



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:37 PM UTC

PDB ID : 1HZX / pdb_00001hzz
Title : CRYSTAL STRUCTURE OF BOVINE RHODOPSIN
Authors : Teller, D.C.; Okada, T.; Behnke, C.A.; Palczewski, K.; Stenkamp, R.E.
Deposited on : 2001-01-26
Resolution : 2.80 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

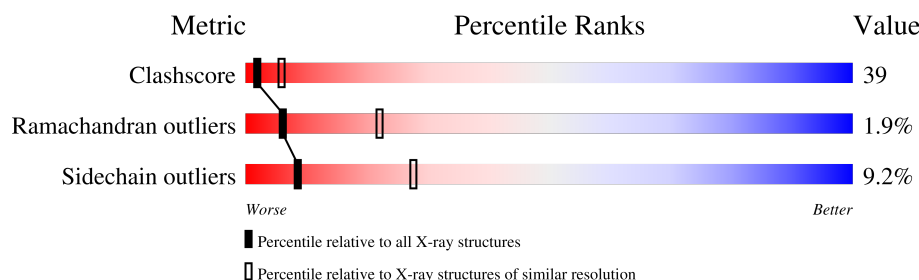
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)

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	349	38% 48% 11% ..
1	B	349	34% 44% 8% 13%
2	C	3	33% 67%
3	D	2	50% 50%
3	F	2	100%
4	E	4	50% 25% 25%

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
5	BNG	B	1506	-	-	X	-

2 Entry composition [i](#)

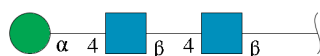
There are 11 unique types of molecules in this entry. The entry contains 5551 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 RHODOPSIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	0	0	0
			2685	1783	413	463	26			
1	B	302	Total	C	N	O	S	0	0	0
			2398	1603	366	404	25			

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



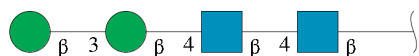
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



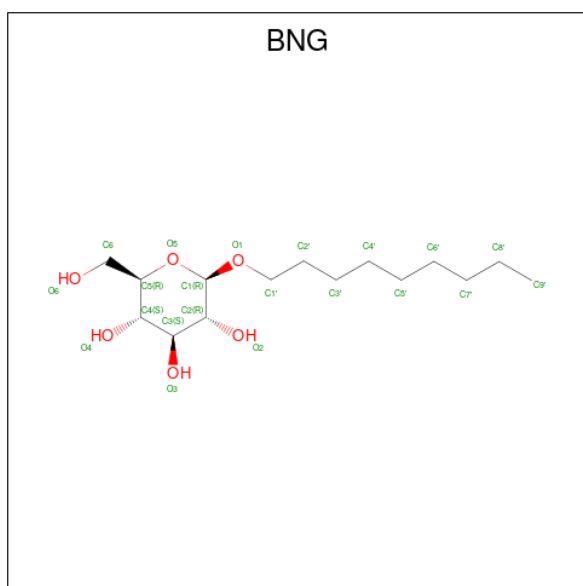
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	F	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is an oligosaccharide called beta-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 5 is nonyl beta-D-glucopyranoside (CCD ID: BNG) (formula: C₁₅H₃₀O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			21	15	6		
5	A	1	Total	C	O	0	0
			21	15	6		
5	A	1	Total	C	O	0	0
			21	15	6		
5	A	1	Total	C	O	0	0
			21	15	6		
5	A	1	Total	C	O	0	0
			21	15	6		
5	B	1	Total	C	O	0	0
			21	15	6		
5	B	1	Total	C	O	0	0
			21	15	6		

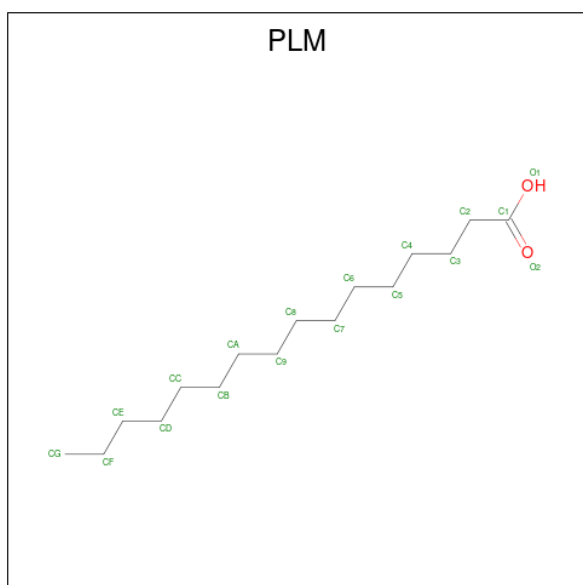
- Molecule 6 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	3	Total Hg 3 3	0	0
6	B	3	Total Hg 3 3	0	0

- Molecule 7 is ZINC ION (CCD ID: ZN) (formula: Zn).

