



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 12:40 pm BST

PDB ID : 6J1L
Title : Crystal Structure Analysis of the ROR gamma(C455E)
Authors : zhang, Y.; Li, C.C.; wu, X.S.
Deposited on : 2018-12-28
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

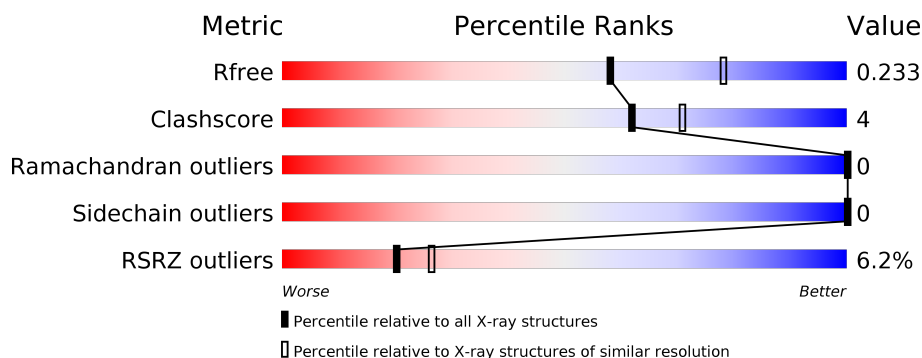
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	267	<div> <div>3%</div> <div> <div></div> <div>74%</div> <div>9%</div> <div>17%</div> </div> </div>
1	B	267	<div> <div>2%</div> <div> <div></div> <div>77%</div> <div>6%</div> <div>17%</div> </div> </div>
1	C	267	<div> <div>11%</div> <div> <div></div> <div>70%</div> <div>12%</div> <div>18%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5609 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 Nuclear receptor ROR-gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	222	Total	C	N	O	S	0	0	0
			1820	1152	328	327	13			
1	B	222	Total	C	N	O	S	0	0	0
			1816	1150	328	325	13			
1	C	219	Total	C	N	O	S	0	0	0
			1788	1130	323	322	13			

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	241	GLU	-	expression tag	UNP P51449
A	242	GLY	-	expression tag	UNP P51449
A	243	ASP	-	expression tag	UNP P51449
A	244	ILE	-	expression tag	UNP P51449
A	245	HIS	-	expression tag	UNP P51449
A	246	MET	-	expression tag	UNP P51449
A	247	LYS	-	expression tag	UNP P51449
A	248	LYS	-	expression tag	UNP P51449
A	249	GLY	-	expression tag	UNP P51449
A	250	TYR	-	expression tag	UNP P51449
A	251	HIS	-	expression tag	UNP P51449
A	252	HIS	-	expression tag	UNP P51449
A	253	HIS	-	expression tag	UNP P51449
A	254	HIS	-	expression tag	UNP P51449
A	255	HIS	-	expression tag	UNP P51449
A	256	LEU	-	expression tag	UNP P51449
A	257	VAL	-	expression tag	UNP P51449
A	258	PRO	-	expression tag	UNP P51449
A	259	ARG	-	expression tag	UNP P51449
A	260	GLY	-	expression tag	UNP P51449
A	261	SER	-	expression tag	UNP P51449
A	455	GLU	CYS	engineered mutation	UNP P51449
B	241	GLU	-	expression tag	UNP P51449

Continued on next page...

Continued from previous page...

