8	C17-C16-N15	-3.89	106.79	112.73
2	D	401	B97	C17-C16-C11	3.89	121.21	117.96
4	D	403	MW8	C34-N29-C30	3.79	120.08	111.57
4	B	403	MW8	C34-N29-C30	3.63	119.74	111.57
4	E	403	MW8	C9-C8-C5	3.61	132.15	128.77
4	A	403	MW8	C9-C8-C5	3.55	132.09	128.77
4	A	403	MW8	C23-N15-C13	3.40	127.79	122.73
2	C	401	B97	C17-C16-C11	3.39	120.79	117.96
4	F	403	MW8	C34-N29-C30	3.24	118.86	111.57
4	C	403	MW8	C34-N29-C30	3.14	118.63	111.57
4	B	403	MW8	C17-C16-N15	-3.08	108.03	112.73
4	E	403	MW8	C34-N29-C30	3.06	118.46	111.57
4	A	403	MW8	C34-N29-C30	2.80	117.87	111.57
4	C	403	MW8	C9-C8-C5	2.65	131.25	128.77
2	B	401	B97	C16-C11-N12	-2.65	122.02	123.56
2	F	401	B97	C10-C11-N12	-2.56	115.82	119.10
4	A	403	MW8	C4-C5-C8	-2.52	117.30	120.44
4	F	403	MW8	C23-N15-C13	2.46	126.39	122.73
4	D	403	MW8	C4-C5-C8	-2.45	117.39	120.44
4	E	403	MW8	C27-C26-N29	-2.40	118.06	121.39
2	E	401	B97	C17-C16-C11	2.37	119.94	117.96
2	A	401	B97	C9-C8-C17	-2.32	122.19	123.41
4	F	403	MW8	C9-C8-C5	2.30	130.93	128.77
4	D	403	MW8	BR1-C2-C7	2.30	122.43	119.30
4	C	403	MW8	C25-C26-N29	-2.29	118.21	121.39
4	D	403	MW8	C17-C16-N15	-2.20	109.38	112.73
4	C	403	MW8	C6-C5-C4	2.19	121.60	117.68
4	F	403	MW8	C4-C5-C8	-2.18	117.73	120.44
2	B	401	B97	C15-C14-C13	-2.16	117.07	120.73
4	C	403	MW8	C7-C6-C5	-2.14	118.37	121.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	403	MW8	C28-C23-N15	-2.13	117.12	120.18
4	C	403	MW8	C4-C5-C8	-2.04	117.89	120.44
4	A	403	MW8	C33-N32-C31	2.04	116.18	110.40
4	B	403	MW8	C28-C23-N15	2.03	123.10	120.18
2	F	401	B97	C17-C16-C11	2.02	119.65	117.96

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	F	403	MW8	C27-C26-N29-C34
4	F	403	MW8	C25-C26-N29-C34
4	B	403	MW8	C27-C26-N29-C34
4	B	403	MW8	C25-C26-N29-C34
4	D	403	MW8	C27-C26-N29-C30

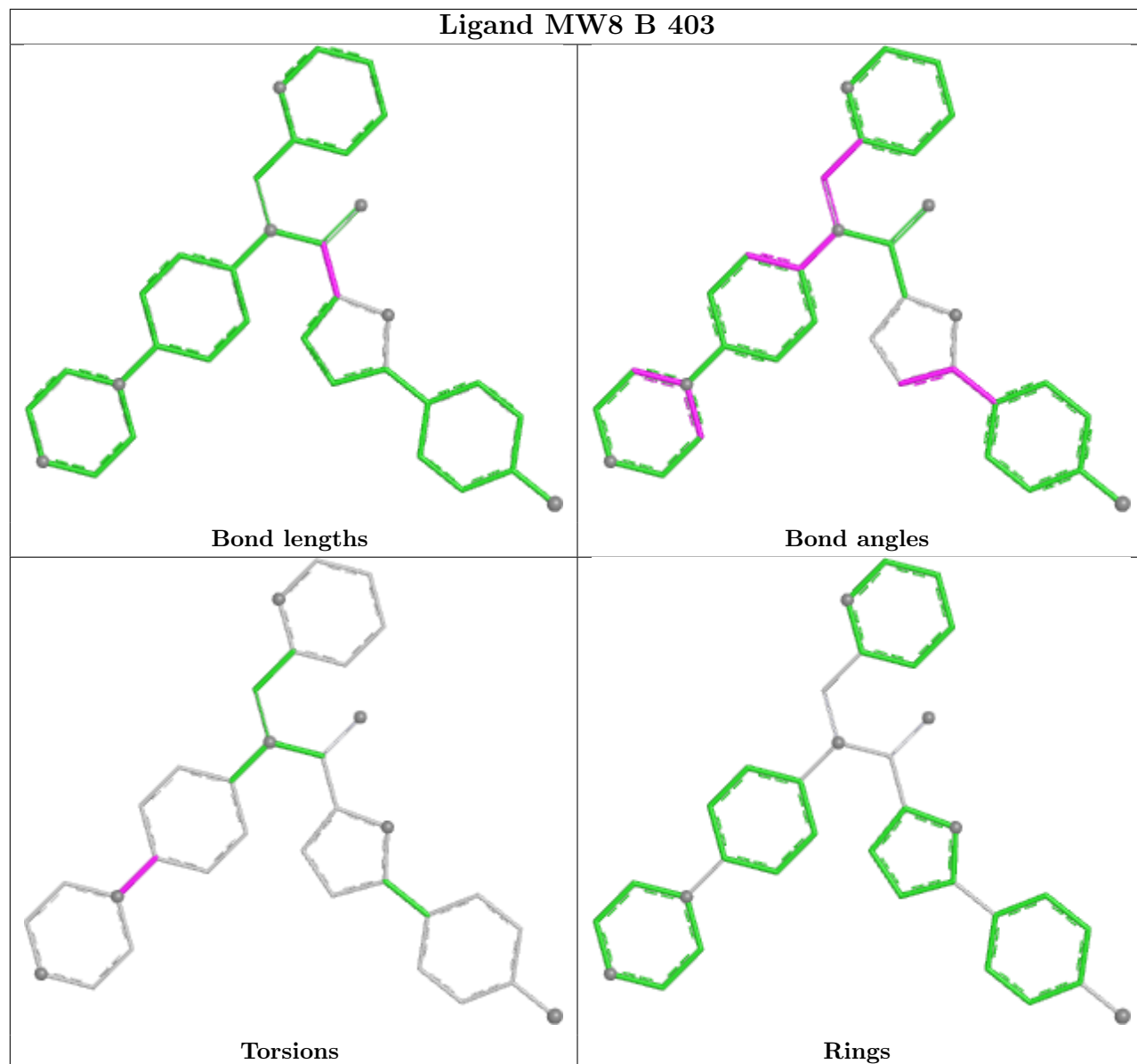
There are no ring outliers.

10 monomers are involved in 14 short contacts:

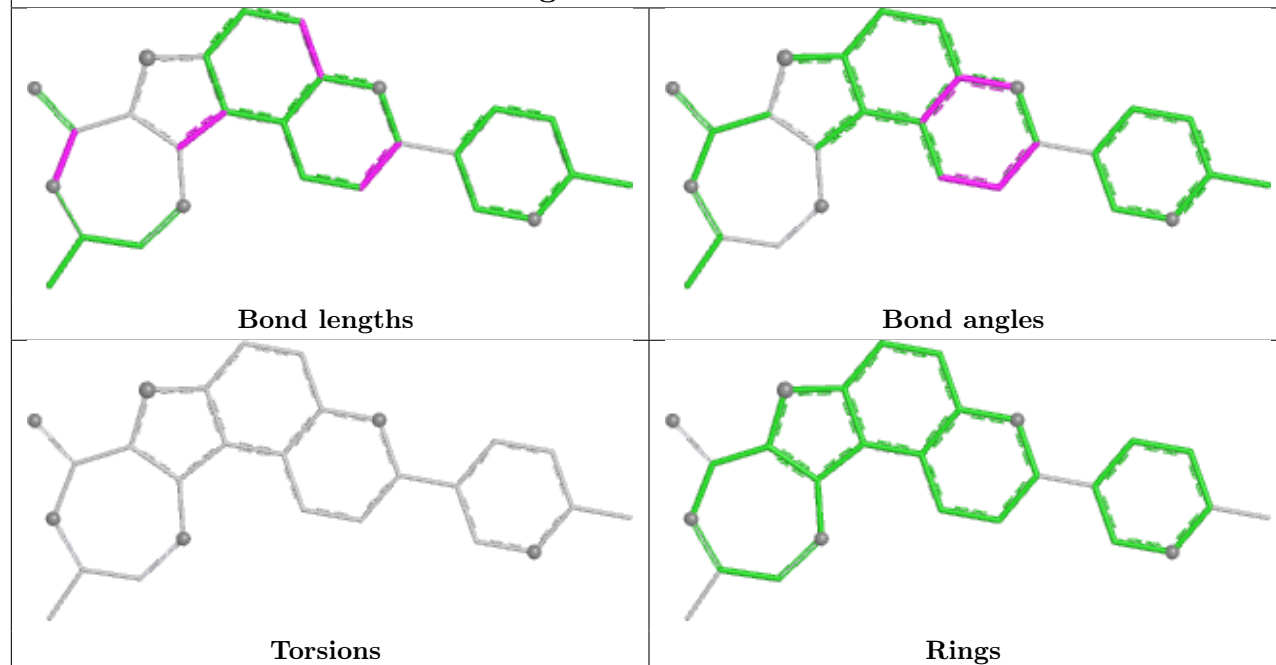
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	B97	1	0
2	A	401	B97	3	0
2	E	401	B97	2	0
2	C	401	B97	1	0
4	E	403	MW8	1	0
4	A	403	MW8	1	0
2	F	401	B97	2	0
4	C	403	MW8	1	0
4	F	403	MW8	1	0
2	D	401	B97	1	0

