



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 2, 2025 – 01:16 am BST

PDB ID : 1HBN / pdb_00001hbn
Title : METHYL-COENZYME M REDUCTASE
Authors : Ermler, U.; Grabarse, W.
Deposited on : 2001-04-20
Resolution : 1.16 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 : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.42

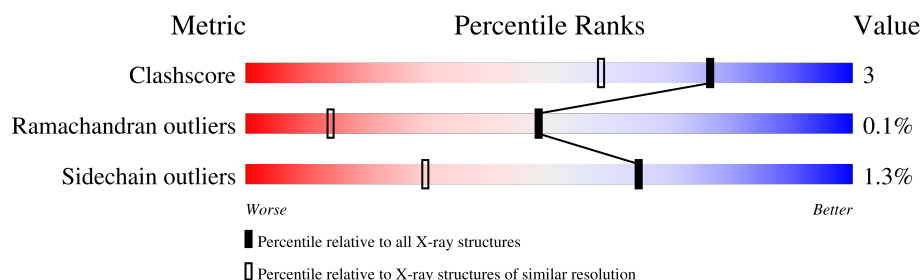
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.16 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))
Clashscore	180529	1146 (1.18-1.14)
Ramachandran outliers	177936	1120 (1.18-1.14)
Sidechain outliers	177891	1120 (1.18-1.14)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	549	
1	D	549	
2	B	442	
2	E	442	
3	C	248	
3	F	248	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	GOL	D	1556	-	X	-	-

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 21958 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 METHYL-COENZYME M REDUCTASE I ALPHA SUB-UNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	548	Total	C	N	O	S	23	13	0
			4291	2712	723	836	20			
1	D	548	Total	C	N	O	S	23	16	0
			4296	2717	721	838	20			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	257	MHS	HIS	modified residue	UNP P11558
A	271	AGM	ARG	modified residue	UNP P11558
A	400	MGN	GLN	modified residue	UNP P11558
A	445	GL3	GLY	modified residue	UNP P11558
A	452	SMC	CYS	modified residue	UNP P11558
D	257	MHS	HIS	modified residue	UNP P11558
D	271	AGM	ARG	modified residue	UNP P11558
D	400	MGN	GLN	modified residue	UNP P11558
D	445	GL3	GLY	modified residue	UNP P11558
D	452	SMC	CYS	modified residue	UNP P11558

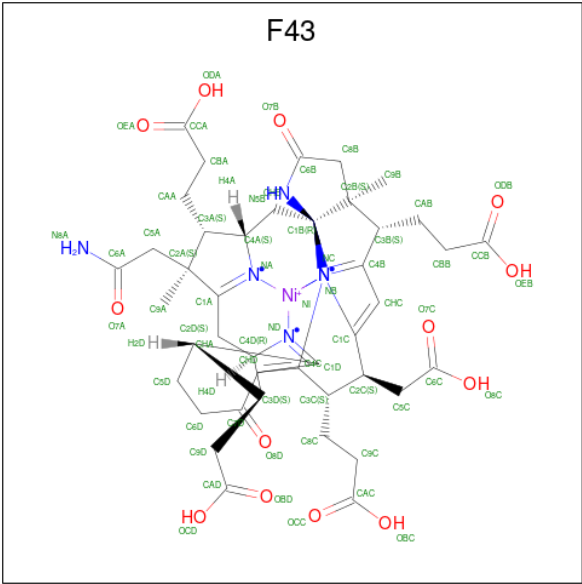
- Molecule 2 is a protein called METHYL-COENZYME M REDUCTASE I BETA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	442	Total	C	N	O	S	26	15	0
			3348	2115	557	654	22			
2	E	442	Total	C	N	O	S	41	12	0
			3340	2115	551	653	21			

- Molecule 3 is a protein called METHYL-COENZYME M REDUCTASE I GAMMA SUB-UNIT.

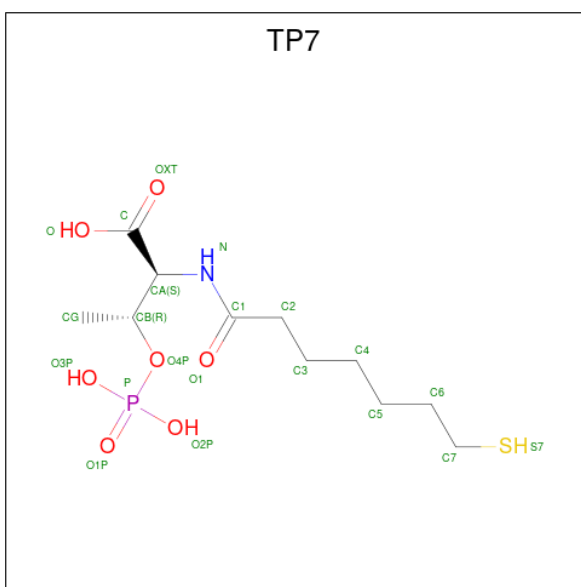
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	247	Total	C	N	O	S	39	5	0
			2007	1242	355	398	12			
3	F	247	Total	C	N	O	S	44	5	0
			2006	1242	354	398	12			

- Molecule 4 is FACTOR 430 (CCD ID: F43) (formula: C₄₂H₅₁N₆NiO₁₃).