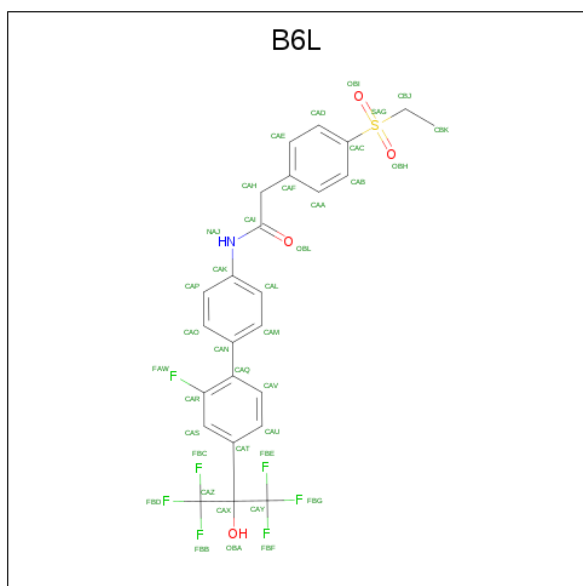
Chain	Residue	Modelled	Actual	Comment	Reference
B	242	GLY	-	expression tag	UNP P51449
B	243	ASP	-	expression tag	UNP P51449
B	244	ILE	-	expression tag	UNP P51449
B	245	HIS	-	expression tag	UNP P51449
B	246	MET	-	expression tag	UNP P51449
B	247	LYS	-	expression tag	UNP P51449
B	248	LYS	-	expression tag	UNP P51449
B	249	GLY	-	expression tag	UNP P51449
B	250	TYR	-	expression tag	UNP P51449
B	251	HIS	-	expression tag	UNP P51449
B	252	HIS	-	expression tag	UNP P51449
B	253	HIS	-	expression tag	UNP P51449
B	254	HIS	-	expression tag	UNP P51449
B	255	HIS	-	expression tag	UNP P51449
B	256	LEU	-	expression tag	UNP P51449
B	257	VAL	-	expression tag	UNP P51449
B	258	PRO	-	expression tag	UNP P51449
B	259	ARG	-	expression tag	UNP P51449
B	260	GLY	-	expression tag	UNP P51449
B	261	SER	-	expression tag	UNP P51449
B	455	GLU	CYS	engineered mutation	UNP P51449
C	241	GLU	-	expression tag	UNP P51449
C	242	GLY	-	expression tag	UNP P51449
C	243	ASP	-	expression tag	UNP P51449
C	244	ILE	-	expression tag	UNP P51449
C	245	HIS	-	expression tag	UNP P51449
C	246	MET	-	expression tag	UNP P51449
C	247	LYS	-	expression tag	UNP P51449
C	248	LYS	-	expression tag	UNP P51449
C	249	GLY	-	expression tag	UNP P51449
C	250	TYR	-	expression tag	UNP P51449
C	251	HIS	-	expression tag	UNP P51449
C	252	HIS	-	expression tag	UNP P51449
C	253	HIS	-	expression tag	UNP P51449
C	254	HIS	-	expression tag	UNP P51449
C	255	HIS	-	expression tag	UNP P51449
C	256	LEU	-	expression tag	UNP P51449
C	257	VAL	-	expression tag	UNP P51449
C	258	PRO	-	expression tag	UNP P51449
C	259	ARG	-	expression tag	UNP P51449
C	260	GLY	-	expression tag	UNP P51449
C	261	SER	-	expression tag	UNP P51449

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	455	GLU	CYS	engineered mutation	UNP P51449

- Molecule 2 is 2-[4-(ethylsulfonyl)phenyl]-N-[2'-fluoro-4'-(1,1,1,3,3,3-hexafluoro-2-hydroxypropyl)-[1,1'-biphenyl]-4-yl]acetamide (three-letter code: B6L) (formula: C₂₅H₂₀F₇NO₄S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	S	0	0
			38	25	7	1	4	1		
2	B	1	Total	C	F	N	O	S	0	0
			38	25	7	1	4	1		
2	C	1	Total	C	F	N	O	S	0	0
			38	25	7	1	4	1		

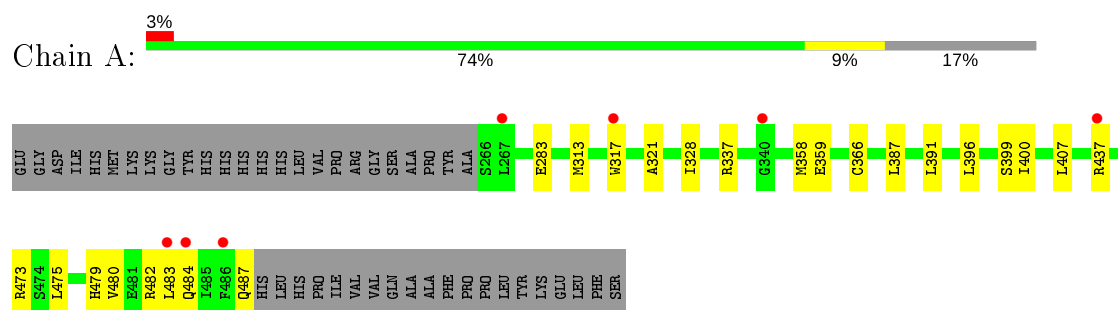
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	29	Total	O	0	0
			29	29		
3	B	36	Total	O	0	0
			36	36		
3	C	6	Total	O	0	0
			6	6		

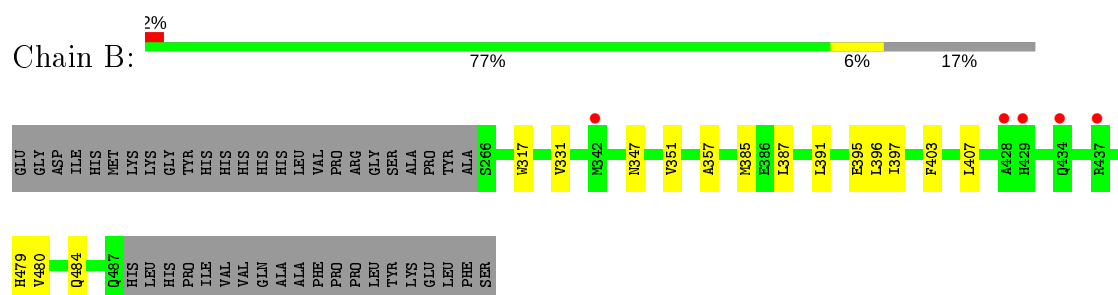
3 Residue-property plots [i](#)