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

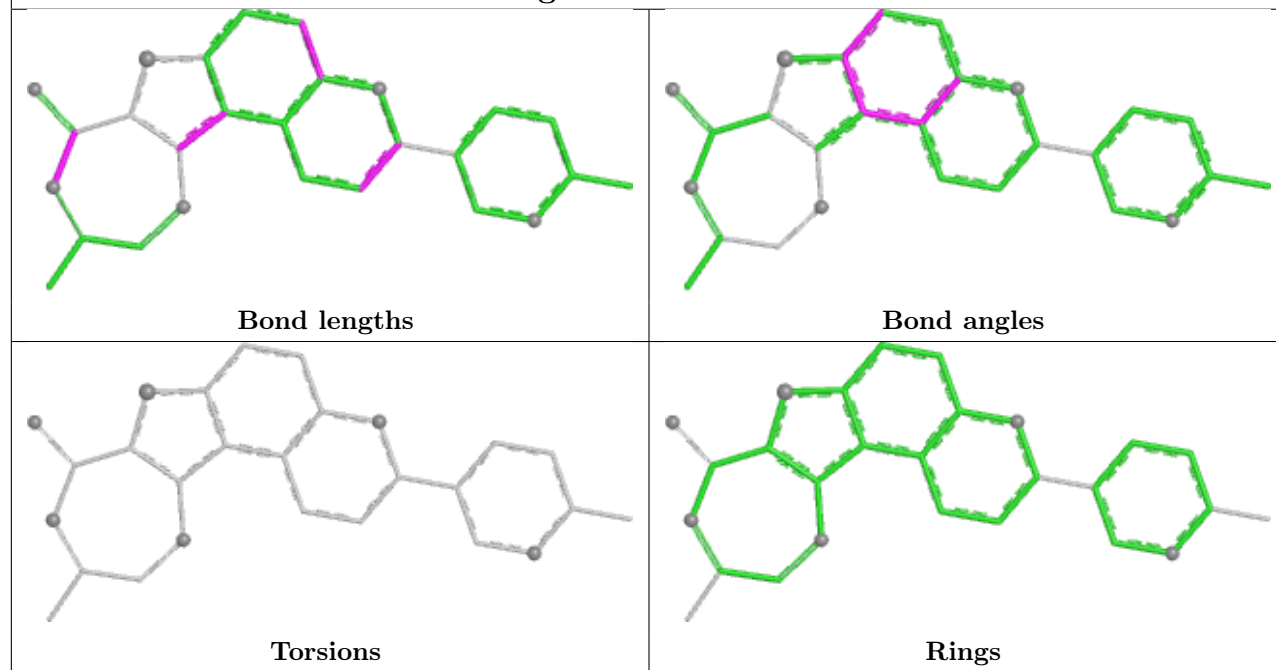
equivalents in the CSD to analyse the geometry.



