



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 17, 2025 – 02:16 PM EDT

PDB ID : 9NSC / pdb_00009nsc
Title : Bacterial Pictet-Spenglerase KslB in complex with L-Trp
Authors : Kim, K.; Kim, W.
Deposited on : 2025-03-16
Resolution : 3.10 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 : 4.02b-467
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) : 2.42

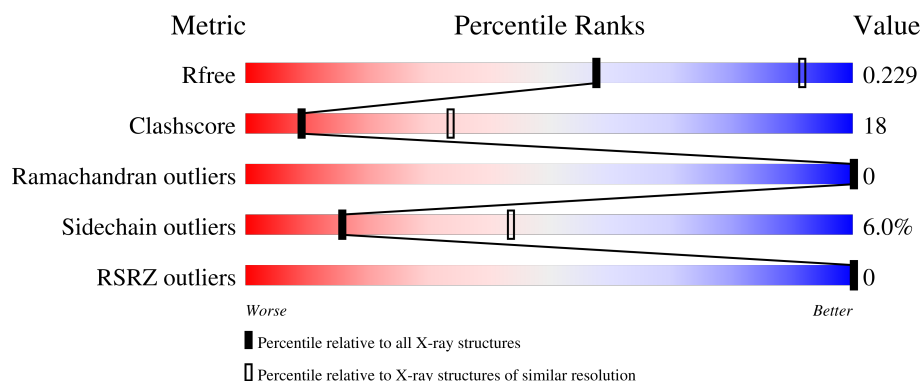
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1351 (3.10-3.10)
Clashscore	180529	1454 (3.10-3.10)
Ramachandran outliers	177936	1391 (3.10-3.10)
Sidechain outliers	177891	1391 (3.10-3.10)
RSRZ outliers	164620	1351 (3.10-3.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	317	
1	B	317	
1	C	317	
1	D	317	
1	E	317	

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Mol	Chain	Length	Quality of chain
1	F	317	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: green (57%), yellow (32%), and orange (9%). The segments are labeled with their respective percentages: 57%, 32%, and 9%.

2 Entry composition [i](#)

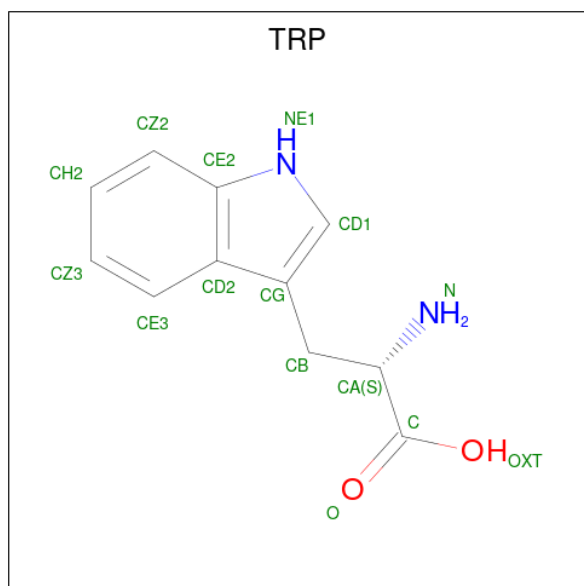
There are 2 unique types of molecules in this entry. The entry contains 14024 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

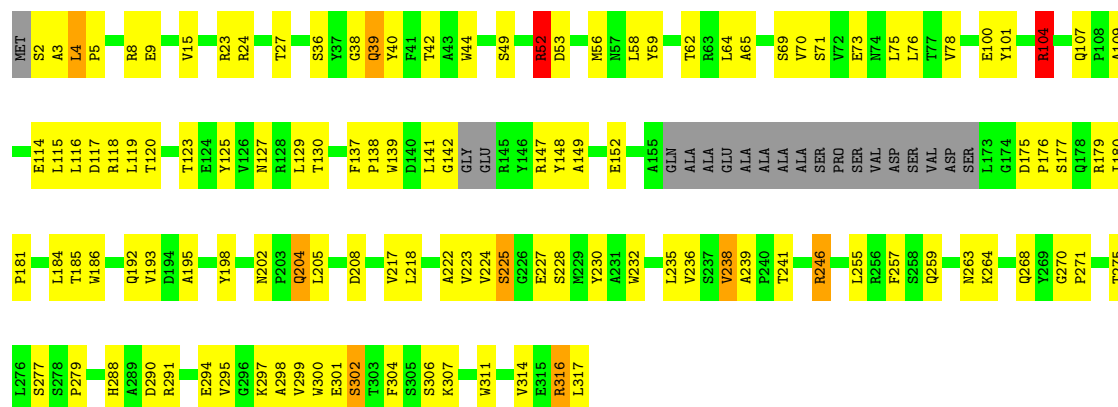
- Molecule 1 is a protein called Pictet-Spenglerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	283	Total	C	N	O	S	0	0	0
			2255	1450	381	422	2			
1	B	295	Total	C	N	O	S	0	0	0
			2337	1496	393	444	4			
1	C	297	Total	C	N	O	S	0	0	0
			2356	1509	398	446	3			
1	D	301	Total	C	N	O	S	0	0	0
			2386	1526	403	454	3			
1	E	288	Total	C	N	O	S	0	0	0
			2299	1474	389	433	3			
1	F	290	Total	C	N	O	S	0	0	0
			2301	1475	387	436	3			

