



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 23, 2024 – 10:02 PM EDT

PDB ID : 1LGH
Title : CRYSTAL STRUCTURE OF THE LIGHT-HARVESTING COMPLEX II
(B800-850) FROM RHODOSPIRILLUM MOLISCHIANUM
Authors : Koepke, J.; Hu, X.; Schulten, K.; Michel, H.
Deposited on : 1996-03-20
Resolution : 2.40 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.1

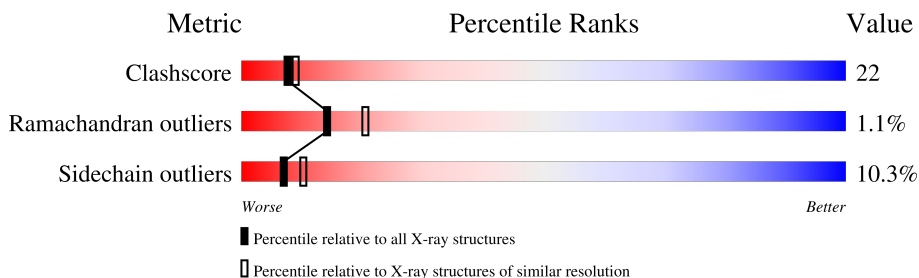
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)

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	56	
1	D	56	
1	G	56	
1	J	56	
2	B	45	
2	E	45	
2	H	45	
2	K	45	

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	HTO	D	100	-	-	X	-
6	HTO	D	98	-	-	X	-
6	HTO	E	61	-	-	X	-
6	HTO	K	61	-	-	X	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 5436 atoms, of which 948 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.

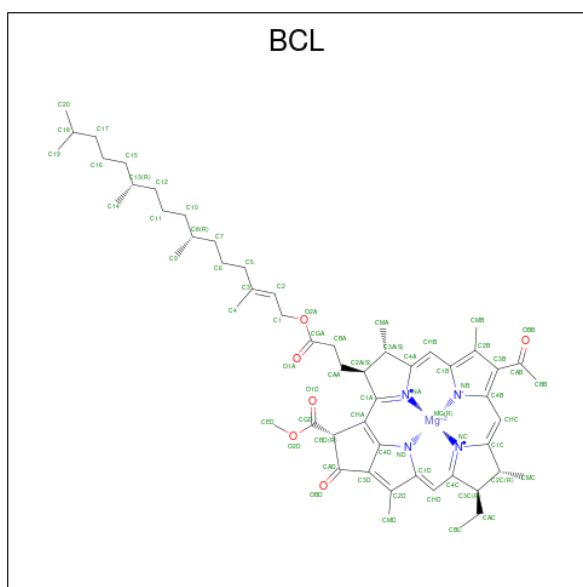
- Molecule 1 is a protein called LIGHT HARVESTING COMPLEX II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	56	Total	C	H	N	O	0	0	0
			507	284	84	69	70			
1	D	56	Total	C	H	N	O	0	0	0
			507	284	84	69	70			
1	G	56	Total	C	H	N	O	0	0	0
			507	284	84	69	70			
1	J	56	Total	C	H	N	O	0	0	0
			507	284	84	69	70			

- Molecule 2 is a protein called LIGHT HARVESTING COMPLEX II.

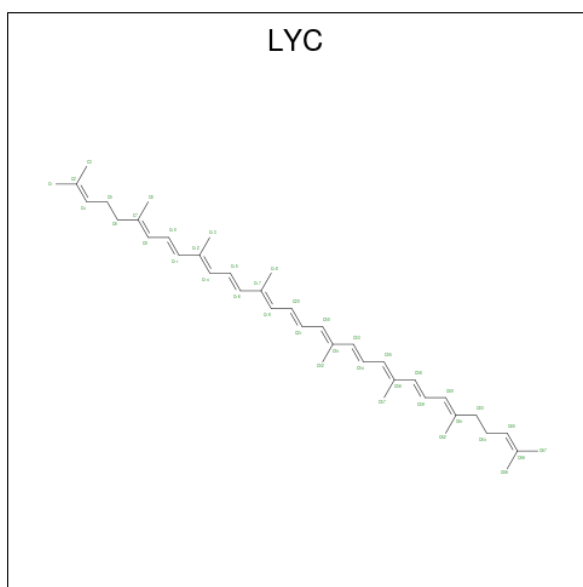
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	43	Total	C	H	N	O	0	0	0
			418	236	67	56	59			
2	E	43	Total	C	H	N	O	0	0	0
			418	236	67	56	59			
2	H	43	Total	C	H	N	O	0	0	0
			418	236	67	56	59			
2	K	43	Total	C	H	N	O	0	0	0
			418	236	67	56	59			

- Molecule 3 is BACTERIOCHLOROPHYLL A (three-letter code: BCL) (formula: C₅₅H₇₄MgN₄O₆).



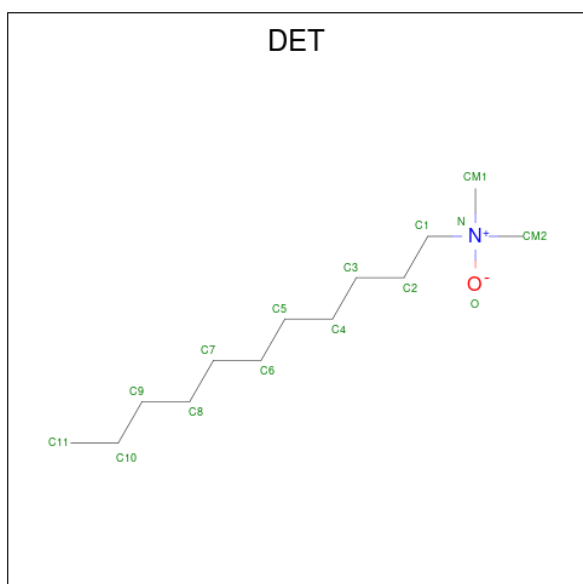
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	A	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
3	A	1	Total	C	Mg	N	O	0	0
			55	44	1	4	6		
3	B	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
3	D	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
3	D	1	Total	C	Mg	N	O	0	0
			55	44	1	4	6		
3	E	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
3	G	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
3	G	1	Total	C	Mg	N	O	0	0
			55	44	1	4	6		
3	H	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
3	J	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
3	J	1	Total	C	Mg	N	O	0	0
			55	44	1	4	6		
3	K	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

- Molecule 4 is LYCOPENE (three-letter code: LYC) (formula: C₄₀H₅₆).



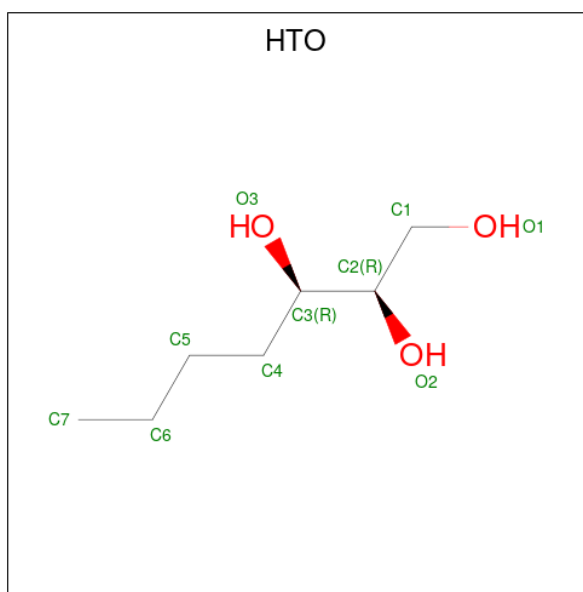
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	H	0	0
			96	40	56		
4	D	1	Total	C	H	0	0
			96	40	56		
4	G	1	Total	C	H	0	0
			96	40	56		
4	J	1	Total	C	H	0	0
			96	40	56		

- Molecule 5 is UNDECYLAMINE-N,N-DIMETHYL-N-OXIDE (three-letter code: DET) (formula: $C_{13}H_{29}NO$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			15	13	1	1		
5	B	1	Total	C	N	O	0	0
			15	13	1	1		
5	B	1	Total	C	N	O	0	0
			15	13	1	1		
5	D	1	Total	C	N	O	0	0
			15	13	1	1		
5	E	1	Total	C	N	O	0	0
			15	13	1	1		
5	E	1	Total	C	N	O	0	0
			15	13	1	1		
5	G	1	Total	C	N	O	0	0
			15	13	1	1		
5	H	1	Total	C	N	O	0	0
			15	13	1	1		
5	H	1	Total	C	N	O	0	0
			15	13	1	1		
5	J	1	Total	C	N	O	0	0
			15	13	1	1		
5	K	1	Total	C	N	O	0	0
			15	13	1	1		
5	K	1	Total	C	N	O	0	0
			15	13	1	1		

- Molecule 6 is HEPTANE-1,2,3-TRIOL (three-letter code: HTO) (formula: C₇H₁₆O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 10 7 3	0	0
6	A	1	Total C O 10 7 3	0	0
6	B	1	Total C O 10 7 3	0	0
6	D	1	Total C O 10 7 3	0	0
6	D	1	Total C O 10 7 3	0	0
6	D	1	Total C O 10 7 3	0	0
6	D	1	Total C O 10 7 3	0	0
6	D	1	Total C O 10 7 3	0	0
6	D	1	Total C O 10 7 3	0	0
6	D	1	Total C O 10 7 3	0	0
6	E	1	Total C O 10 7 3	0	0
6	E	1	Total C O 10 7 3	0	0
6	E	1	Total C O 10 7 3	0	0
6	G	1	Total C O 10 7 3	0	0
6	G	1	Total C O 10 7 3	0	0
6	H	1	Total C O 10 7 3	0	0
6	J	1	Total C O 10 7 3	0	0
6	J	1	Total C O 10 7 3	0	0
6	J	1	Total C O 10 7 3	0	0
6	J	1	Total C O 10 7 3	0	0
6	J	1	Total C O 10 7 3	0	0
6	K	1	Total C O 10 7 3	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	K	1	Total	C	O	0	0
			10	7	3		
6	K	1	Total	C	O	0	0
			10	7	3		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	4	Total	H	O	0	0
			12	8	4		
7	B	9	Total	H	O	0	0
			25	16	9		
7	D	8	Total	H	O	0	0
			24	16	8		
7	E	11	Total	H	O	0	0
			31	20	11		
7	G	4	Total	H	O	0	0
			12	8	4		
7	H	10	Total	H	O	0	0
			28	18	10		
7	J	7	Total	H	O	0	0
			21	14	7		
7	K	11	Total	H	O	0	0
			31	20	11		

3 Residue-property plots [i](#)

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.

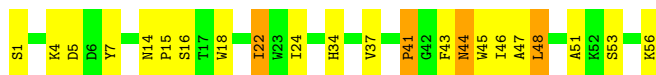
- Molecule 1: LIGHT HARVESTING COMPLEX II

Chain A: 62% 32% 5%



- Molecule 1: LIGHT HARVESTING COMPLEX II

Chain D: 61% 32% 7%



- Molecule 1: LIGHT HARVESTING COMPLEX II

Chain G: 64% 29% 7%



- Molecule 1: LIGHT HARVESTING COMPLEX II

Chain J: 62% 30% 7%



- Molecule 2: LIGHT HARVESTING COMPLEX II

Chain B: 64% 29% . .



- Molecule 2: LIGHT HARVESTING COMPLEX II

Chain E: 67% 24% . .



- Molecule 2: LIGHT HARVESTING COMPLEX II



- Molecule 2: LIGHT HARVESTING COMPLEX II



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	91.60Å 91.60Å 209.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.40	Depositor
% Data completeness (in resolution range)	87.2 (8.00-2.40)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.211 , 0.232	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	5436	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

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: HTO, DET, BCL, LYC

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	0.63	0/437	0.71	0/603
1	D	0.63	0/437	0.71	0/603
1	G	0.63	0/437	0.71	0/603
1	J	0.63	0/437	0.71	0/603
2	B	0.59	0/363	0.60	0/495
2	E	0.59	0/363	0.60	0/495
2	H	0.59	0/363	0.60	0/495
2	K	0.59	0/363	0.60	0/495
All	All	0.61	0/3200	0.66	0/4392

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	423	84	441	16	0
1	D	423	84	441	20	0
1	G	423	84	441	16	0
1	J	423	84	441	20	0
2	B	351	67	346	14	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	351	67	346	15	0
2	H	351	67	346	13	1
2	K	351	67	346	13	0
3	A	121	0	124	16	0
3	B	66	0	74	13	0
3	D	121	0	124	19	1
3	E	66	0	74	9	0
3	G	121	0	124	11	0
3	H	66	0	74	11	0
3	J	121	0	124	17	1
3	K	66	0	74	9	0
4	A	40	56	56	2	0
4	D	40	56	56	3	0
4	G	40	56	56	1	0
4	J	40	56	56	1	0
5	A	15	0	29	1	0
5	B	30	0	58	2	0
5	D	15	0	29	0	0
5	E	30	0	58	2	0
5	G	15	0	29	1	0
5	H	30	0	58	2	0
5	J	15	0	29	0	0
5	K	30	0	58	2	0
6	A	20	0	32	1	0
6	B	10	0	16	0	0
6	D	70	0	111	17	0
6	E	30	0	48	11	0
6	G	20	0	32	0	0
6	H	10	0	16	0	0
6	J	50	0	79	5	0
6	K	30	0	48	13	0
7	A	4	8	0	0	0
7	B	9	16	0	0	1
7	D	8	16	0	0	0
7	E	11	20	0	0	0
7	G	4	8	0	0	0
7	H	10	18	0	0	0
7	J	7	14	0	0	0
7	K	11	20	0	0	0
All	All	4488	948	4894	205	3

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 22.