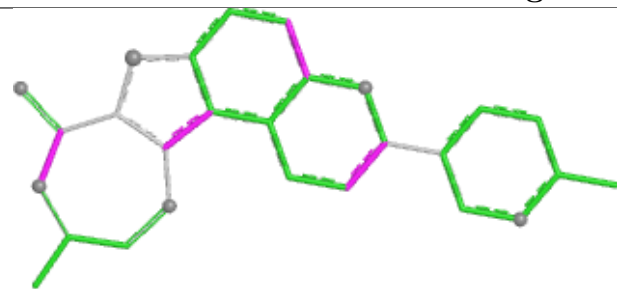
Ligand B97 B 401



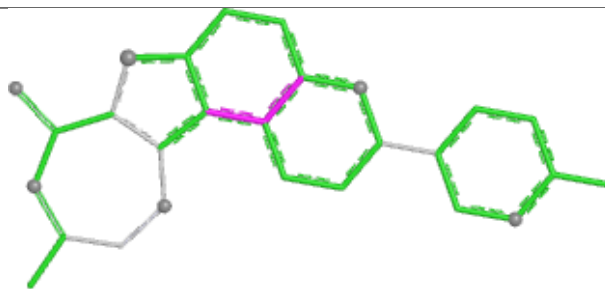
Ligand B97 A 401



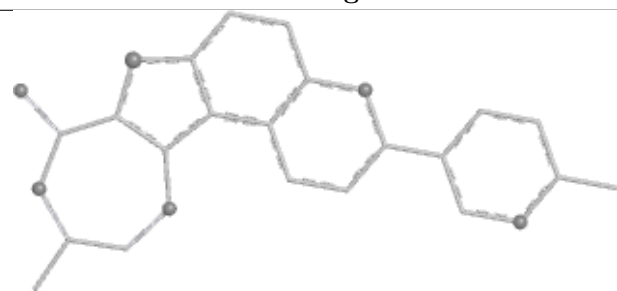
Ligand B97 E 401



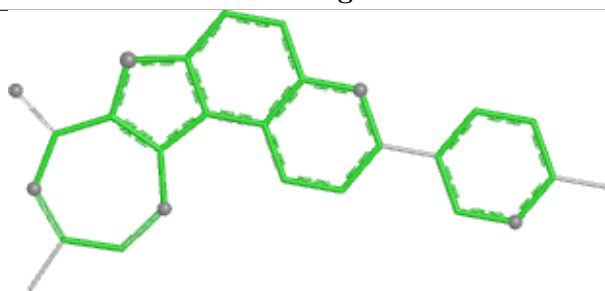
Bond lengths



Bond angles

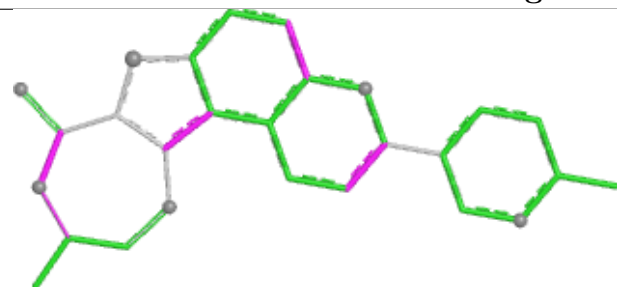


Torsions

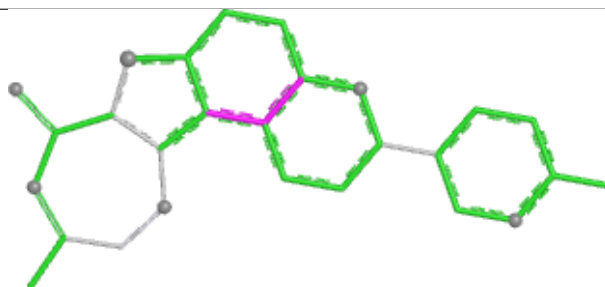


Rings

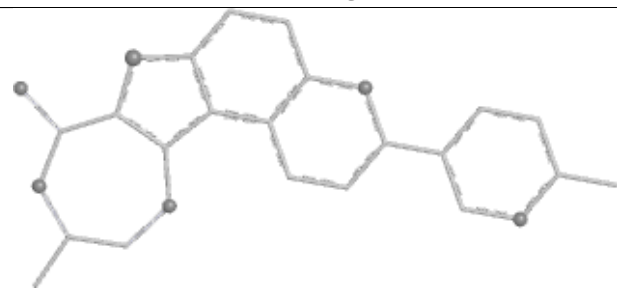
Ligand B97 C 401



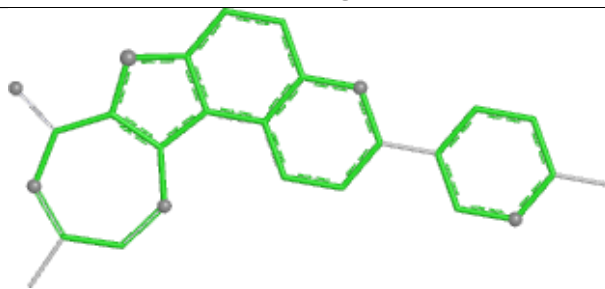
Bond lengths



Bond angles

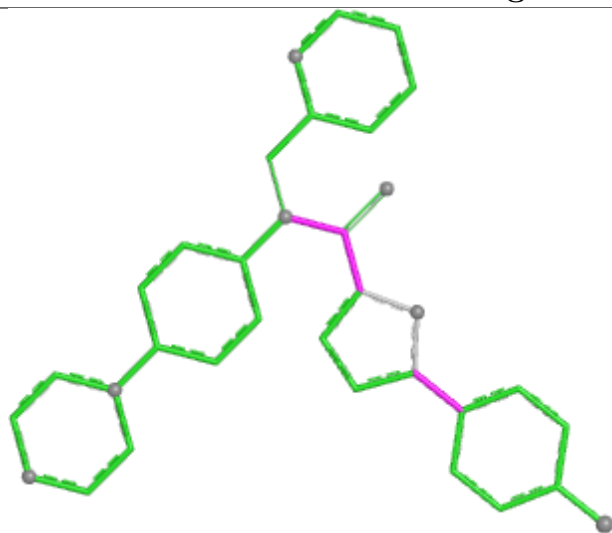


Torsions

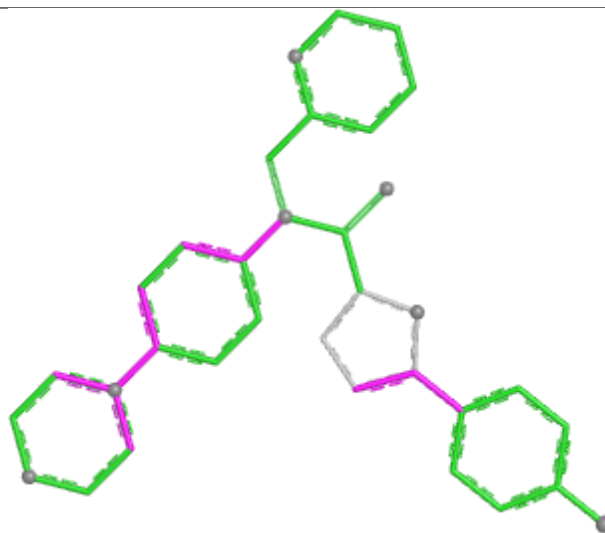


Rings

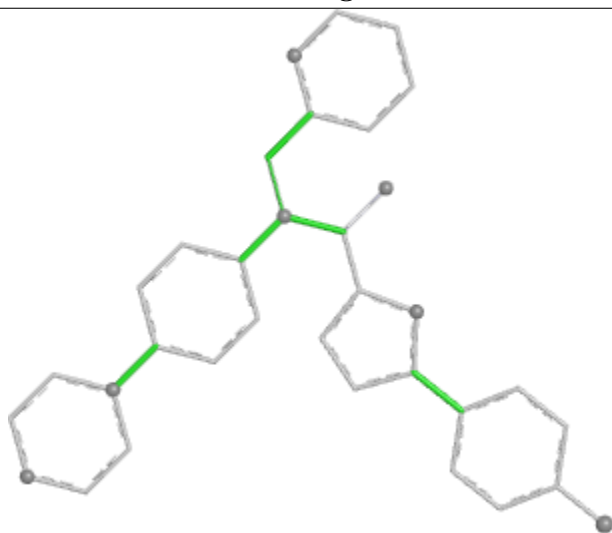
Ligand MW8 E 403



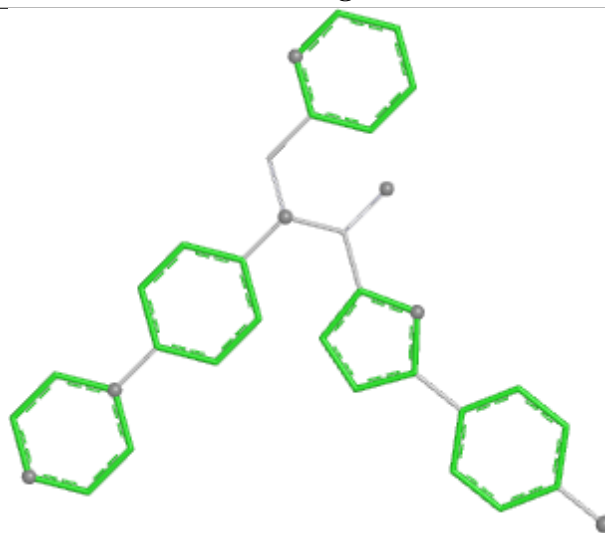
Bond lengths



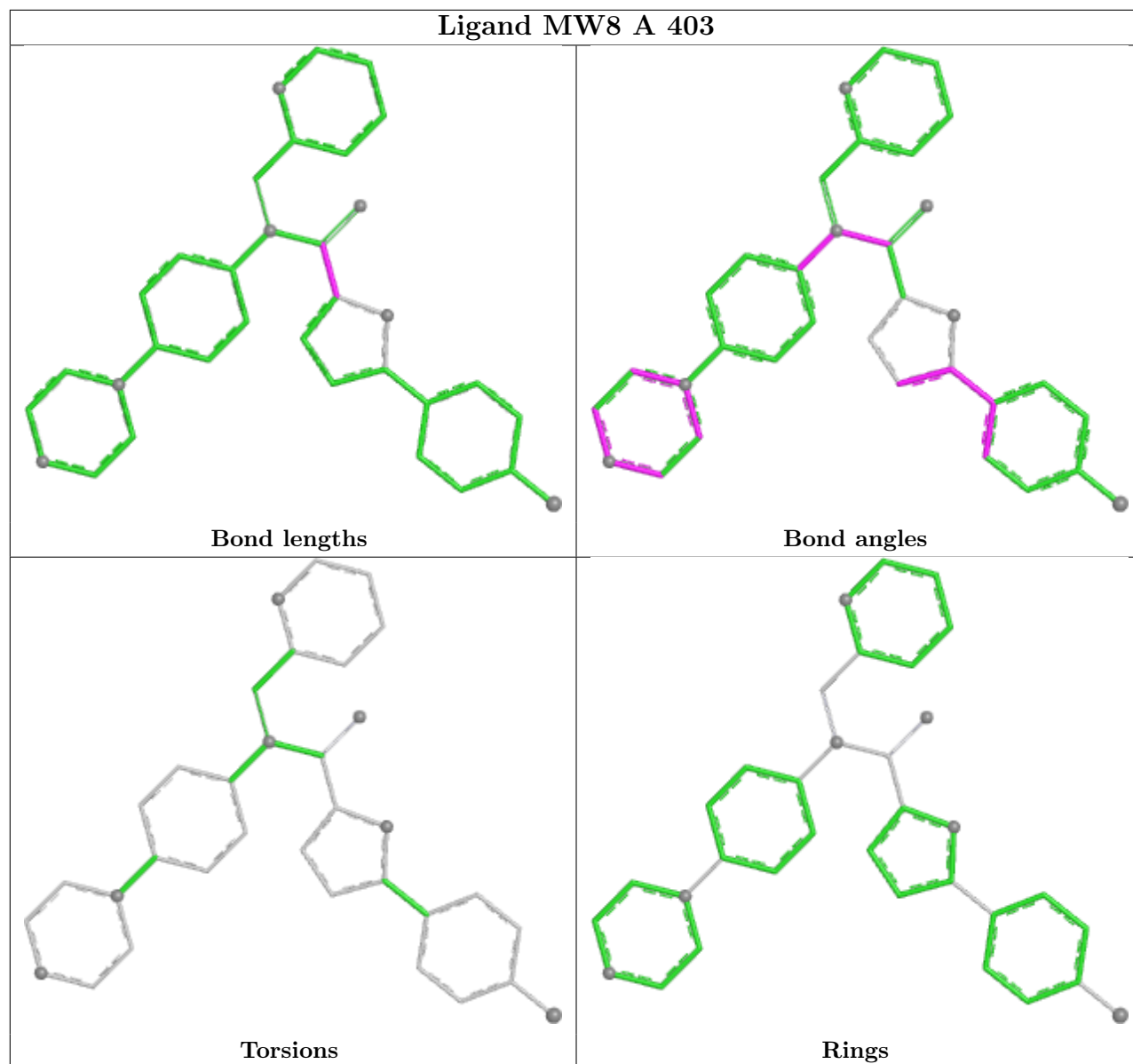
Bond angles

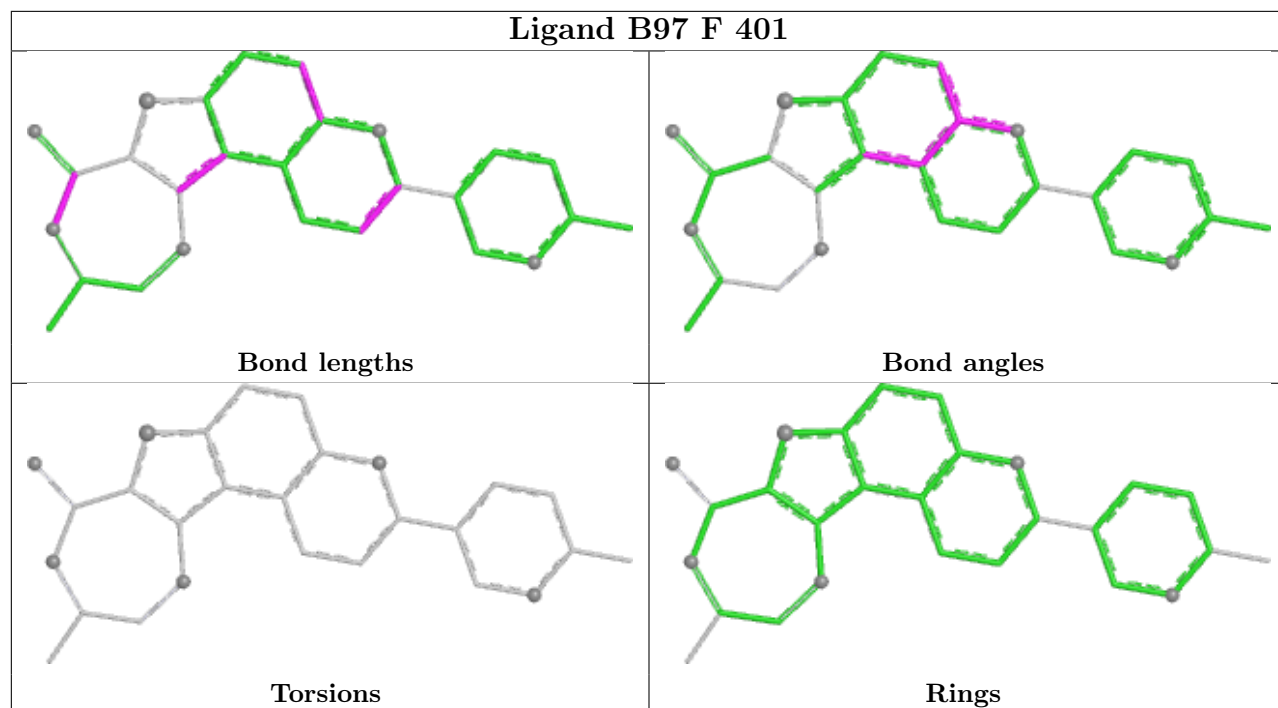


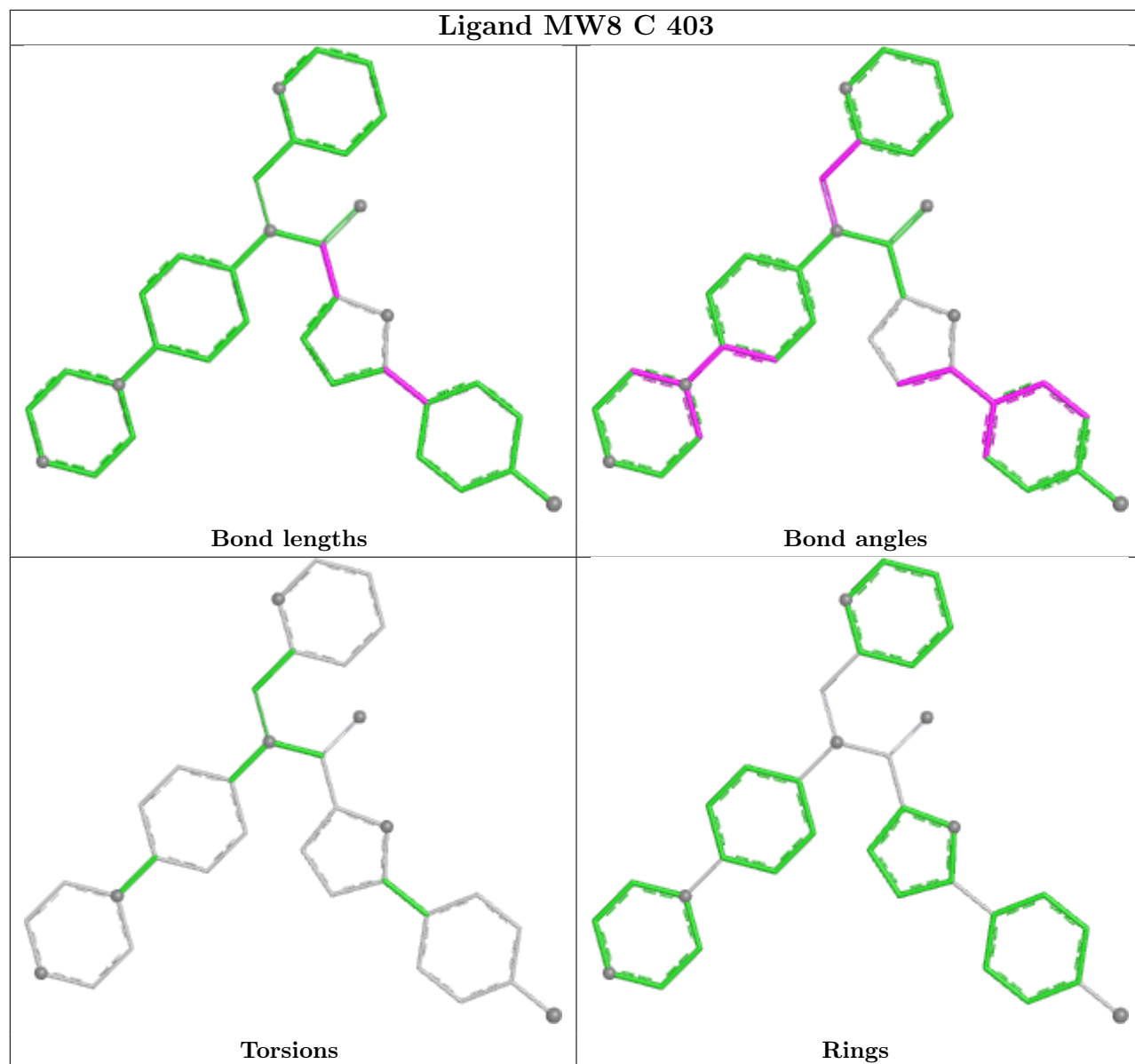
Torsions



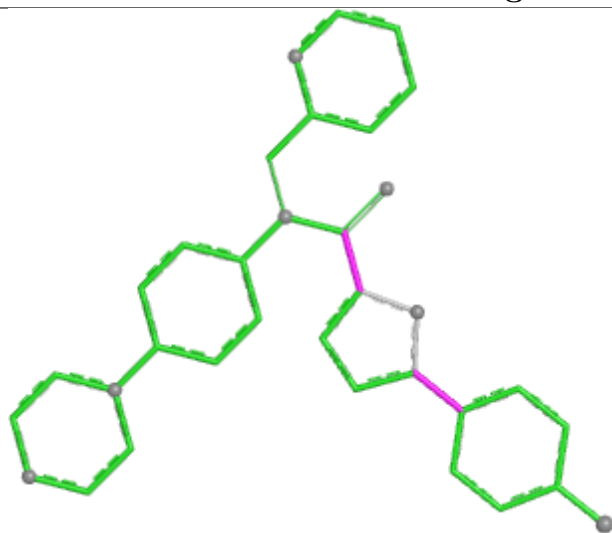
Rings



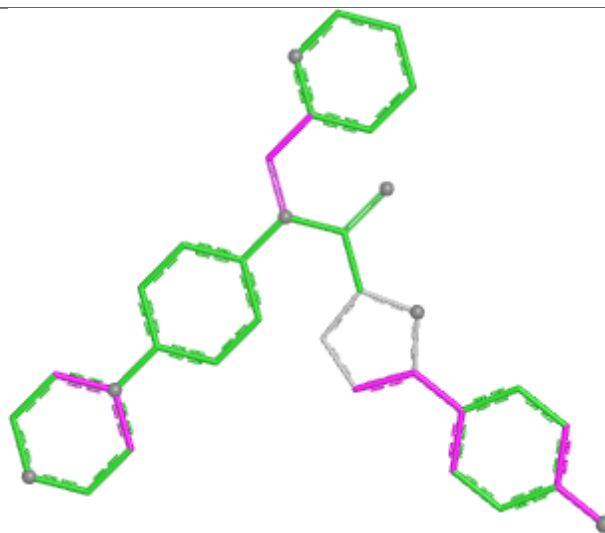




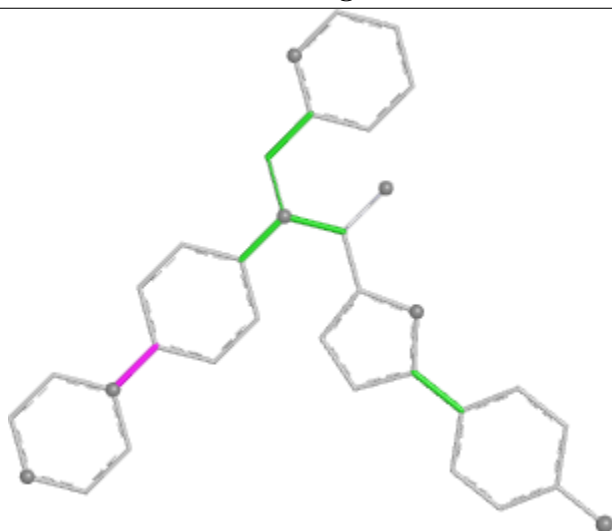
Ligand MW8 D 403



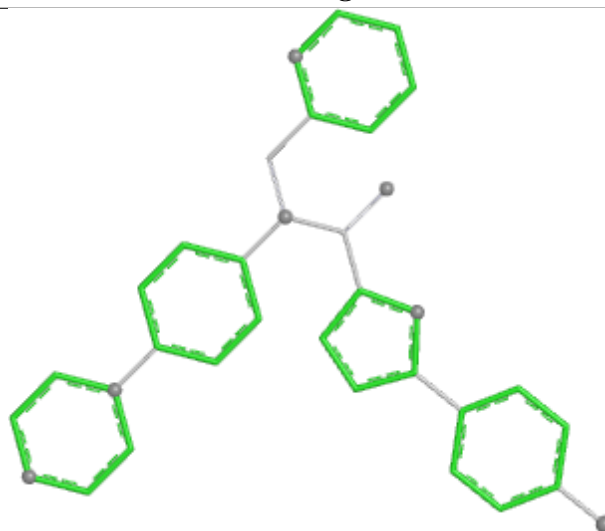