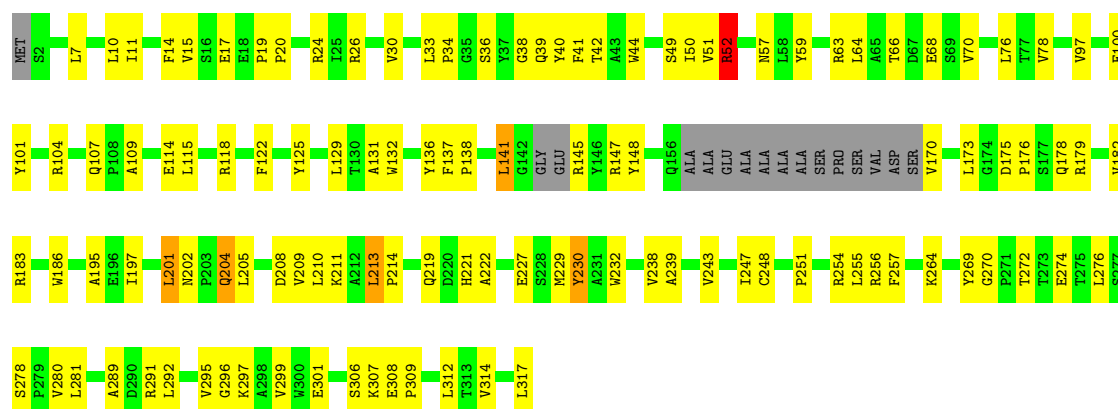
- Molecule 2 is TRYPTOPHAN (CCD ID: TRP) (formula: $C_{11}H_{12}N_2O_2$) (labeled as "Ligand of Interest" by depositor).



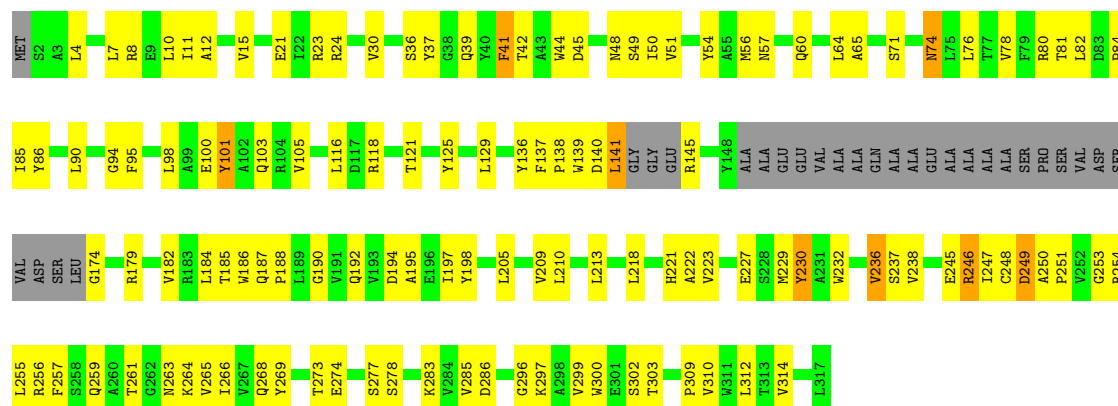
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			15	11	2	2		
2	B	1	Total	C	N	O	0	0
			15	11	2	2		
2	C	1	Total	C	N	O	0	0
			15	11	2	2		
2	D	1	Total	C	N	O	0	0
			15	11	2	2		
2	E	1	Total	C	N	O	0	0
			15	11	2	2		
2	F	1	Total	C	N	O	0	0
			15	11	2	2		



• Molecule 1: Pictet-Spenglerase



• Molecule 1: Pictet-Spenglerase



• Molecule 1: Pictet-Spenglerase





4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	94.53Å 94.53Å 193.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.27 – 3.10 47.27 – 3.10	Depositor EDS
% Data completeness (in resolution range)	98.0 (47.27-3.10) 97.9 (47.27-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.176 , 0.234 0.177 , 0.229	Depositor DCC
R_{free} test set	32294 reflections (5.88%)	wwPDB-VP
Wilson B-factor (Å ²)	52.9	Xtriage
Anisotropy	0.398	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 25.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.067 for -h,-k,l 0.417 for h,-h-k,-l 0.076 for -k,-h,-l	Xtriage
Reported twinning fraction	0.480 for h,-h-k,-l	Depositor
Outliers	0 of 34328 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	14024	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% 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](#)

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	0.62	0/2314	0.72	0/3161
1	B	0.63	0/2397	0.72	0/3279
1	C	0.63	0/2417	0.73	0/3306
1	D	0.63	0/2447	0.73	0/3347
1	E	0.62	0/2360	0.70	0/3228
1	F	0.63	0/2361	0.70	0/3231
All	All	0.63	0/14296	0.72	0/19552

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	1
1	C	0	2
1	D	0	1
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	147	ARG	Sidechain
1	C	104	ARG	Sidechain
1	C	52	ARG	Sidechain
1	D	52	ARG	Sidechain

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	2255	0	2211	97	0
1	B	2337	0	2291	92	0
1	C	2356	0	2309	87	0
1	D	2386	0	2335	84	0
1	E	2299	0	2254	91	0
1	F	2301	0	2255	77	0
2	A	15	0	9	0	0
2	B	15	0	9	0	0
2	C	15	0	9	0	0
2	D	15	0	9	0	0
2	E	15	0	9	2	0
2	F	15	0	9	0	0
All	All	14024	0	13709	496	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 18.

The worst 5 of 496 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:39:GLN:HE21	1:E:41:PHE:HB3	1.37	0.89
1:B:64:LEU:HB3	1:B:75:LEU:HD21	1.60	0.83
1:D:173:LEU:HB3	1:D:201:LEU:HD11	1.61	0.82
1:B:175:ASP:HB3	1:B:178:GLN:HG3	1.62	0.81
1:D:104:ARG:HA	1:D:107:GLN:HG3	1.64	0.80