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

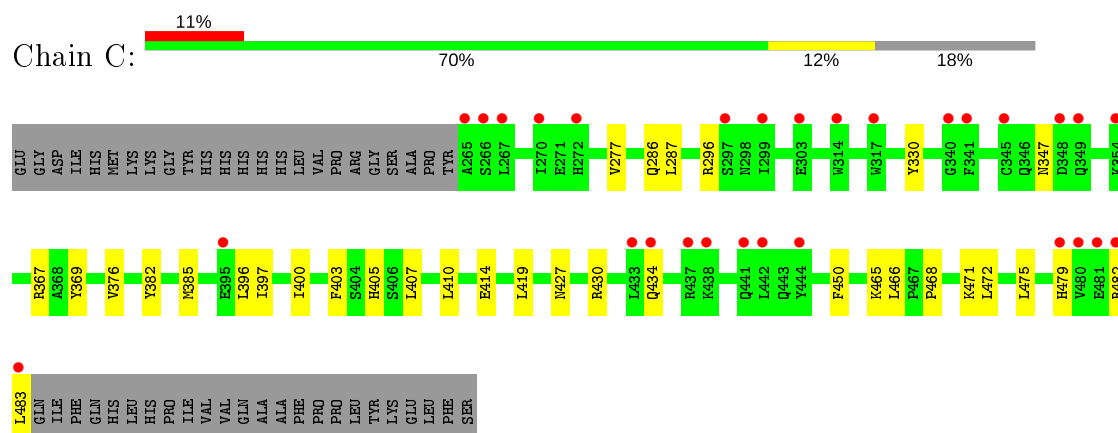
- Molecule 1: Nuclear receptor ROR-gamma



- Molecule 1: Nuclear receptor ROR-gamma



- Molecule 1: Nuclear receptor ROR-gamma



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	110.14Å 110.14Å 114.23Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.61 – 2.30 49.61 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.61-2.30) 100.0 (49.61-2.30)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.202 , 0.226 0.212 , 0.233	Depositor DCC
R_{free} test set	2548 reflections (7.30%)	wwPDB-VP
Wilson B-factor (Å ²)	52.8	Xtriage
Anisotropy	0.131	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.041 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5609	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% 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: B6L

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	0.38	0/1855	0.59	1/2493 (0.0%)
1	B	0.35	0/1851	0.59	1/2488 (0.0%)
1	C	0.31	0/1822	0.55	1/2449 (0.0%)
All	All	0.35	0/5528	0.58	3/7430 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	483	LEU	CA-CB-CG	-6.12	101.22	115.30
1	A	387	LEU	CA-CB-CG	6.08	129.29	115.30
1	B	387	LEU	CA-CB-CG	5.07	126.97	115.30

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	1820	0	1820	17	0
1	B	1816	0	1816	8	0
1	C	1788	0	1789	26	0
2	A	38	0	0	0	0
2	B	38	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	38	0	0	1	0
3	A	29	0	0	0	0
3	B	36	0	0	0	0
3	C	6	0	0	0	0
All	All	5609	0	5425	48	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 4.