Bond lengths



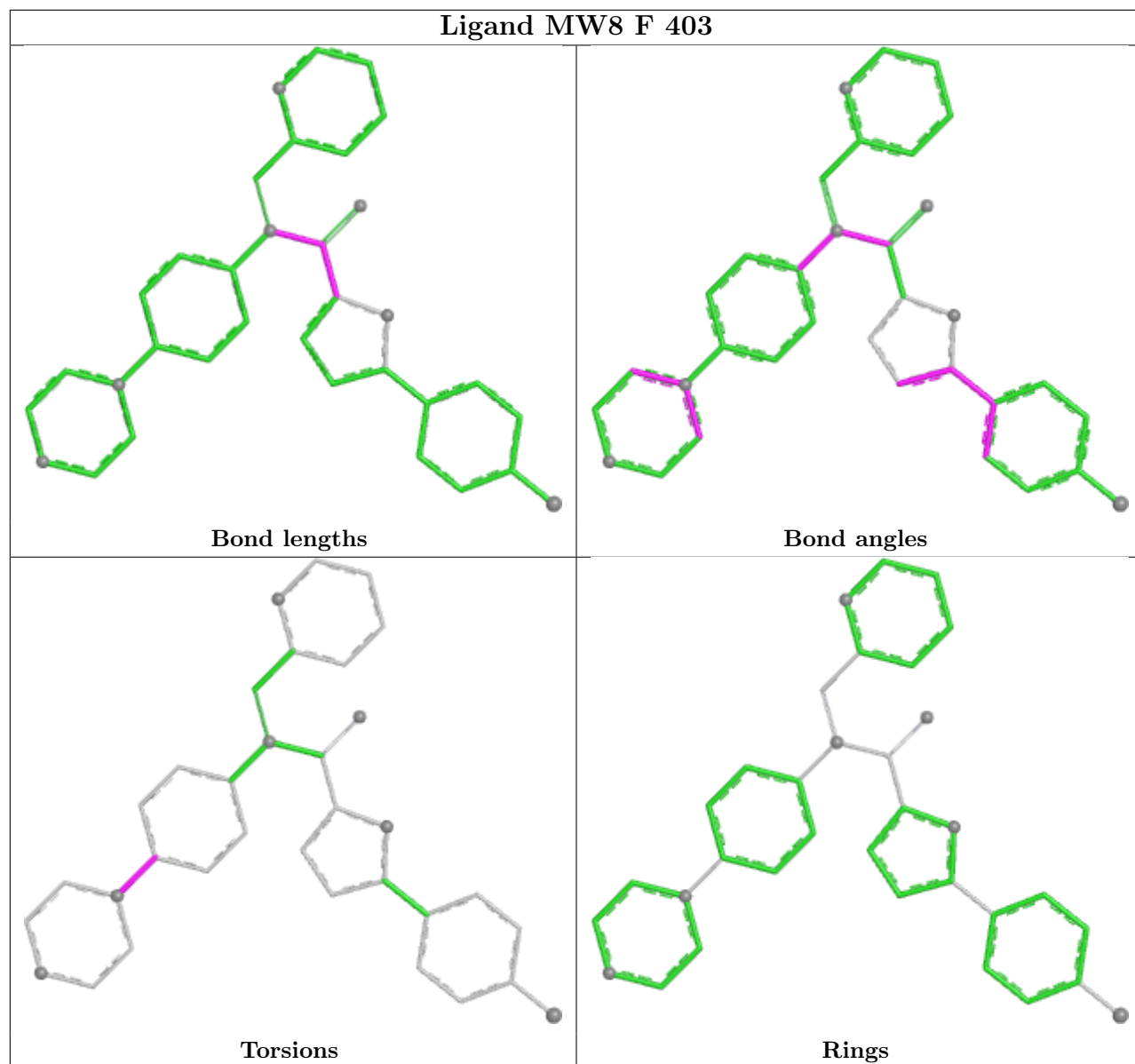
Bond angles

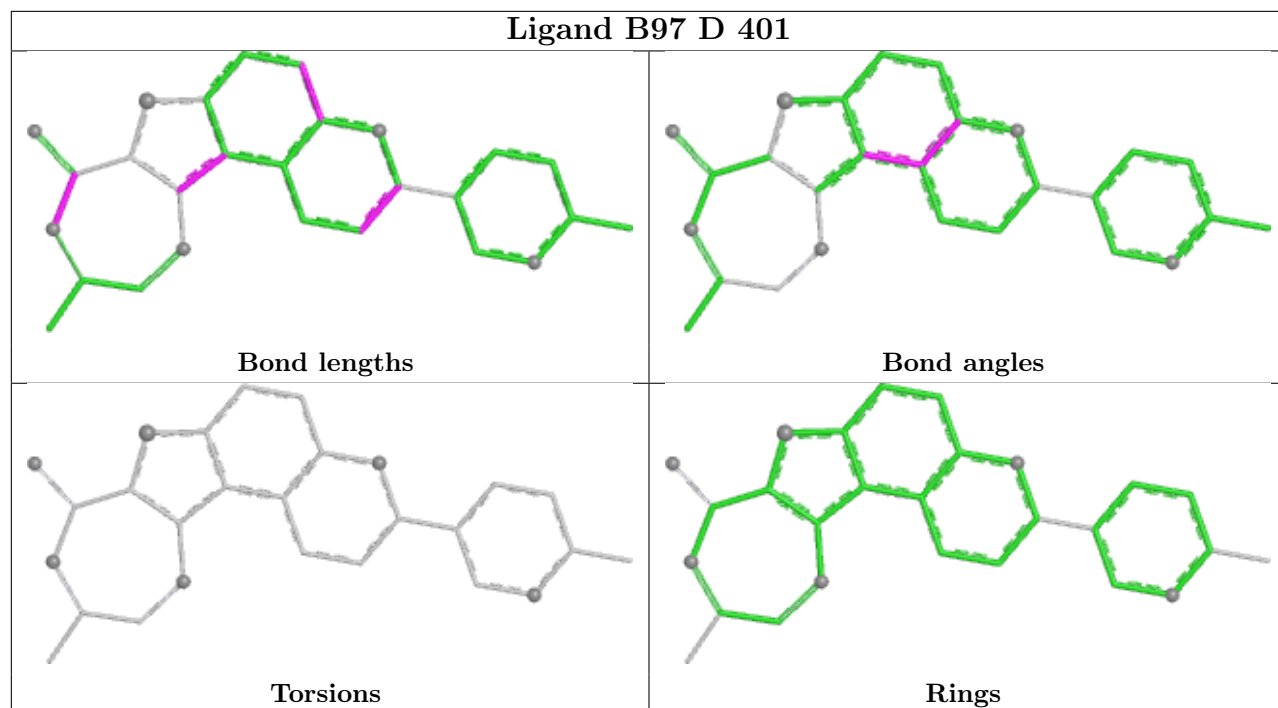


Torsions



Rings





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	279/319 (87%)	0.65	7 (2%) 57 47	66, 86, 114, 127	0
1	B	276/319 (86%)	0.72	8 (2%) 51 41	68, 89, 114, 127	0
1	C	276/319 (86%)	0.58	9 (3%) 46 36	69, 87, 110, 126	0
1	D	280/319 (87%)	0.76	12 (4%) 35 25	65, 90, 120, 133	0
1	E	277/319 (86%)	0.75	10 (3%) 42 32	74, 92, 119, 134	0
1	F	276/319 (86%)	0.73	13 (4%) 31 22	75, 100, 126, 141	0
All	All	1664/1914 (86%)	0.70	59 (3%) 44 34	65, 90, 120, 141	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	267	HIS	6.2
1	F	225	TYR	4.8
1	E	225	TYR	4.0
1	B	329	THR	3.6
1	E	240	TYR	3.6
1	D	150	ILE	3.2
1	B	226	THR	3.2
1	C	240	TYR	3.2
1	B	240	TYR	3.1
1	A	225	TYR	3.1
1	F	324	TRP	3.0
1	D	59	ILE	2.9
1	D	240	TYR	2.9
1	F	325	ILE	2.9
1	E	46	PHE	2.9
1	A	94	MET	2.8
1	C	226	THR	2.8
1	F	240	TYR	2.8
1	D	44	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	F	94	MET	2.7
1	B	59	ILE	2.7
1	E	269	LEU	2.6
