



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 01:44 am BST

PDB ID : 5LAY
Title : Discovery of New Natural-product-inspired Spiro-oxindole Compounds as Orally Active Inhibitors of the MDM2-p53 Interaction: HDM2 (MDM2) IN COMPLEX WITH COMPOUND 6g
Authors : Kessler, D.; Gollner, A.
Deposited on : 2016-06-15
Resolution : 2.71 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

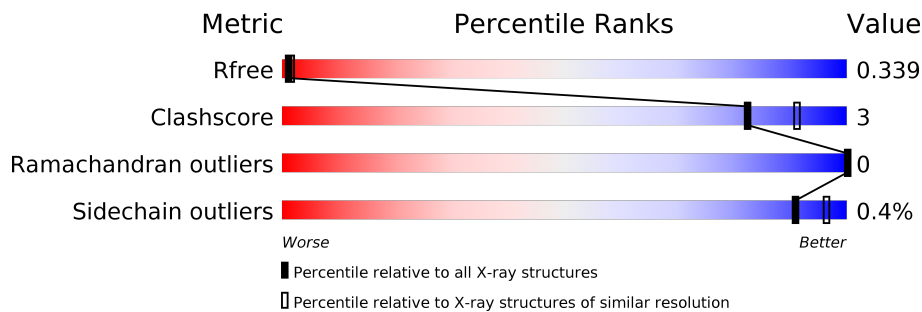
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.71 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	3359 (2.74-2.70)
Clashscore	141614	3686 (2.74-2.70)
Ramachandran outliers	138981	3622 (2.74-2.70)
Sidechain outliers	138945	3623 (2.74-2.70)

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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	95	94% 5% .
1	B	95	95% . .
1	C	95	94% 5% .
1	D	95	86% 13% .
1	E	95	93% 6% .
1	F	95	79% 6% 15%

2 Entry composition [i](#)

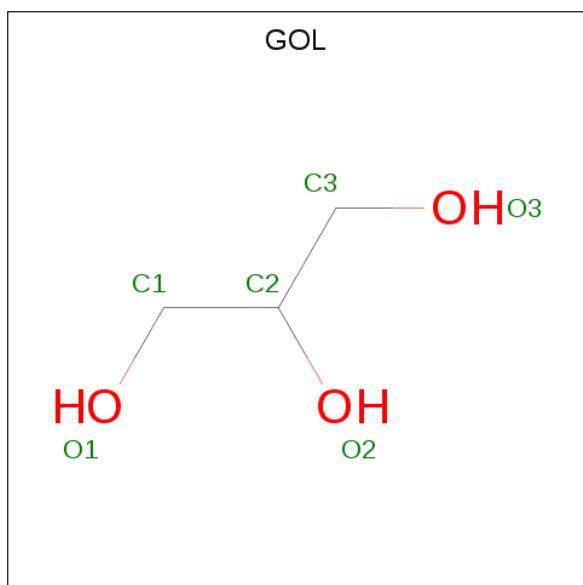
There are 5 unique types of molecules in this entry. The entry contains 4950 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 E3 ubiquitin-protein ligase Mdm2.

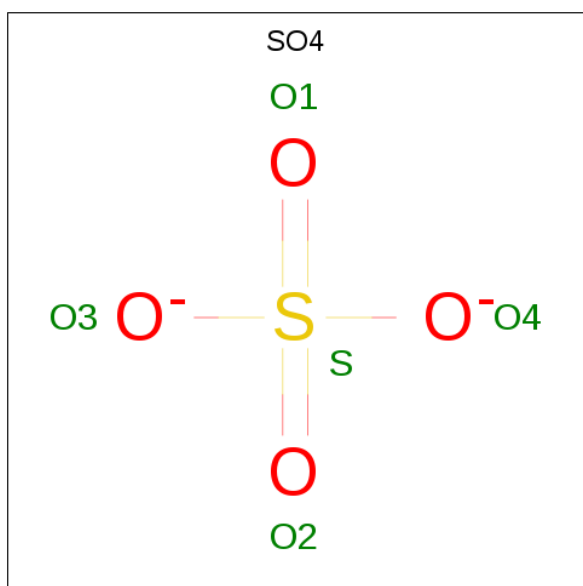
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	95	Total 801	C 518	N 137	O 142	S 4	11	1	0
1	B	94	Total 796	C 517	N 132	O 143	S 4	12	3	0
1	C	94	Total 773	C 503	N 128	O 138	S 4	31	0	0
1	D	94	Total 771	C 502	N 127	O 138	S 4	11	0	0
1	E	94	Total 774	C 503	N 128	O 139	S 4	46	0	0
1	F	81	Total 677	C 441	N 111	O 121	S 4	108	0	0

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



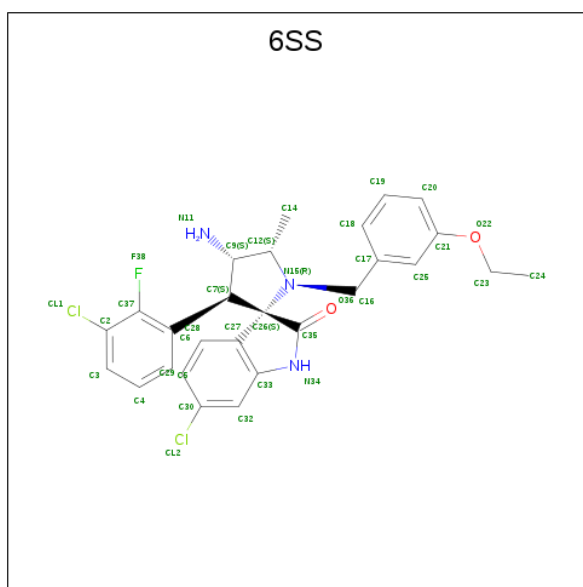
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	E	1	Total C O 6 3 3	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	E	1	Total O S 5 4 1	0	0

- Molecule 4 is (3 {S},3' {S},4' {S},5' {S})-4'-azanyl-6-chloranyl-3'-(3-chloranyl-2-fluoranyl-phenyl)-1'-[(3-ethoxyphenyl)methyl]-5'-methyl-spiro[1 {H}-indole-3,2'-pyrrolidine]-2-one (three-letter code: 6SS) (formula: C₂₇H₂₆Cl₂FN₃O₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	Cl	F	N			O
4	A	1	Total 35	C 27	Cl 2	F 1	N 3	O 2	0	0
4	B	1	Total 35	C 27	Cl 2	F 1	N 3	O 2	0	0
4	C	1	Total 35	C 27	Cl 2	F 1	N 3	O 2	0	0
4	D	1	Total 35	C 27	Cl 2	F 1	N 3	O 2	0	0
4	E	1	Total 35	C 27	Cl 2	F 1	N 3	O 2	0	0
4	F	1	Total 35	C 27	Cl 2	F 1	N 3	O 2	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	26	Total 26	O 26	0	0
5	B	23	Total 23	O 23	0	0
5	C	14	Total 14	O 14	0	0
5	D	26	Total 26	O 26	0	0
5	E	6	Total 6	O 6	0	0
5	F	3	Total 3	O 3	0	0

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. 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. 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.