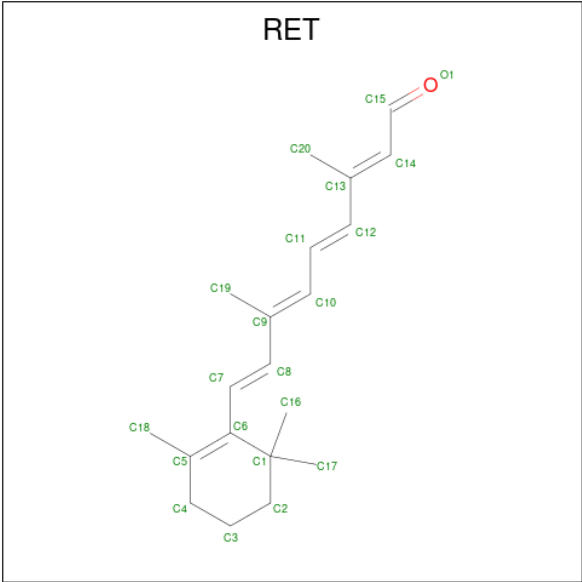
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	4	Total Zn 4 4	0	0
7	B	3	Total Zn 3 3	0	0

- Molecule 8 is PALMITIC ACID (CCD ID: PLM) (formula: C₁₆H₃₂O₂).



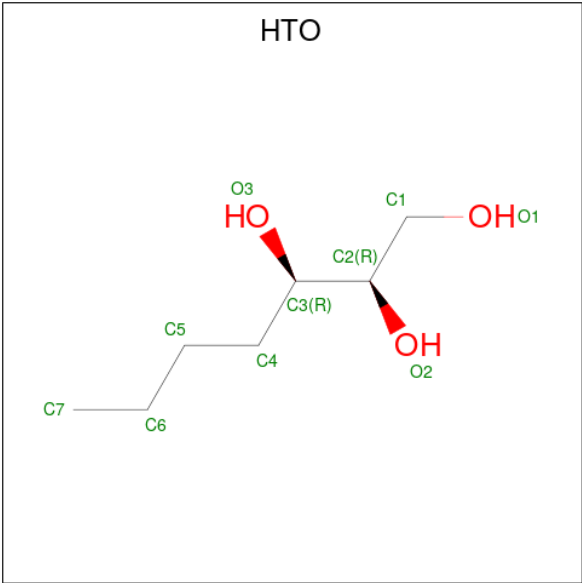
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 17 16 1	0	0
8	A	1	Total C O 17 16 1	0	0
8	B	1	Total C O 17 16 1	0	0

- Molecule 9 is RETINAL (CCD ID: RET) (formula: C₂₀H₂₈O).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	1	Total	C	0	0
			20	20		
9	B	1	Total	C	0	0
			20	20		

- Molecule 10 is HEPTANE-1,2,3-TRIOL (CCD ID: HTO) (formula: C₇H₁₆O₃).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			10	7	3		
10	A	1	Total	C	O	0	0
			10	7	3		
10	B	1	Total	C	O	0	0
			10	7	3		
10	B	1	Total	C	O	0	0
			10	7	3		

- Molecule 11 is water.

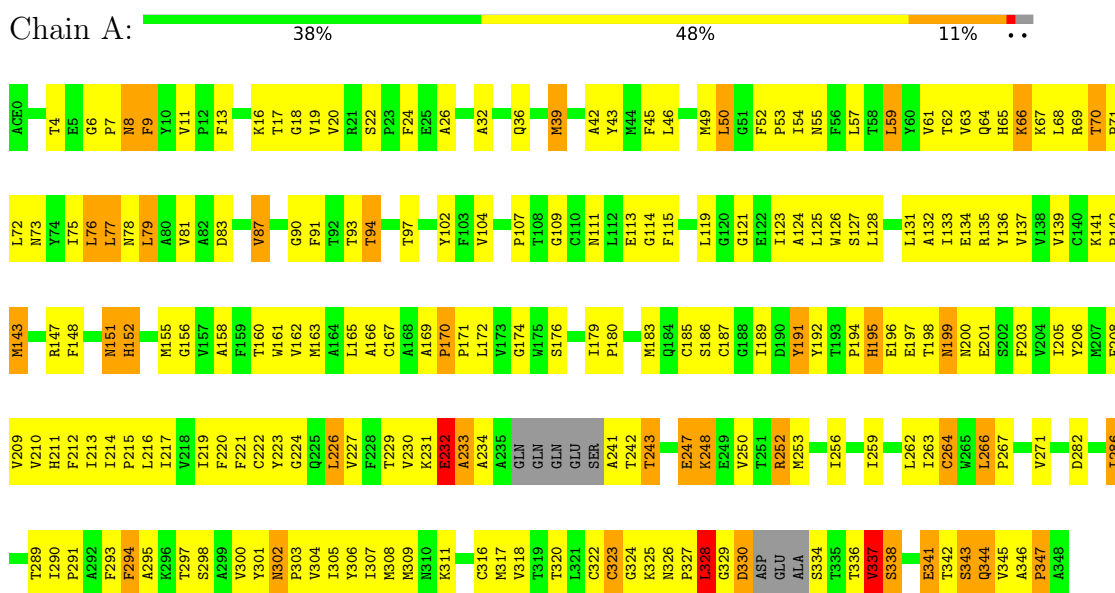
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	7	Total	O	0	0
			7	7		
11	B	5	Total	O	0	0
			5	5		