1	A	240	TYR	2.6
1	A	133	CYS	2.6
1	C	298	VAL	2.5
1	A	267	HIS	2.5
1	E	226	THR	2.5
1	C	174	ILE	2.5
1	F	150	ILE	2.4
1	F	266	ASN	2.4
1	A	202	ILE	2.4
1	E	203	LEU	2.4
1	C	44	PRO	2.4
1	B	239	LYS	2.4
1	F	146	LEU	2.4
1	D	75	ASN	2.4
1	D	225	TYR	2.3
1	F	298	VAL	2.2
1	E	163	ALA	2.2
1	E	85	ARG	2.2
1	F	199	PRO	2.2
1	B	78	VAL	2.2
1	F	158	PHE	2.2
1	D	271	ILE	2.2
1	D	165	GLU	2.1
1	D	267	HIS	2.1
1	D	200	ASN	2.1
1	D	146	LEU	2.1
1	E	271	ILE	2.1
1	D	164	SER	2.1
1	C	187	VAL	2.1
1	B	46	PHE	2.1
1	C	215	THR	2.0
1	C	118	VAL	2.0
1	B	64	LYS	2.0
1	F	226	THR	2.0
1	C	46	PHE	2.0
1	A	249	LEU	2.0
1	E	134	LEU	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 monosaccharides in this entry.

6.4 Ligands [i](#)

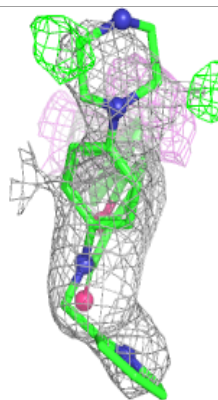
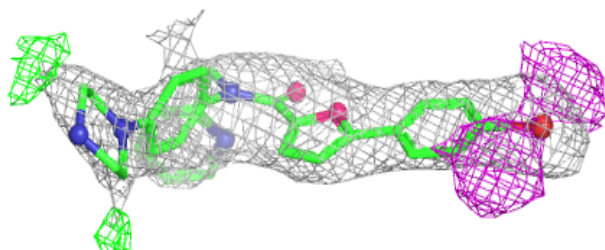
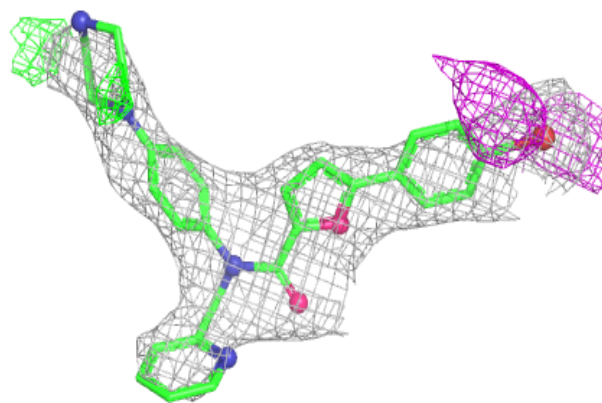
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	CL	E	402	1/1	0.49	0.11	127,127,127,127	0
3	CL	C	402	1/1	0.50	0.11	115,115,115,115	0
3	CL	F	402	1/1	0.76	0.09	122,122,122,122	0
3	CL	B	402	1/1	0.79	0.18	123,123,123,123	0