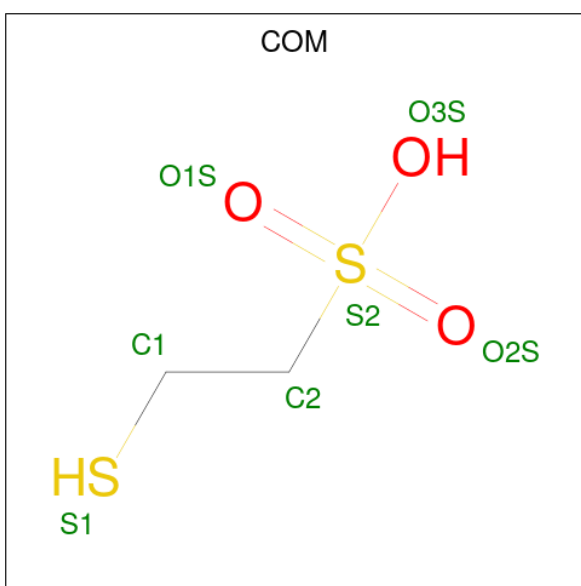
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	Ni	O	0	0
			62	42	6	1	13		
4	D	1	Total	C	N	Ni	O	0	0
			62	42	6	1	13		

- Molecule 5 is Coenzyme B (CCD ID: TP7) (formula: C₁₁H₂₂NO₇PS).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	S	0	0
			21	11	1	7	1	1		
5	D	1	Total	C	N	O	P	S	0	0
			21	11	1	7	1	1		

- Molecule 6 is 1-THIOETHANESULFONIC ACID (CCD ID: COM) (formula: $C_2H_6O_3S_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	O	S	0	0
			7	2	3	2		
6	D	1	Total	C	O	S	0	0
			7	2	3	2		

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	B	1	Total	C	O	0	0
			6	3	3		
7	C	1	Total	C	O	0	0
			6	3	3		
7	D	1	Total	C	O	0	0
			6	3	3		
7	D	1	Total	C	O	0	0
			6	3	3		
7	D	1	Total	C	O	0	0
			6	3	3		
7	E	1	Total	C	O	0	0
			6	3	3		

- Molecule 8 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total 1	Zn 1	0	1

- Molecule 9 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	5	Total 5	Mg 5	0	0
9	B	2	Total 2	Mg 2	0	0
9	C	2	Total 2	Mg 2	0	0
9	D	2	Total 2	Mg 2	0	0
9	E	5	Total 5	Mg 5	0	0
9	F	2	Total 2	Mg 2	0	0

- Molecule 10 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	5	Total 5	Na 5	0	0
10	B	2	Total 2	Na 2	0	0
10	D	3	Total 3	Na 3	0	0

- Molecule 11 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	B	1	Total 1	Cl 1	0	0
11	E	1	Total 1	Cl 1	0	0

- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	489	Total 489	O 489	0	29

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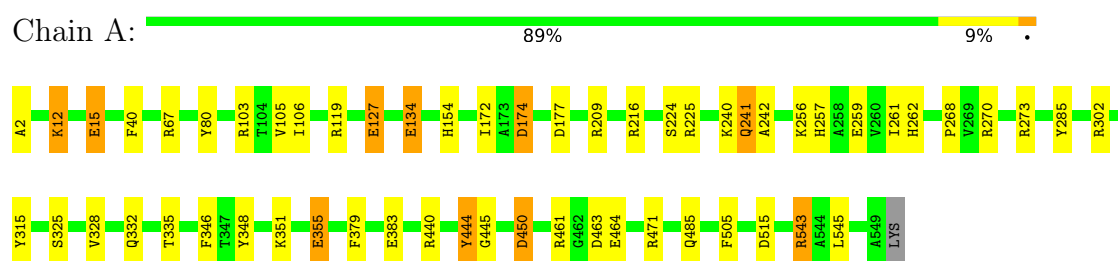
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	B	423	Total 423	O 423	0	31
12	C	294	Total 294	O 294	0	25
12	D	510	Total 510	O 510	0	37
12	E	402	Total 402	O 402	0	33
12	F	275	Total 275	O 275	0	15