3 Residue-property plots

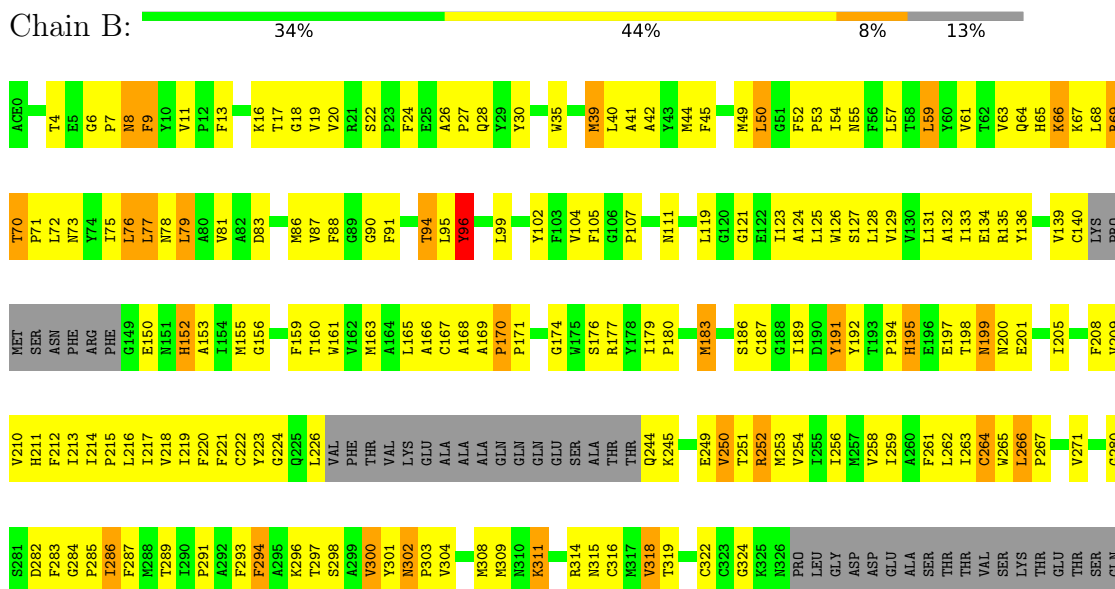
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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



• Molecule 1: RHODOPSIN



VAL
ALA
PRO
ALA

- Molecule 2: alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:  33% 67%

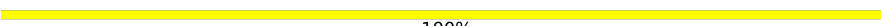
NAG1
NAG2
NAG3

- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  50% 50%

NAG1
NAG2

- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  100%

NAG1
NAG2

- Molecule 4: beta-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  50% 25% 25%

NAG1
NAG2
BMA3
BMA4

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	97.25Å 97.25Å 149.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80	Depositor
% Data completeness (in resolution range)	69.8 (30.00-2.80)	Depositor
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
Refinement program		Depositor
R, R_{free}	0.175 , 0.212	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5551	wwPDB-VP
Average B, all atoms (Å ²)	45.0	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: BMA, HTO, HG, NAG, BNG, MAN, ACE, ZN, PLM, RET

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	0.59	0/2765	1.14	22/3767 (0.6%)
1	B	0.58	0/2472	1.11	17/3368 (0.5%)
All	All	0.59	0/5237	1.13	39/7135 (0.5%)

There are no bond length outliers.