All (205) 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)
3:A:57:BCL:OBB	3:A:57:BCL:HHC	1.42	1.18
3:D:57:BCL:HHC	3:D:57:BCL:OBB	1.42	1.12
3:J:57:BCL:OBB	3:J:57:BCL:HHC	1.42	1.12
3:G:57:BCL:HHC	3:G:57:BCL:OBB	1.42	1.10
1:D:48:LEU:HD12	6:D:100:HTO:H42	1.29	1.10
3:B:59:BCL:HMC3	3:D:57:BCL:HBB1	1.36	1.06
3:J:58:BCL:OBB	3:J:58:BCL:HHC	1.51	1.05
3:H:59:BCL:OBB	3:H:59:BCL:HHC	1.53	1.05
3:A:58:BCL:HHC	3:A:58:BCL:OBB	1.51	1.04
6:D:98:HTO:H42	1:J:48:LEU:HD12	1.32	1.04
3:G:58:BCL:OBB	3:G:58:BCL:HHC	1.51	1.04
3:B:59:BCL:HHC	3:B:59:BCL:OBB	1.53	1.04
3:D:58:BCL:HHC	3:D:58:BCL:OBB	1.51	1.03
3:E:59:BCL:HHC	3:E:59:BCL:OBB	1.53	1.03
3:K:59:BCL:OBB	3:K:59:BCL:HHC	1.53	1.02
1:D:48:LEU:CD1	6:D:100:HTO:H42	1.92	0.97
6:D:98:HTO:H42	1:J:48:LEU:CD1	1.95	0.95
3:G:57:BCL:OBB	3:G:57:BCL:CHC	2.10	0.94
3:H:59:BCL:HMC3	3:J:57:BCL:HBB1	1.47	0.93
3:D:57:BCL:OBB	3:D:57:BCL:CHC	2.10	0.93
3:A:57:BCL:OBB	3:A:57:BCL:CHC	2.10	0.90
3:J:57:BCL:OBB	3:J:57:BCL:CHC	2.10	0.86
3:E:59:BCL:OBB	3:E:59:BCL:CHC	2.11	0.86
3:B:59:BCL:OBB	3:B:59:BCL:CHC	2.11	0.83
3:K:59:BCL:OBB	3:K:59:BCL:CHC	2.11	0.81
3:H:59:BCL:OBB	3:H:59:BCL:CHC	2.11	0.79
3:D:58:BCL:OBB	3:D:58:BCL:CHC	2.21	0.77
1:D:48:LEU:CD1	6:D:100:HTO:C4	2.65	0.75
1:G:1:SER:HB2	1:G:4:LYS:HE2	1.70	0.74
1:D:1:SER:HB2	1:D:4:LYS:HE2	1.70	0.74
1:A:1:SER:HB2	1:A:4:LYS:HE2	1.70	0.74
1:J:1:SER:HB2	1:J:4:LYS:HE2	1.70	0.74
3:E:59:BCL:H51	6:E:79:HTO:H42	1.70	0.73
3:K:59:BCL:H51	6:K:79:HTO:H42	1.70	0.73
3:B:59:BCL:H51	6:E:61:HTO:H42	1.70	0.73
3:H:59:BCL:H51	6:K:61:HTO:H42	1.70	0.72
6:D:98:HTO:C4	1:J:48:LEU:CD1	2.68	0.70
1:G:34:HIS:CE1	3:H:59:BCL:HMD1	2.27	0.69
1:J:34:HIS:CE1	3:K:59:BCL:HMD1	2.27	0.69
1:D:34:HIS:CE1	3:E:59:BCL:HMD1	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:53:SER:HA	1:D:56:LYS:HE2	1.76	0.68
1:J:53:SER:HA	1:J:56:LYS:HE2	1.76	0.68
1:A:34:HIS:CE1	3:B:59:BCL:HMD1	2.27	0.68
1:A:53:SER:HA	1:A:56:LYS:HE2	1.76	0.67
1:G:53:SER:HA	1:G:56:LYS:HE2	1.76	0.67
5:G:75:DET:HM13	6:J:62:HTO:H3	1.75	0.67
3:B:59:BCL:CHB	3:D:57:BCL:HMB3	2.25	0.66
1:D:44:ASN:ND2	1:D:47:ALA:H	1.94	0.65
1:J:44:ASN:ND2	1:J:47:ALA:H	1.94	0.65
1:A:44:ASN:ND2	1:A:47:ALA:H	1.94	0.65
1:G:44:ASN:ND2	1:G:47:ALA:H	1.94	0.65
5:A:75:DET:HM13	6:D:62:HTO:H3	1.80	0.63
2:K:45:PHE:CE2	3:K:59:BCL:HBB1	2.35	0.61
2:E:45:PHE:CE2	3:E:59:BCL:HBB1	2.35	0.61
2:B:45:PHE:HE2	3:B:59:BCL:HBB1	1.66	0.61
2:H:45:PHE:HE2	3:H:59:BCL:HBB1	1.66	0.61
2:H:45:PHE:CE2	3:H:59:BCL:HBB1	2.35	0.61
3:G:58:BCL:OBB	3:G:58:BCL:CHC	2.21	0.60
2:K:45:PHE:HE2	3:K:59:BCL:HBB1	1.66	0.60
2:E:45:PHE:HE2	3:E:59:BCL:HBB1	1.66	0.60
2:B:45:PHE:CE2	3:B:59:BCL:HBB1	2.35	0.60
3:B:59:BCL:CMC	3:D:57:BCL:HBB1	2.22	0.59
3:K:59:BCL:H51	6:K:79:HTO:C4	2.33	0.59
3:B:59:BCL:H51	6:E:61:HTO:C4	2.33	0.59
3:E:59:BCL:H51	6:E:79:HTO:C4	2.33	0.59
3:H:59:BCL:H51	6:K:61:HTO:C4	2.33	0.59
3:A:58:BCL:HBC1	3:D:57:BCL:H91	1.86	0.58
3:J:57:BCL:H202	6:K:61:HTO:H41	1.86	0.57
2:B:3:ARG:HA	2:B:3:ARG:NE	2.19	0.57
3:A:58:BCL:O1A	3:D:57:BCL:H172	2.04	0.57
1:D:37:VAL:HG22	1:D:45:TRP:HZ3	1.69	0.57
2:H:3:ARG:HA	2:H:3:ARG:NE	2.19	0.57
3:J:57:BCL:HMD1	2:K:35:HIS:CE1	2.40	0.57
1:A:37:VAL:HG22	1:A:45:TRP:HZ3	1.69	0.57
2:K:3:ARG:NE	2:K:3:ARG:HA	2.19	0.57
3:A:57:BCL:HMD1	2:B:35:HIS:CE1	2.40	0.57
1:G:37:VAL:HG22	1:G:45:TRP:HZ3	1.69	0.57
2:E:3:ARG:HA	2:E:3:ARG:NE	2.19	0.56
1:J:37:VAL:HG22	1:J:45:TRP:HZ3	1.70	0.56
1:J:44:ASN:C	1:J:44:ASN:HD22	2.08	0.56
3:D:57:BCL:HMD1	2:E:35:HIS:CE1	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:44:ASN:HD22	1:D:44:ASN:C	2.08	0.56
1:G:14:ASN:ND2	1:G:16:SER:H	2.04	0.56
3:G:57:BCL:HMD1	2:H:35:HIS:CE1	2.40	0.56
1:J:14:ASN:ND2	1:J:16:SER:H	2.04	0.56
1:G:53:SER:HA	1:G:56:LYS:CE	2.36	0.56
1:A:14:ASN:ND2	1:A:16:SER:H	2.04	0.56
1:A:53:SER:HA	1:A:56:LYS:CE	2.36	0.56
1:D:14:ASN:ND2	1:D:16:SER:H	2.04	0.56
3:J:58:BCL:OBB	3:J:58:BCL:CHC	2.21	0.55
1:G:44:ASN:C	1:G:44:ASN:HD22	2.08	0.55
3:A:58:BCL:OBB	3:A:58:BCL:CHC	2.21	0.55
1:A:44:ASN:C	1:A:44:ASN:HD22	2.08	0.55
3:D:57:BCL:H102	6:E:61:HTO:H72	1.88	0.55
1:J:53:SER:HA	1:J:56:LYS:CE	2.36	0.55
1:D:53:SER:HA	1:D:56:LYS:CE	2.36	0.55
2:K:39:TRP:HB2	5:K:78:DET:HM13	1.89	0.55
2:E:39:TRP:HB2	5:E:78:DET:HM13	1.89	0.55
2:H:39:TRP:HB2	5:H:78:DET:HM13	1.89	0.54
2:B:39:TRP:HB2	5:B:78:DET:HM13	1.89	0.54
6:D:64:HTO:H52	6:D:100:HTO:H52	1.89	0.54
3:D:57:BCL:HMD1	2:E:35:HIS:ND1	2.23	0.54
3:D:57:BCL:H202	6:E:61:HTO:H41	1.90	0.54
3:J:57:BCL:HMD1	2:K:35:HIS:ND1	2.23	0.53
3:G:57:BCL:HMD1	2:H:35:HIS:ND1	2.23	0.53
3:A:57:BCL:HMD1	2:B:35:HIS:ND1	2.23	0.53
6:D:98:HTO:H2	6:D:98:HTO:H62	1.92	0.52
6:D:64:HTO:H62	6:D:64:HTO:H2	1.92	0.51
6:D:100:HTO:H2	6:D:100:HTO:H62	1.92	0.51
6:J:64:HTO:H2	6:J:64:HTO:H62	1.92	0.51
3:J:57:BCL:H202	6:K:61:HTO:C4	2.40	0.51
3:G:58:BCL:HBC1	3:J:57:BCL:H91	1.93	0.50
1:J:4:LYS:HE3	1:J:5:ASP:OD1	2.12	0.49
1:D:4:LYS:HE3	1:D:5:ASP:OD1	2.12	0.49
2:K:39:TRP:HH2	2:K:45:PHE:HB2	1.77	0.49
2:H:25:SER:O	2:H:29:ILE:HG23	2.13	0.49
3:A:58:BCL:H2C	4:D:97:LYC:H133	1.93	0.49
2:B:39:TRP:HH2	2:B:45:PHE:HB2	1.77	0.49
2:B:25:SER:O	2:B:29:ILE:HG23	2.12	0.49
4:A:97:LYC:H61	2:B:16:VAL:HA	1.95	0.49
4:D:97:LYC:H61	2:E:16:VAL:HA	1.95	0.49
1:G:4:LYS:HE3	1:G:5:ASP:OD1	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:59:BCL:CHB	3:J:57:BCL:HMB3	2.42	0.49
4:J:97:LYC:H61	2:K:16:VAL:HA	1.95	0.49
2:E:39:TRP:HH2	2:E:45:PHE:HB2	1.77	0.49
4:G:97:LYC:H61	2:H:16:VAL:HA	1.95	0.49
1:A:4:LYS:HE3	1:A:5:ASP:OD1	2.12	0.48
6:D:98:HTO:H52	6:J:64:HTO:H52	1.96	0.48
2:E:25:SER:O	2:E:29:ILE:HG23	2.13	0.48
2:K:25:SER:O	2:K:29:ILE:HG23	2.13	0.48
3:J:57:BCL:H171	6:K:61:HTO:H52	1.95	0.48
2:H:39:TRP:HH2	2:H:45:PHE:HB2	1.77	0.48
1:J:51:ALA:HA	6:J:64:HTO:H3	1.43	0.48
1:J:1:SER:O	1:J:4:LYS:HB3	2.14	0.48
3:A:58:BCL:HMD1	2:B:20:PHE:CD1	2.49	0.47
1:G:1:SER:O	1:G:4:LYS:HB3	2.14	0.47
1:A:1:SER:O	1:A:4:LYS:HB3	2.14	0.47
3:J:58:BCL:H62	3:J:58:BCL:H41	1.65	0.47
3:J:58:BCL:HMD1	2:K:20:PHE:CD1	2.49	0.47
1:D:1:SER:O	1:D:4:LYS:HB3	2.14	0.47
3:J:57:BCL:C20	6:K:61:HTO:H41	2.44	0.47
3:G:58:BCL:HMD1	2:H:20:PHE:CD1	2.49	0.47
3:D:58:BCL:HMD1	2:E:20:PHE:CD1	2.49	0.47
1:G:18:TRP:O	1:G:22:ILE:HG23	2.16	0.46
1:A:18:TRP:O	1:A:22:ILE:HG23	2.16	0.46
1:D:18:TRP:O	1:D:22:ILE:HG23	2.16	0.46
3:K:59:BCL:H41	6:K:79:HTO:H73	1.97	0.46
2:B:39:TRP:CH2	2:B:45:PHE:HB2	2.51	0.46
3:E:59:BCL:H41	6:E:79:HTO:H73	1.97	0.46
3:H:59:BCL:H41	6:K:61:HTO:H73	1.97	0.46
1:J:18:TRP:O	1:J:22:ILE:HG23	2.16	0.46
3:B:59:BCL:C4A	3:D:57:BCL:HMB3	2.45	0.46
3:B:59:BCL:H41	6:E:61:HTO:H73	1.97	0.46
5:B:78:DET:H21	5:B:78:DET:H51	1.44	0.46
2:H:39:TRP:CH2	2:H:45:PHE:HB2	2.51	0.46
2:K:39:TRP:CH2	2:K:45:PHE:HB2	2.51	0.46
3:A:58:BCL:H2C	4:D:97:LYC:C13	2.46	0.45
5:K:78:DET:H21	5:K:78:DET:H51	1.44	0.45
2:E:39:TRP:CH2	2:E:45:PHE:HB2	2.51	0.45
3:J:57:BCL:H102	6:K:61:HTO:H72	1.99	0.45
3:A:58:BCL:H62	3:A:58:BCL:H41	1.65	0.44
5:E:78:DET:H21	5:E:78:DET:H51	1.44	0.44
3:K:59:BCL:C5	6:K:79:HTO:H11	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:58:BCL:H62	3:G:58:BCL:H41	1.65	0.44
6:A:66:HTO:H3	6:D:66:HTO:H52	1.98	0.44
2:B:5:LEU:CD2	2:E:8:LEU:HD22	2.48	0.44
1:D:51:ALA:HA	6:D:64:HTO:H3	1.55	0.44
3:B:59:BCL:C5	6:E:61:HTO:H11	2.48	0.44
3:E:59:BCL:C5	6:E:79:HTO:H11	2.48	0.44
4:A:97:LYC:H54	4:A:97:LYC:H521	1.91	0.43
3:A:58:BCL:HMD3	2:B:17:HIS:CE1	2.54	0.43
5:H:78:DET:H21	5:H:78:DET:H51	1.44	0.43
3:H:59:BCL:C5	6:K:61:HTO:H11	2.48	0.43
3:A:57:BCL:H41	3:A:57:BCL:H62	1.74	0.43
6:D:100:HTO:H72	6:D:100:HTO:H12	2.01	0.43
6:D:64:HTO:H12	6:D:64:HTO:H72	2.01	0.43
6:D:98:HTO:H12	6:D:98:HTO:H72	2.01	0.43
6:J:64:HTO:H72	6:J:64:HTO:H12	2.01	0.43
3:D:58:BCL:HMD3	2:E:17:HIS:CE1	2.54	0.43
1:J:7:TYR:CD2	2:K:14:ILE:HG13	2.54	0.42
1:G:7:TYR:CD2	2:H:14:ILE:HG13	2.54	0.42
3:G:58:BCL:HMD3	2:H:17:HIS:CE1	2.54	0.42
3:A:58:BCL:CBC	3:D:57:BCL:H91	2.50	0.42
3:D:58:BCL:H62	3:D:58:BCL:H41	1.65	0.42
3:J:58:BCL:HMD3	2:K:17:HIS:CE1	2.54	0.42
1:A:44:ASN:HD21	1:A:46:ILE:HB	1.84	0.42
1:D:44:ASN:HD21	1:D:46:ILE:HB	1.84	0.42
1:J:48:LEU:HD12	1:J:48:LEU:HA	1.93	0.42
1:D:7:TYR:CD2	2:E:14:ILE:HG13	2.54	0.42
1:G:44:ASN:HD21	1:G:46:ILE:HB	1.84	0.42
1:A:7:TYR:CD2	2:B:14:ILE:HG13	2.54	0.42
1:D:14:ASN:HA	1:D:15:PRO:HD2	1.84	0.42
1:J:44:ASN:HD21	1:J:46:ILE:HB	1.84	0.41
1:D:14:ASN:HD22	1:D:15:PRO:CD	2.33	0.41
3:G:57:BCL:H41	3:G:57:BCL:H62	1.74	0.41
1:G:48:LEU:HD12	1:G:48:LEU:HA	1.93	0.41
1:J:14:ASN:HD22	1:J:15:PRO:CD	2.33	0.41
1:A:14:ASN:HD22	1:A:15:PRO:CD	2.33	0.41
1:G:14:ASN:HD22	1:G:15:PRO:CD	2.33	0.41
1:G:14:ASN:HA	1:G:15:PRO:HD2	1.84	0.41
3:A:58:BCL:OBB	2:E:23:THR:OG1	2.24	0.41
1:A:25:VAL:HG21	1:D:24:ILE:HG12	2.03	0.40
1:J:14:ASN:HA	1:J:15:PRO:HD2	1.84	0.40
1:A:14:ASN:HA	1:A:15:PRO:HD2	1.84	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:57:BCL:H202	6:E:61:HTO:C4	2.50	0.40