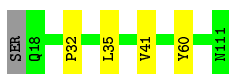
- Molecule 1: E3 ubiquitin-protein ligase Mdm2

Chain A:  94% 5%



- Molecule 1: E3 ubiquitin-protein ligase Mdm2

Chain B:  95%




- Molecule 1: E3 ubiquitin-protein ligase Mdm2

Chain C:  94% 5%



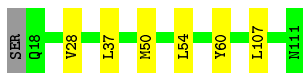
- Molecule 1: E3 ubiquitin-protein ligase Mdm2

Chain D:  86% 13%




- Molecule 1: E3 ubiquitin-protein ligase Mdm2

Chain E:  93% 6%



- Molecule 1: E3 ubiquitin-protein ligase Mdm2

Chain F:  79% 6% 15%

SER	GLN	ILE	PRO	ALA	SER	E23	Q24	K39	A43	Y48	Y56	S92	VAL	LYS	E96	Y104	R105	ASN	LEU	VAL	VAL	VAL	ASN
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4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	71.45Å 71.45Å 520.98Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	87.45 – 2.71 44.86 – 2.71	Depositor EDS
% Data completeness (in resolution range)	99.0 (87.45-2.71) 99.0 (44.86-2.71)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.94 (at 2.73Å)	Xtrriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.251 , 0.292 0.319 , 0.339	Depositor DCC
R_{free} test set	806 reflections (3.55%)	wwPDB-VP
Wilson B-factor (Å ²)	25.4	Xtrriage
Anisotropy	0.435	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 32.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.79	EDS
Total number of atoms	4950	wwPDB-VP
Average B, all atoms (Å ²)	8.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% 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: GOL, 6SS, SO4

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	0.42	0/816	0.54	0/1101
1	B	0.42	0/811	0.58	0/1095
1	C	0.48	0/788	0.56	0/1063
1	D	0.41	0/786	0.56	0/1060
1	E	0.43	0/789	0.54	0/1063
1	F	0.65	2/690 (0.3%)	0.80	3/926 (0.3%)
All	All	0.47	2/4680 (0.0%)	0.60	3/6308 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	104	TYR	CG-CD1	8.86	1.50	1.39
1	F	104	TYR	CG-CD2	-7.22	1.29	1.39

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	104	TYR	CB-CG-CD1	-11.13	114.32	121.00
1	F	104	TYR	CB-CG-CD2	11.09	127.65	121.00
1	F	56	TYR	CB-CG-CD1	5.86	124.52	121.00

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	801	0	836	5	0
1	B	796	0	826	4	0
1	C	773	0	805	5	0
1	D	771	0	804	8	0
1	E	774	0	805	3	0
1	F	677	0	696	2	0
2	A	18	0	24	0	0
2	C	6	0	8	1	0
2	E	6	0	8	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	D	5	0	0	0	0
3	E	5	0	0	0	0
4	A	35	0	0	0	0
4	B	35	0	0	0	0
4	C	35	0	0	0	0
4	D	35	0	0	0	0
4	E	35	0	0	0	0
4	F	35	0	0	0	0
5	A	26	0	0	0	0
5	B	23	0	0	0	0
5	C	14	0	0	0	0
5	D	26	0	0	0	0
5	E	6	0	0	0	0
5	F	3	0	0	0	0
All	All	4950	0	4812	23	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 (23) 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:18:GLN:N	2:C:201:GOL:HO2	1.95	0.64
1:C:37:LEU:HD13	1:C:85:LEU:HD21	1.88	0.56
1:D:41:VAL:HG11	1:D:60:TYR:HA	1.89	0.54
1:F:39:LYS:HA	1:F:43:ALA:HB3	1.90	0.53
1:D:38:LEU:HD22	1:D:56:TYR:HB2	1.92	0.52
1:F:43:ALA:HB1	1:F:48:TYR:OH	2.10	0.51
1:B:41:VAL:HG11	1:B:60:TYR:HA	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:43:ALA:HB1	1:D:48:TYR:OH	2.11	0.49
1:D:19:ILE:HG21	1:D:100:TYR:CZ	2.47	0.49
1:E:50:MET:O	1:E:54:LEU:HD23	2.14	0.48
1:D:54:LEU:HD11	1:D:100:TYR:CE1	2.49	0.47
1:D:54:LEU:HD11	1:D:100:TYR:CZ	2.50	0.47
1:A:78:SER:HB3	1:B:35:LEU:HD23	1.97	0.46
1:C:78:SER:HB2	1:D:35:LEU:HD23	1.99	0.44
1:E:28:VAL:HB	1:E:107:LEU:HD13	2.00	0.44
1:E:37:LEU:HD21	1:E:60:TYR:CG	2.53	0.43
1:C:37:LEU:CD1	1:C:85:LEU:HD21	2.48	0.43
1:A:66:LEU:HD22	1:B:32:PRO:HG2	2.00	0.42
1:A:28:VAL:HB	1:A:107:LEU:HD13	2.01	0.42
1:C:85:LEU:HD12	1:C:86:PHE:HD1	1.84	0.42
1:A:79:ASN:HB2	1:B:35:LEU:HD22	2.02	0.41
1:D:28:VAL:HB	1:D:107:LEU:HD13	2.03	0.40
1:A:37:LEU:C	1:A:37:LEU:HD23	2.41	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	95/95 (100%)	92 (97%)	3 (3%)	0	100	100
1	B	95/95 (100%)	93 (98%)	2 (2%)	0	100	100
1	C	92/95 (97%)	89 (97%)	3 (3%)	0	100	100
1	D	92/95 (97%)	90 (98%)	2 (2%)	0	100	100
1	E	92/95 (97%)	88 (96%)	4 (4%)	0	100	100
1	F	77/95 (81%)	72 (94%)	5 (6%)	0	100	100
All	All	543/570 (95%)	524 (96%)	19 (4%)	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	91/89 (102%)	90 (99%)	1 (1%)	73	89
1	B	91/89 (102%)	91 (100%)	0	100	100
1	C	88/89 (99%)	88 (100%)	0	100	100
1	D	88/89 (99%)	88 (100%)	0	100	100
1	E	88/89 (99%)	88 (100%)	0	100	100
1	F	76/89 (85%)	75 (99%)	1 (1%)	69	86
All	All	522/534 (98%)	520 (100%)	2 (0%)	91	96