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	273/317 (86%)	260 (95%)	13 (5%)	0	100	100
1	B	289/317 (91%)	278 (96%)	11 (4%)	0	100	100
1	C	291/317 (92%)	276 (95%)	15 (5%)	0	100	100
1	D	295/317 (93%)	281 (95%)	14 (5%)	0	100	100
1	E	282/317 (89%)	268 (95%)	14 (5%)	0	100	100
1	F	284/317 (90%)	272 (96%)	12 (4%)	0	100	100
All	All	1714/1902 (90%)	1635 (95%)	79 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	245/268 (91%)	238 (97%)	7 (3%)	37	65
1	B	255/268 (95%)	243 (95%)	12 (5%)	22	52
1	C	255/268 (95%)	236 (92%)	19 (8%)	11	36
1	D	259/268 (97%)	244 (94%)	15 (6%)	17	45
1	E	251/268 (94%)	233 (93%)	18 (7%)	12	38
1	F	252/268 (94%)	232 (92%)	20 (8%)	10	34
All	All	1517/1608 (94%)	1426 (94%)	91 (6%)	16	44

5 of 91 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	101	TYR
1	F	29	THR
1	E	141	LEU
1	E	245	GLU
1	F	67	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	39	GLN
1	E	74	ASN
1	F	74	ASN
1	E	259	GLN
1	C	202	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](#)

6 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	TRP	E	401	-	14,16,16	0.93	1 (7%)	13,22,22	1.04	1 (7%)
2	TRP	A	401	-	14,16,16	0.77	0	13,22,22	0.91	1 (7%)
2	TRP	F	401	-	14,16,16	0.75	0	13,22,22	0.87	1 (7%)
2	TRP	B	401	-	14,16,16	0.79	1 (7%)	13,22,22	0.88	0
2	TRP	C	401	-	14,16,16	0.78	0	13,22,22	0.94	1 (7%)
2	TRP	D	401	-	14,16,16	0.79	0	13,22,22	0.92	1 (7%)

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	TRP	E	401	-	-	5/7/8/8	0/2/2/2
2	TRP	A	401	-	-	2/7/8/8	0/2/2/2
2	TRP	F	401	-	-	1/7/8/8	0/2/2/2
2	TRP	B	401	-	-	5/7/8/8	0/2/2/2
2	TRP	C	401	-	-	4/7/8/8	0/2/2/2
2	TRP	D	401	-	-	6/7/8/8	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	401	TRP	OXT-C	-2.10	1.24	1.30
2	B	401	TRP	OXT-C	-2.02	1.24	1.30

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	401	TRP	CH2-CZ2-CE2	-2.28	116.97	120.09
2	C	401	TRP	CH2-CZ2-CE2	-2.12	117.19	120.09
2	D	401	TRP	CH2-CZ2-CE2	-2.11	117.20	120.09
2	A	401	TRP	CH2-CZ2-CE2	-2.06	117.28	120.09
2	F	401	TRP	CH2-CZ2-CE2	-2.04	117.30	120.09

There are no chirality outliers.

5 of 23 torsion outliers are listed below:

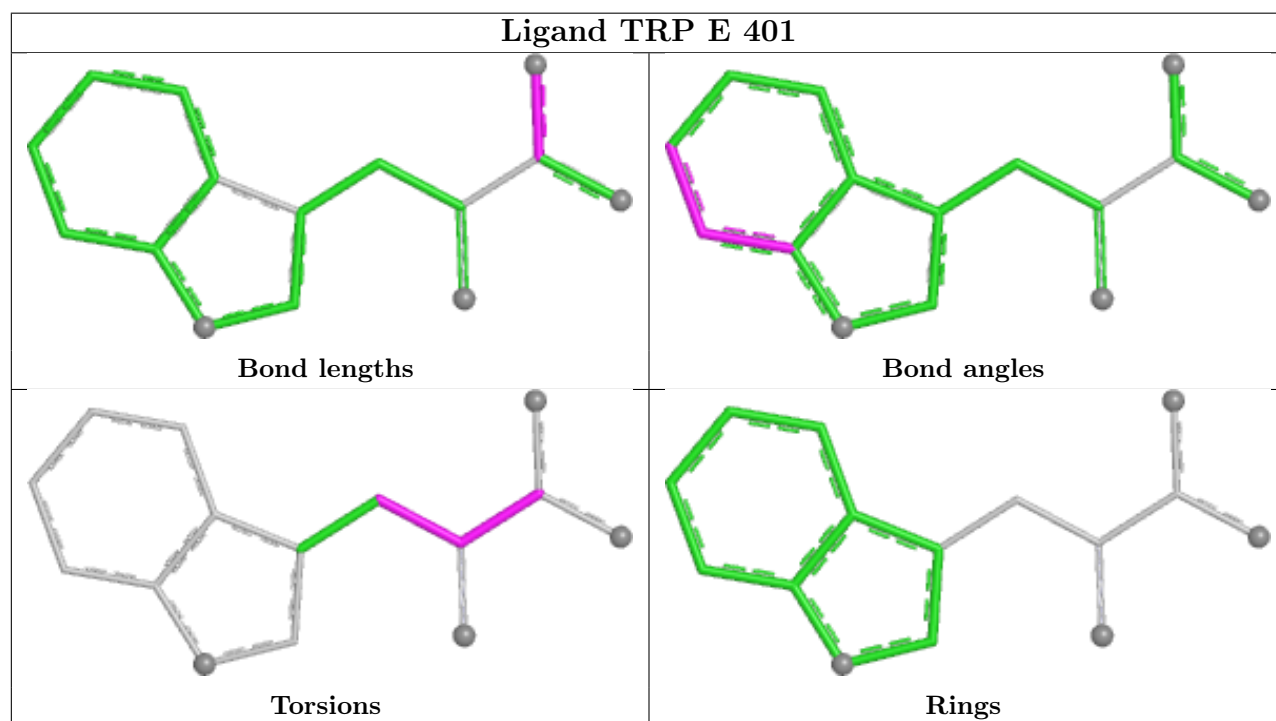
Mol	Chain	Res	Type	Atoms
2	C	401	TRP	O-C-CA-N
2	D	401	TRP	N-CA-CB-CG
2	D	401	TRP	C-CA-CB-CG
2	E	401	TRP	O-C-CA-N
2	E	401	TRP	N-CA-CB-CG