All (3) 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)
7:B:79:HOH:O	7:B:79:HOH:O[7_646]	1.27	0.93
2:H:23:THR:HG1	3:J:58:BCL:OBB[3_545]	1.37	0.23
2:B:23:THR:HG1	3:D:58:BCL:OBB[4_555]	1.43	0.17

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	54/56 (96%)	52 (96%)	1 (2%)	1 (2%)	8	10
1	D	54/56 (96%)	52 (96%)	1 (2%)	1 (2%)	8	10
1	G	54/56 (96%)	52 (96%)	1 (2%)	1 (2%)	8	10
1	J	54/56 (96%)	52 (96%)	1 (2%)	1 (2%)	8	10
2	B	41/45 (91%)	40 (98%)	1 (2%)	0	100	100
2	E	41/45 (91%)	40 (98%)	1 (2%)	0	100	100
2	H	41/45 (91%)	40 (98%)	1 (2%)	0	100	100
2	K	41/45 (91%)	40 (98%)	1 (2%)	0	100	100
All	All	380/404 (94%)	368 (97%)	8 (2%)	4 (1%)	14	20

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	41	PRO
1	D	41	PRO

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Mol	Chain	Res	Type
1	G	41	PRO
1	J	41	PRO

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	42/42 (100%)	37 (88%)	5 (12%)	5	6
1	D	42/42 (100%)	37 (88%)	5 (12%)	5	6
1	G	42/42 (100%)	37 (88%)	5 (12%)	5	6
1	J	42/42 (100%)	37 (88%)	5 (12%)	5	6
2	B	36/37 (97%)	33 (92%)	3 (8%)	11	17
2	E	36/37 (97%)	33 (92%)	3 (8%)	11	17
2	H	36/37 (97%)	33 (92%)	3 (8%)	11	17
2	K	36/37 (97%)	33 (92%)	3 (8%)	11	17
All	All	312/316 (99%)	280 (90%)	32 (10%)	7	10

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ILE
1	A	41	PRO
1	A	43	PHE
1	A	44	ASN
1	A	48	LEU
2	B	8	LEU
2	B	29	ILE
2	B	40	VAL
1	D	22	ILE
1	D	41	PRO
1	D	43	PHE
1	D	44	ASN
1	D	48	LEU

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Mol	Chain	Res	Type
2	E	8	LEU
2	E	29	ILE
2	E	40	VAL
1	G	22	ILE
1	G	41	PRO
1	G	43	PHE
1	G	44	ASN
1	G	48	LEU
2	H	8	LEU
2	H	29	ILE
2	H	40	VAL
1	J	22	ILE
1	J	41	PRO
1	J	43	PHE
1	J	44	ASN
1	J	48	LEU
2	K	8	LEU
2	K	29	ILE
2	K	40	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	44	ASN
1	D	14	ASN
1	D	44	ASN
1	G	14	ASN
1	G	44	ASN
1	J	14	ASN
1	J	44	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

52 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
6	HTO	B	63	-	9,9,9	0.85	0	10,10,10	0.77	0
3	BCL	K	59	2	64,74,74	1.43	11 (17%)	78,115,115	2.34	12 (15%)
6	HTO	D	64	1	9,9,9	0.95	0	10,10,10	0.86	0
6	HTO	K	61	-	9,9,9	1.14	1 (11%)	10,10,10	0.45	0
3	BCL	J	57	1	64,74,74	1.28	7 (10%)	78,115,115	2.43	14 (17%)
3	BCL	B	59	2	64,74,74	1.43	11 (17%)	78,115,115	2.34	12 (15%)
5	DET	E	77	-	11,14,14	0.44	0	13,16,16	0.55	0
4	LYC	J	97	-	39,39,39	0.99	1 (2%)	44,46,46	2.06	11 (25%)
3	BCL	A	57	1	64,74,74	1.28	7 (10%)	78,115,115	2.43	14 (17%)
3	BCL	H	59	2	64,74,74	1.43	11 (17%)	78,115,115	2.34	11 (14%)
6	HTO	D	62	-	9,9,9	0.76	0	10,10,10	0.65	0
6	HTO	D	100	1	9,9,9	0.95	0	10,10,10	0.86	0
6	HTO	G	62	-	9,9,9	0.76	0	10,10,10	0.65	0
6	HTO	J	66	-	9,9,9	0.65	0	10,10,10	0.59	0
6	HTO	J	64	1	9,9,9	0.95	0	10,10,10	0.87	0
6	HTO	G	66	-	9,9,9	0.65	0	10,10,10	0.59	0
5	DET	D	75	-	11,14,14	0.56	0	13,16,16	0.47	0
3	BCL	E	59	2	64,74,74	1.43	11 (17%)	78,115,115	2.34	11 (14%)
6	HTO	J	65	-	9,9,9	0.82	0	10,10,10	0.24	0
6	HTO	E	61	-	9,9,9	1.14	1 (11%)	10,10,10	0.45	0
6	HTO	A	62	-	9,9,9	0.76	0	10,10,10	0.65	0
6	HTO	H	63	-	9,9,9	0.85	0	10,10,10	0.77	0
6	HTO	J	62	-	9,9,9	0.75	0	10,10,10	0.65	0
5	DET	B	77	-	11,14,14	0.45	0	13,16,16	0.55	0
5	DET	J	75	-	11,14,14	0.56	0	13,16,16	0.48	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	HTO	D	66	-	9,9,9	0.65	0	10,10,10	0.60	0
4	LYC	A	97	-	39,39,39	0.99	1 (2%)	44,46,46	2.06	11 (25%)
3	BCL	G	57	1	64,74,74	1.28	7 (10%)	78,115,115	2.42	14 (17%)
6	HTO	E	63	-	9,9,9	0.85	0	10,10,10	0.77	0
3	BCL	G	58	1	53,63,74	1.57	9 (16%)	64,101,115	2.51	14 (21%)
5	DET	K	77	-	11,14,14	0.45	0	13,16,16	0.55	0
6	HTO	K	79	-	9,9,9	1.14	1 (11%)	10,10,10	0.45	0
6	HTO	K	63	-	9,9,9	0.86	0	10,10,10	0.77	0
5	DET	K	78	-	11,14,14	0.52	0	13,16,16	0.64	0
3	BCL	J	58	1	53,63,74	1.56	9 (16%)	64,101,115	2.51	14 (21%)
6	HTO	J	98	-	9,9,9	0.82	0	10,10,10	0.24	0
3	BCL	A	58	1	53,63,74	1.56	9 (16%)	64,101,115	2.51	15 (23%)
5	DET	H	78	-	11,14,14	0.52	0	13,16,16	0.64	0
5	DET	G	75	-	11,14,14	0.57	0	13,16,16	0.48	0
4	LYC	G	97	-	39,39,39	0.99	1 (2%)	44,46,46	2.05	11 (25%)
5	DET	B	78	-	11,14,14	0.53	0	13,16,16	0.64	0
5	DET	E	78	-	11,14,14	0.53	0	13,16,16	0.64	0
3	BCL	D	57	1	64,74,74	1.28	7 (10%)	78,115,115	2.43	14 (17%)
3	BCL	D	58	1	53,63,74	1.56	9 (16%)	64,101,115	2.50	15 (23%)
6	HTO	A	66	-	9,9,9	0.65	0	10,10,10	0.60	0
6	HTO	D	98	1	9,9,9	0.95	0	10,10,10	0.86	0
6	HTO	D	65	-	9,9,9	0.82	0	10,10,10	0.23	0
5	DET	A	75	-	11,14,14	0.56	0	13,16,16	0.47	0
5	DET	H	77	-	11,14,14	0.45	0	13,16,16	0.55	0
6	HTO	D	99	-	9,9,9	0.82	0	10,10,10	0.23	0
6	HTO	E	79	-	9,9,9	1.14	1 (11%)	10,10,10	0.45	0
4	LYC	D	97	-	39,39,39	0.99	1 (2%)	44,46,46	2.06	11 (25%)

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
6	HTO	B	63	-	-	1/10/10/10	-
3	BCL	K	59	2	-	3/37/137/137	-
6	HTO	D	64	1	-	0/10/10/10	-
6	HTO	K	61	-	-	1/10/10/10	-
3	BCL	J	57	1	-	12/37/137/137	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	BCL	B	59	2	-	3/37/137/137	-
5	DET	E	77	-	-	2/12/12/12	-
4	LYC	J	97	-	-	7/43/43/43	-
3	BCL	A	57	1	-	13/37/137/137	-
3	BCL	H	59	2	-	3/37/137/137	-
6	HTO	D	62	-	-	1/10/10/10	-
6	HTO	D	100	1	-	0/10/10/10	-
6	HTO	G	62	-	-	1/10/10/10	-
6	HTO	J	66	-	-	0/10/10/10	-
6	HTO	J	64	1	-	0/10/10/10	-
6	HTO	G	66	-	-	0/10/10/10	-
5	DET	D	75	-	-	1/12/12/12	-
3	BCL	E	59	2	-	3/37/137/137	-
6	HTO	J	65	-	-	0/10/10/10	-
6	HTO	E	61	-	-	1/10/10/10	-
6	HTO	A	62	-	-	1/10/10/10	-
6	HTO	H	63	-	-	1/10/10/10	-
6	HTO	J	62	-	-	1/10/10/10	-
5	DET	B	77	-	-	2/12/12/12	-
5	DET	J	75	-	-	1/12/12/12	-
6	HTO	D	66	-	-	0/10/10/10	-
4	LYC	A	97	-	-	7/43/43/43	-
3	BCL	G	57	1	-	12/37/137/137	-
6	HTO	E	63	-	-	1/10/10/10	-
3	BCL	G	58	1	-	7/24/124/137	-
5	DET	K	77	-	-	2/12/12/12	-
6	HTO	K	79	-	-	1/10/10/10	-
6	HTO	K	63	-	-	1/10/10/10	-
5	DET	K	78	-	-	6/12/12/12	-
3	BCL	J	58	1	-	7/24/124/137	-
6	HTO	J	98	-	-	0/10/10/10	-
3	BCL	A	58	1	-	7/24/124/137	-
5	DET	H	78	-	-	6/12/12/12	-
5	DET	G	75	-	-	1/12/12/12	-
4	LYC	G	97	-	-	7/43/43/43	-
5	DET	B	78	-	-	6/12/12/12	-
5	DET	E	78	-	-	6/12/12/12	-
3	BCL	D	57	1	-	12/37/137/137	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	BCL	D	58	1	-	7/24/124/137	-
6	HTO	A	66	-	-	0/10/10/10	-
6	HTO	D	98	1	-	0/10/10/10	-
6	HTO	D	65	-	-	0/10/10/10	-
5	DET	A	75	-	-	1/12/12/12	-
5	DET	H	77	-	-	2/12/12/12	-
6	HTO	D	99	-	-	0/10/10/10	-
6	HTO	E	79	-	-	1/10/10/10	-
4	LYC	D	97	-	-	7/43/43/43	-