3 Residue-property plots

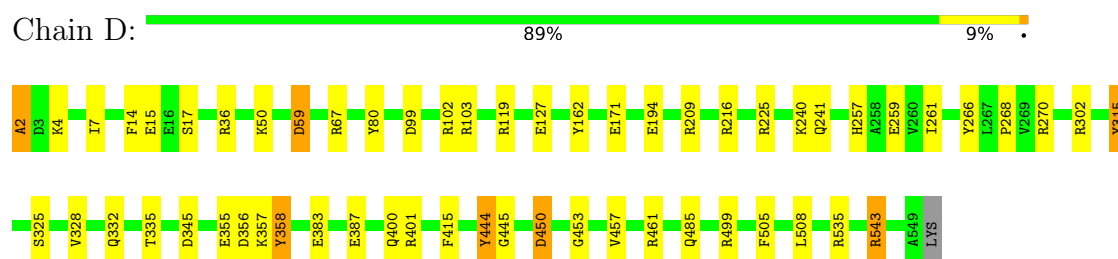
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

Note EDS was not executed.

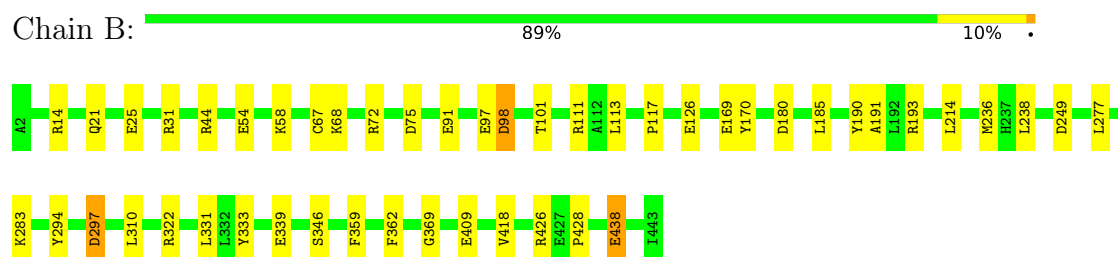
• Molecule 1: METHYL-COENZYME M REDUCTASE I ALPHA SUBUNIT



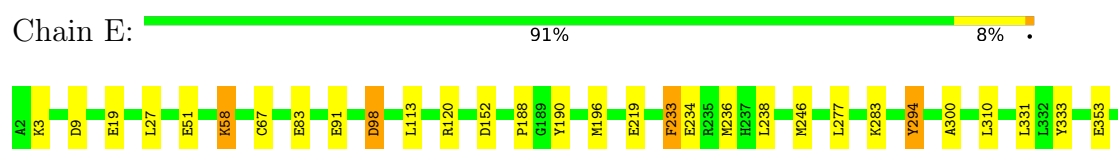
• Molecule 1: METHYL-COENZYME M REDUCTASE I ALPHA SUBUNIT



• Molecule 2: METHYL-COENZYME M REDUCTASE I BETA SUBUNIT



• Molecule 2: METHYL-COENZYME M REDUCTASE I BETA SUBUNIT





- Molecule 3: METHYL-COENZYME M REDUCTASE I GAMMA SUBUNIT

Chain C: 88% 10% .



- Molecule 3: METHYL-COENZYME M REDUCTASE I GAMMA SUBUNIT

Chain F: 89% 9% .



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.72Å 116.88Å 122.58Å 90.00° 92.02° 90.00°	Depositor
Resolution (Å)	30.00 – 1.16	Depositor
% Data completeness (in resolution range)	93.1 (30.00-1.16)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.124 , 0.190	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	21958	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: COM, SMC, ZN, GOL, MHS, TP7, AGM, MGN, GL3, MG, F43, NA, CL

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	0.78	3/4411 (0.1%)	1.25	45/5986 (0.8%)
1	D	0.93	7/4432 (0.2%)	1.23	44/6015 (0.7%)
2	B	0.76	3/3493 (0.1%)	1.22	31/4724 (0.7%)
2	E	1.65	6/3459 (0.2%)	1.19	20/4682 (0.4%)
3	C	0.98	8/2072 (0.4%)	1.47	31/2790 (1.1%)
3	F	0.78	7/2070 (0.3%)	1.37	25/2789 (0.9%)
All	All	1.03	34/19937 (0.2%)	1.27	196/26986 (0.7%)

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	3
1	D	0	3
2	B	0	1
2	E	0	1
3	C	0	1
All	All	0	9

The worst 5 of 34 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	438	GLU	CG-CD	83.79	2.77	1.51
2	E	83	GLU	CG-CD	28.40	1.94	1.51
1	D	355	GLU	CG-CD	26.78	1.92	1.51
1	D	127	GLU	CG-CD	-25.44	1.13	1.51
3	C	179	LYS	CD-CE	16.63	1.92	1.51

The worst 5 of 196 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	60	ASP	CB-CG-OD1	-22.90	97.69	118.30
2	E	438	GLU	CB-CG-CD	-20.86	57.87	114.20
3	F	196	ASP	CB-CG-OD2	18.44	134.90	118.30
2	B	98	ASP	CA-CB-CG	-16.64	76.78	113.40
3	C	120	ARG	NE-CZ-NH2	-15.87	112.36	120.30

There are no chirality outliers.