All (48) 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:C:468:PRO:HG2	1:C:471:LYS:HB2	1.61	0.83
1:C:286:GLN:HG2	1:C:330:TYR:CD2	2.20	0.77
1:C:403:PHE:CE2	1:C:407:LEU:HD11	2.20	0.77
1:A:480:VAL:O	1:A:484:GLN:HG2	1.93	0.69
1:C:396:LEU:HG	1:C:479:HIS:CE1	2.28	0.69
1:A:399:SER:HB3	1:A:475:LEU:HD11	1.76	0.68
1:B:403:PHE:CE2	1:B:407:LEU:HD11	2.29	0.67
1:A:313:MET:O	1:A:317:TRP:HD1	1.84	0.60
1:C:367:ARG:NH1	1:C:414:GLU:OE1	2.36	0.59
1:B:395:GLU:HG2	1:B:396:LEU:N	2.18	0.58
1:A:317:TRP:CE2	1:A:391:LEU:HD22	2.38	0.58
1:C:347:ASN:ND2	1:C:434:GLN:OE1	2.37	0.58
1:A:396:LEU:HD22	1:A:482:ARG:HG3	1.86	0.57
1:C:400:ILE:HD11	1:C:479:HIS:NE2	2.20	0.56
1:A:313:MET:O	1:A:317:TRP:CD1	2.58	0.56
1:C:286:GLN:HG2	1:C:330:TYR:CG	2.41	0.55
1:B:317:TRP:CE2	1:B:391:LEU:HD22	2.43	0.54
1:B:480:VAL:O	1:B:484:GLN:HG2	2.09	0.53
1:A:359:GLU:OE2	1:A:473:ARG:HD3	2.10	0.52
1:A:437:ARG:NH1	1:C:466:LEU:O	2.44	0.51
1:C:287:LEU:O	1:C:367:ARG:NH2	2.37	0.51
1:C:397:ILE:HG22	2:C:601:B6L:FBF	2.00	0.51
1:A:400:ILE:HD11	1:A:479:HIS:NE2	2.27	0.50
1:C:410:LEU:HD21	1:C:465:LYS:HB3	1.94	0.50
1:A:321:ALA:HB2	1:A:483:LEU:HD11	1.94	0.49
1:C:434:GLN:N	1:C:434:GLN:OE1	2.44	0.49
1:C:475:LEU:O	1:C:479:HIS:CD2	2.66	0.48
1:C:385:MET:HB2	1:C:397:ILE:HD13	1.96	0.48
1:B:385:MET:HG2	1:B:397:ILE:HG22	1.96	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:427:ASN:HD22	1:C:430:ARG:NH1	2.13	0.46
1:C:396:LEU:HD22	1:C:482:ARG:HE	1.80	0.46
1:C:296:ARG:HA	1:C:382:TYR:CE1	2.52	0.45
1:A:483:LEU:O	1:A:487:GLN:N	2.50	0.45
1:B:347:ASN:O	1:B:351:VAL:HG23	2.17	0.45
1:A:400:ILE:HD11	1:A:479:HIS:CE1	2.52	0.45
1:B:331:VAL:HG11	1:B:357:ALA:HA	1.99	0.45
1:A:366:CYS:HB3	1:A:407:LEU:HD23	2.00	0.44
1:A:437:ARG:HH12	1:C:468:PRO:HD3	1.82	0.44
1:A:328:ILE:HD11	1:A:358:MET:HG2	2.00	0.43
1:C:396:LEU:HD21	1:C:482:ARG:HG3	2.00	0.43
1:B:396:LEU:HD22	1:B:479:HIS:CE1	2.53	0.43
1:A:437:ARG:NH1	1:C:468:PRO:HD3	2.34	0.42
1:C:419:LEU:HD13	1:C:450:PHE:HA	2.02	0.42
1:A:283:GLU:OE1	1:A:337:ARG:NH1	2.53	0.41
1:C:277:VAL:HG11	1:C:419:LEU:HD23	2.02	0.41
1:C:376:VAL:O	1:C:382:TYR:HA	2.21	0.41
1:C:369:TYR:CE1	1:C:405:HIS:HA	2.56	0.40
1:C:403:PHE:CZ	1:C:472:LEU:HD13	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	220/267 (82%)	217 (99%)	3 (1%)	0	100	100
1	B	220/267 (82%)	217 (99%)	3 (1%)	0	100	100
1	C	217/267 (81%)	215 (99%)	2 (1%)	0	100	100
All	All	657/801 (82%)	649 (99%)	8 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	199/237 (84%)	199 (100%)	0	100	100
1	B	198/237 (84%)	198 (100%)	0	100	100
1	C	195/237 (82%)	195 (100%)	0	100	100
All	All	592/711 (83%)	592 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	347	ASN
1	C	479	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

3 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	B6L	A	601	-	40,40,40	3.29	6 (15%)	61,62,62	1.80	16 (26%)
2	B6L	B	601	-	40,40,40	3.36	6 (15%)	61,62,62	1.91	15 (24%)
2	B6L	C	601	-	40,40,40	3.42	7 (17%)	61,62,62	1.69	11 (18%)

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	B6L	A	601	-	-	2/45/45/45	0/3/3/3
2	B6L	B	601	-	-	0/45/45/45	0/3/3/3
2	B6L	C	601	-	-	0/45/45/45	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	B6L	OBH-SAG	14.14	1.65	1.44
2	C	601	B6L	OBI-SAG	13.99	1.64	1.44
2	B	601	B6L	OBI-SAG	13.78	1.64	1.44
2	B	601	B6L	OBH-SAG	13.75	1.64	1.44
2	A	601	B6L	OBI-SAG	13.66	1.64	1.44
2	A	601	B6L	OBH-SAG	13.47	1.64	1.44
2	C	601	B6L	CAI-NAJ	4.37	1.45	1.35
2	B	601	B6L	CAI-NAJ	4.26	1.45	1.35
2	A	601	B6L	CAI-NAJ	4.20	1.44	1.35
2	C	601	B6L	CBJ-SAG	3.81	1.82	1.77
2	B	601	B6L	CBJ-SAG	3.46	1.82	1.77
2	C	601	B6L	CAC-SAG	3.29	1.82	1.76
2	A	601	B6L	CBJ-SAG	3.25	1.81	1.77
2	B	601	B6L	CAC-SAG	2.94	1.81	1.76
2	A	601	B6L	CAC-SAG	2.56	1.81	1.76
2	B	601	B6L	CAY-CAX	-2.31	1.50	1.54
2	A	601	B6L	OBL-CAI	-2.20	1.18	1.23
2	C	601	B6L	CAK-NAJ	2.14	1.46	1.41
2	C	601	B6L	OBL-CAI	-2.07	1.19	1.23