All (116) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	K	59	BCL	MG-NA	5.30	2.18	2.06
3	B	59	BCL	MG-NA	5.30	2.18	2.06
3	H	59	BCL	MG-NA	5.30	2.18	2.06
3	E	59	BCL	MG-NA	5.30	2.18	2.06
3	G	58	BCL	O2A-CGA	4.63	1.46	1.33
3	D	58	BCL	O2A-CGA	4.63	1.46	1.33
3	A	58	BCL	O2A-CGA	4.63	1.46	1.33
3	J	58	BCL	O2A-CGA	4.63	1.46	1.33
3	D	57	BCL	MG-NC	4.36	2.16	2.06
3	G	57	BCL	MG-NC	4.36	2.16	2.06
3	A	57	BCL	MG-NC	4.36	2.16	2.06
3	J	57	BCL	MG-NC	4.36	2.16	2.06
3	B	59	BCL	C1B-NB	3.85	1.38	1.35
3	H	59	BCL	C1B-NB	3.85	1.38	1.35
3	K	59	BCL	C1B-NB	3.84	1.38	1.35
3	E	59	BCL	C1B-NB	3.79	1.38	1.35
3	D	57	BCL	C1-C2	-3.71	1.38	1.49
3	A	57	BCL	C1-C2	-3.71	1.38	1.49
3	J	57	BCL	C1-C2	-3.70	1.38	1.49
3	G	57	BCL	C1-C2	-3.68	1.38	1.49
3	D	57	BCL	O2D-CED	-3.54	1.37	1.45
3	G	57	BCL	O2D-CED	-3.52	1.37	1.45
3	A	57	BCL	O2D-CED	-3.52	1.37	1.45
3	J	57	BCL	O2D-CED	-3.50	1.37	1.45
3	A	57	BCL	C4B-NB	3.43	1.38	1.35
3	J	57	BCL	C4B-NB	3.43	1.38	1.35
3	G	57	BCL	C4B-NB	3.43	1.38	1.35
3	D	57	BCL	C4B-NB	3.42	1.38	1.35
3	K	59	BCL	C4B-NB	3.34	1.38	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	59	BCL	C4B-NB	3.33	1.38	1.35
3	E	59	BCL	C4B-NB	3.33	1.38	1.35
3	H	59	BCL	C4B-NB	3.32	1.38	1.35
3	G	58	BCL	O2D-CED	-3.30	1.37	1.45
3	A	58	BCL	O2D-CED	-3.26	1.37	1.45
3	D	58	BCL	O2D-CED	-3.23	1.37	1.45
3	J	58	BCL	O2D-CED	-3.23	1.37	1.45
3	B	59	BCL	MG-ND	-3.21	1.99	2.05
3	D	58	BCL	C2-C3	3.20	1.40	1.33
3	H	59	BCL	MG-ND	-3.20	1.99	2.05
3	J	58	BCL	C2-C3	3.20	1.40	1.33
3	K	59	BCL	MG-ND	-3.20	1.99	2.05
3	A	58	BCL	C2-C3	3.19	1.40	1.33
3	E	59	BCL	MG-ND	-3.19	1.99	2.05
3	G	58	BCL	C2-C3	3.18	1.40	1.33
3	G	58	BCL	MG-NA	3.15	2.13	2.06
3	J	57	BCL	MG-NA	3.14	2.13	2.06
3	G	57	BCL	MG-NA	3.14	2.13	2.06
3	A	58	BCL	MG-NA	3.14	2.13	2.06
3	A	57	BCL	MG-NA	3.14	2.13	2.06
3	J	58	BCL	MG-NA	3.14	2.13	2.06
3	D	57	BCL	MG-NA	3.14	2.13	2.06
3	D	58	BCL	MG-NA	3.14	2.13	2.06
3	D	58	BCL	O2D-CGD	3.12	1.40	1.33
3	J	58	BCL	O2D-CGD	3.12	1.40	1.33
3	G	58	BCL	MG-ND	-3.12	1.99	2.05
3	A	58	BCL	O2D-CGD	3.12	1.40	1.33
3	G	58	BCL	O2D-CGD	3.12	1.40	1.33
3	J	58	BCL	MG-ND	-3.10	1.99	2.05
3	A	58	BCL	MG-ND	-3.09	1.99	2.05
3	D	58	BCL	MG-ND	-3.07	1.99	2.05
3	J	57	BCL	C1B-NB	2.97	1.37	1.35
3	A	57	BCL	C1B-NB	2.96	1.37	1.35
3	D	57	BCL	C1B-NB	2.96	1.37	1.35
3	D	58	BCL	C4B-NB	2.94	1.37	1.35
3	G	58	BCL	C4B-NB	2.94	1.37	1.35
3	G	58	BCL	C1B-NB	2.92	1.37	1.35
3	G	57	BCL	C1B-NB	2.92	1.37	1.35
3	E	59	BCL	O2D-CED	-2.91	1.38	1.45
3	B	59	BCL	O2D-CED	-2.90	1.38	1.45
3	H	59	BCL	O2D-CED	-2.90	1.38	1.45
3	J	58	BCL	C4B-NB	2.90	1.37	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	58	BCL	C1B-NB	2.89	1.37	1.35
3	A	58	BCL	C4B-NB	2.89	1.37	1.35
3	K	59	BCL	O2D-CED	-2.89	1.38	1.45
3	J	58	BCL	C1B-NB	2.89	1.37	1.35
3	D	58	BCL	C1B-NB	2.89	1.37	1.35
3	H	59	BCL	O2D-CGD	2.81	1.40	1.33
3	K	59	BCL	O2D-CGD	2.80	1.40	1.33
3	B	59	BCL	O2D-CGD	2.80	1.40	1.33
3	E	59	BCL	O2D-CGD	2.80	1.40	1.33
3	G	58	BCL	C5-C3	2.76	1.57	1.51
3	J	58	BCL	C5-C3	2.75	1.57	1.51
3	A	58	BCL	C5-C3	2.74	1.57	1.51
3	D	58	BCL	C5-C3	2.71	1.56	1.51
6	K	79	HTO	C3-C2	2.63	1.59	1.52
6	K	61	HTO	C3-C2	2.63	1.59	1.52
3	E	59	BCL	MG-NC	2.63	2.12	2.06
3	H	59	BCL	MG-NC	2.62	2.12	2.06
3	K	59	BCL	MG-NC	2.62	2.12	2.06
3	B	59	BCL	MG-NC	2.62	2.12	2.06
3	E	59	BCL	C2-C3	2.61	1.39	1.33
3	H	59	BCL	C2-C3	2.61	1.39	1.33
6	E	61	HTO	C3-C2	2.61	1.59	1.52
3	B	59	BCL	C2-C3	2.61	1.39	1.33
6	E	79	HTO	C3-C2	2.60	1.59	1.52
3	K	59	BCL	C2-C3	2.59	1.39	1.33
3	H	59	BCL	C1-C2	-2.32	1.42	1.49
3	B	59	BCL	C1-C2	-2.30	1.42	1.49
3	K	59	BCL	C1-C2	-2.29	1.42	1.49
3	E	59	BCL	C1-C2	-2.27	1.42	1.49
4	D	97	LYC	C6-C7	2.24	1.55	1.51
4	G	97	LYC	C6-C7	2.23	1.55	1.51
4	A	97	LYC	C6-C7	2.23	1.55	1.51
4	J	97	LYC	C6-C7	2.21	1.55	1.51
3	D	57	BCL	C2-C3	2.17	1.38	1.33
3	G	57	BCL	C2-C3	2.16	1.38	1.33
3	J	57	BCL	C2-C3	2.16	1.38	1.33
3	A	57	BCL	C2-C3	2.16	1.38	1.33
3	H	59	BCL	C4-C3	2.10	1.56	1.50
3	B	59	BCL	C4-C3	2.09	1.56	1.50
3	E	59	BCL	C4-C3	2.09	1.56	1.50
3	K	59	BCL	C4-C3	2.09	1.56	1.50
3	E	59	BCL	C5-C3	2.07	1.55	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	K	59	BCL	C5-C3	2.05	1.55	1.51
3	B	59	BCL	C5-C3	2.04	1.55	1.51
3	H	59	BCL	C5-C3	2.01	1.55	1.51