5 of 9 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	134[A]	GLU	Sidechain
1	A	134[B]	GLU	Sidechain
1	A	241	GLN	Peptide
2	B	333	TYR	Sidechain
3	C	60	ASP	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	4291	0	4093	24	0
1	D	4296	0	4108	16	1
2	B	3348	0	3331	31	0
2	E	3340	0	3325	28	0
3	C	2007	0	1937	9	0
3	F	2006	0	1937	3	0
4	A	62	0	43	1	0
4	D	62	0	43	1	0
5	A	21	0	19	0	0
5	D	21	0	19	0	0
6	A	7	0	4	0	0
6	D	7	0	5	0	0
7	A	24	0	31	7	0
7	B	6	0	7	1	0
7	C	6	0	8	0	0
7	D	24	0	31	2	0
7	E	6	0	8	0	0
8	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	A	5	0	0	0	0
9	B	2	0	0	0	0
9	C	2	0	0	0	0
9	D	2	0	0	0	0
9	E	5	0	0	0	0
9	F	2	0	0	0	0
10	A	5	0	0	0	0
10	B	2	0	0	0	0
10	D	3	0	0	0	0
11	B	1	0	0	0	0
11	E	1	0	0	0	0
12	A	489	0	0	10	0
12	B	423	0	0	10	0
12	C	294	0	0	3	0
12	D	510	0	0	6	0
12	E	402	0	0	6	0
12	F	275	0	0	0	0
All	All	21958	0	18949	106	1

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.

The worst 5 of 106 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)
2:B:21[B]:GLN:OE1	12:B:2048[B]:HOH:O	1.58	1.20
2:E:196[B]:MET:HE1	2:E:374:ILE:HG22	1.39	1.01
2:B:236[B]:MET:CE	2:B:236[B]:MET:SD	2.53	0.97
2:B:236[B]:MET:HB2	3:C:248:LEU:HD11	1.51	0.92
2:B:236[A]:MET:HB2	3:C:248:LEU:HD11	1.53	0.91

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:59:ASP:OD2	1:D:357:LYS:NZ[2_644]	2.19	0.01

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	554/549 (101%)	535 (97%)	18 (3%)	1 (0%)	44	13
1	D	557/549 (102%)	539 (97%)	17 (3%)	1 (0%)	44	13
2	B	456/442 (103%)	447 (98%)	9 (2%)	0	100	100
2	E	452/442 (102%)	445 (98%)	7 (2%)	0	100	100
3	C	250/248 (101%)	245 (98%)	5 (2%)	0	100	100
3	F	250/248 (101%)	242 (97%)	8 (3%)	0	100	100
All	All	2519/2478 (102%)	2453 (97%)	64 (2%)	2 (0%)	48	14

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	325	SER
1	D	325	SER

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	446/434 (103%)	440 (99%)	6 (1%)	65	29
1	D	449/434 (104%)	445 (99%)	4 (1%)	75	46
2	B	357/341 (105%)	353 (99%)	4 (1%)	70	37
2	E	353/341 (104%)	350 (99%)	3 (1%)	79	51
3	C	220/216 (102%)	216 (98%)	4 (2%)	54	17

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	F	220/216 (102%)	215 (98%)	5 (2%)	45 9
All	All	2045/1982 (103%)	2019 (99%)	26 (1%)	65 29

5 of 26 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	17	SER
1	D	543	ARG
3	F	196	ASP
1	D	450	ASP
2	E	98	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	365	ASN
1	D	485	GLN
2	E	40	GLN
1	D	42	ASN
1	A	485	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

10 non-standard protein/DNA/RNA residues 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
1	MHS	D	257	1	7,11,12	1.57	1 (14%)	6,14,16	2.01	2 (33%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	AGM	D	271	1	10,11,12	0.90	0	6,13,15	0.94	0
1	GL3	D	445	1	2,3,4	3.10	1 (50%)	1,2,4	0.21	0
1	SMC	A	452	1	5,6,7	0.63	0	2,6,8	0.96	0
1	MGN	D	400	1	6,9,10	1.21	1 (16%)	5,12,14	0.72	0
1	AGM	A	271	1	10,11,12	0.77	0	6,13,15	1.00	0
1	MHS	A	257	1	7,11,12	1.72	1 (14%)	6,14,16	1.37	1 (16%)
1	MGN	A	400	1	6,9,10	0.72	0	5,12,14	0.80	0
1	GL3	A	445	1	2,3,4	2.45	1 (50%)	1,2,4	0.03	0
1	SMC	D	452	1	5,6,7	0.95	0	2,6,8	1.41	0