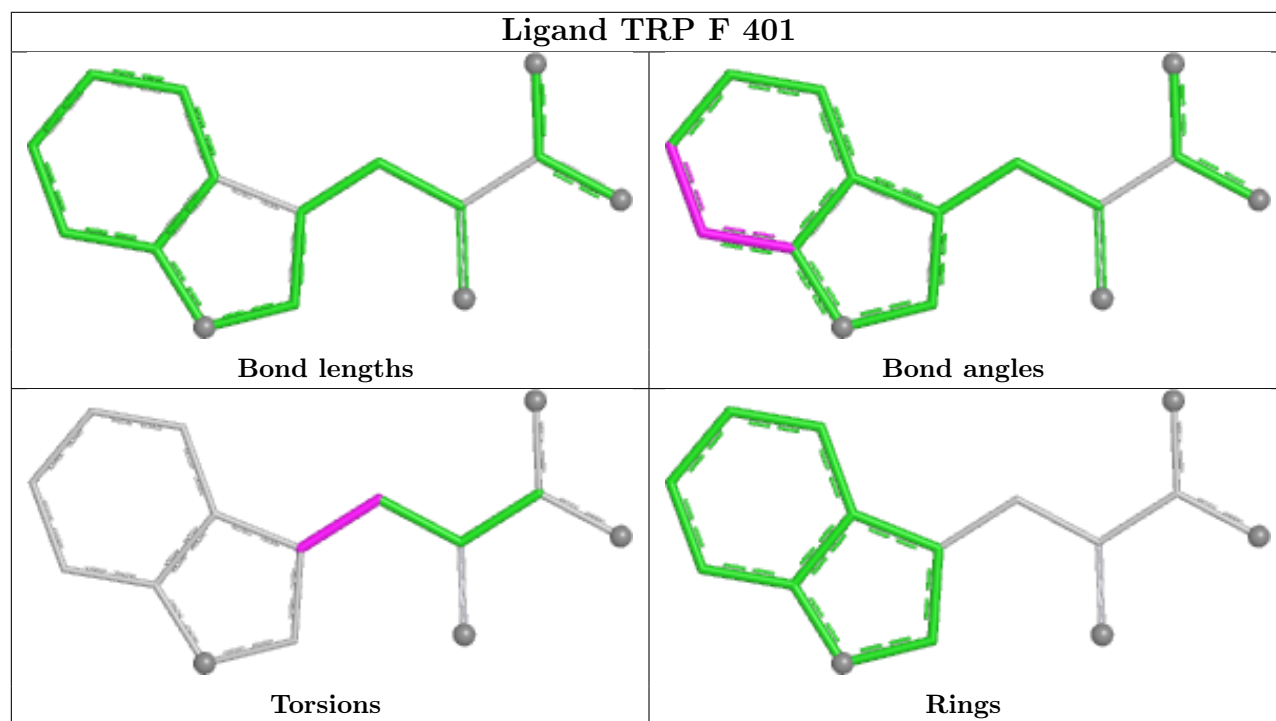
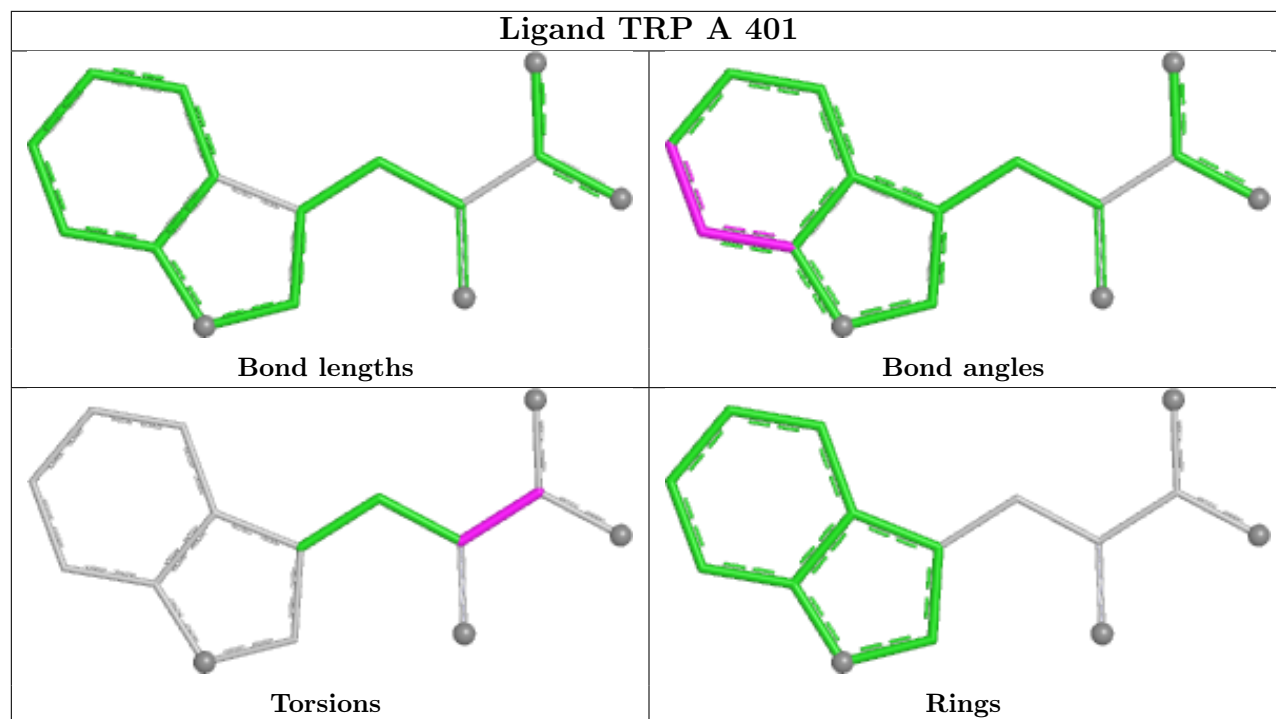
There are no ring outliers.