All (204) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	57	BCL	OBB-CAB-C3B	-9.79	102.62	119.99
3	J	57	BCL	OBB-CAB-C3B	-9.79	102.62	119.99
3	A	57	BCL	OBB-CAB-C3B	-9.79	102.63	119.99
3	G	57	BCL	OBB-CAB-C3B	-9.78	102.63	119.99
3	E	59	BCL	OBB-CAB-C3B	-9.10	103.85	119.99
3	H	59	BCL	OBB-CAB-C3B	-9.09	103.86	119.99
3	B	59	BCL	OBB-CAB-C3B	-9.08	103.87	119.99
3	K	59	BCL	OBB-CAB-C3B	-9.06	103.92	119.99
3	K	59	BCL	C4B-C3B-CAB	-8.99	109.78	127.13
3	D	57	BCL	C4A-NA-C1A	8.98	110.74	106.71
3	B	59	BCL	C4B-C3B-CAB	-8.98	109.79	127.13
3	H	59	BCL	C4B-C3B-CAB	-8.98	109.79	127.13
3	E	59	BCL	C4B-C3B-CAB	-8.97	109.80	127.13
3	A	57	BCL	C4A-NA-C1A	8.97	110.74	106.71
3	G	57	BCL	C4A-NA-C1A	8.97	110.74	106.71
3	J	57	BCL	C4A-NA-C1A	8.97	110.74	106.71
3	G	58	BCL	C4A-NA-C1A	8.33	110.45	106.71
3	J	58	BCL	C4A-NA-C1A	8.33	110.45	106.71
3	A	58	BCL	C4A-NA-C1A	8.32	110.44	106.71
3	D	58	BCL	C4A-NA-C1A	8.28	110.43	106.71
3	G	58	BCL	C4B-C3B-CAB	-7.64	112.38	127.13
3	D	58	BCL	C4B-C3B-CAB	-7.63	112.39	127.13
3	A	58	BCL	C4B-C3B-CAB	-7.63	112.40	127.13
3	J	58	BCL	C4B-C3B-CAB	-7.62	112.41	127.13
3	G	58	BCL	OBB-CAB-C3B	-7.51	106.67	119.99
3	J	58	BCL	OBB-CAB-C3B	-7.51	106.67	119.99
3	A	58	BCL	OBB-CAB-C3B	-7.50	106.67	119.99
3	D	58	BCL	OBB-CAB-C3B	-7.50	106.69	119.99
3	G	58	BCL	C1C-NC-C4C	7.48	110.07	106.71
3	J	58	BCL	C1C-NC-C4C	7.46	110.06	106.71
3	D	57	BCL	C1C-NC-C4C	7.46	110.06	106.71
3	A	57	BCL	C1C-NC-C4C	7.45	110.06	106.71
3	J	57	BCL	C1C-NC-C4C	7.45	110.05	106.71
3	A	58	BCL	C1C-NC-C4C	7.44	110.05	106.71
3	G	57	BCL	C1C-NC-C4C	7.44	110.05	106.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	58	BCL	C1C-NC-C4C	7.42	110.04	106.71
3	H	59	BCL	C4A-NA-C1A	7.41	110.04	106.71
3	E	59	BCL	C4A-NA-C1A	7.40	110.03	106.71
3	B	59	BCL	C4A-NA-C1A	7.39	110.03	106.71
3	K	59	BCL	C4A-NA-C1A	7.39	110.03	106.71
3	J	57	BCL	C4B-C3B-CAB	-7.02	113.57	127.13
3	G	57	BCL	C4B-C3B-CAB	-7.02	113.58	127.13
3	D	57	BCL	C4B-C3B-CAB	-7.01	113.59	127.13
3	A	57	BCL	C4B-C3B-CAB	-7.01	113.59	127.13
4	J	97	LYC	C5-C6-C7	7.00	136.01	112.98
4	A	97	LYC	C5-C6-C7	6.98	135.95	112.98
4	G	97	LYC	C5-C6-C7	6.98	135.95	112.98
4	D	97	LYC	C5-C6-C7	6.98	135.93	112.98
3	B	59	BCL	C1C-NC-C4C	6.89	109.80	106.71
3	K	59	BCL	C1C-NC-C4C	6.88	109.80	106.71
3	H	59	BCL	C1C-NC-C4C	6.88	109.80	106.71
3	E	59	BCL	C1C-NC-C4C	6.87	109.80	106.71
3	J	57	BCL	O2D-CGD-CBD	6.84	123.42	111.27
3	A	57	BCL	O2D-CGD-CBD	6.84	123.42	111.27
3	G	57	BCL	O2D-CGD-CBD	6.84	123.42	111.27
3	D	57	BCL	O2D-CGD-CBD	6.83	123.41	111.27
3	E	59	BCL	O2D-CGD-CBD	6.60	122.99	111.27
3	B	59	BCL	O2D-CGD-CBD	6.58	122.96	111.27
3	K	59	BCL	O2D-CGD-CBD	6.56	122.93	111.27
3	H	59	BCL	O2D-CGD-CBD	6.56	122.92	111.27
4	D	97	LYC	C8-C7-C6	5.77	124.98	115.27
4	A	97	LYC	C8-C7-C6	5.77	124.98	115.27
4	G	97	LYC	C8-C7-C6	5.76	124.96	115.27
4	J	97	LYC	C8-C7-C6	5.76	124.95	115.27
3	D	58	BCL	O2D-CGD-CBD	5.51	121.06	111.27
3	A	58	BCL	O2D-CGD-CBD	5.51	121.06	111.27
3	G	58	BCL	O2D-CGD-CBD	5.51	121.06	111.27
3	J	58	BCL	O2D-CGD-CBD	5.48	121.01	111.27
3	A	57	BCL	O1D-CGD-CBD	-4.42	115.43	124.48
3	D	57	BCL	O1D-CGD-CBD	-4.42	115.44	124.48
3	G	57	BCL	O1D-CGD-CBD	-4.42	115.44	124.48
3	J	57	BCL	O1D-CGD-CBD	-4.41	115.46	124.48
3	K	59	BCL	C1-C2-C3	4.36	133.59	126.04
3	B	59	BCL	C1-C2-C3	4.35	133.57	126.04
3	H	59	BCL	C1-C2-C3	4.35	133.57	126.04
3	E	59	BCL	C1-C2-C3	4.34	133.56	126.04
3	G	58	BCL	O2A-CGA-CBA	4.11	124.81	111.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	58	BCL	O2A-CGA-CBA	4.11	124.80	111.91
3	J	58	BCL	O2A-CGA-CBA	4.10	124.78	111.91
3	D	58	BCL	O2A-CGA-CBA	4.10	124.78	111.91
3	G	58	BCL	O1D-CGD-CBD	-4.05	116.19	124.48
3	A	58	BCL	O1D-CGD-CBD	-4.04	116.22	124.48
3	D	58	BCL	O1D-CGD-CBD	-4.03	116.24	124.48
3	J	58	BCL	O1D-CGD-CBD	-4.02	116.26	124.48
3	G	57	BCL	O2A-CGA-CBA	3.57	123.11	111.91
3	D	57	BCL	O2A-CGA-CBA	3.55	123.05	111.91
3	J	57	BCL	O2A-CGA-CBA	3.55	123.05	111.91
3	A	57	BCL	O2A-CGA-CBA	3.55	123.05	111.91
3	E	59	BCL	O1D-CGD-CBD	-3.45	117.43	124.48
3	B	59	BCL	O1D-CGD-CBD	-3.45	117.44	124.48
3	H	59	BCL	O1D-CGD-CBD	-3.44	117.44	124.48
3	K	59	BCL	O1D-CGD-CBD	-3.44	117.44	124.48
4	J	97	LYC	C52-C51-C53	-3.26	112.94	118.08
4	A	97	LYC	C52-C51-C53	-3.26	112.94	118.08
4	D	97	LYC	C52-C51-C53	-3.25	112.95	118.08
4	G	97	LYC	C52-C51-C53	-3.24	112.97	118.08
4	D	97	LYC	C6-C7-C9	-3.22	112.23	121.98
4	A	97	LYC	C6-C7-C9	-3.21	112.25	121.98
4	J	97	LYC	C6-C7-C9	-3.21	112.27	121.98
4	G	97	LYC	C6-C7-C9	-3.20	112.29	121.98
3	J	58	BCL	O2A-CGA-O1A	-3.16	115.61	123.59
3	A	58	BCL	O2A-CGA-O1A	-3.16	115.62	123.59
3	G	58	BCL	O2A-CGA-O1A	-3.16	115.63	123.59
3	D	58	BCL	O2A-CGA-O1A	-3.16	115.63	123.59
4	A	97	LYC	C13-C12-C11	-3.07	113.23	118.08
4	G	97	LYC	C13-C12-C11	-3.07	113.24	118.08
4	J	97	LYC	C13-C12-C11	-3.05	113.27	118.08
4	D	97	LYC	C13-C12-C11	-3.05	113.28	118.08
3	D	57	BCL	C1-C2-C3	2.97	131.17	126.04
3	A	57	BCL	C1-C2-C3	2.95	131.15	126.04
3	J	57	BCL	C1-C2-C3	2.95	131.14	126.04
3	G	57	BCL	C1-C2-C3	2.92	131.09	126.04
4	G	97	LYC	C6-C5-C4	2.91	121.44	111.88
3	J	57	BCL	OBB-CAB-CBB	-2.90	113.63	120.17
3	G	57	BCL	OBB-CAB-CBB	-2.90	113.64	120.17
3	A	57	BCL	OBB-CAB-CBB	-2.90	113.64	120.17
3	D	57	BCL	OBB-CAB-CBB	-2.90	113.65	120.17
4	A	97	LYC	C6-C5-C4	2.89	121.38	111.88
4	J	97	LYC	C6-C5-C4	2.89	121.37	111.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	97	LYC	C6-C5-C4	2.89	121.37	111.88
3	J	58	BCL	CHD-C1D-ND	-2.83	121.85	124.45
3	G	58	BCL	CHD-C1D-ND	-2.82	121.86	124.45
3	K	59	BCL	O2A-CGA-CBA	2.82	120.75	111.91
3	H	59	BCL	O2A-CGA-CBA	2.82	120.74	111.91
3	E	59	BCL	O2A-CGA-CBA	2.81	120.74	111.91
3	B	59	BCL	O2A-CGA-CBA	2.81	120.72	111.91
3	A	58	BCL	CHD-C1D-ND	-2.81	121.87	124.45
3	J	57	BCL	CMB-C2B-C1B	-2.80	124.16	128.46
3	D	57	BCL	CMB-C2B-C1B	-2.80	124.16	128.46
3	A	57	BCL	CMB-C2B-C1B	-2.80	124.16	128.46
3	G	57	BCL	CMB-C2B-C1B	-2.80	124.17	128.46
3	D	58	BCL	CHD-C1D-ND	-2.79	121.89	124.45
3	G	58	BCL	C1-C2-C3	2.77	130.83	126.04
3	A	58	BCL	C1-C2-C3	2.77	130.83	126.04
3	J	58	BCL	C1-C2-C3	2.75	130.80	126.04
3	D	58	BCL	C1-C2-C3	2.75	130.80	126.04
3	E	59	BCL	OBB-CAB-CBB	-2.67	114.16	120.17
3	B	59	BCL	OBB-CAB-CBB	-2.67	114.17	120.17
3	H	59	BCL	OBB-CAB-CBB	-2.67	114.17	120.17
3	K	59	BCL	OBB-CAB-CBB	-2.66	114.17	120.17
3	G	57	BCL	O2A-CGA-O1A	-2.65	116.91	123.59
3	J	57	BCL	O2A-CGA-O1A	-2.63	116.94	123.59
3	A	57	BCL	O2A-CGA-O1A	-2.63	116.95	123.59
3	D	57	BCL	O2A-CGA-O1A	-2.63	116.96	123.59
3	H	59	BCL	C2C-C3C-C4C	2.52	105.12	101.34
4	A	97	LYC	C13-C12-C14	2.52	126.45	122.92
4	D	97	LYC	C21-C20-C19	-2.52	118.31	123.47
4	A	97	LYC	C21-C20-C19	-2.52	118.31	123.47
3	E	59	BCL	C2C-C3C-C4C	2.52	105.11	101.34
4	J	97	LYC	C21-C20-C19	-2.52	118.32	123.47
4	G	97	LYC	C13-C12-C14	2.52	126.45	122.92
4	D	97	LYC	C13-C12-C14	2.52	126.45	122.92
4	J	97	LYC	C13-C12-C14	2.51	126.44	122.92
3	A	57	BCL	CBB-CAB-C3B	2.51	127.80	120.34
3	D	57	BCL	CBB-CAB-C3B	2.51	127.78	120.34
3	G	57	BCL	CBB-CAB-C3B	2.51	127.78	120.34
4	G	97	LYC	C21-C20-C19	-2.51	118.34	123.47
3	J	57	BCL	CBB-CAB-C3B	2.50	127.78	120.34
3	B	59	BCL	C2C-C3C-C4C	2.50	105.09	101.34
3	K	59	BCL	C2C-C3C-C4C	2.49	105.08	101.34
4	J	97	LYC	C3-C2-C4	2.40	129.58	122.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	97	LYC	C3-C2-C4	2.39	129.56	122.65
4	D	97	LYC	C3-C2-C4	2.39	129.56	122.65
3	G	58	BCL	C1D-ND-C4D	-2.39	104.64	106.33
4	A	97	LYC	C3-C2-C4	2.39	129.54	122.65
3	J	57	BCL	C3A-C2A-C1A	2.38	104.90	101.34
3	J	58	BCL	C1D-ND-C4D	-2.38	104.64	106.33
3	D	57	BCL	C3A-C2A-C1A	2.38	104.90	101.34
3	A	57	BCL	C3A-C2A-C1A	2.37	104.89	101.34
3	G	57	BCL	C3A-C2A-C1A	2.36	104.88	101.34
3	A	58	BCL	C4D-CHA-C1A	2.36	124.12	121.25
3	J	58	BCL	C4D-CHA-C1A	2.36	124.12	121.25
3	A	58	BCL	C1D-ND-C4D	-2.35	104.66	106.33
3	D	57	BCL	C2C-C3C-C4C	2.35	104.86	101.34
3	J	57	BCL	C2C-C3C-C4C	2.35	104.86	101.34
3	G	58	BCL	C4D-CHA-C1A	2.35	124.10	121.25
3	D	58	BCL	C1D-ND-C4D	-2.34	104.67	106.33
3	D	58	BCL	C4D-CHA-C1A	2.34	124.10	121.25
3	A	57	BCL	C2C-C3C-C4C	2.34	104.84	101.34
3	G	57	BCL	C2C-C3C-C4C	2.33	104.84	101.34
4	J	97	LYC	C57-C56-C58	-2.20	114.62	118.08
4	A	97	LYC	C57-C56-C55	2.19	126.00	122.92
4	D	97	LYC	C57-C56-C55	2.19	125.99	122.92
3	E	59	BCL	O2D-CGD-O1D	-2.19	119.56	123.84
4	A	97	LYC	C57-C56-C58	-2.19	114.63	118.08
4	J	97	LYC	C57-C56-C55	2.19	125.99	122.92
4	D	97	LYC	C57-C56-C58	-2.19	114.63	118.08
4	G	97	LYC	C57-C56-C58	-2.18	114.64	118.08
4	G	97	LYC	C57-C56-C55	2.18	125.98	122.92
3	B	59	BCL	O2D-CGD-O1D	-2.17	119.59	123.84
3	K	59	BCL	O2D-CGD-O1D	-2.16	119.61	123.84
3	H	59	BCL	O2D-CGD-O1D	-2.16	119.62	123.84
3	D	58	BCL	CED-O2D-CGD	2.08	120.65	115.94
3	G	58	BCL	CED-O2D-CGD	2.08	120.64	115.94
3	A	58	BCL	CED-O2D-CGD	2.08	120.63	115.94
3	J	58	BCL	CBB-CAB-C3B	2.07	126.49	120.34
3	A	58	BCL	CBB-CAB-C3B	2.06	126.47	120.34
3	J	58	BCL	CED-O2D-CGD	2.06	120.59	115.94
3	D	58	BCL	CBB-CAB-C3B	2.05	126.44	120.34
3	G	58	BCL	CBB-CAB-C3B	2.05	126.43	120.34
3	D	58	BCL	CBC-CAC-C3C	2.01	117.94	113.47
3	A	58	BCL	CBC-CAC-C3C	2.00	117.92	113.47
3	B	59	BCL	C4D-CHA-C1A	2.00	123.68	121.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	K	59	BCL	C4D-CHA-C1A	2.00	123.68	121.25

There are no chirality outliers.

All (165) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	58	BCL	C2-C3-C5-C6
3	A	58	BCL	C4-C3-C5-C6
3	D	58	BCL	C2-C3-C5-C6
3	D	58	BCL	C4-C3-C5-C6
3	G	58	BCL	C2-C3-C5-C6
3	G	58	BCL	C4-C3-C5-C6
3	J	58	BCL	C2-C3-C5-C6
3	J	58	BCL	C4-C3-C5-C6
4	A	97	LYC	C1-C2-C4-C5
4	A	97	LYC	C57-C56-C58-C59
4	D	97	LYC	C1-C2-C4-C5
4	D	97	LYC	C57-C56-C58-C59
4	G	97	LYC	C1-C2-C4-C5
4	G	97	LYC	C57-C56-C58-C59
4	J	97	LYC	C1-C2-C4-C5
4	J	97	LYC	C57-C56-C58-C59
5	B	78	DET	C2-C1-N-O
5	B	78	DET	C2-C1-N-CM1
5	E	78	DET	C2-C1-N-O
5	E	78	DET	C2-C1-N-CM1
5	H	78	DET	C2-C1-N-O
5	H	78	DET	C2-C1-N-CM1
5	K	78	DET	C2-C1-N-O
5	K	78	DET	C2-C1-N-CM1
3	A	58	BCL	CBA-CGA-O2A-C1
3	D	58	BCL	CBA-CGA-O2A-C1
3	G	58	BCL	CBA-CGA-O2A-C1
3	J	58	BCL	CBA-CGA-O2A-C1
3	A	57	BCL	CBD-CGD-O2D-CED
3	D	57	BCL	CBD-CGD-O2D-CED
3	G	57	BCL	CBD-CGD-O2D-CED
3	J	57	BCL	CBD-CGD-O2D-CED
3	A	58	BCL	O1A-CGA-O2A-C1
3	D	58	BCL	O1A-CGA-O2A-C1
3	G	58	BCL	O1A-CGA-O2A-C1
3	J	58	BCL	O1A-CGA-O2A-C1