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	SER
1	F	24	GLN

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

Mol	Chain	Res	Type
1	B	18	GLN
1	F	79	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

15 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
3	SO4	B	201	-	4,4,4	0.34	0	6,6,6	0.08	0
2	GOL	A	203	-	5,5,5	0.34	0	5,5,5	0.23	0
3	SO4	D	201	-	4,4,4	0.34	0	6,6,6	0.09	0
3	SO4	E	202	-	4,4,4	0.36	0	6,6,6	0.18	0
4	6SS	A	205	-	39,39,39	1.24	2 (5%)	48,59,59	1.88	9 (18%)
4	6SS	B	202	-	39,39,39	1.21	2 (5%)	48,59,59	1.95	11 (22%)
3	SO4	A	204	-	4,4,4	0.33	0	6,6,6	0.07	0
4	6SS	D	202	-	39,39,39	1.17	2 (5%)	48,59,59	1.75	9 (18%)
2	GOL	A	202	-	5,5,5	0.28	0	5,5,5	0.18	0
2	GOL	A	201	-	5,5,5	0.43	0	5,5,5	0.43	0
2	GOL	C	201	-	5,5,5	0.30	0	5,5,5	0.29	0
2	GOL	E	201	-	5,5,5	0.31	0	5,5,5	0.35	0
4	6SS	F	201	-	39,39,39	1.17	3 (7%)	48,59,59	1.60	7 (14%)
4	6SS	E	203	-	39,39,39	1.09	2 (5%)	48,59,59	1.69	9 (18%)
4	6SS	C	202	-	39,39,39	1.07	2 (5%)	48,59,59	1.66	11 (22%)

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	GOL	A	203	-	-	2/4/4/4	-
4	6SS	A	205	-	-	2/11/51/51	0/5/5/5
4	6SS	B	202	-	-	2/11/51/51	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	6SS	F	201	-	-	2/11/51/51	0/5/5/5
4	6SS	D	202	-	-	2/11/51/51	0/5/5/5
2	GOL	A	202	-	-	2/4/4/4	-
2	GOL	A	201	-	-	2/4/4/4	-
2	GOL	C	201	-	-	2/4/4/4	-
2	GOL	E	201	-	-	0/4/4/4	-
4	6SS	E	203	-	-	3/11/51/51	0/5/5/5
4	6SS	C	202	-	-	2/11/51/51	0/5/5/5

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	205	6SS	C26-C35	-5.87	1.51	1.55
4	B	202	6SS	C26-C35	-5.31	1.51	1.55
4	E	203	6SS	C26-C35	-5.14	1.51	1.55
4	D	202	6SS	C26-C35	-5.01	1.51	1.55
4	C	202	6SS	C26-C35	-4.54	1.52	1.55
4	F	201	6SS	C26-C35	-4.37	1.52	1.55
4	F	201	6SS	C12-N15	3.38	1.51	1.47
4	B	202	6SS	C12-N15	3.23	1.51	1.47
4	C	202	6SS	C12-N15	2.69	1.50	1.47
4	A	205	6SS	C12-N15	2.66	1.50	1.47
4	E	203	6SS	C12-N15	2.56	1.50	1.47
4	D	202	6SS	C12-N15	2.30	1.50	1.47
4	F	201	6SS	C6-C37	2.05	1.41	1.38

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	205	6SS	C27-C26-C35	7.71	104.92	101.59
4	B	202	6SS	C27-C26-C35	7.40	104.79	101.59
4	E	203	6SS	C27-C26-C35	6.87	104.56	101.59
4	D	202	6SS	C27-C26-C35	6.56	104.43	101.59
4	C	202	6SS	C27-C26-C35	5.56	103.99	101.59
4	F	201	6SS	C27-C26-C35	5.47	103.95	101.59
4	B	202	6SS	C26-C27-C33	-5.43	105.74	108.66
4	A	205	6SS	C26-C27-C33	-5.15	105.89	108.66
4	D	202	6SS	C26-C27-C33	-4.55	106.21	108.66
4	E	203	6SS	C26-C27-C33	-4.41	106.29	108.66
4	B	202	6SS	O36-C35-C26	4.00	128.49	125.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	201	6SS	C26-C27-C33	-3.71	106.67	108.66
4	F	201	6SS	C7-C9-C12	-3.39	100.16	105.13
4	C	202	6SS	C26-C27-C33	-3.39	106.84	108.66
4	F	201	6SS	O36-C35-C26	3.23	127.92	125.48
4	D	202	6SS	C37-C2-CL1	3.15	120.71	119.50
4	D	202	6SS	O36-C35-C26	3.09	127.81	125.48
4	A	205	6SS	O36-C35-C26	3.06	127.78	125.48
4	A	205	6SS	C3-C2-C37	3.00	120.94	119.29
4	C	202	6SS	C37-C2-CL1	-3.00	118.34	119.50
4	C	202	6SS	C27-C26-N15	-2.99	108.99	116.44
4	C	202	6SS	O36-C35-C26	2.87	127.64	125.48
4	A	205	6SS	C26-C35-N34	-2.75	106.51	108.04
4	E	203	6SS	O36-C35-C26	2.73	127.54	125.48
4	B	202	6SS	C28-C27-C26	2.66	134.71	131.33
4	B	202	6SS	C37-C2-CL1	-2.64	118.48	119.50
4	C	202	6SS	C7-C9-C12	-2.62	101.30	105.13
4	B	202	6SS	C26-C35-N34	-2.55	106.62	108.04
4	A	205	6SS	C27-C26-N15	-2.54	110.12	116.44
4	F	201	6SS	C7-C26-C35	2.42	115.90	110.60
4	E	203	6SS	C3-C2-C37	2.41	120.61	119.29
4	A	205	6SS	C17-C16-N15	-2.37	108.98	113.78
4	D	202	6SS	C26-C35-N34	-2.37	106.72	108.04
4	D	202	6SS	C3-C2-C37	2.34	120.58	119.29
4	C	202	6SS	F38-C37-C6	2.34	121.96	119.56
4	A	205	6SS	C7-C9-C12	-2.31	101.75	105.13
4	E	203	6SS	C7-C9-C12	-2.30	101.76	105.13
4	E	203	6SS	C7-C26-C35	2.29	115.61	110.60
4	D	202	6SS	C7-C9-C12	-2.27	101.81	105.13
4	B	202	6SS	C27-C26-N15	-2.26	110.82	116.44
4	B	202	6SS	C7-C9-C12	-2.26	101.82	105.13
4	D	202	6SS	C17-C16-N15	-2.25	109.22	113.78
4	D	202	6SS	C28-C27-C26	2.24	134.18	131.33
4	C	202	6SS	C17-C16-N15	-2.22	109.29	113.78
4	E	203	6SS	C17-C16-N15	-2.21	109.31	113.78
4	C	202	6SS	C7-C26-C35	2.19	115.39	110.60
4	E	203	6SS	C26-C35-N34	-2.18	106.83	108.04
4	B	202	6SS	C17-C16-N15	-2.18	109.38	113.78
4	E	203	6SS	C27-C26-N15	-2.15	111.08	116.44
4	F	201	6SS	C17-C16-N15	-2.13	109.47	113.78
4	B	202	6SS	C7-C26-C35	2.13	115.27	110.60
4	C	202	6SS	C35-C26-N15	2.13	115.24	108.45
4	F	201	6SS	F38-C37-C6	2.12	121.73	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	202	6SS	C3-C2-C37	2.08	120.44	119.29
4	B	202	6SS	F38-C37-C6	2.08	121.69	119.56
4	A	205	6SS	C35-C26-N15	2.01	114.87	108.45