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
1	MHS	D	257	1	-	0/5/6/8	0/1/1/1
1	AGM	D	271	1	-	3/10/11/13	-
1	GL3	D	445	1	-	0/1/1/2	-
1	SMC	A	452	1	-	1/3/5/7	-
1	MGN	D	400	1	-	0/7/9/12	-
1	AGM	A	271	1	-	3/10/11/13	-
1	MHS	A	257	1	-	0/5/6/8	0/1/1/1
1	MGN	A	400	1	-	0/7/9/12	-
1	GL3	A	445	1	-	1/1/1/2	-
1	SMC	D	452	1	-	1/3/5/7	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	445	GL3	C-S	-4.38	1.65	1.80
1	A	257	MHS	CM-ND1	3.49	1.55	1.47
1	A	445	GL3	C-S	-3.43	1.68	1.80
1	D	257	MHS	CM-ND1	3.14	1.54	1.47
1	D	400	MGN	CB1-CA	-2.29	1.52	1.55

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	257	MHS	NE2-CE1-ND1	-3.34	107.30	112.26
1	D	257	MHS	CD2-NE2-CE1	2.96	110.40	105.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	257	MHS	CD2-NE2-CE1	2.02	108.93	105.78

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	452	SMC	CA-CB-SG-CS
1	D	452	SMC	CA-CB-SG-CS
1	A	271	AGM	CE2-CD-NE1-CZ
1	D	271	AGM	CE2-CD-NE1-CZ
1	A	445	GL3	S-C-CA-N

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 48 ligands modelled in this entry, 31 are monoatomic - leaving 17 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	$\# Z > 2$
7	GOL	D	1554	10	5,5,5	0.55	0	5,5,5	0.58	0
6	COM	D	1552	4	6,6,6	1.35	1 (16%)	7,8,8	0.56	0
4	F43	D	1550	6,1	61,71,71	2.77	11 (18%)	64,118,118	1.50	15 (23%)
6	COM	A	1552	4	6,6,6	1.43	2 (33%)	7,8,8	1.23	1 (14%)
7	GOL	A	1554	-	5,5,5	0.62	0	5,5,5	1.28	1 (20%)
5	TP7	A	1551	-	19,20,20	1.57	3 (15%)	24,26,26	1.11	2 (8%)
4	F43	A	1550	6,1	61,71,71	2.34	12 (19%)	64,118,118	1.63	11 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	GOL	D	1553	-	5,5,5	0.77	0	5,5,5	0.85	0
7	GOL	B	1444	-	5,5,5	1.21	1 (20%)	5,5,5	1.24	0
7	GOL	D	1555	-	5,5,5	0.66	0	5,5,5	0.99	0
7	GOL	D	1556	-	5,5,5	1.07	1 (20%)	5,5,5	1.68	2 (40%)
7	GOL	A	1555	-	5,5,5	0.88	0	5,5,5	0.98	0
7	GOL	E	1444	-	5,5,5	0.81	0	5,5,5	0.84	0
7	GOL	A	1553	10	5,5,5	0.35	0	5,5,5	1.96	2 (40%)
7	GOL	C	1249	-	5,5,5	0.50	0	5,5,5	0.37	0
5	TP7	D	1551	-	19,20,20	1.84	5 (26%)	24,26,26	1.10	2 (8%)
7	GOL	A	1556	-	5,5,5	1.05	0	5,5,5	1.00	1 (20%)

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
7	GOL	D	1554	10	-	0/4/4/4	-
6	COM	D	1552	4	-	0/4/4/4	-
4	F43	D	1550	6,1	-	9/28/185/185	-
6	COM	A	1552	4	-	0/4/4/4	-
7	GOL	A	1554	-	-	0/4/4/4	-
5	TP7	A	1551	-	-	0/24/24/24	-
4	F43	A	1550	6,1	-	9/28/185/185	-
7	GOL	D	1553	-	-	0/4/4/4	-
7	GOL	B	1444	-	-	2/4/4/4	-
7	GOL	D	1555	-	-	0/4/4/4	-
7	GOL	D	1556	-	-	4/4/4/4	-
7	GOL	A	1555	-	-	2/4/4/4	-
7	GOL	E	1444	-	-	0/4/4/4	-
7	GOL	A	1553	10	-	2/4/4/4	-
7	GOL	C	1249	-	-	0/4/4/4	-
5	TP7	D	1551	-	-	0/24/24/24	-
7	GOL	A	1556	-	-	0/4/4/4	-

The worst 5 of 36 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1550	F43	NI-NA	12.74	2.17	1.89
4	D	1550	F43	NI-NB	11.20	2.13	1.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1550	F43	NI-NA	9.86	2.10	1.89
4	A	1550	F43	NI-NB	8.52	2.07	1.89
4	D	1550	F43	NI-ND	7.29	2.05	1.89

The worst 5 of 37 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1550	F43	C3D-C4D-ND	5.02	110.16	102.34
4	A	1550	F43	C5D-C2D-C1D	4.78	116.88	110.45
4	A	1550	F43	CAB-C3B-C2B	-3.99	110.64	119.09
4	D	1550	F43	CAB-C3B-C2B	-3.79	111.08	119.09
4	D	1550	F43	C3B-C4B-CHC	-3.52	115.72	123.32

There are no chirality outliers.