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	601	B6L	OBI-SAG-OBH	-7.12	110.12	118.44
2	B	601	B6L	OBI-SAG-OBH	-6.35	111.01	118.44
2	A	601	B6L	CAZ-CAX-CAY	4.78	116.26	110.28
2	B	601	B6L	CAF-CAH-CAI	-4.53	99.13	112.57
2	B	601	B6L	FBB-CAZ-CAX	-4.11	104.00	111.85
2	A	601	B6L	CBJ-SAG-CAC	3.91	110.34	105.02
2	B	601	B6L	FBE-CAY-CAX	-3.87	104.46	111.85
2	A	601	B6L	CAV-CAQ-CAR	3.72	121.47	116.10
2	A	601	B6L	OBI-SAG-OBH	-3.71	114.10	118.44
2	A	601	B6L	CAF-CAH-CAI	-3.52	102.15	112.57
2	C	601	B6L	CAV-CAQ-CAR	3.41	121.02	116.10
2	B	601	B6L	CBJ-SAG-CAC	3.36	109.59	105.02
2	C	601	B6L	CAF-CAH-CAI	-3.34	102.67	112.57
2	B	601	B6L	CAD-CAC-SAG	3.32	123.12	119.49
2	B	601	B6L	OBH-SAG-CAC	3.30	111.72	108.36
2	C	601	B6L	CAS-CAR-CAQ	-3.16	118.66	123.64
2	A	601	B6L	CAO-CAN-CAQ	-3.02	116.02	120.91
2	A	601	B6L	CAS-CAR-CAQ	-3.00	118.92	123.64
2	B	601	B6L	CAB-CAC-SAG	-2.90	116.32	119.49
2	B	601	B6L	CAV-CAQ-CAR	2.78	120.11	116.10
2	A	601	B6L	OBH-SAG-CBJ	-2.74	105.86	108.29
2	B	601	B6L	OBI-SAG-CAC	-2.65	105.67	108.36
2	A	601	B6L	FBD-CAZ-CAX	-2.60	106.88	111.85
2	B	601	B6L	CAO-CAN-CAQ	-2.59	116.71	120.91
2	A	601	B6L	FBB-CAZ-CAX	-2.57	106.94	111.85
2	B	601	B6L	CAS-CAR-CAQ	-2.53	119.66	123.64
2	A	601	B6L	CAB-CAC-SAG	-2.49	116.76	119.49
2	A	601	B6L	FBG-CAY-CAX	-2.36	107.35	111.85
2	C	601	B6L	CAK-NAJ-CAI	-2.35	123.38	127.50
2	A	601	B6L	FBF-CAY-CAX	-2.35	107.36	111.85
2	A	601	B6L	CAD-CAC-SAG	2.33	122.03	119.49
2	C	601	B6L	CAH-CAF-CAE	-2.32	117.57	120.89
2	C	601	B6L	CAZ-CAX-CAY	2.30	113.16	110.28
2	B	601	B6L	FBG-CAY-FBE	2.27	114.48	107.53
2	B	601	B6L	FBF-CAY-CAX	-2.14	107.76	111.85
2	C	601	B6L	FBF-CAY-CAX	-2.11	107.83	111.85
2	A	601	B6L	CAO-CAN-CAM	2.11	121.79	117.59
2	B	601	B6L	CAH-CAF-CAA	-2.07	117.93	120.89
2	A	601	B6L	OBH-SAG-CAC	2.05	110.45	108.36
2	C	601	B6L	CAO-CAN-CAM	2.03	121.64	117.59
2	C	601	B6L	OBI-SAG-CAC	2.02	110.42	108.36
2	C	601	B6L	CBJ-SAG-CAC	2.01	107.75	105.02

There are no chirality outliers.

All (2) torsion outliers are listed below:

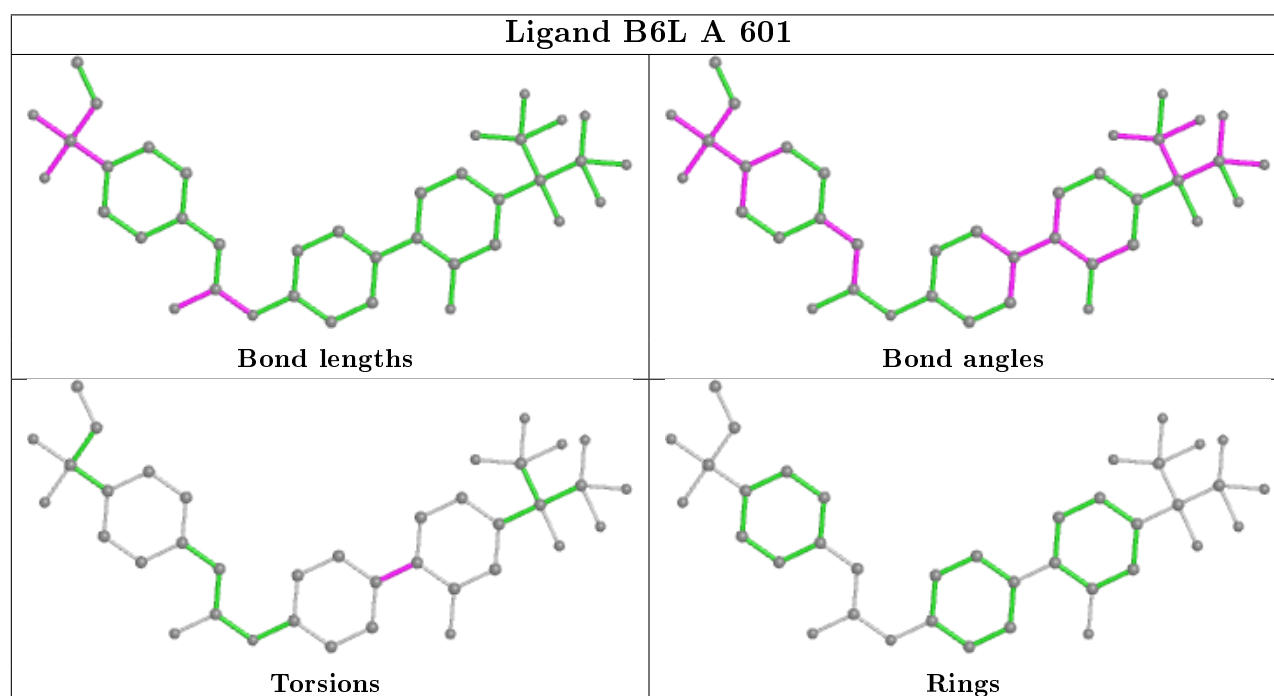
Mol	Chain	Res	Type	Atoms
2	A	601	B6L	CAM-CAN-CAQ-CAR
2	A	601	B6L	CAO-CAN-CAQ-CAR