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	343	SER	N-CA-C	-9.81	101.48	112.57
1	B	7	PRO	N-CA-C	-8.22	105.14	114.92
1	A	16	LYS	N-CA-C	-7.29	104.12	113.16
1	A	7	PRO	N-CA-C	-7.15	106.41	114.92
1	B	250	VAL	N-CA-C	-7.07	103.88	110.53
1	A	227	VAL	N-CA-C	-7.06	105.63	111.91
1	A	66	LYS	N-CA-C	6.99	121.42	112.34
1	A	174	GLY	N-CA-C	6.80	122.57	114.48
1	A	77	LEU	N-CA-C	-6.63	104.06	111.28
1	B	77	LEU	N-CA-C	-6.62	104.06	111.28
1	B	16	LYS	N-CA-C	-6.31	105.57	113.20
1	B	26	ALA	N-CA-C	6.28	116.20	108.11
1	B	282	ASP	N-CA-C	6.06	119.32	107.57
1	A	170	PRO	N-CA-C	5.99	118.01	110.70
1	B	294	PHE	N-CA-C	-5.91	104.84	111.28
1	B	265	TRP	N-CA-C	5.90	121.36	114.04
1	A	26	ALA	N-CA-C	5.89	115.71	108.11
1	B	174	GLY	N-CA-C	5.86	121.46	114.48
1	A	282	ASP	N-CA-C	5.84	118.89	107.57
1	A	242	THR	N-CA-C	-5.79	106.22	113.28
1	A	191	TYR	N-CA-C	-5.72	106.65	113.97
1	A	192	TYR	N-CA-C	5.70	117.49	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	9	PHE	N-CA-C	5.67	115.50	108.19
1	B	168	ALA	N-CA-C	5.66	119.36	112.23
1	B	192	TYR	N-CA-C	5.64	117.43	111.28
1	A	328	LEU	N-CA-C	5.63	122.80	110.80
1	B	170	PRO	N-CA-C	5.60	117.53	110.70
1	A	87	VAL	N-CA-C	5.58	116.21	110.36
1	B	191	TYR	N-CA-C	-5.53	107.08	113.88
1	A	294	PHE	N-CA-C	-5.52	105.16	111.07
1	B	297	THR	N-CA-C	-5.48	106.44	113.01
1	A	243	THR	N-CA-C	-5.37	106.69	113.18
1	A	336	THR	N-CA-C	5.34	117.38	109.69
1	A	232	GLU	N-CA-C	5.30	122.09	110.80
1	B	9	PHE	N-CA-C	5.26	114.97	108.19
1	B	96	TYR	N-CA-C	5.24	117.07	111.36
1	A	231	LYS	N-CA-C	5.21	119.91	111.37
1	B	183	MET	N-CA-C	-5.10	107.22	112.93
1	A	297	THR	N-CA-C	-5.06	107.06	113.18

There are no chirality outliers.

There are no planarity outliers.

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	2685	0	2655	240	1
1	B	2398	0	2368	209	0
2	C	39	0	34	5	0
3	D	28	0	25	2	0
3	F	28	0	25	1	0
4	E	50	0	43	2	0
5	A	105	0	150	9	0
5	B	42	0	60	10	0
6	A	3	0	0	0	0
6	B	3	0	0	0	0
7	A	4	0	0	0	0
7	B	3	0	0	0	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	34	0	62	1	0
8	B	17	0	31	1	0
9	A	20	0	27	1	0
9	B	20	0	27	1	0
10	A	40	0	64	4	0
10	B	20	0	32	2	0
11	A	7	0	0	0	0
11	B	5	0	0	1	0
All	All	5551	0	5603	434	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 39.