5 of 28 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	1553	GOL	C1-C2-C3-O3
7	B	1444	GOL	C1-C2-C3-O3
7	D	1556	GOL	O1-C1-C2-C3
7	D	1556	GOL	C1-C2-C3-O3
7	D	1556	GOL	O2-C2-C3-O3

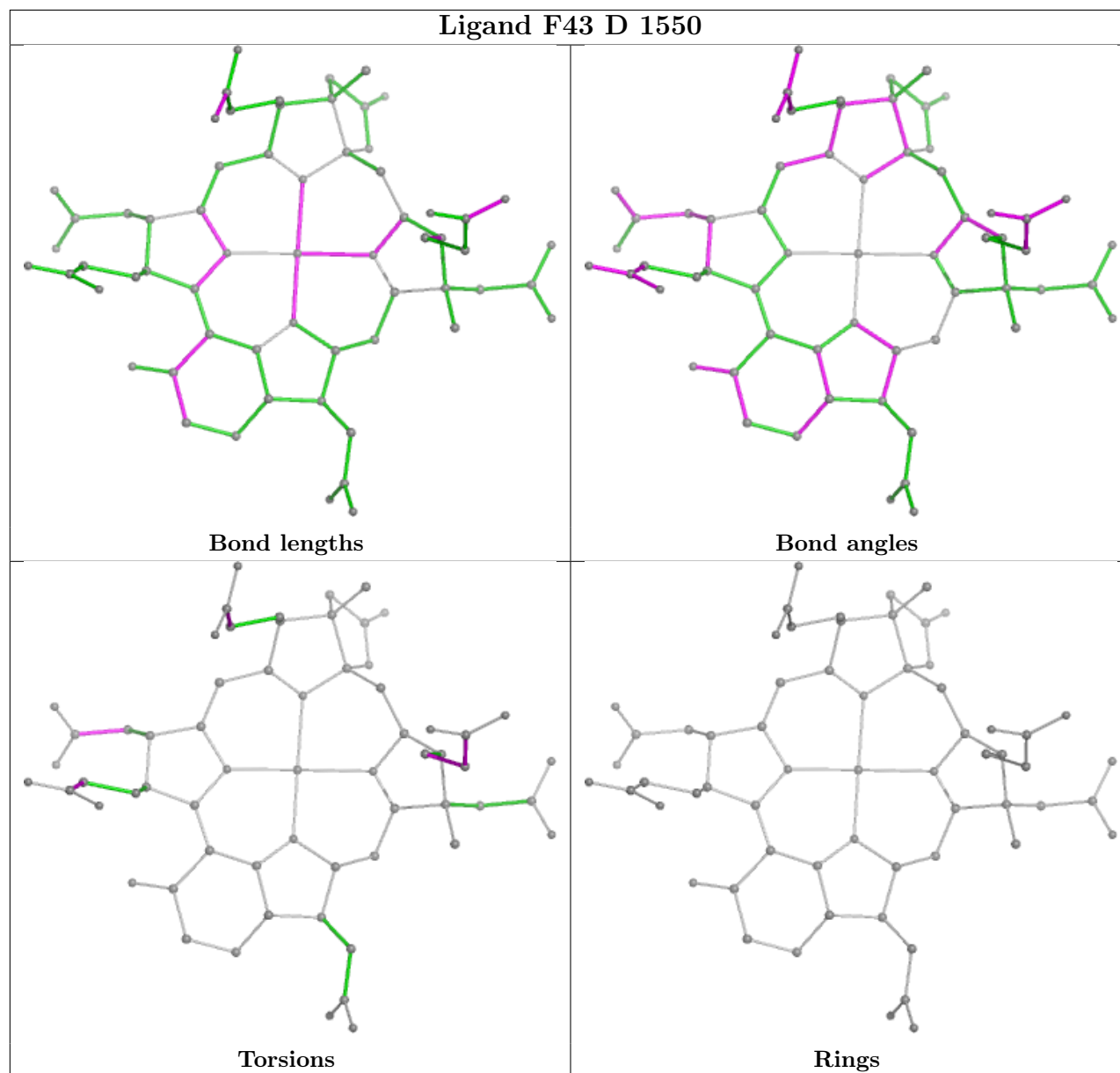
There are no ring outliers.