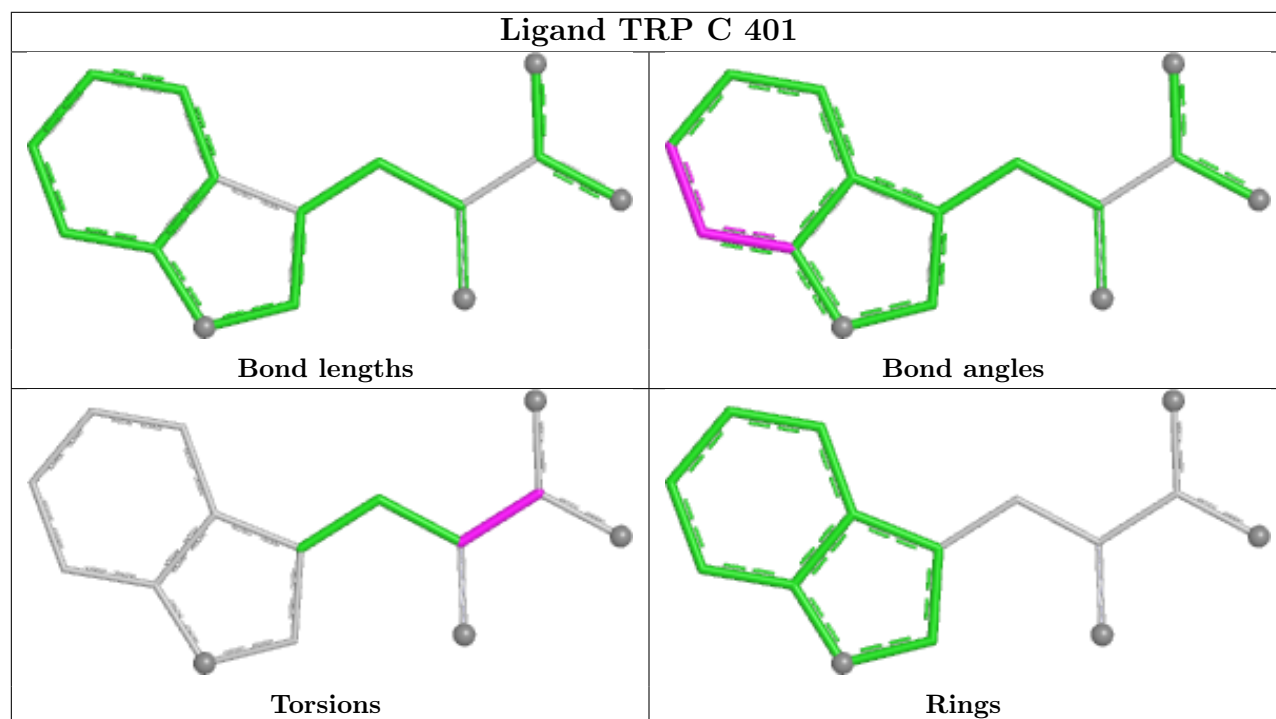
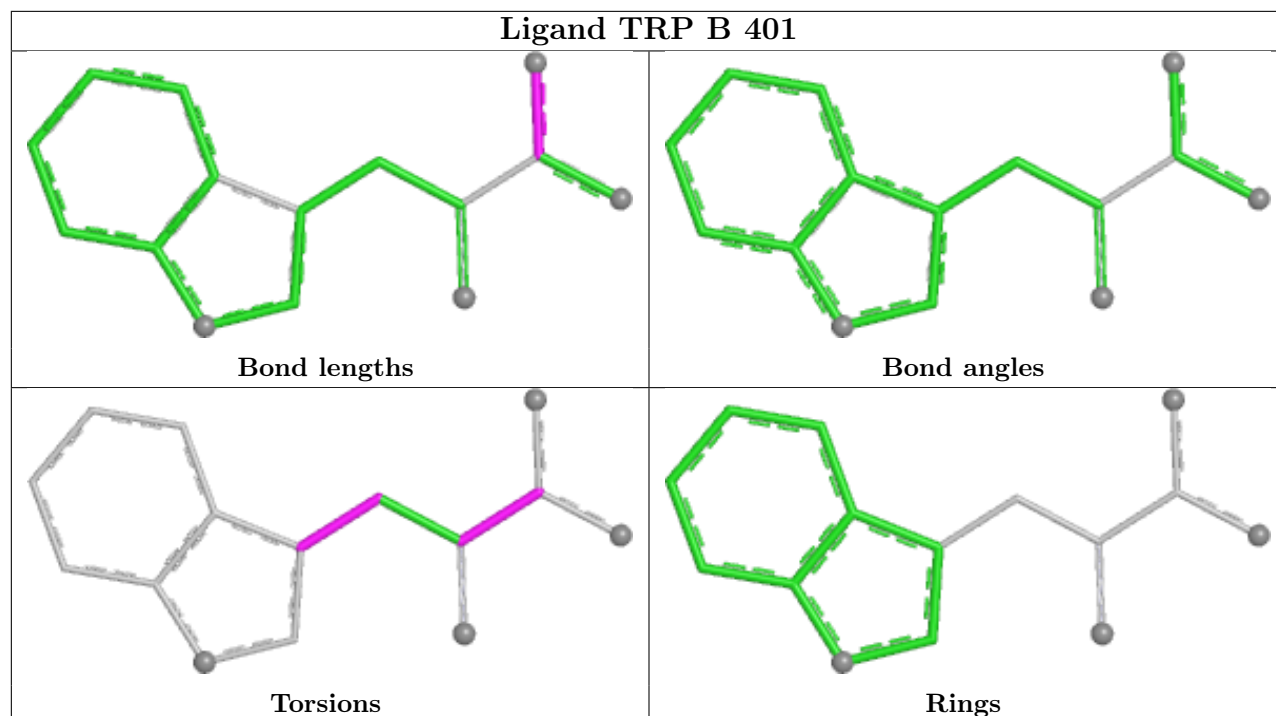
1 monomer is involved in 2 short contacts:

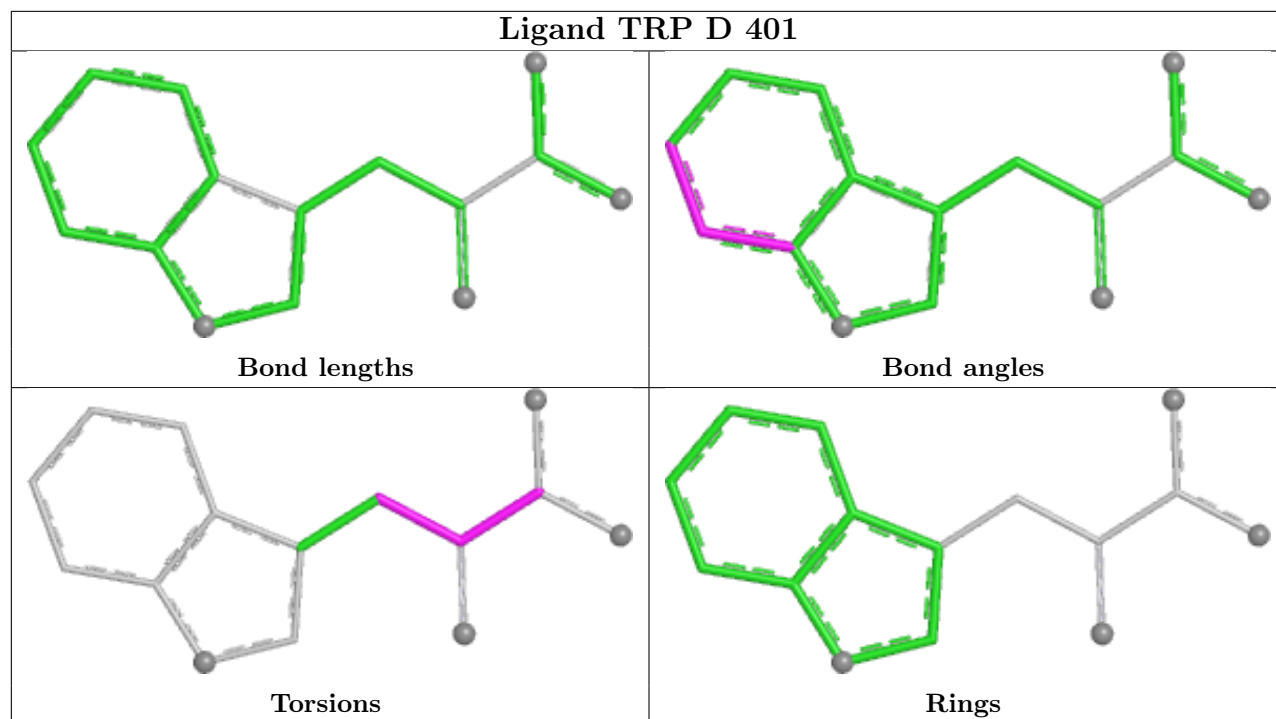
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	401	TRP	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	283/317 (89%)	-1.98	0 100 100	18, 36, 55, 77	0
1	B	295/317 (93%)	-1.96	0 100 100	19, 37, 58, 66	0
1	C	297/317 (93%)	-1.96	0 100 100	20, 37, 56, 77	0
1	D	301/317 (94%)	-1.95	0 100 100	19, 40, 58, 84	0
1	E	288/317 (90%)	-1.94	0 100 100	25, 47, 68, 79	0
1	F	290/317 (91%)	-1.93	0 100 100	24, 50, 67, 80	0
All	All	1754/1902 (92%)	-1.95	0 100 100	18, 41, 62, 84	0

There are no RSRZ outliers to report.

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.

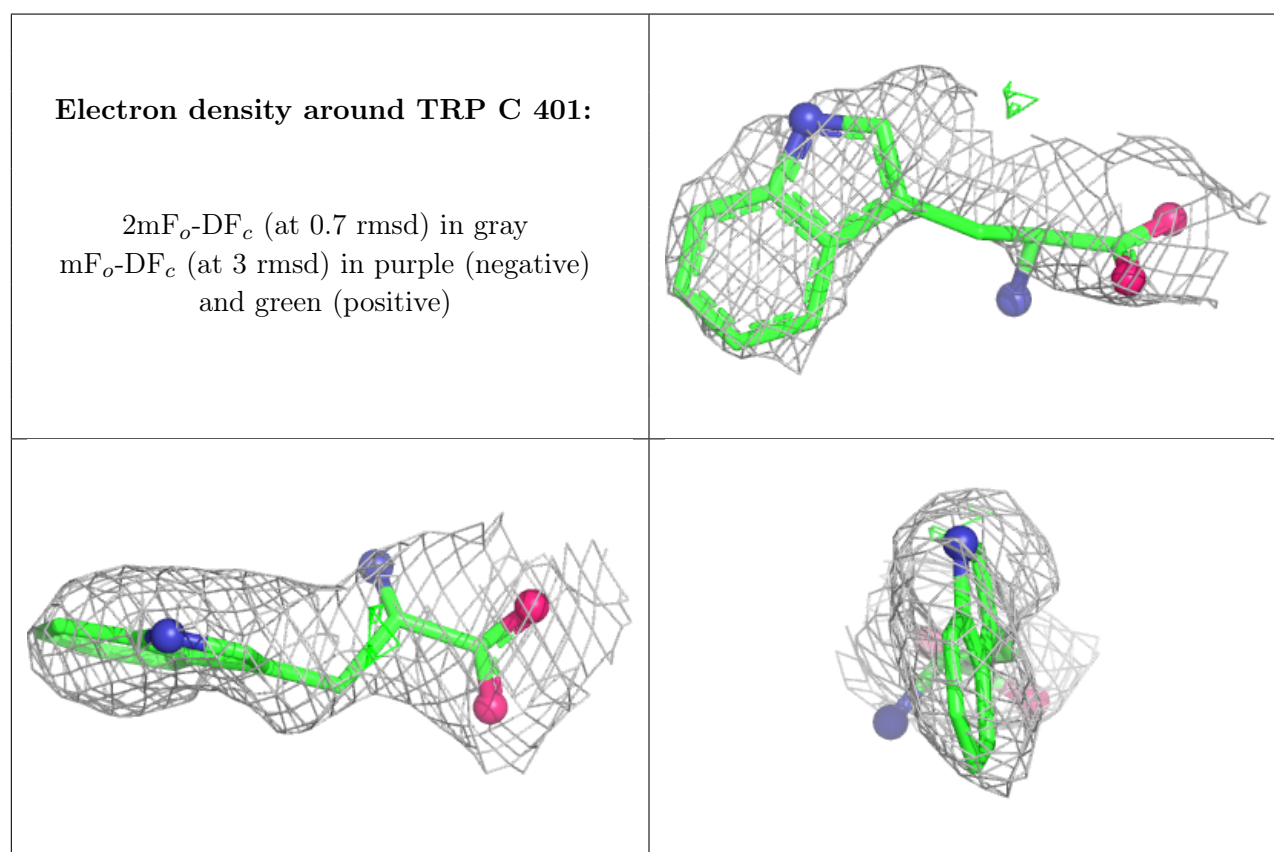
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	TRP	C	401	15/15	0.99	0.04	49,57,67,68	0