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Mol	Chain	Res	Type	Atoms
4	A	97	LYC	C5-C6-C7-C9
4	D	97	LYC	C5-C6-C7-C9
4	G	97	LYC	C5-C6-C7-C9
4	J	97	LYC	C5-C6-C7-C9
5	B	78	DET	C2-C3-C4-C5
5	E	78	DET	C2-C3-C4-C5
5	H	78	DET	C2-C3-C4-C5
5	K	78	DET	C2-C3-C4-C5
4	A	97	LYC	C5-C6-C7-C8
4	D	97	LYC	C5-C6-C7-C8
4	G	97	LYC	C5-C6-C7-C8
4	J	97	LYC	C5-C6-C7-C8
3	A	57	BCL	C13-C15-C16-C17
3	D	57	BCL	C13-C15-C16-C17
3	G	57	BCL	C13-C15-C16-C17
3	J	57	BCL	C13-C15-C16-C17
3	J	57	BCL	O1D-CGD-O2D-CED
3	A	57	BCL	O1D-CGD-O2D-CED
3	D	57	BCL	O1D-CGD-O2D-CED
3	G	57	BCL	O1D-CGD-O2D-CED
3	A	57	BCL	C3-C5-C6-C7
3	D	57	BCL	C3-C5-C6-C7
3	G	57	BCL	C3-C5-C6-C7
3	J	57	BCL	C3-C5-C6-C7
4	A	97	LYC	C3-C2-C4-C5
4	D	97	LYC	C3-C2-C4-C5
4	G	97	LYC	C3-C2-C4-C5
4	J	97	LYC	C3-C2-C4-C5
3	A	57	BCL	C6-C7-C8-C10
3	D	57	BCL	C6-C7-C8-C10
3	G	57	BCL	C6-C7-C8-C10
3	J	57	BCL	C6-C7-C8-C10
5	B	77	DET	C6-C7-C8-C9
5	E	77	DET	C6-C7-C8-C9
5	H	77	DET	C6-C7-C8-C9
5	K	77	DET	C6-C7-C8-C9
4	A	97	LYC	C55-C56-C58-C59
4	D	97	LYC	C55-C56-C58-C59
4	G	97	LYC	C55-C56-C58-C59
4	J	97	LYC	C55-C56-C58-C59
5	B	78	DET	C1-C2-C3-C4
5	H	78	DET	C1-C2-C3-C4

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Mol	Chain	Res	Type	Atoms
5	K	78	DET	C1-C2-C3-C4
5	E	78	DET	C1-C2-C3-C4
3	A	57	BCL	C8-C10-C11-C12
3	D	57	BCL	C8-C10-C11-C12
3	G	57	BCL	C8-C10-C11-C12
3	J	57	BCL	C8-C10-C11-C12
3	B	59	BCL	C3A-C2A-CAA-CBA
3	E	59	BCL	C3A-C2A-CAA-CBA
3	H	59	BCL	C3A-C2A-CAA-CBA
3	K	59	BCL	C3A-C2A-CAA-CBA
3	A	58	BCL	C6-C7-C8-C9
3	D	58	BCL	C6-C7-C8-C9
3	G	58	BCL	C6-C7-C8-C9
3	J	58	BCL	C6-C7-C8-C9
3	A	57	BCL	C6-C7-C8-C9
3	D	57	BCL	C6-C7-C8-C9
3	G	57	BCL	C6-C7-C8-C9
3	J	57	BCL	C6-C7-C8-C9
3	B	59	BCL	C8-C10-C11-C12
3	E	59	BCL	C8-C10-C11-C12
3	H	59	BCL	C8-C10-C11-C12
3	K	59	BCL	C8-C10-C11-C12
6	B	63	HTO	C3-C4-C5-C6
6	E	63	HTO	C3-C4-C5-C6
6	H	63	HTO	C3-C4-C5-C6
6	K	63	HTO	C3-C4-C5-C6
3	J	57	BCL	CBA-CGA-O2A-C1
3	A	57	BCL	CBA-CGA-O2A-C1
3	D	57	BCL	CBA-CGA-O2A-C1
3	G	57	BCL	CBA-CGA-O2A-C1
4	G	97	LYC	C18-C17-C19-C20
5	J	75	DET	C7-C8-C9-C10
5	A	75	DET	C7-C8-C9-C10
5	G	75	DET	C7-C8-C9-C10
5	D	75	DET	C7-C8-C9-C10
5	B	78	DET	C11-C10-C9-C8
5	E	78	DET	C11-C10-C9-C8
5	K	78	DET	C11-C10-C9-C8
5	H	78	DET	C11-C10-C9-C8
4	A	97	LYC	C18-C17-C19-C20
4	D	97	LYC	C18-C17-C19-C20
4	J	97	LYC	C18-C17-C19-C20

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Mol	Chain	Res	Type	Atoms
3	A	57	BCL	O1A-CGA-O2A-C1
3	G	57	BCL	O1A-CGA-O2A-C1
3	J	57	BCL	O1A-CGA-O2A-C1
3	D	57	BCL	O1A-CGA-O2A-C1
3	A	58	BCL	C2-C1-O2A-CGA
3	D	58	BCL	C2-C1-O2A-CGA
3	G	58	BCL	C2-C1-O2A-CGA
3	J	58	BCL	C2-C1-O2A-CGA
3	A	57	BCL	CAD-CBD-CGD-O2D
3	D	57	BCL	CAD-CBD-CGD-O2D
3	G	57	BCL	CAD-CBD-CGD-O2D
3	J	57	BCL	CAD-CBD-CGD-O2D
5	B	77	DET	C2-C1-N-CM2
5	B	78	DET	C2-C1-N-CM2
5	E	77	DET	C2-C1-N-CM2
5	E	78	DET	C2-C1-N-CM2
5	H	77	DET	C2-C1-N-CM2
5	H	78	DET	C2-C1-N-CM2
5	K	77	DET	C2-C1-N-CM2
5	K	78	DET	C2-C1-N-CM2
3	A	57	BCL	CAA-CBA-CGA-O2A
3	D	57	BCL	CAA-CBA-CGA-O2A
3	G	57	BCL	CAA-CBA-CGA-O2A
3	J	57	BCL	CAA-CBA-CGA-O2A
6	A	62	HTO	C3-C4-C5-C6
6	D	62	HTO	C3-C4-C5-C6
6	E	61	HTO	C3-C4-C5-C6
6	E	79	HTO	C3-C4-C5-C6
6	G	62	HTO	C3-C4-C5-C6
6	J	62	HTO	C3-C4-C5-C6
6	K	61	HTO	C3-C4-C5-C6
6	K	79	HTO	C3-C4-C5-C6
3	A	58	BCL	C1A-C2A-CAA-CBA
3	B	59	BCL	C1A-C2A-CAA-CBA
3	D	58	BCL	C1A-C2A-CAA-CBA
3	E	59	BCL	C1A-C2A-CAA-CBA
3	G	58	BCL	C1A-C2A-CAA-CBA
3	H	59	BCL	C1A-C2A-CAA-CBA
3	J	58	BCL	C1A-C2A-CAA-CBA
3	K	59	BCL	C1A-C2A-CAA-CBA
3	A	57	BCL	CAA-CBA-CGA-O1A
3	D	57	BCL	CAA-CBA-CGA-O1A

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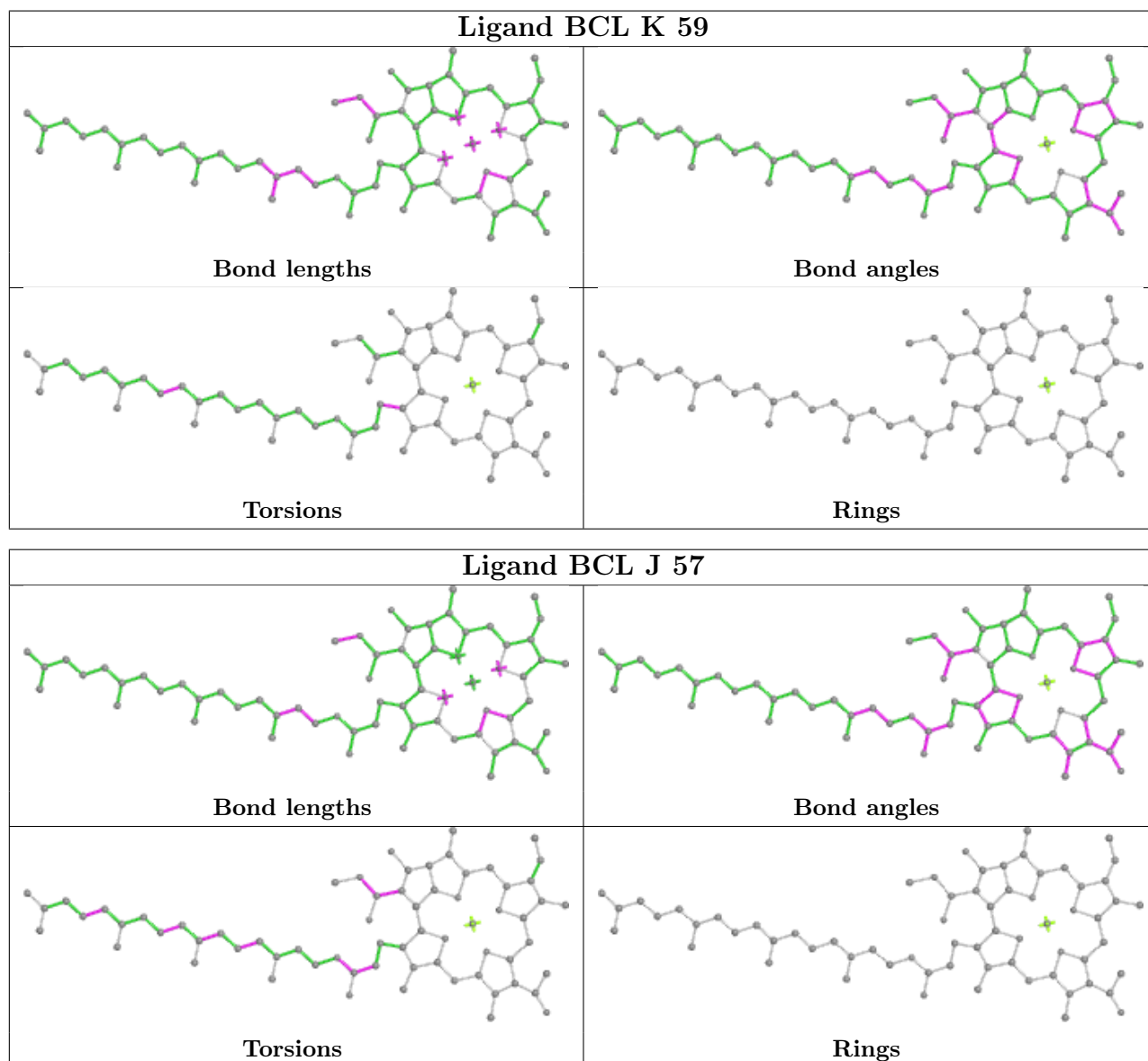
Mol	Chain	Res	Type	Atoms
3	G	57	BCL	CAA-CBA-CGA-O1A
3	J	57	BCL	CAA-CBA-CGA-O1A
3	A	57	BCL	C4-C3-C5-C6

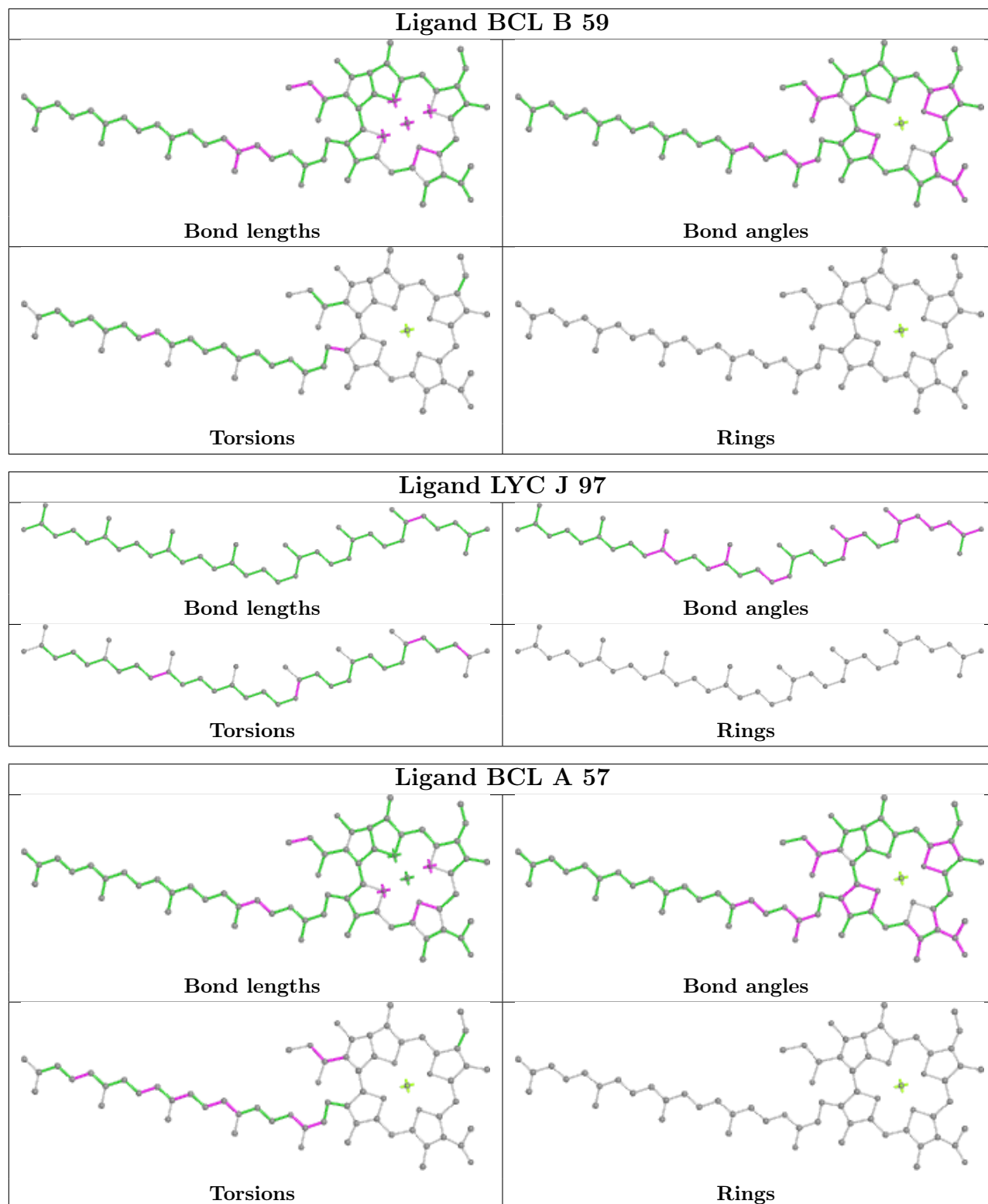
There are no ring outliers.