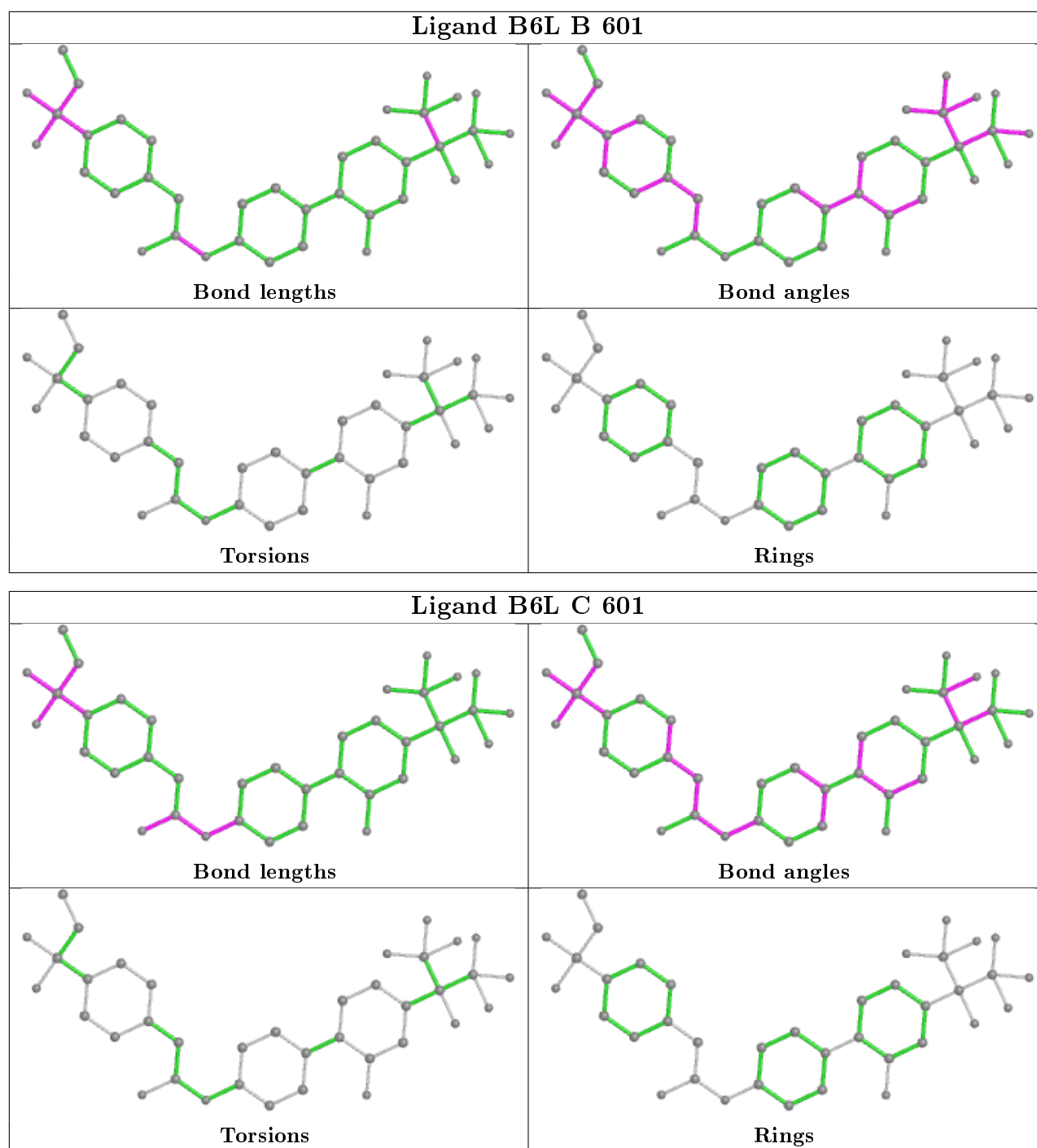
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	601	B6L	1	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	222/267 (83%)	0.21	7 (3%) 47 54	36, 60, 100, 139	0
1	B	222/267 (83%)	0.21	5 (2%) 60 67	37, 56, 92, 119	0
1	C	219/267 (82%)	0.69	29 (13%) 3 4	62, 84, 116, 170	0
All	All	663/801 (82%)	0.37	41 (6%) 20 26	36, 69, 107, 170	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	479	HIS	6.4
1	C	480	VAL	5.1
1	C	340	GLY	4.7
1	C	317	TRP	4.7
1	C	481	GLU	4.3
1	C	441	GLN	4.2
1	A	486	PHE	4.0
1	C	438	LYS	4.0
1	A	317	TRP	3.8
1	C	270	ILE	3.8
1	B	434	GLN	3.6
1	C	272	HIS	3.6
1	A	267	LEU	3.5
1	C	341	PHE	3.5
1	B	428	ALA	3.3
1	C	314	TRP	3.0
1	A	340	GLY	3.0
1	C	437	ARG	2.9
1	C	265	ALA	2.9
1	B	429	HIS	2.8
1	C	267	LEU	2.8
1	C	444	TYR	2.7
1	C	433	LEU	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	484	GLN	2.6
1	C	483	LEU	2.5
1	C	354	LYS	2.5
1	C	266	SER	2.5
1	C	297	SER	2.4
1	B	437	ARG	2.4
1	A	483	LEU	2.4
1	C	482	ARG	2.3
1	A	437	ARG	2.3
1	C	345	CYS	2.3
1	C	348	ASP	2.3
1	C	349	GLN	2.2
1	C	299	ILE	2.2
1	C	442	LEU	2.2
1	C	395	GLU	2.1
1	B	342	MET	2.1
1	C	303	GLU	2.1
1	C	434	GLN	2.1

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 carbohydrates 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.

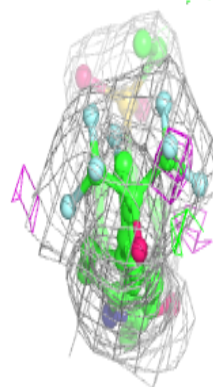
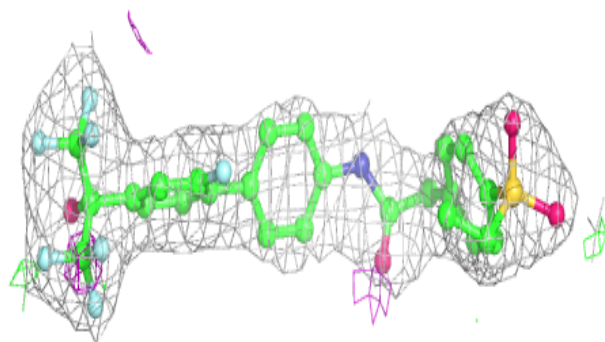
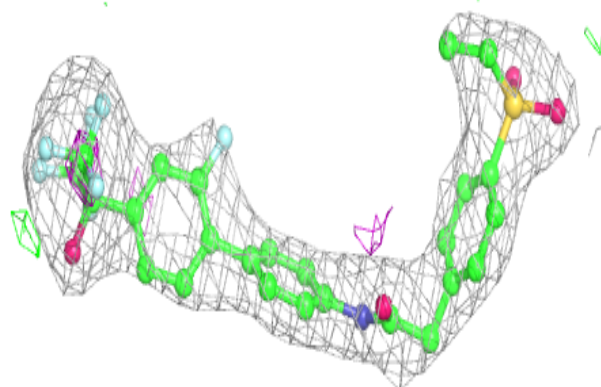
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	B6L	C	601	38/38	0.93	0.15	61,71,85,86	0
2	B6L	A	601	38/38	0.96	0.15	36,48,72,74	0
2	B6L	B	601	38/38	0.97	0.15	33,43,57,61	0

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.