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	202	GOL	O1-C1-C2-C3
2	A	201	GOL	O1-C1-C2-C3
2	C	201	GOL	C1-C2-C3-O3
4	E	203	6SS	C25-C21-O22-C23
2	C	201	GOL	O2-C2-C3-O3
4	E	203	6SS	C20-C21-O22-C23
2	A	203	GOL	O1-C1-C2-C3
4	F	201	6SS	C25-C21-O22-C23
4	F	201	6SS	C20-C21-O22-C23
4	C	202	6SS	C25-C21-O22-C23
4	C	202	6SS	C20-C21-O22-C23
2	A	201	GOL	O1-C1-C2-O2
4	A	205	6SS	C25-C21-O22-C23
4	A	205	6SS	C20-C21-O22-C23
2	A	202	GOL	O1-C1-C2-O2
2	A	203	GOL	O1-C1-C2-O2
4	D	202	6SS	C20-C21-O22-C23
4	E	203	6SS	C24-C23-O22-C21
4	D	202	6SS	C25-C21-O22-C23
4	B	202	6SS	C20-C21-O22-C23
4	B	202	6SS	C25-C21-O22-C23

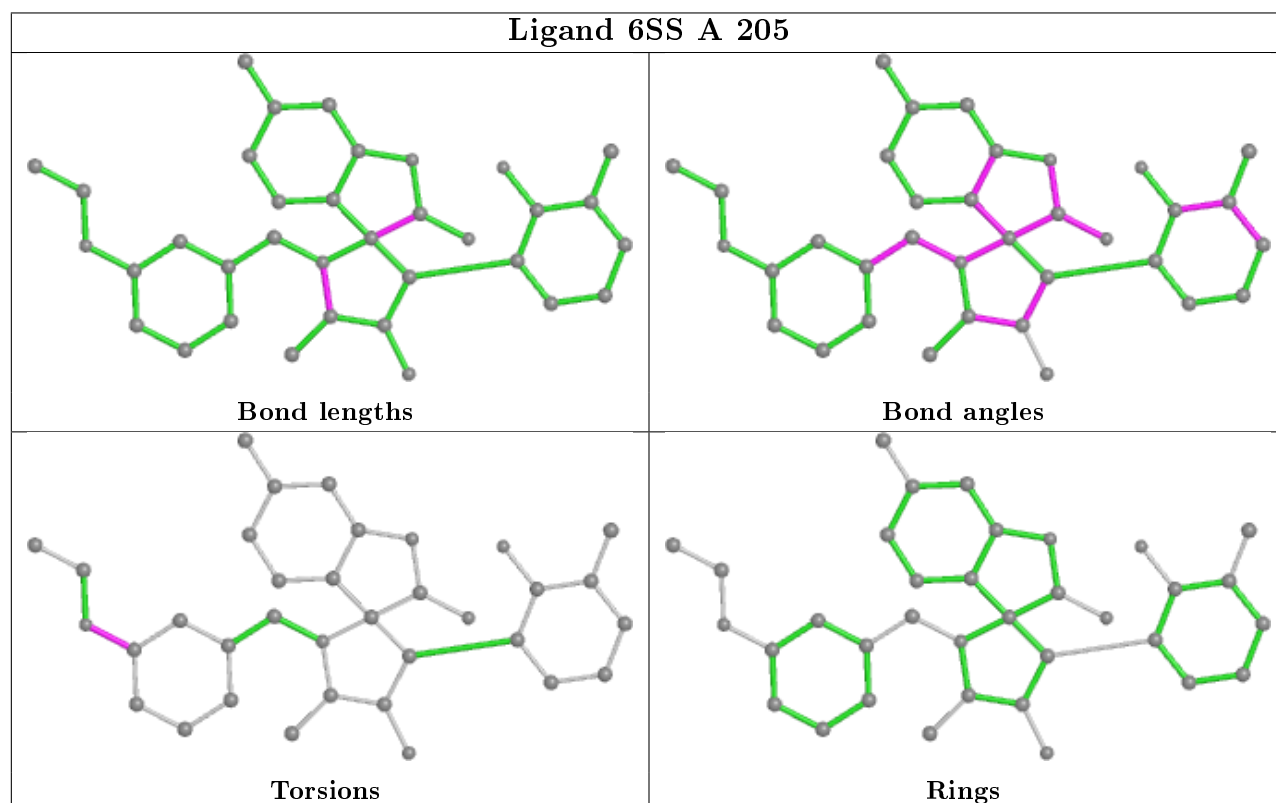
There are no ring outliers.