All (434) 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:A:328:LEU:HB2	1:B:96:TYR:CE2	1.67	1.29
1:A:50:LEU:HD12	1:B:50:LEU:HD12	1.11	1.09
1:A:50:LEU:HD12	1:B:50:LEU:CD1	1.83	1.09
1:A:65:HIS:HB3	1:A:337:VAL:HG22	1.38	1.06
1:A:328:LEU:HB2	1:B:96:TYR:HE2	1.04	1.02
1:B:136:TYR:HA	1:B:226:LEU:HD11	1.42	1.01
1:B:72:LEU:HD22	1:B:250:VAL:HG13	1.39	1.01
1:A:54:ILE:HG12	1:B:49:MET:HE2	1.47	0.96
1:A:142:PRO:HB2	1:A:143:MET:HE2	1.46	0.95
1:A:230:VAL:HG23	1:A:248:LYS:HG3	1.48	0.94
1:A:75:ILE:HG21	1:A:131:LEU:HD21	1.49	0.93
1:A:67:LYS:H	1:A:337:VAL:HG23	1.34	0.92
1:A:50:LEU:CD1	1:B:50:LEU:HD12	1.99	0.91
1:A:241:ALA:HB3	1:A:243:THR:HG22	1.54	0.90
1:A:65:HIS:ND1	1:A:338:SER:HA	1.85	0.89
1:B:75:ILE:HG21	1:B:131:LEU:HD21	1.55	0.88
1:A:9:PHE:HA	1:A:179:ILE:HD11	1.55	0.86
1:B:136:TYR:CA	1:B:226:LEU:HD11	2.07	0.84
1:A:67:LYS:HB2	1:A:337:VAL:HB	1.58	0.84
1:B:212:PHE:O	1:B:216:LEU:HD23	1.77	0.83
1:A:328:LEU:N	1:A:328:LEU:HD22	1.93	0.83
1:A:91:PHE:HA	1:A:94:THR:CG2	2.08	0.82
1:A:212:PHE:O	1:A:216:LEU:HD23	1.79	0.82
1:A:325:LYS:HG2	1:A:326:ASN:H	1.43	0.81
1:A:346:ALA:N	1:A:347:PRO:HD3	1.95	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:MET:HE3	1:B:53:PRO:HB2	1.60	0.81
1:B:311:LYS:HG2	1:B:314:ARG:HH21	1.45	0.81
1:A:328:LEU:HB3	1:B:105:PHE:HZ	1.48	0.77
1:A:68:LEU:C	1:A:69:ARG:HD2	2.09	0.77
1:B:91:PHE:HA	1:B:94:THR:CG2	2.14	0.77
1:B:50:LEU:HD23	1:B:54:ILE:HD12	1.66	0.77
1:B:322:CYS:C	1:B:324:GLY:H	1.93	0.77
1:B:68:LEU:O	1:B:69:ARG:HD2	1.86	0.76
1:B:214:ILE:HB	1:B:215:PRO:HD3	1.67	0.76
1:A:136:TYR:HA	1:A:226:LEU:HD11	1.67	0.76
1:B:314:ARG:O	1:B:318:VAL:HG12	1.85	0.76
1:B:161:TRP:O	1:B:165:LEU:HD23	1.87	0.75
1:B:139:VAL:HB	1:B:226:LEU:CD2	2.17	0.75
1:B:9:PHE:HA	1:B:179:ILE:HD11	1.68	0.74
1:A:230:VAL:HG23	1:A:248:LYS:CG	2.17	0.74
1:A:65:HIS:ND1	1:A:337:VAL:O	2.20	0.74
1:A:214:ILE:HB	1:A:215:PRO:HD3	1.69	0.73
1:A:90:GLY:O	1:A:94:THR:HG22	1.89	0.73
1:A:155:MET:HA	1:A:155:MET:HE2	1.68	0.73
1:B:66:LYS:HZ2	1:B:67:LYS:HG2	1.51	0.73
1:B:209:VAL:HA	1:B:213:ILE:HD12	1.71	0.73
1:A:316:CYS:SG	1:A:337:VAL:HG13	2.29	0.73
1:B:304:VAL:O	1:B:308:MET:HG2	1.90	0.72
2:C:1:NAG:H62	2:C:2:NAG:H83	1.70	0.72
1:A:304:VAL:O	1:A:308:MET:HG2	1.90	0.72
1:B:39:MET:HA	1:B:39:MET:HE3	1.70	0.72
1:B:126:TRP:CH2	1:B:215:PRO:HG3	2.25	0.72
1:B:155:MET:HE2	1:B:155:MET:HA	1.70	0.72
1:B:90:GLY:O	1:B:94:THR:HG22	1.89	0.72
1:B:139:VAL:HB	1:B:226:LEU:HD23	1.72	0.72
1:B:244:GLN:HB2	1:B:245:LYS:HD2	1.72	0.71
1:B:311:LYS:HG2	1:B:314:ARG:NH2	2.06	0.70