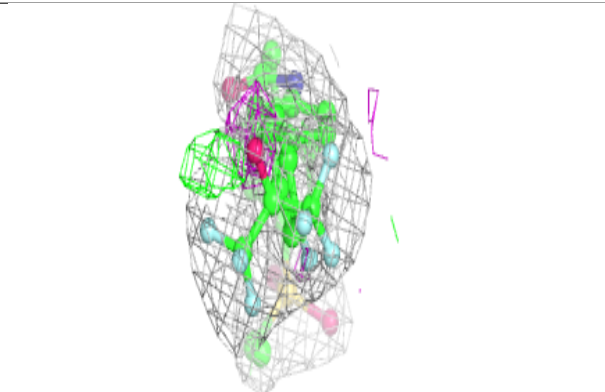
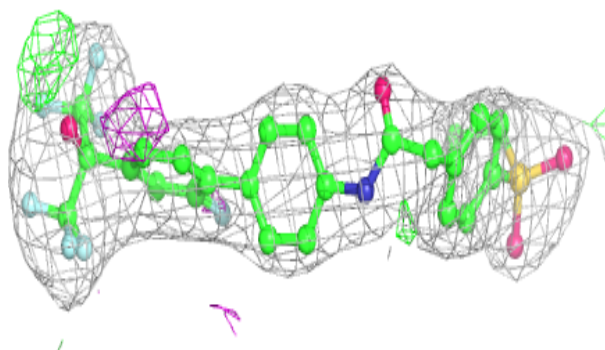
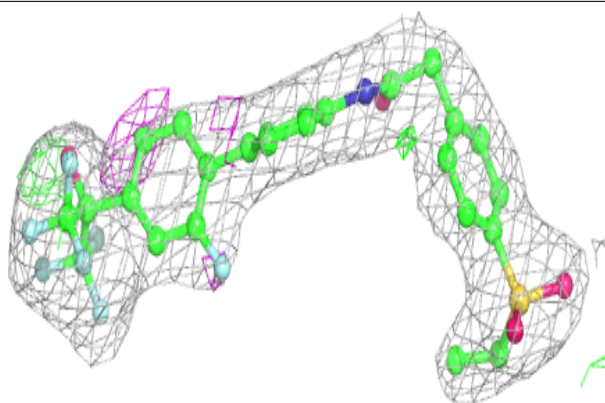
Electron density around B6L C 601:

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

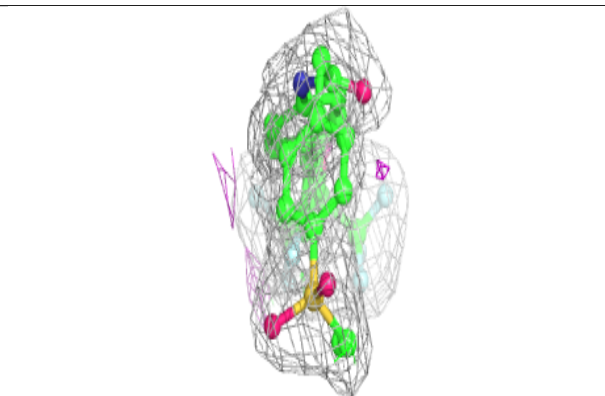
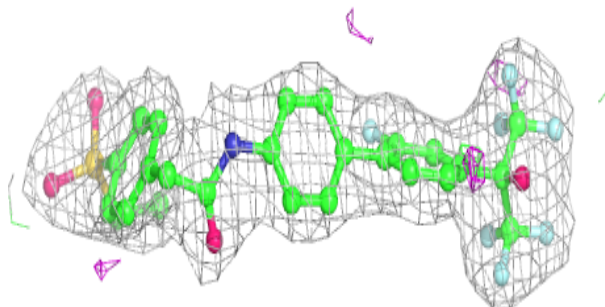
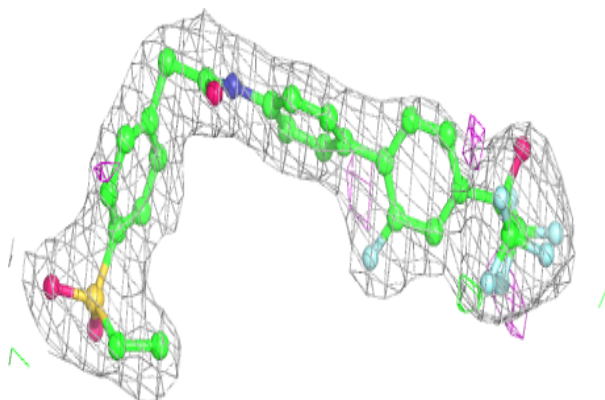


Electron density around B6L A 601:

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

**Electron density around B6L B 601:**

$2mF_o-DF_c$ (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.