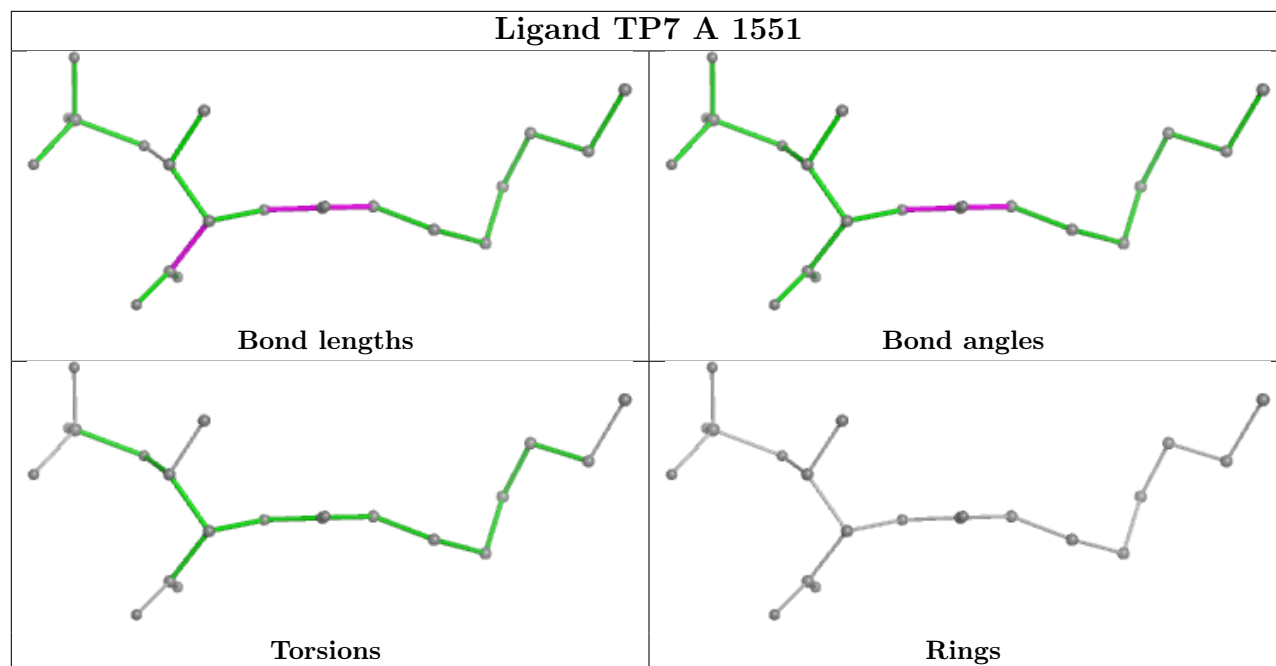
8 monomers are involved in 12 short contacts:

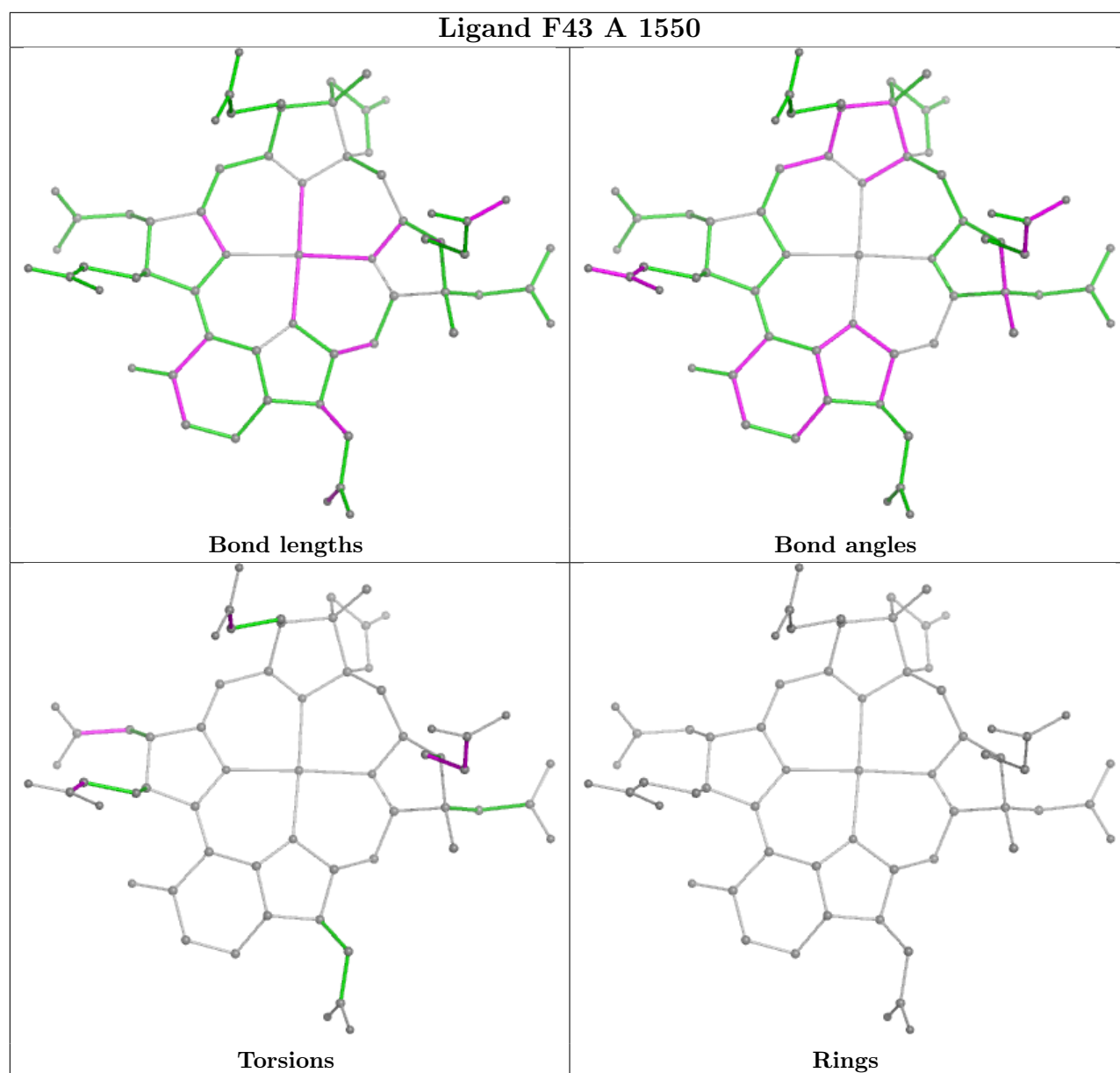
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	1550	F43	1	0
7	A	1554	GOL	1	0
4	A	1550	F43	1	0
7	B	1444	GOL	1	0
7	D	1556	GOL	2	0
7	A	1555	GOL	1	0
7	A	1553	GOL	2	0
7	A	1556	GOL	3	0