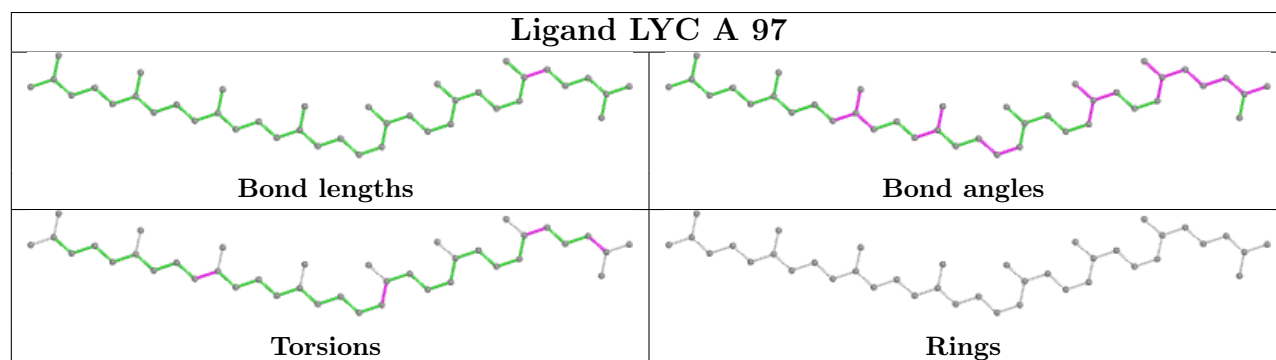
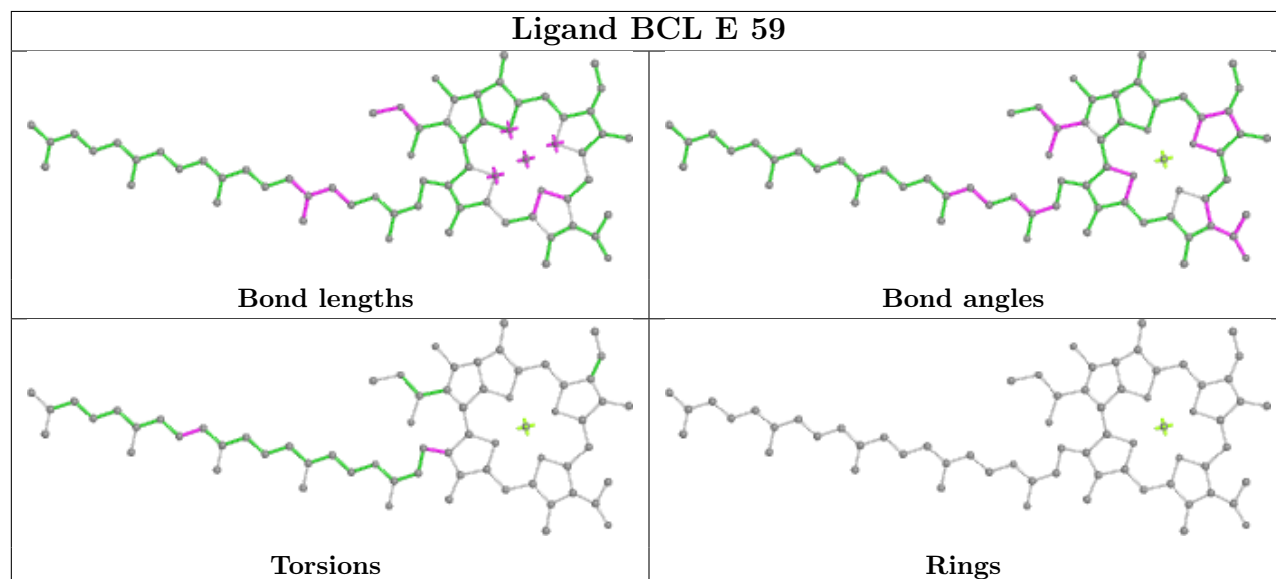
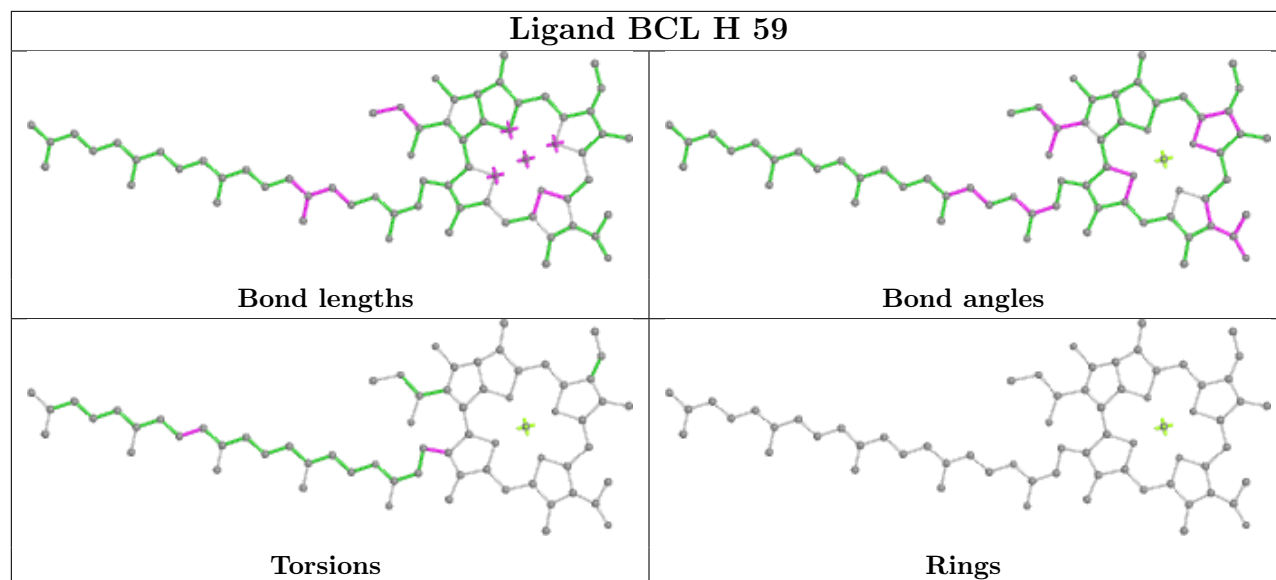
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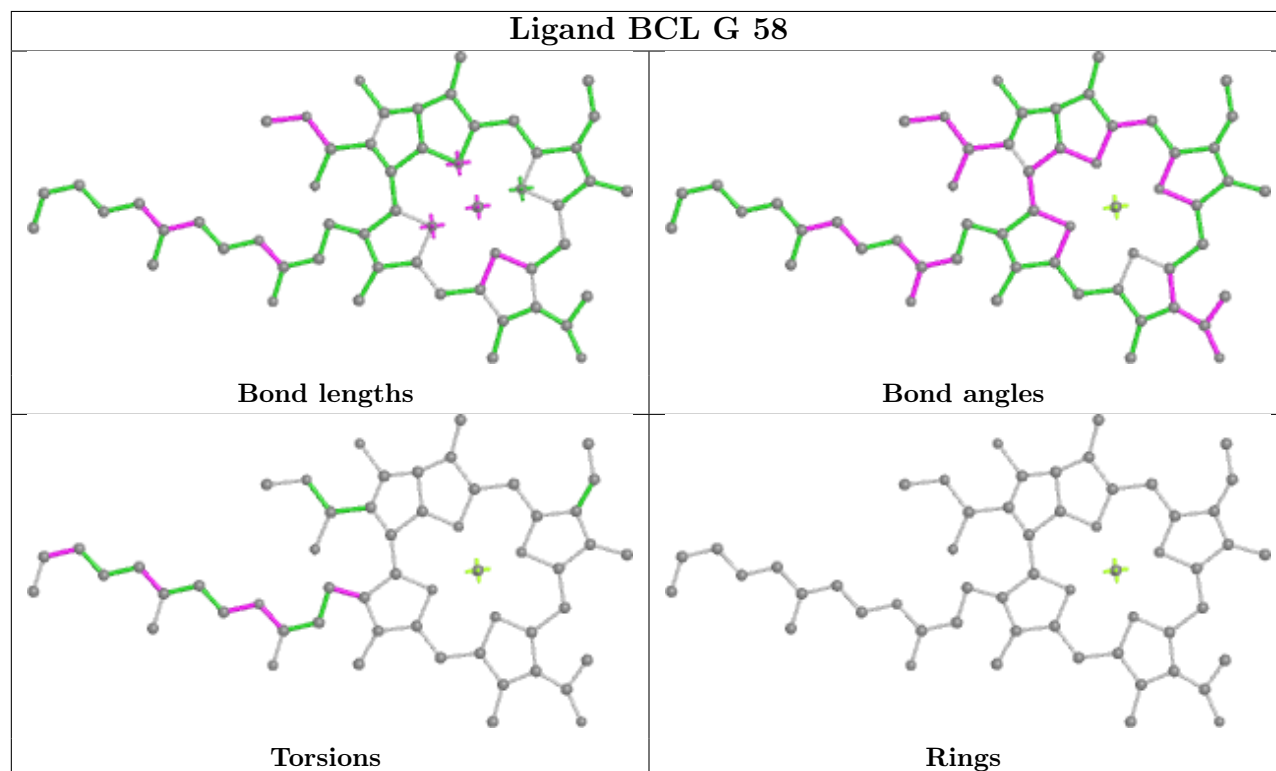
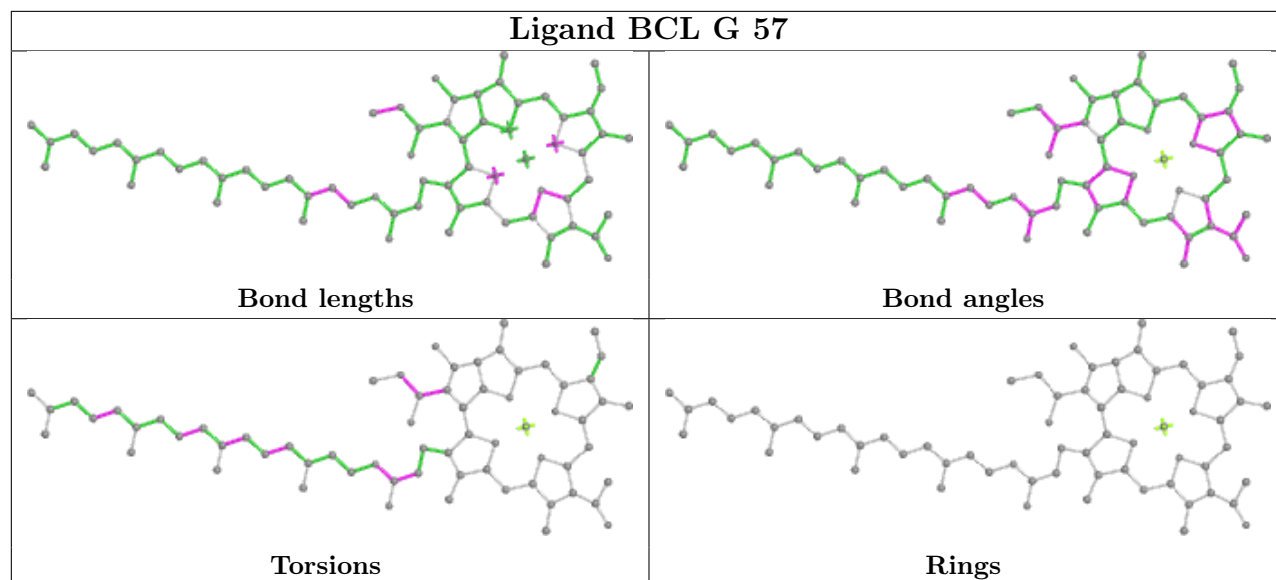
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	K	59	BCL	9	0
6	D	64	HTO	4	0
6	K	61	HTO	9	0
3	J	57	BCL	12	0
3	B	59	BCL	13	0
4	J	97	LYC	1	0
3	A	57	BCL	5	0
3	H	59	BCL	11	0
6	D	62	HTO	1	0
6	D	100	HTO	6	0
6	J	64	HTO	4	0
3	E	59	BCL	9	0
6	E	61	HTO	7	0
6	J	62	HTO	1	0
6	D	66	HTO	1	0
4	A	97	LYC	2	0
3	G	57	BCL	5	0
3	G	58	BCL	6	0
6	K	79	HTO	4	0
5	K	78	DET	2	0
3	J	58	BCL	5	1
3	A	58	BCL	11	0
5	H	78	DET	2	0
5	G	75	DET	1	0
4	G	97	LYC	1	0
5	B	78	DET	2	0
5	E	78	DET	2	0
3	D	57	BCL	14	0
3	D	58	BCL	5	1
6	A	66	HTO	1	0
6	D	98	HTO	6	0
5	A	75	DET	1	0
6	E	79	HTO	4	0
4	D	97	LYC	3	0