1:A:161:TRP:O	1:A:165:LEU:HD23	1.90	0.70
1:B:314:ARG:NH1	10:B:1401:HTO:H51	2.06	0.70
1:B:68:LEU:C	1:B:69:ARG:HD2	2.16	0.70
1:A:65:HIS:HD1	1:A:338:SER:HA	1.53	0.70
1:A:9:PHE:HA	1:A:179:ILE:CD1	2.22	0.69
1:A:68:LEU:O	1:A:69:ARG:HD2	1.93	0.69
1:B:68:LEU:HB3	1:B:73:ASN:HD22	1.57	0.69
1:B:135:ARG:HD2	1:B:251:THR:OG1	1.93	0.69
1:A:230:VAL:CG2	1:A:248:LYS:HG3	2.20	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:SER:O	1:B:131:LEU:HD23	1.92	0.69
1:A:54:ILE:CG1	1:B:49:MET:HE2	2.22	0.68
1:A:346:ALA:H	1:A:347:PRO:HD3	1.58	0.68
1:B:128:LEU:HD22	1:B:258:VAL:HG22	1.76	0.68
1:A:50:LEU:CD2	1:A:54:ILE:HD11	2.23	0.68
1:A:209:VAL:HA	1:A:213:ILE:HD12	1.73	0.68
1:A:127:SER:O	1:A:131:LEU:HD23	1.93	0.68
1:B:91:PHE:CZ	1:B:296:LYS:HB3	2.30	0.67
1:A:126:TRP:CE2	1:A:163:MET:HB3	2.30	0.67
1:A:199:ASN:N	1:A:199:ASN:HD22	1.92	0.67
1:B:285:PRO:HD2	5:B:1506:BNG:H8'1	1.76	0.67
1:A:267:PRO:HA	5:A:1503:BNG:H7'1	1.76	0.67
1:B:66:LYS:HD2	1:B:67:LYS:H	1.60	0.66
1:A:134:GLU:HA	1:A:148:PHE:CE2	2.30	0.66
1:A:151:ASN:OD1	1:A:152:HIS:CE1	2.49	0.66
1:A:50:LEU:CD2	1:A:54:ILE:CD1	2.73	0.66
1:A:39:MET:HE2	1:A:39:MET:HA	1.78	0.66
1:A:65:HIS:HD1	1:A:337:VAL:C	2.05	0.65
1:A:213:ILE:O	1:A:217:ILE:HD13	1.95	0.65
1:B:75:ILE:HG13	1:B:131:LEU:HD21	1.77	0.65
1:A:151:ASN:OD1	1:A:152:HIS:ND1	2.30	0.65
1:A:266:LEU:N	1:A:267:PRO:HD2	2.10	0.65
1:B:266:LEU:N	1:B:267:PRO:HD2	2.12	0.65
1:B:55:ASN:ND2	1:B:83:ASP:HB3	2.12	0.65
1:A:65:HIS:HB3	1:A:337:VAL:CG2	2.22	0.64
1:A:142:PRO:HB2	1:A:143:MET:CE	2.25	0.64
1:A:75:ILE:CG2	1:A:131:LEU:HD21	2.25	0.64
1:A:50:LEU:HD21	1:A:54:ILE:HD11	1.78	0.64
1:B:208:PHE:O	1:B:213:ILE:HG13	1.98	0.63
1:A:208:PHE:O	1:A:213:ILE:HG13	1.97	0.63
1:A:75:ILE:O	1:A:79:LEU:HD22	1.99	0.63
1:A:308:MET:HE1	1:B:41:ALA:HB1	1.81	0.63
1:A:300:VAL:O	1:A:303:PRO:HD2	1.99	0.63
1:B:213:ILE:O	1:B:217:ILE:HD13	1.97	0.63
1:A:68:LEU:HB3	1:A:73:ASN:HD22	1.63	0.63
1:B:139:VAL:CB	1:B:226:LEU:HD23	2.28	0.63
1:B:284:GLY:H	5:B:1506:BNG:H5'2	1.63	0.63
1:B:75:ILE:O	1:B:78:ASN:HB3	1.98	0.62
1:A:68:LEU:H	1:A:337:VAL:HG21	1.63	0.62
1:A:308:MET:HE2	1:B:41:ALA:C	2.24	0.62
1:B:300:VAL:O	1:B:303:PRO:HD2	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:THR:HA	1:A:344:GLN:CD	2.24	0.61
1:A:201:GLU:O	1:A:205:ILE:HG13	2.01	0.61
1:A:328:LEU:N	1:A:328:LEU:CD2	2.63	0.61
1:A:197:GLU:O	1:A:197:GLU:HG2	2.01	0.61
1:B:284:GLY:H	5:B:1506:BNG:H3'1	1.64	0.61
1:A:75:ILE:O	1:A:78:ASN:HB3	2.00	0.61
1:A:102:TYR:CE2	1:A:104:VAL:HA	2.35	0.60
1:B:75:ILE:CG2	1:B:131:LEU:HD21	2.30	0.60