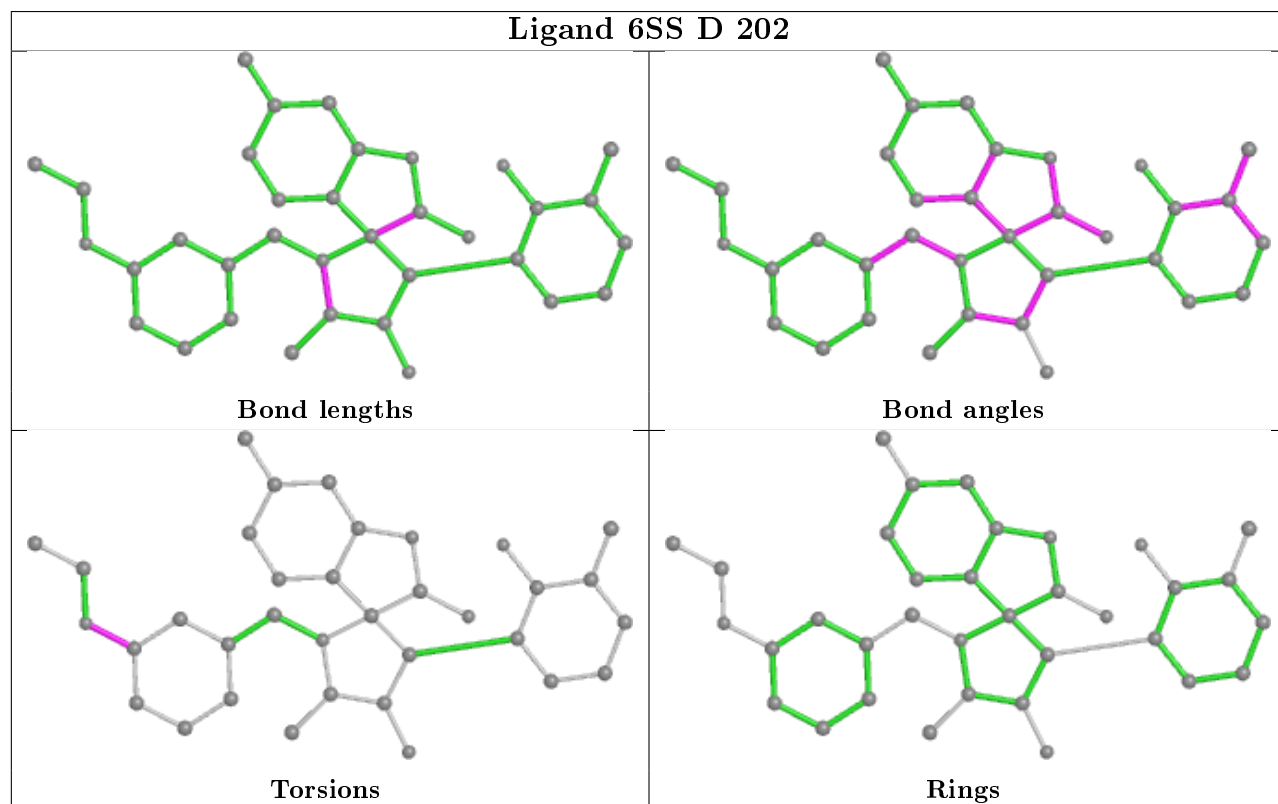
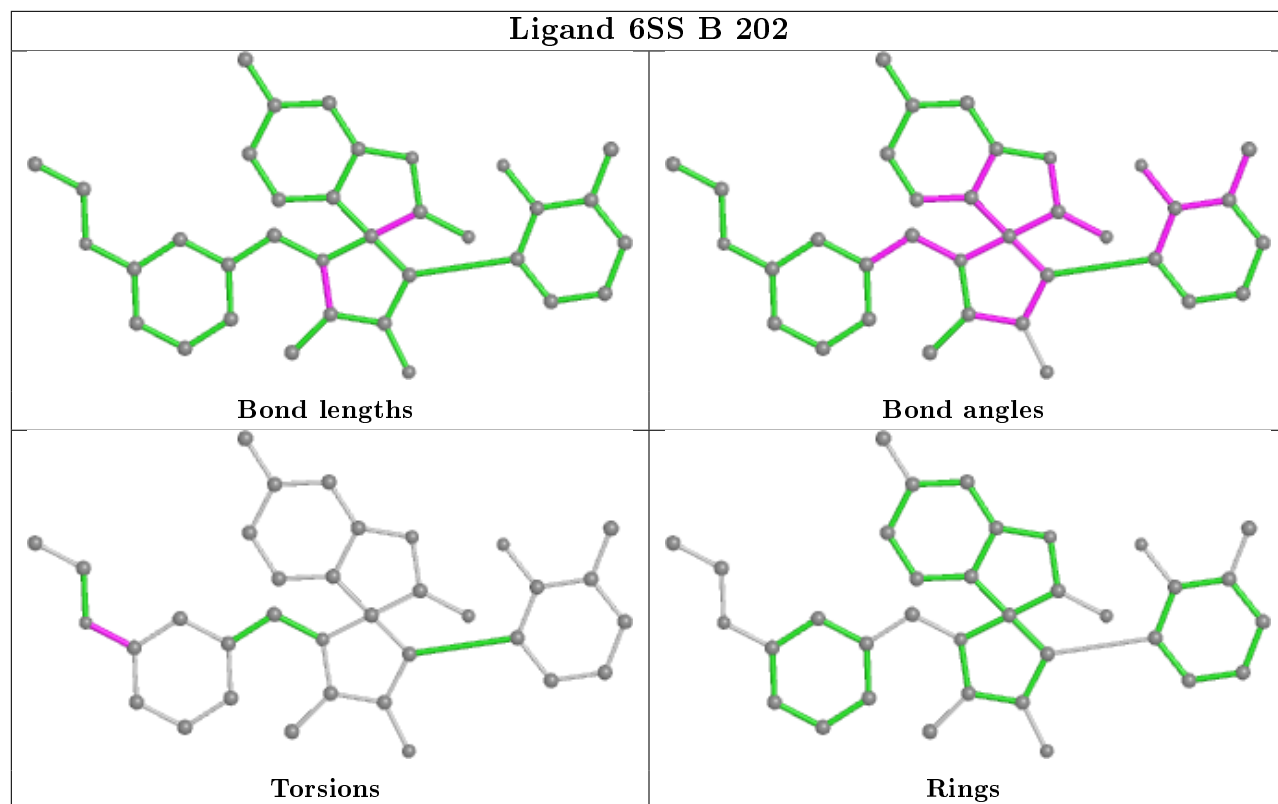
1 monomer is involved in 1 short contact:

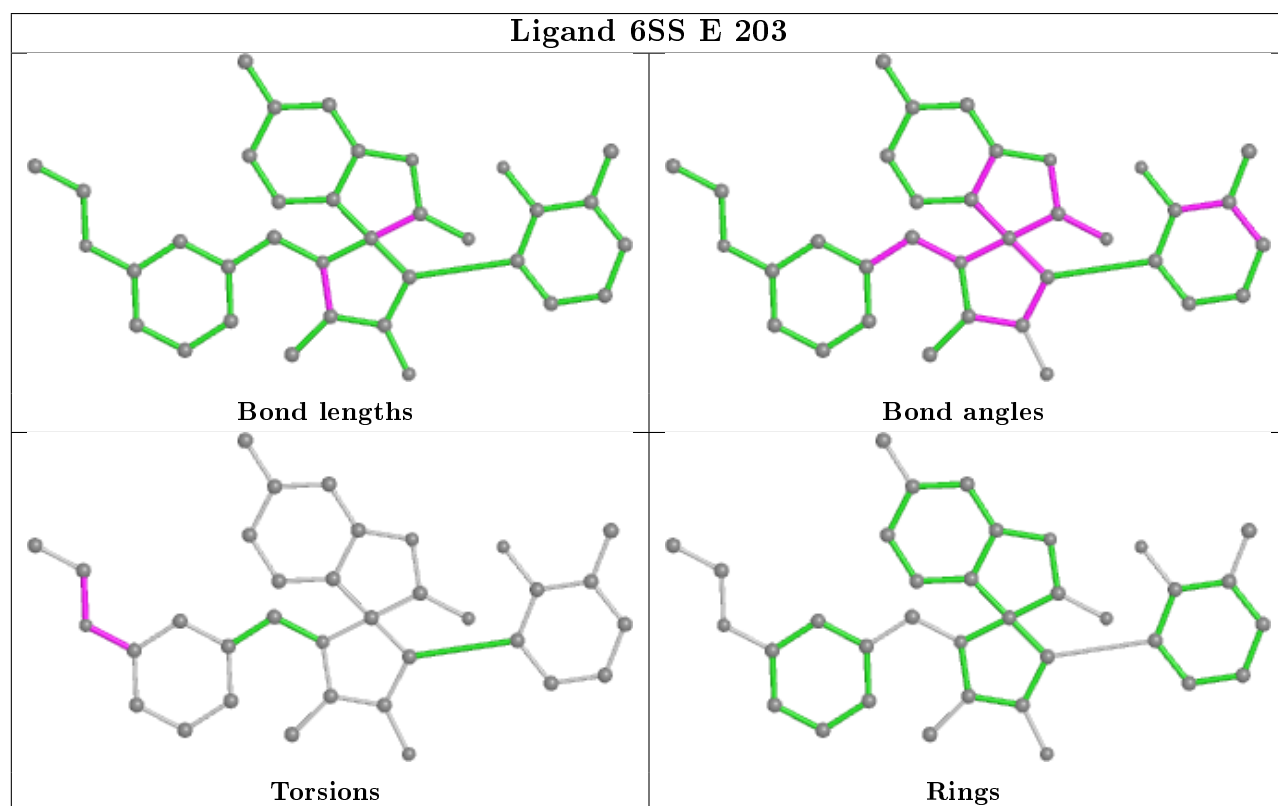
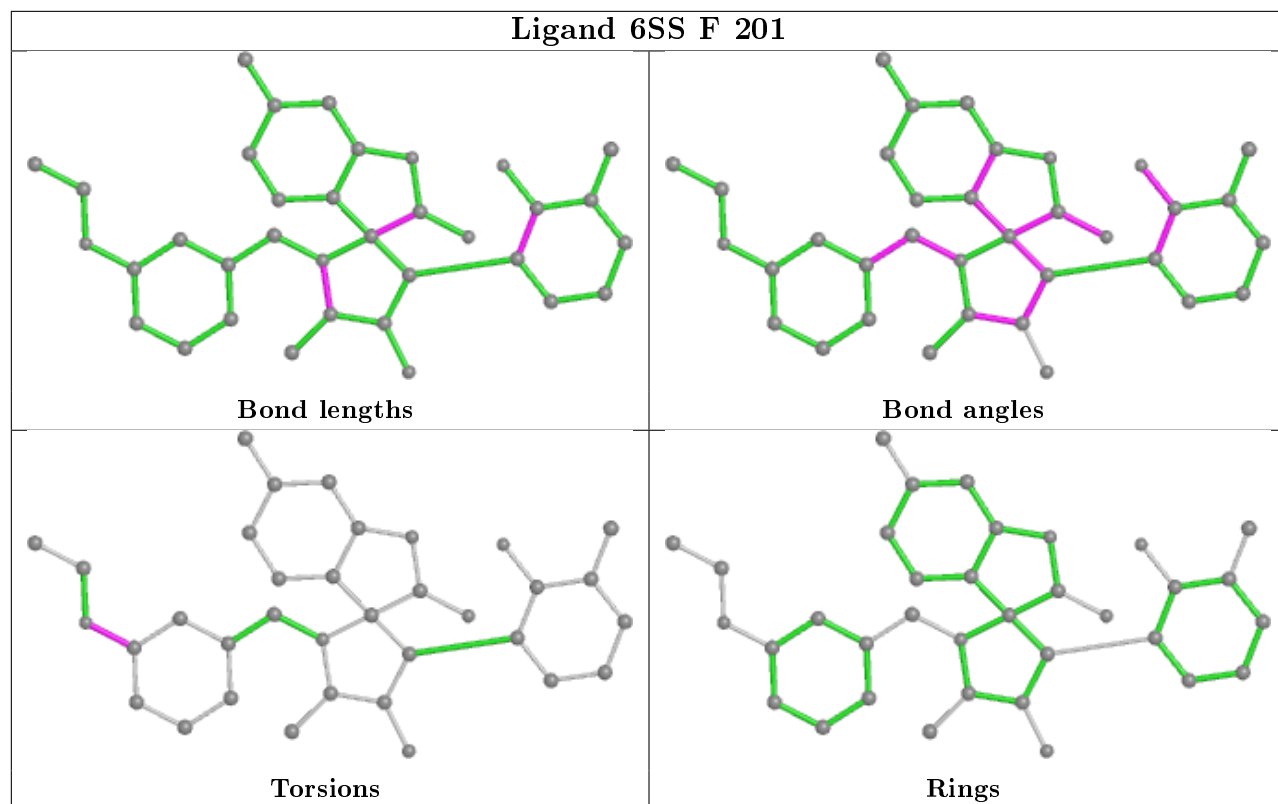
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	201	GOL	1	0