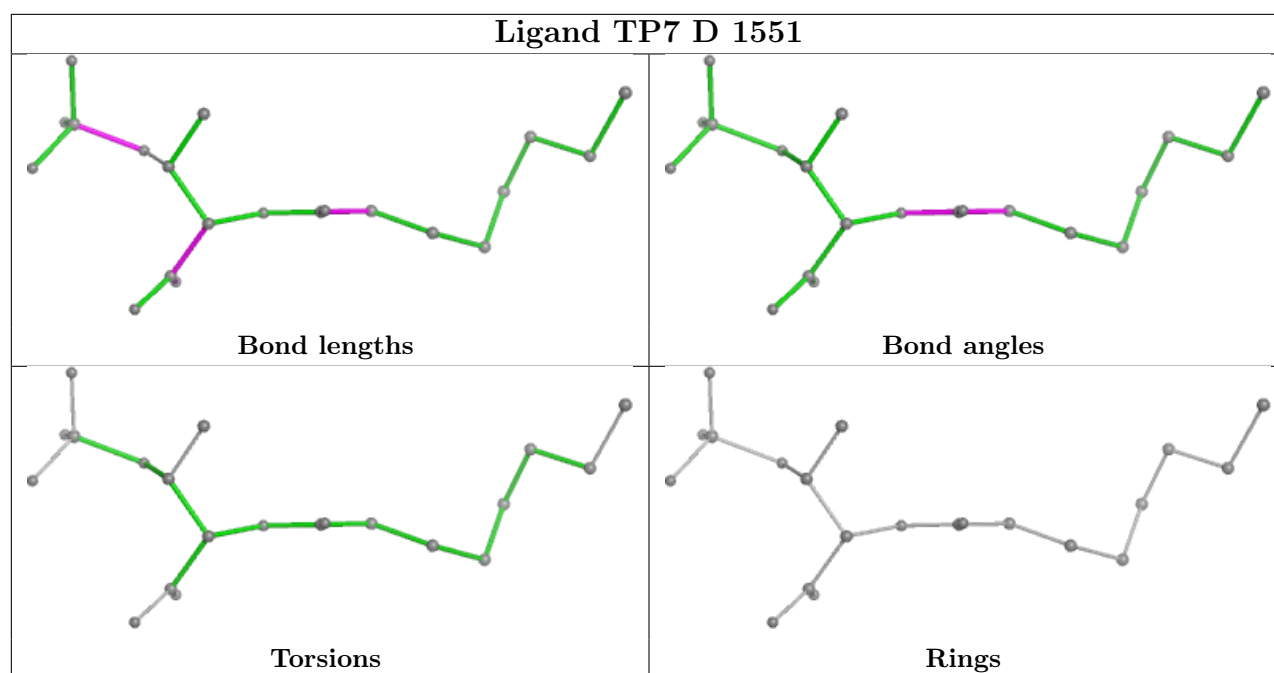
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier.

Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.