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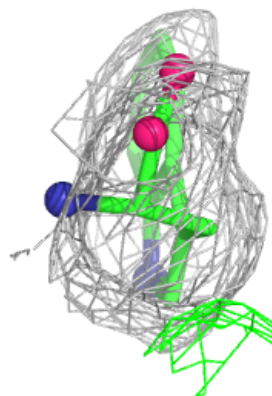
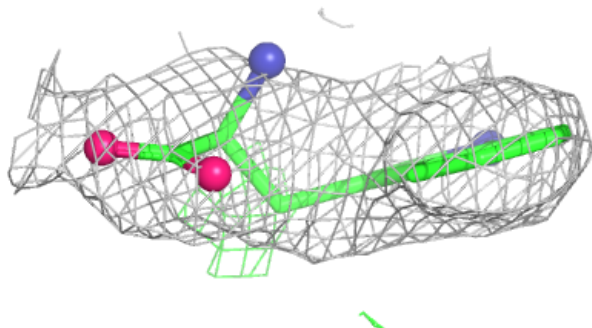
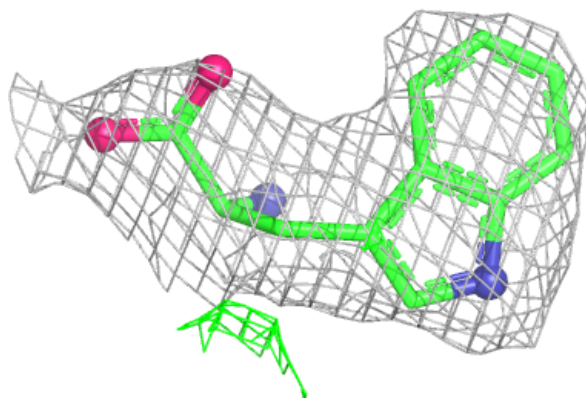
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	TRP	D	401	15/15	0.99	0.03	49,54,57,59	0
2	TRP	E	401	15/15	0.99	0.04	55,59,70,70	0
2	TRP	B	401	15/15	1.00	0.03	35,43,48,50	0
2	TRP	A	401	15/15	1.00	0.04	44,52,59,64	0
2	TRP	F	401	15/15	1.00	0.02	52,56,65,67	0

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.

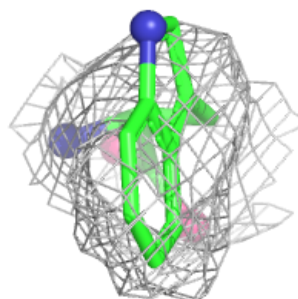
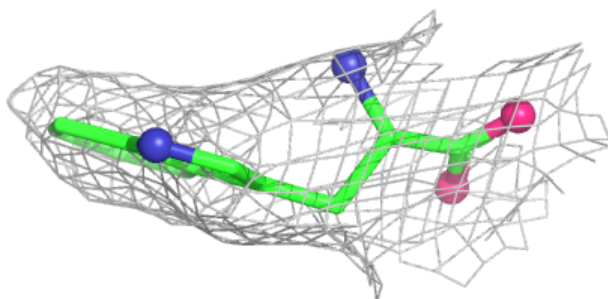
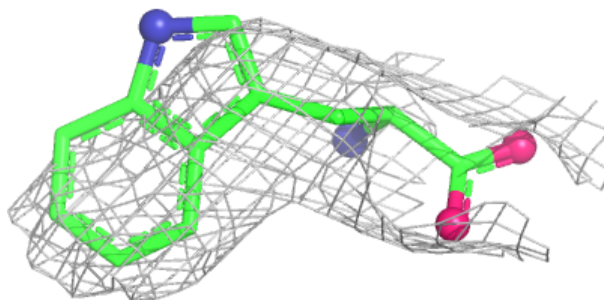


Electron density around TRP D 401:

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

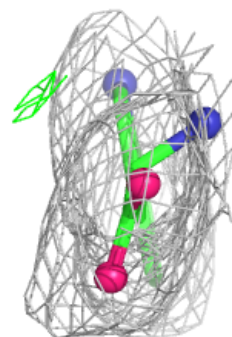
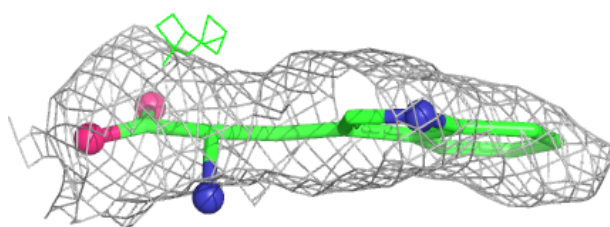
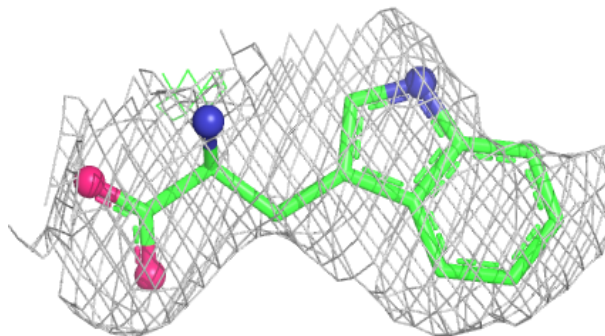
**Electron density around TRP E 401:**

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

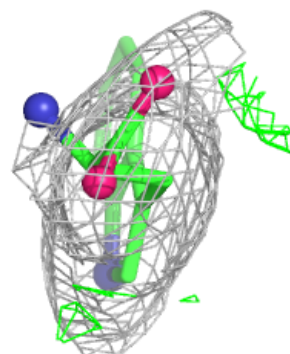
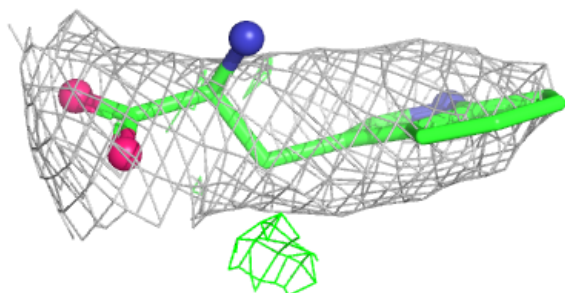
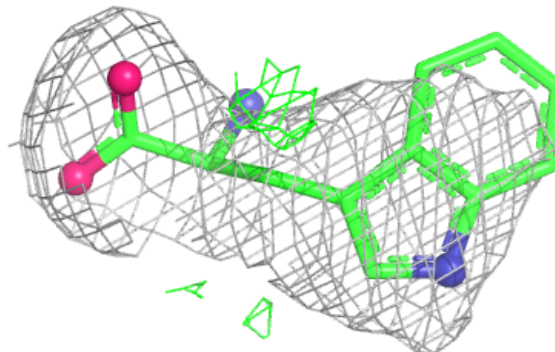


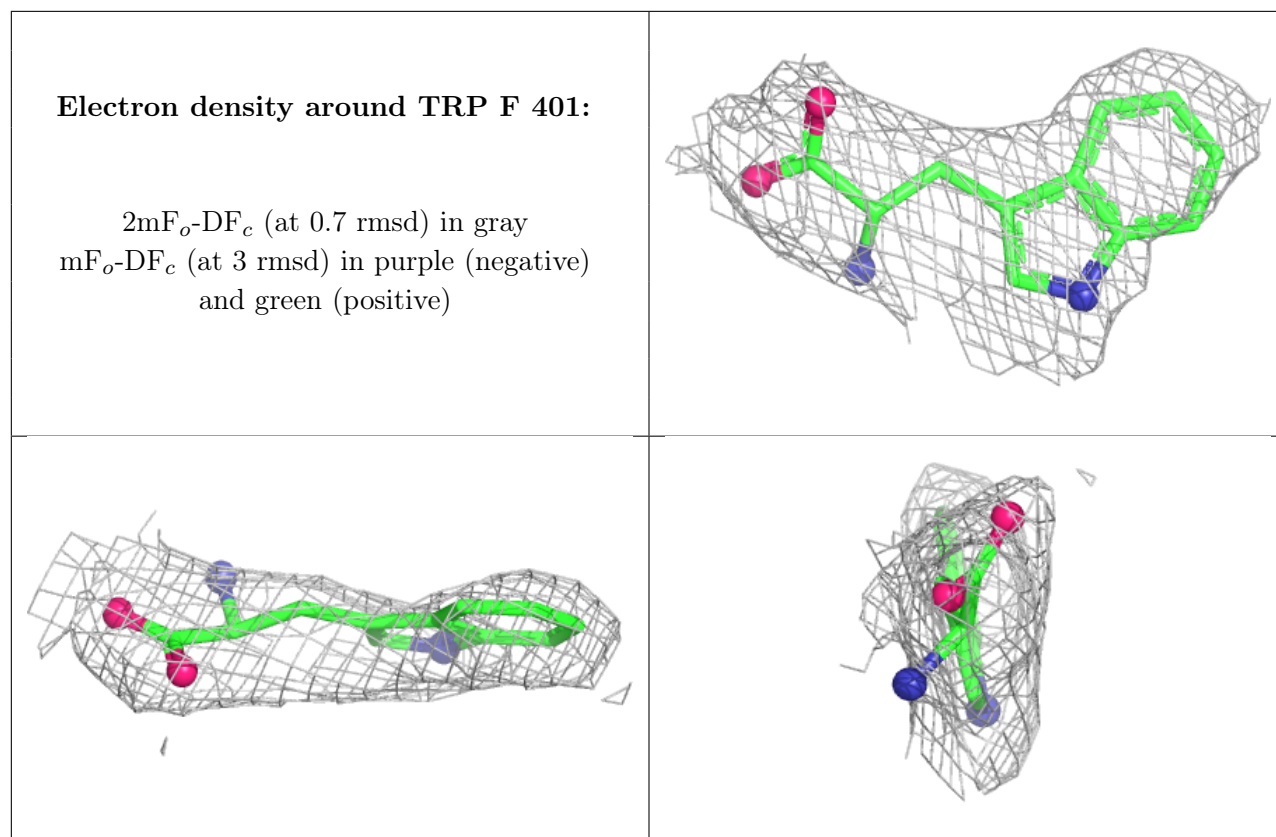
Electron density around TRP B 401:

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

**Electron density around TRP 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.