1:A:49:MET:CE	1:B:53:PRO:HB2	2.32	0.60
1:A:68:LEU:N	1:A:337:VAL:HG21	2.15	0.60
1:B:102:TYR:CE2	1:B:104:VAL:HA	2.36	0.60
1:B:199:ASN:N	1:B:199:ASN:HD22	1.97	0.60
1:B:201:GLU:O	1:B:205:ILE:HG13	2.01	0.60
1:B:152:HIS:N	1:B:152:HIS:ND1	2.47	0.60
1:B:132:ALA:O	1:B:222:CYS:SG	2.60	0.60
1:A:267:PRO:O	1:A:271:VAL:HG23	2.01	0.60
1:A:50:LEU:HD23	1:A:54:ILE:CD1	2.32	0.60
1:A:65:HIS:HD1	1:A:338:SER:CA	2.15	0.60
1:A:252:ARG:HH21	5:A:1501:BNG:C6	2.14	0.59
1:B:210:VAL:HA	1:B:214:ILE:HD12	1.83	0.59
1:B:298:SER:HA	1:B:301:TYR:CE2	2.37	0.59
1:A:328:LEU:HB3	1:B:105:PHE:CZ	2.34	0.59
1:B:9:PHE:HA	1:B:179:ILE:CD1	2.33	0.59
1:A:75:ILE:HG22	1:A:79:LEU:CD2	2.32	0.59
1:A:328:LEU:CB	1:B:96:TYR:CE2	2.62	0.59
1:A:308:MET:HE2	1:B:42:ALA:N	2.18	0.59
1:B:59:LEU:HD12	1:B:77:LEU:HD11	1.84	0.59
1:B:126:TRP:CZ3	1:B:215:PRO:HG3	2.37	0.59
1:A:134:GLU:HG2	1:A:148:PHE:CD2	2.37	0.59
1:A:308:MET:CE	1:B:41:ALA:HB1	2.33	0.59
1:B:170:PRO:HB2	1:B:171:PRO:HD3	1.85	0.59
1:B:222:CYS:O	1:B:226:LEU:HD12	2.02	0.58
1:A:135:ARG:HA	1:A:135:ARG:NE	2.18	0.58
1:A:9:PHE:CA	1:A:179:ILE:HD11	2.29	0.58
1:A:230:VAL:HA	1:A:248:LYS:HE3	1.84	0.58
1:B:267:PRO:O	1:B:271:VAL:HG23	2.02	0.58
1:A:75:ILE:HG13	1:A:131:LEU:HD21	1.85	0.58
1:A:342:THR:O	1:A:342:THR:HG23	2.04	0.58
1:A:294:PHE:HB2	10:A:1403:HTO:H41	1.86	0.58
1:B:72:LEU:HD11	1:B:253:MET:HE2	1.86	0.58
1:B:125:LEU:O	1:B:128:LEU:HB2	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:LYS:N	1:A:337:VAL:HG23	2.13	0.57
1:B:271:VAL:HG21	1:B:291:PRO:HG3	1.85	0.57
1:A:67:LYS:HB2	1:A:337:VAL:CB	2.33	0.57
1:A:125:LEU:O	1:A:128:LEU:HB2	2.05	0.57
1:A:262:LEU:HB3	1:A:266:LEU:HD22	1.87	0.57
1:B:284:GLY:N	5:B:1506:BNG:H5'2	2.20	0.57
3:D:1:NAG:O6	3:D:2:NAG:H82	2.04	0.57
1:B:322:CYS:C	1:B:324:GLY:N	2.56	0.56
1:A:65:HIS:HB2	1:A:68:LEU:HD12	1.87	0.56
1:B:136:TYR:HA	1:B:226:LEU:CD1	2.28	0.56
1:A:325:LYS:HG2	1:A:326:ASN:N	2.17	0.56
1:B:256:ILE:HA	1:B:259:ILE:HD12	1.87	0.56
1:A:290:ILE:HG23	10:A:1403:HTO:H42	1.86	0.56
1:A:298:SER:HA	1:A:301:TYR:CE2	2.41	0.56
1:B:50:LEU:CD2	1:B:54:ILE:HD12	2.34	0.56
1:A:76:LEU:HD22	1:A:306:TYR:CD2	2.40	0.56
1:A:135:ARG:NH1	1:A:247:GLU:HG3	2.20	0.56
1:A:67:LYS:HD3	1:A:343:SER:O	2.06	0.56
1:B:75:ILE:HG13	1:B:131:LEU:CD2	2.36	0.56
1:B:50:LEU:CD2	1:B:54:ILE:CD1	2.84	0.55
1:B:75:ILE:HG22	1:B:79:LEU:CD2	2.35	0.55
1:B:189:ILE:HB	1:B:191:TYR:CE1	2.41	0.55
1:A:139:VAL:HB	1:A:226:LEU:HD21	1.87	0.55
1:B:245:LYS:HD2	1:B:245:LYS:N	2.20	0.55
1:A:91:PHE:HA	1:A:94:THR:HG23	1.89	0.55