3	CL	A	402	1/1	0.79	0.16	119,119,119,119	0
4	MW8	D	403	34/34	0.82	0.32	133,134,135,136	0
4	MW8	A	403	34/34	0.86	0.26	121,121,123,124	0
4	MW8	F	403	34/34	0.86	0.28	137,138,140,142	0
3	CL	D	402	1/1	0.90	0.18	123,123,123,123	0
4	MW8	C	403	34/34	0.90	0.26	114,116,117,117	0
4	MW8	E	403	34/34	0.91	0.26	108,110,112,113	0
4	MW8	B	403	34/34	0.92	0.22	104,109,113,113	0
2	B97	F	401	27/27	0.93	0.22	90,94,102,102	0
2	B97	C	401	27/27	0.94	0.24	75,81,84,84	0
2	B97	D	401	27/27	0.95	0.27	87,88,97,97	0
2	B97	B	401	27/27	0.95	0.24	73,75,83,83	0
2	B97	A	401	27/27	0.96	0.23	84,89,94,94	0
2	B97	E	401	27/27	0.97	0.25	92,95,99,100	0

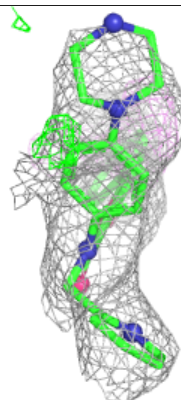
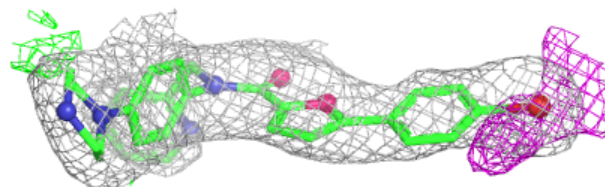
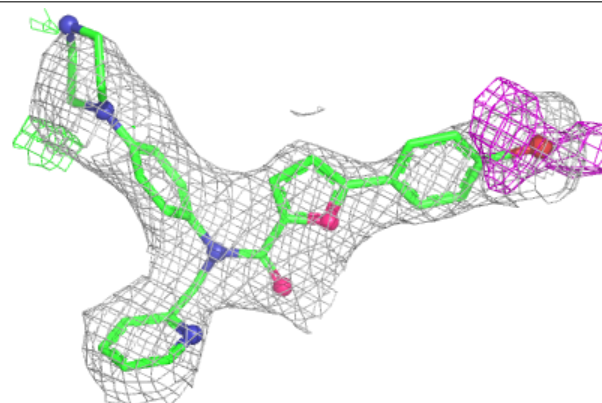
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around MW8 D 403:

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

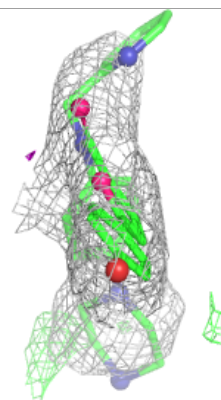
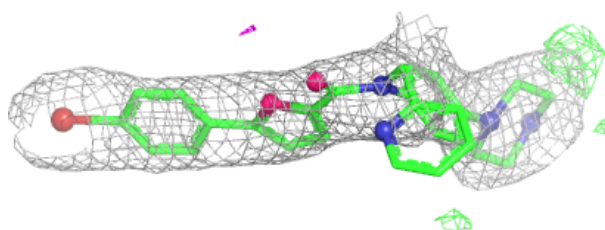
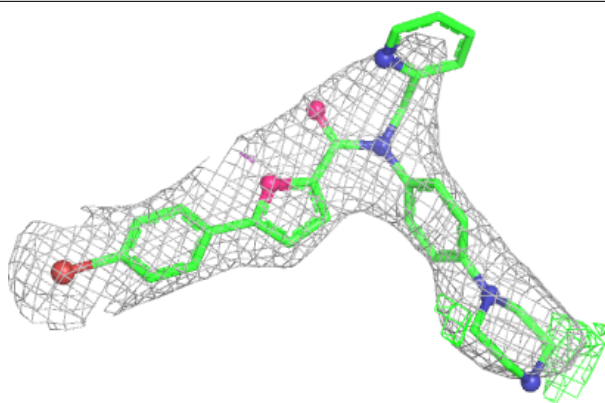
**Electron density around MW8 A 403:**