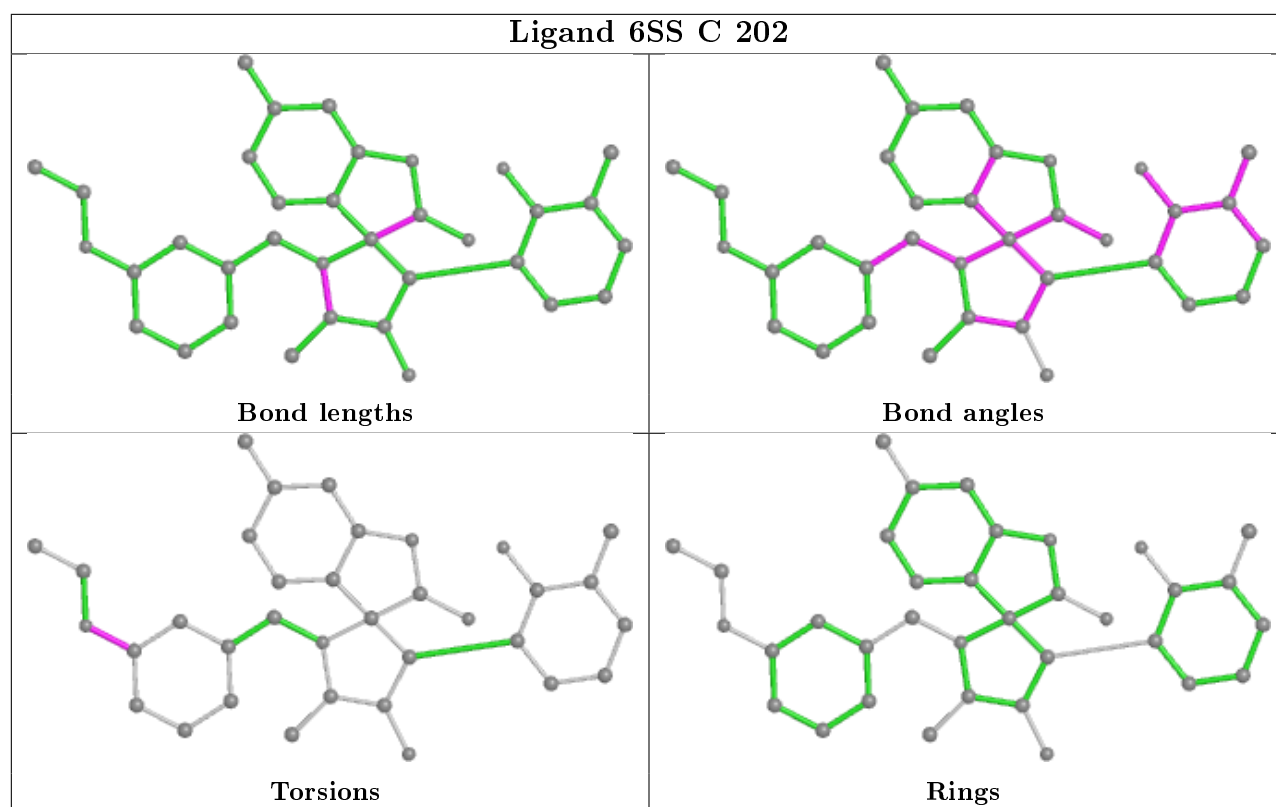
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

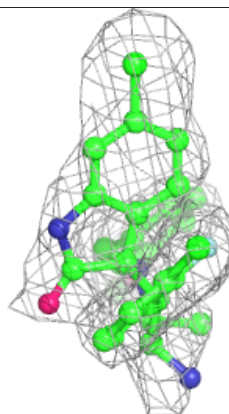
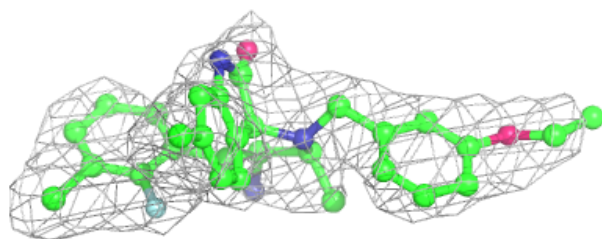
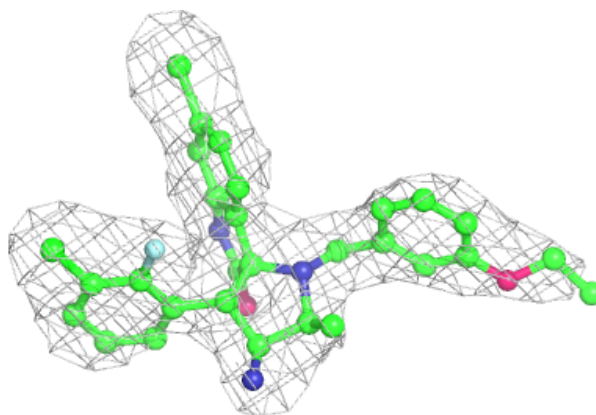
6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

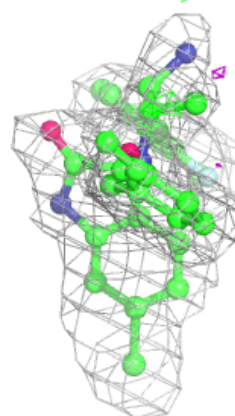
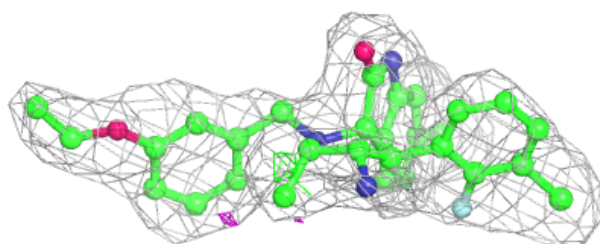
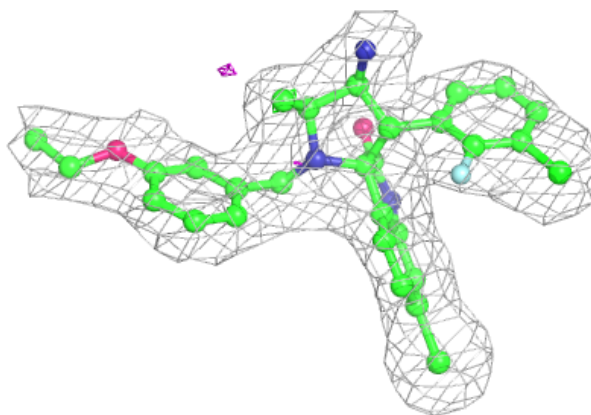
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 6SS A 205:

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

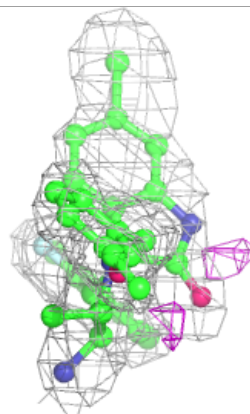
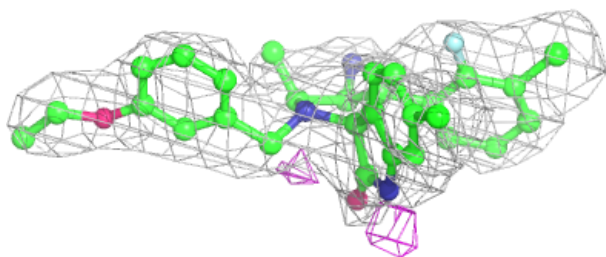
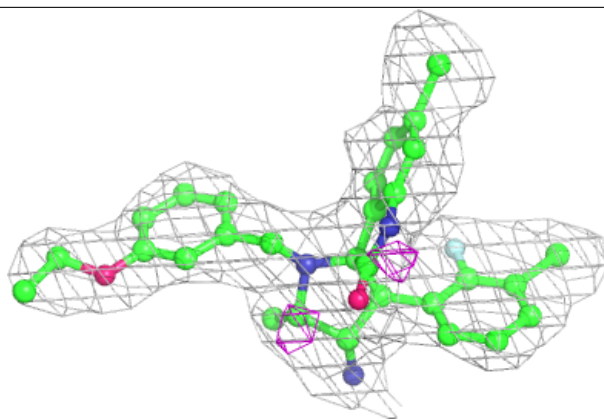
**Electron density around 6SS B 202:**

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

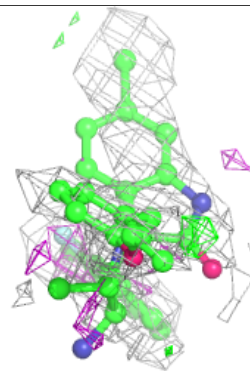
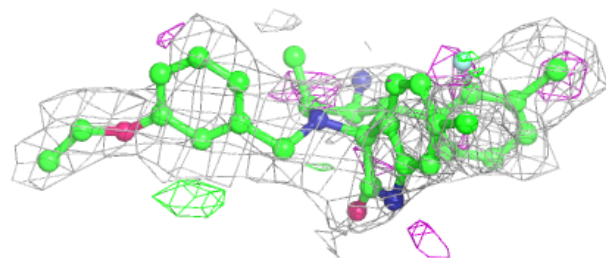
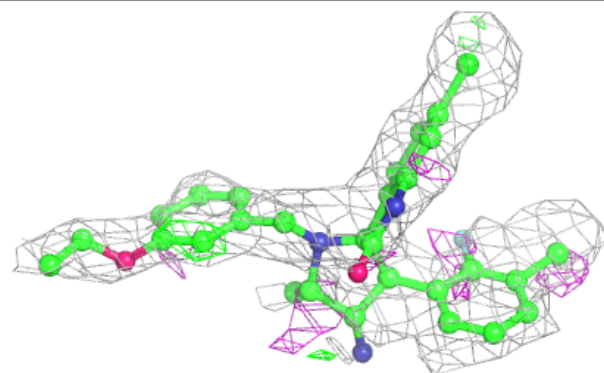


Electron density around 6SS D 202:

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

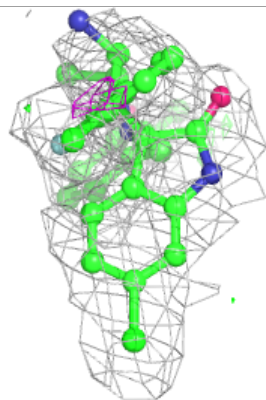
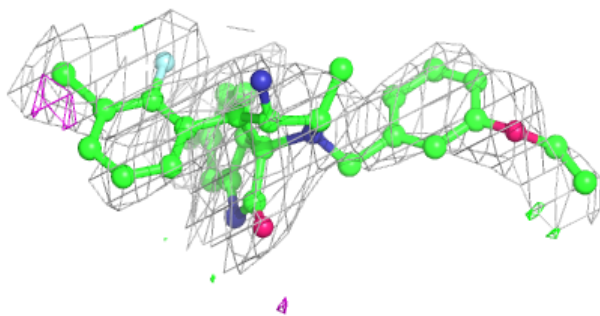
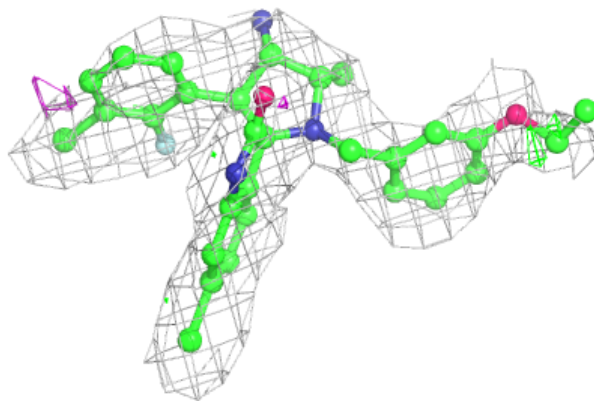
**Electron density around 6SS F 201:**

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

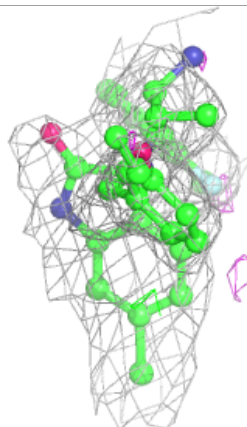
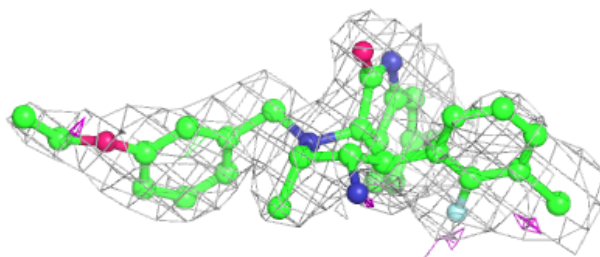
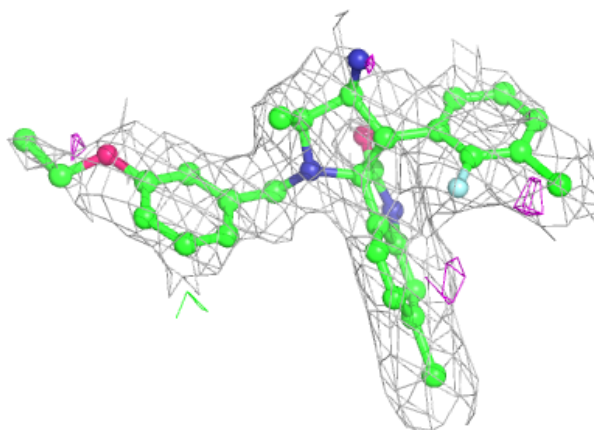


Electron density around 6SS E 203:

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

**Electron density around 6SS C 202:**

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



6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.