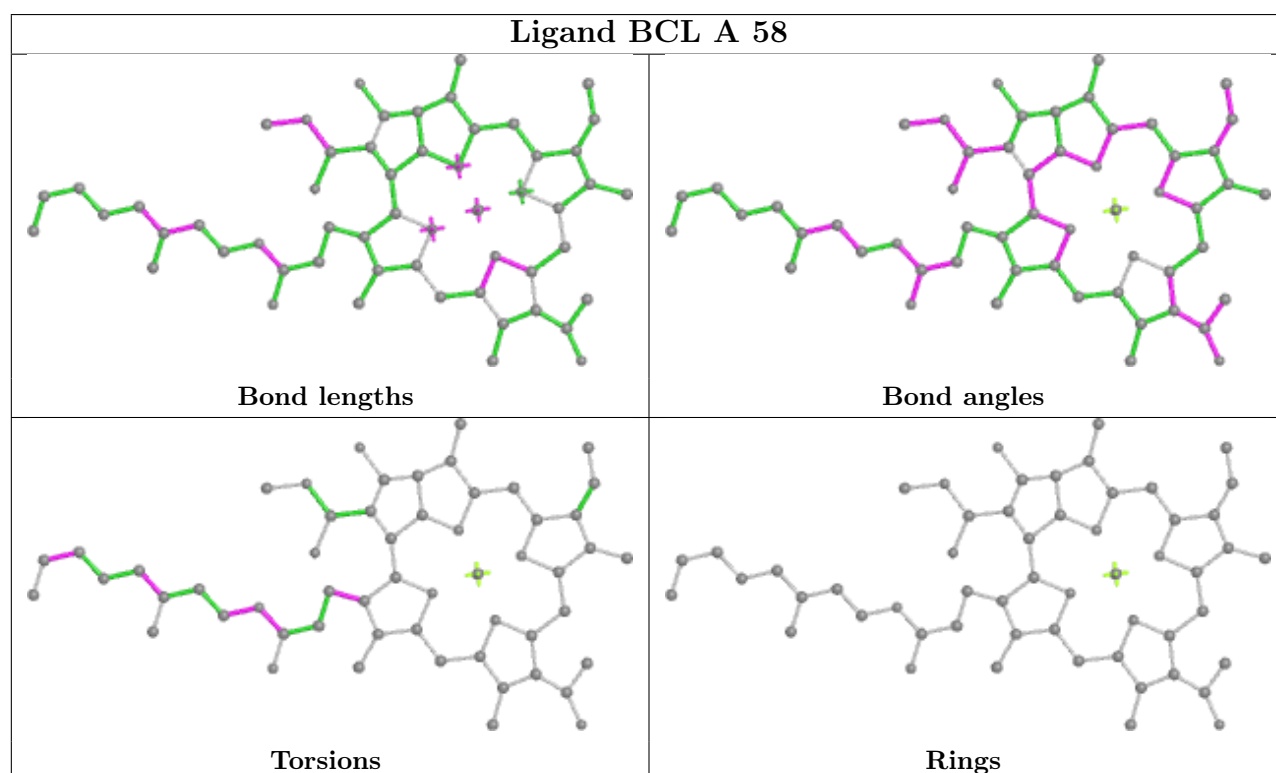
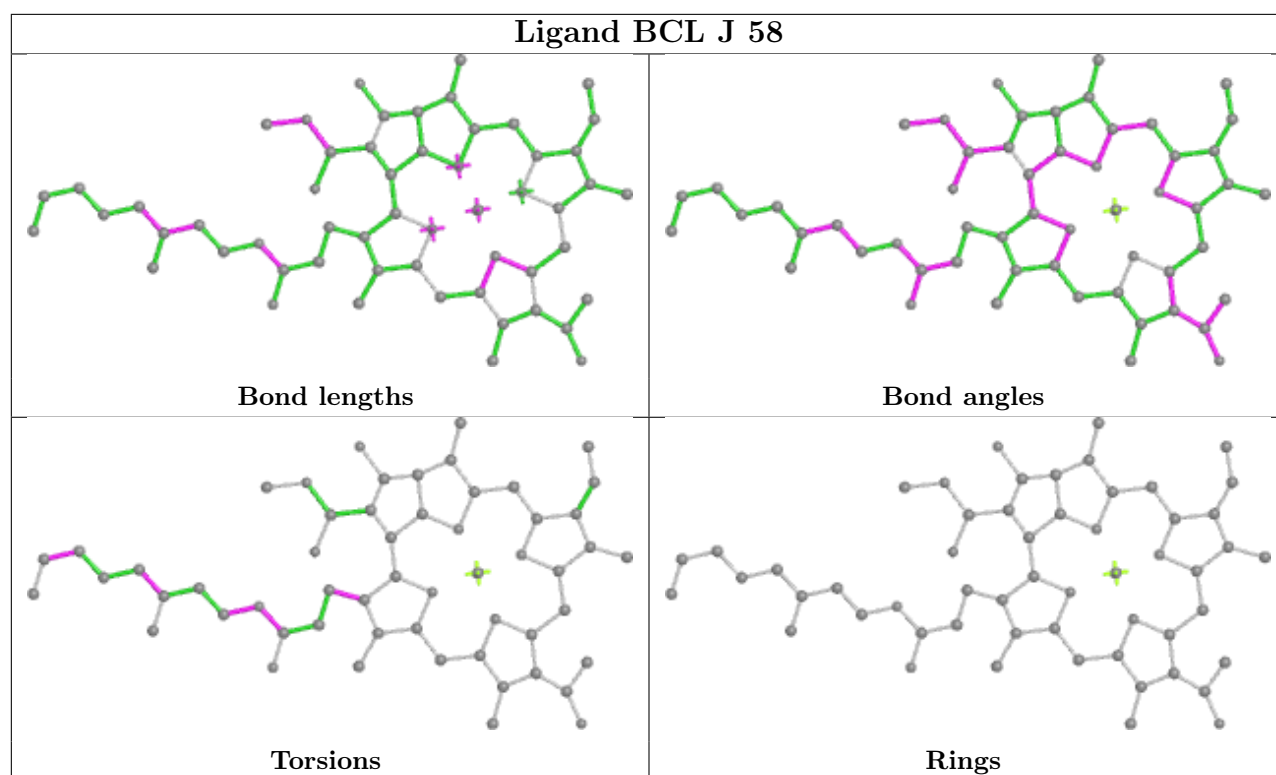
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

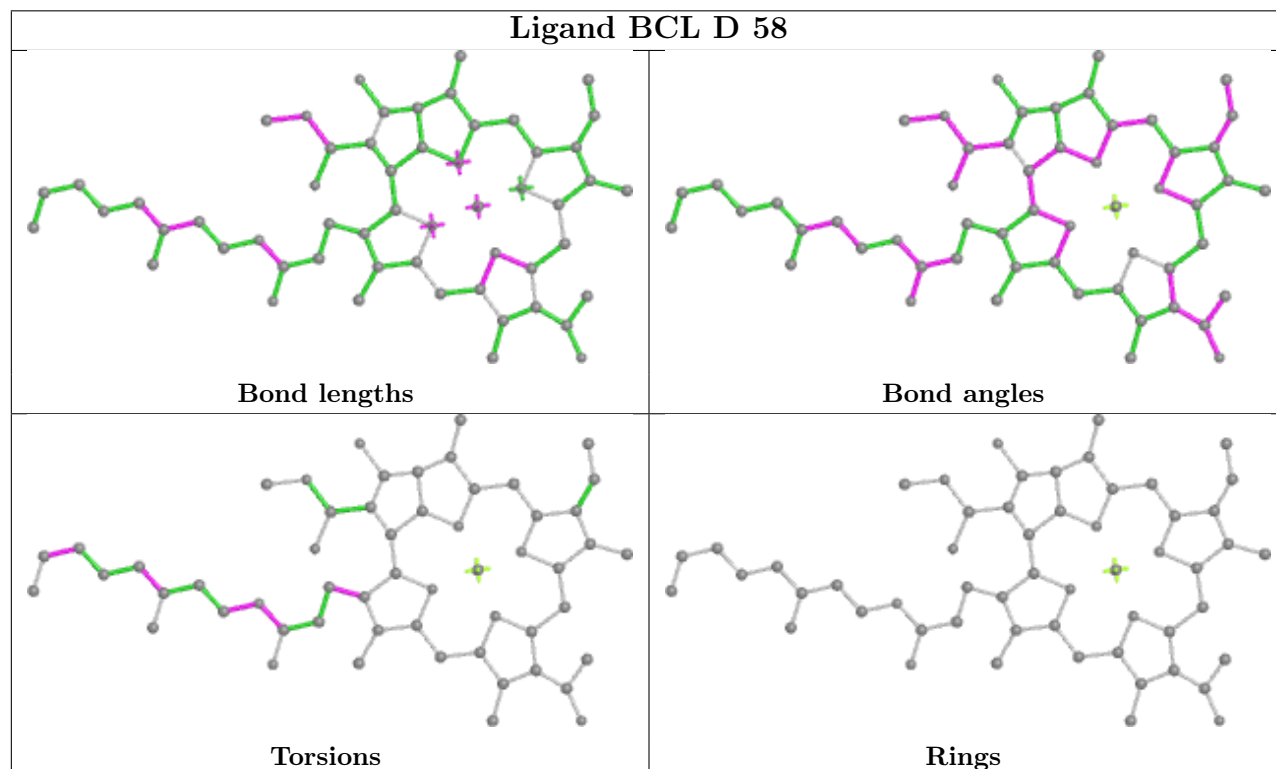
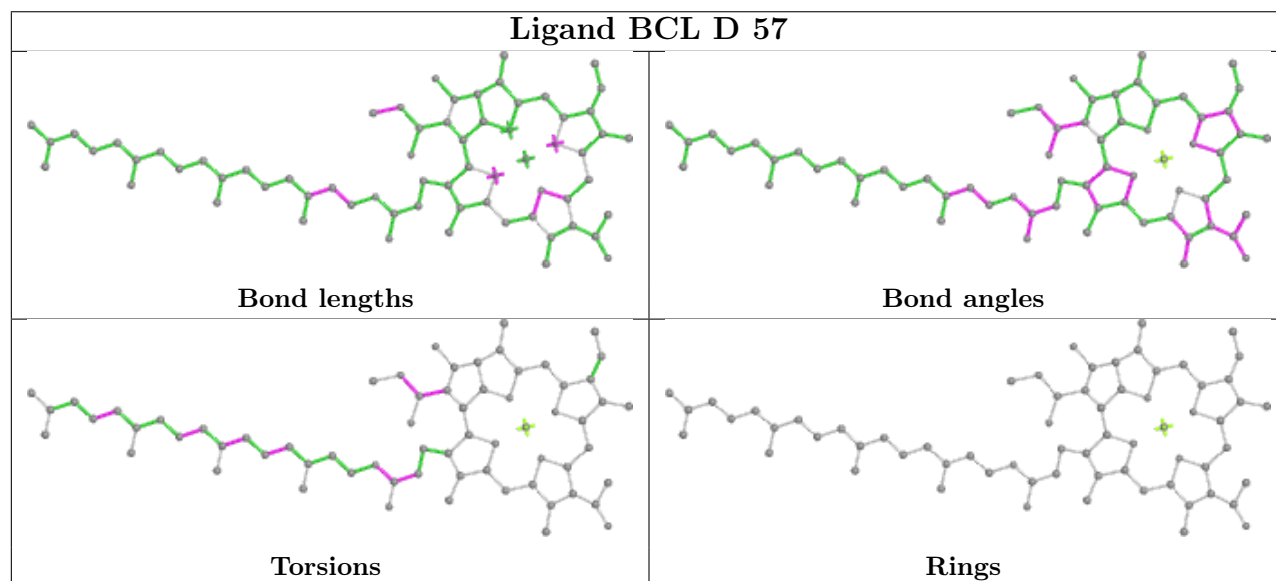
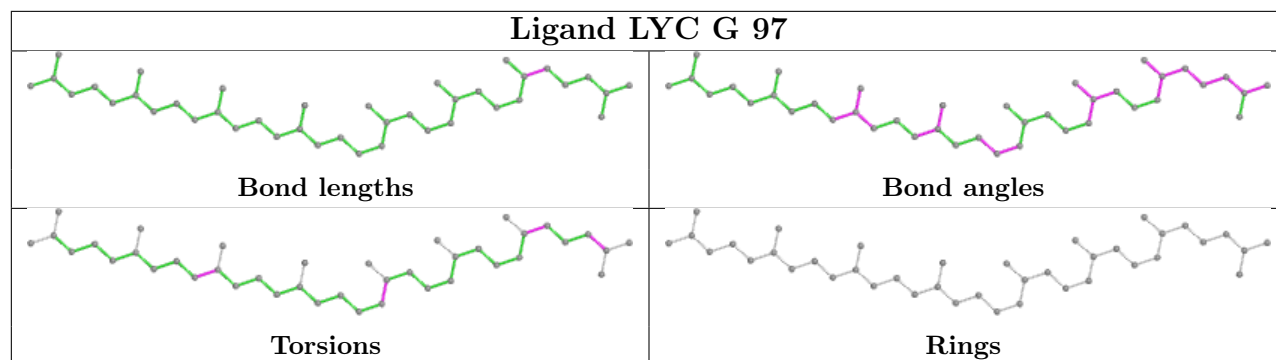


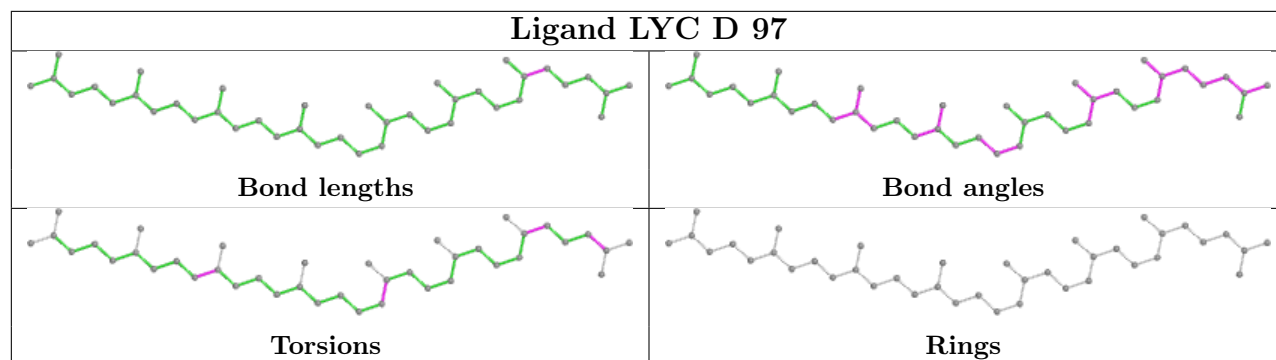












5.7 Other polymers [i](#)

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.