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

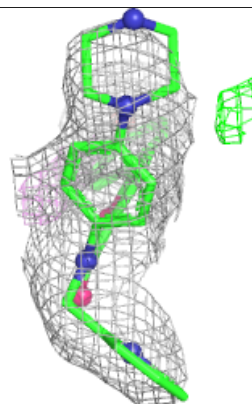
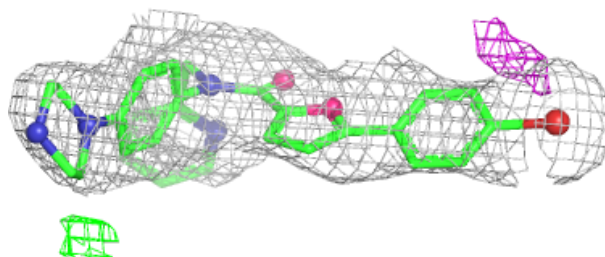
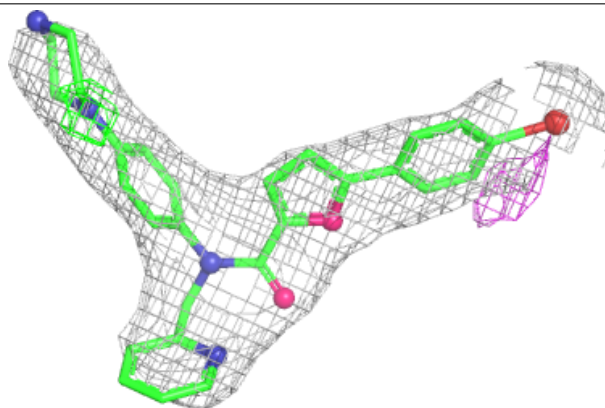


Electron density around MW8 F 403:

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

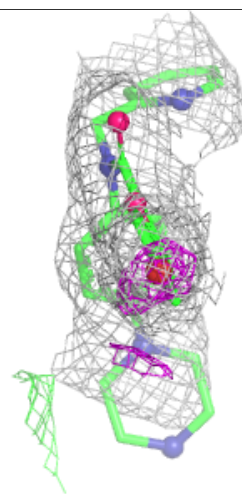
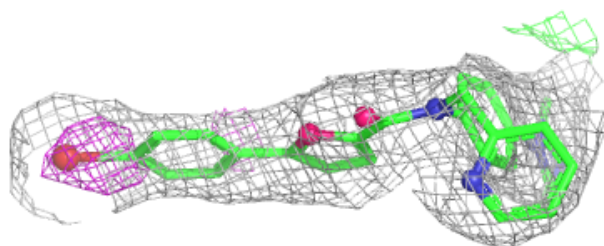
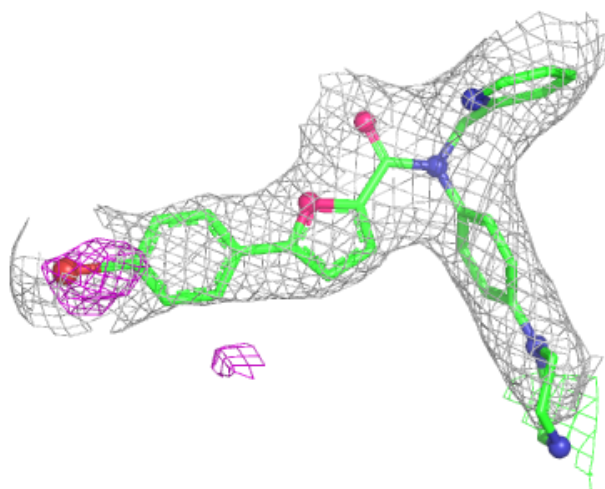
**Electron density around MW8 C 403:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



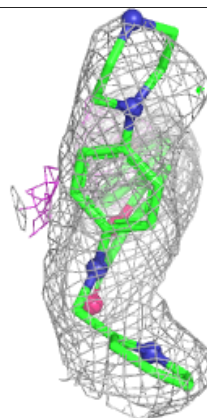
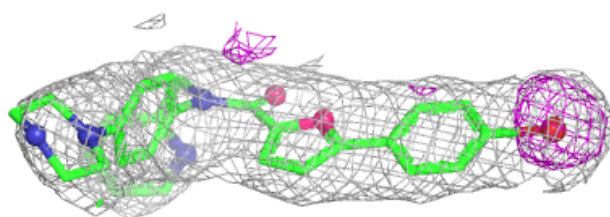
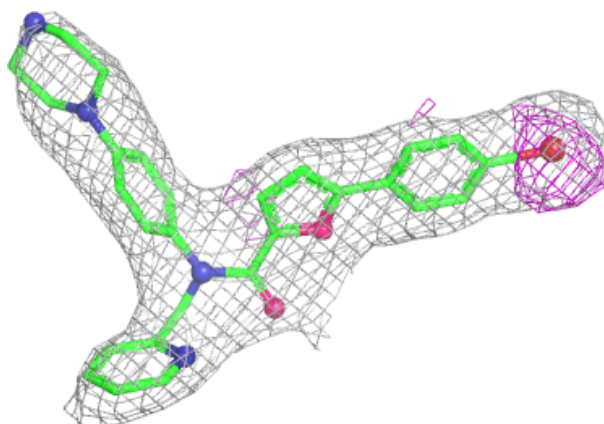
Electron density around MW8 E 403:

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

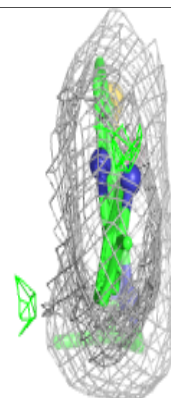
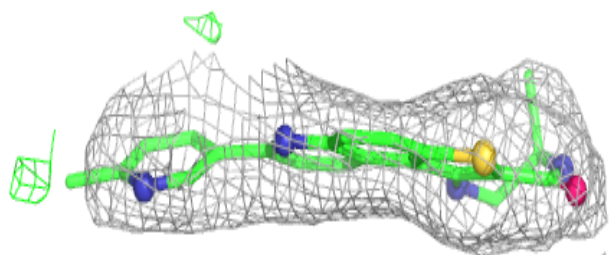
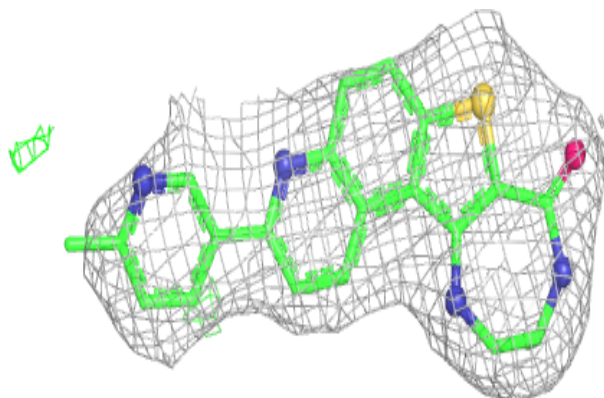


Electron density around MW8 B 403:

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

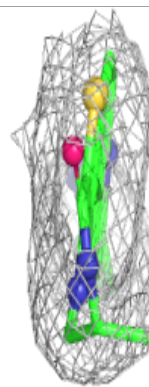
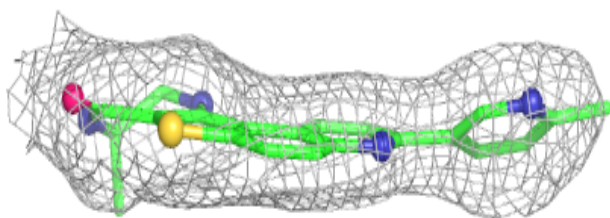
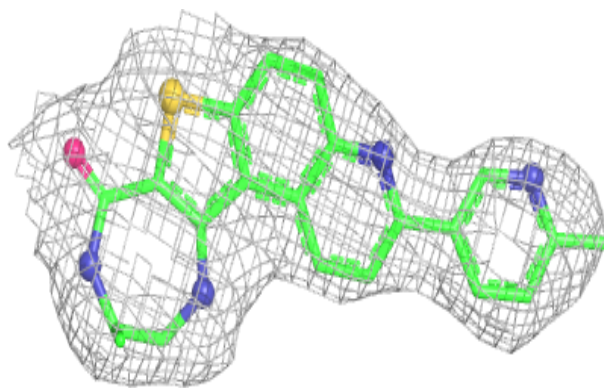
**Electron density around B97 F 401:**

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

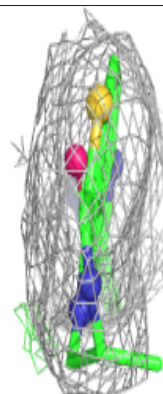
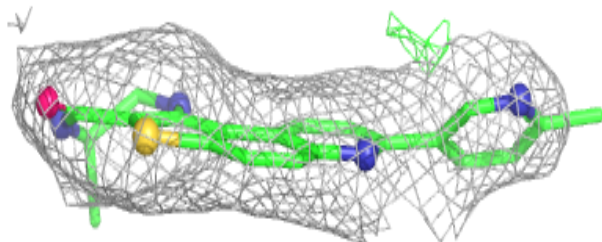
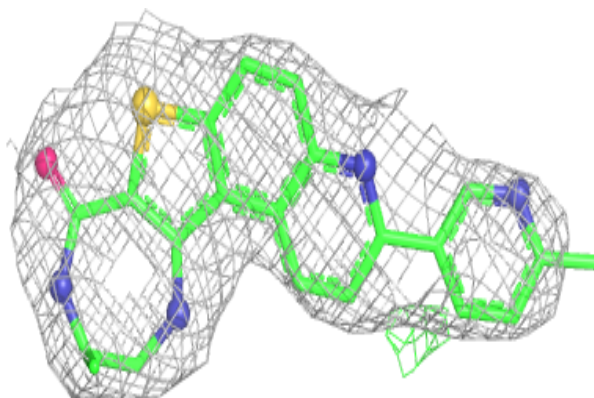


Electron density around B97 C 401:

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

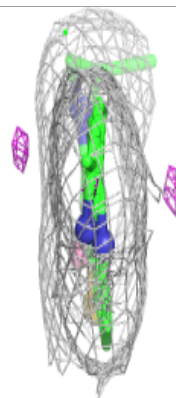
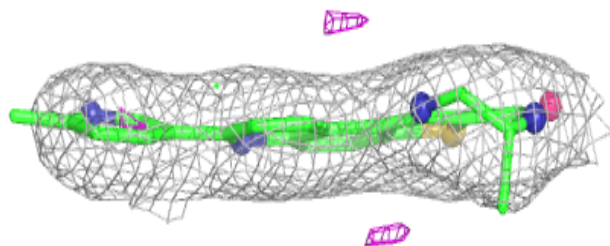
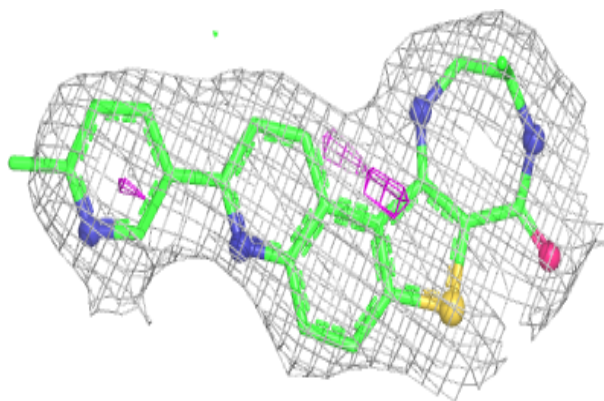
**Electron density around B97 D 401:**

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

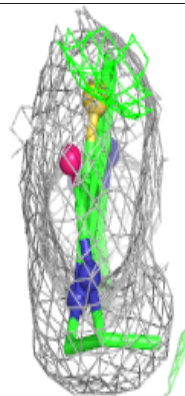
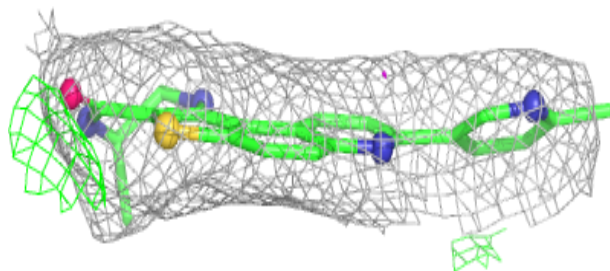
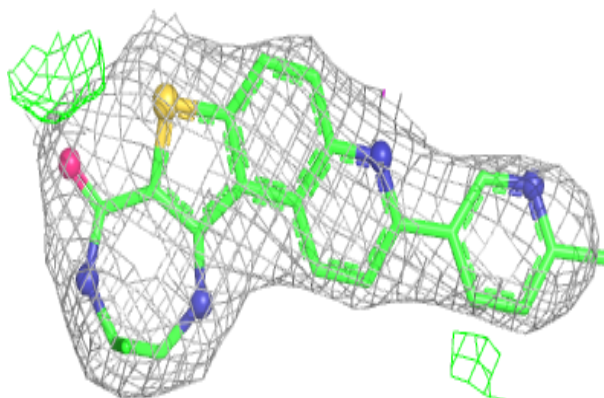


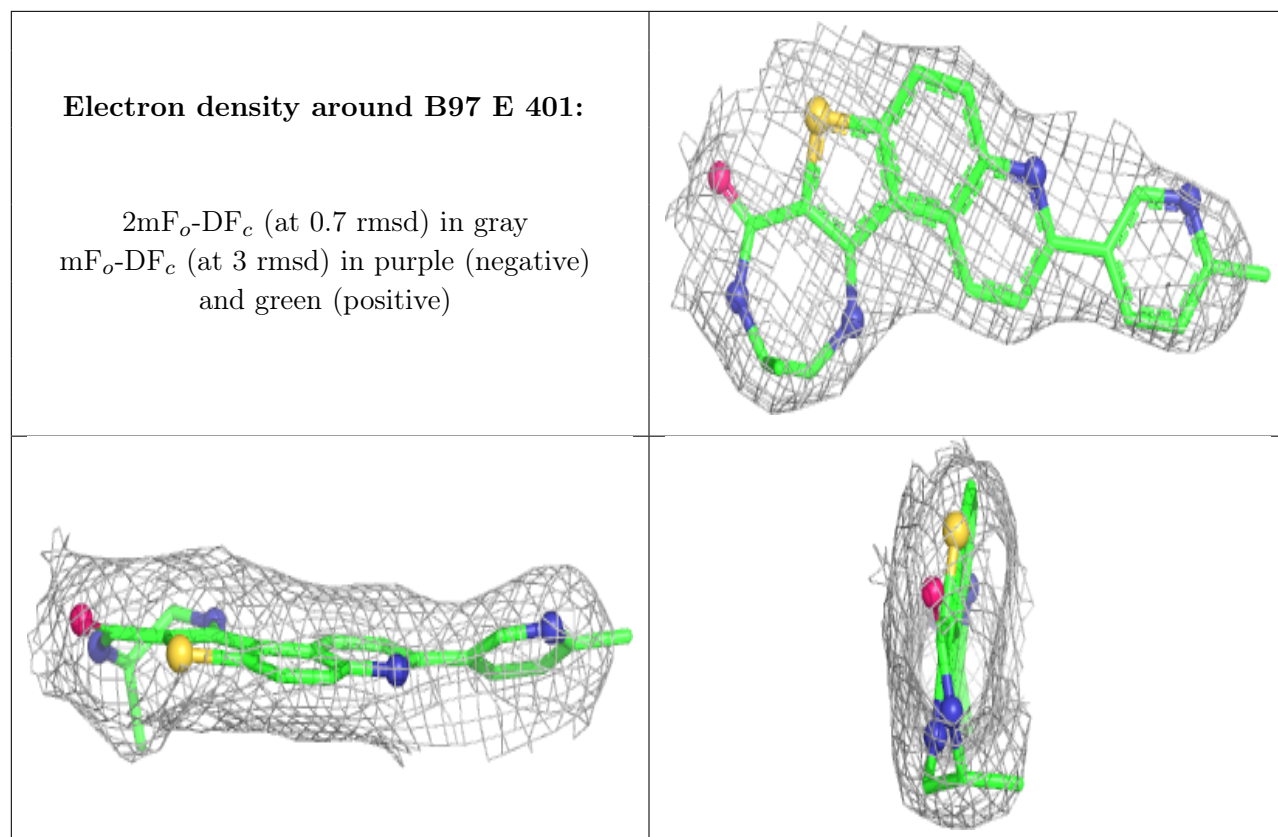
Electron density around B97 B 401:

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

**Electron density around B97 A 401:**

$2mF_o-DF_c$ (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.