1:A:195:HIS:HB3	1:A:200:ASN:HD21	1.71	0.55
1:A:136:TYR:CA	1:A:226:LEU:HD11	2.34	0.55
1:A:119:LEU:O	1:A:123:ILE:HG13	2.07	0.55
5:A:1500:BNG:H3'1	1:B:35:TRP:HB2	1.89	0.55
1:B:50:LEU:HD21	1:B:54:ILE:HD11	1.89	0.55
1:A:4:THR:HG21	1:A:20:VAL:HG11	1.88	0.55
1:B:139:VAL:HB	1:B:226:LEU:HD21	1.87	0.55
1:B:68:LEU:CB	1:B:73:ASN:HD22	2.20	0.55
1:A:187:CYS:O	9:A:1296:RET:H12	2.06	0.54
1:B:128:LEU:HB3	1:B:219:ILE:HD13	1.89	0.54
1:A:210:VAL:HA	1:A:214:ILE:HD12	1.88	0.54
1:A:195:HIS:HB3	1:A:200:ASN:ND2	2.23	0.54
1:A:346:ALA:N	1:A:347:PRO:CD	2.70	0.54
1:B:72:LEU:HD22	1:B:250:VAL:CG1	2.26	0.54
1:B:197:GLU:CG	1:B:197:GLU:O	2.55	0.54
1:B:283:PHE:HA	5:B:1506:BNG:H3'1	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:LYS:HD2	1:B:67:LYS:N	2.23	0.53
1:A:345:VAL:HB	1:A:347:PRO:HD3	1.89	0.53
1:B:75:ILE:O	1:B:79:LEU:HD22	2.09	0.53
1:A:308:MET:CE	1:B:41:ALA:CB	2.87	0.53
1:B:18:GLY:HA2	4:E:1:NAG:C1	2.38	0.53
1:B:179:ILE:HG13	1:B:180:PRO:HD2	1.90	0.53
1:B:284:GLY:H	5:B:1506:BNG:C3'	2.22	0.53
1:A:139:VAL:HB	1:A:226:LEU:CD2	2.38	0.53
1:B:4:THR:HG21	1:B:20:VAL:HG11	1.90	0.52
1:B:135:ARG:NE	1:B:135:ARG:HA	2.22	0.52
1:A:137:VAL:O	1:A:141:LYS:HA	2.08	0.52
1:A:256:ILE:HA	1:A:259:ILE:HD12	1.91	0.52
1:A:342:THR:OG1	1:A:344:GLN:HG2	2.09	0.52
1:B:17:THR:OG1	1:B:19:VAL:HG12	2.09	0.52
1:B:19:VAL:HG22	1:B:19:VAL:O	2.09	0.52
1:B:252:ARG:O	1:B:256:ILE:HG12	2.10	0.52
1:B:199:ASN:N	1:B:199:ASN:ND2	2.58	0.52
1:B:280:GLY:O	3:F:1:NAG:H82	2.09	0.52
1:A:57:LEU:HD23	1:A:61:VAL:HG23	1.91	0.52
1:A:318:VAL:HG11	1:A:328:LEU:CD2	2.40	0.52
1:B:57:LEU:C	1:B:57:LEU:HD23	2.34	0.52
1:A:59:LEU:O	1:A:63:VAL:HG12	2.11	0.51
1:B:152:HIS:O	1:B:155:MET:HB2	2.09	0.51
1:B:221:PHE:O	1:B:224:GLY:N	2.43	0.51
1:A:126:TRP:CH2	1:A:215:PRO:HG3	2.44	0.51
1:A:222:CYS:O	1:A:226:LEU:HD12	2.10	0.51
1:A:328:LEU:HD22	1:A:328:LEU:H	1.73	0.51
1:B:220:PHE:O	1:B:223:TYR:HB3	2.11	0.51
1:A:57:LEU:HD23	1:A:57:LEU:C	2.35	0.51
1:A:322:CYS:O	1:A:323:CYS:SG	2.69	0.51
1:A:62:THR:HG21	1:A:77:LEU:HB2	1.92	0.51
1:A:75:ILE:HG22	1:A:79:LEU:HD21	1.92	0.51
1:A:65:HIS:CE1	1:A:338:SER:OG	2.63	0.50
1:A:322:CYS:C	1:A:323:CYS:SG	2.94	0.50
1:A:327:PRO:O	1:A:329:GLY:N	2.43	0.50
1:A:183:MET:HG3	1:A:289:THR:OG1	2.11	0.50
1:A:43:TYR:OH	1:A:293:PHE:O	2.20	0.50
1:A:57:LEU:HD23	1:A:57:LEU:O	2.11	0.50
1:A:327:PRO:C	1:A:328:LEU:HD22	2.37	0.50
1:B:86:MET:HE1	1:B:298:SER:OG	2.11	0.50
1:B:133:ILE:O	1:B:134:GLU:C	2.55	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:HIS:O	1:A:155:MET:HB2	2.10	0.50
1:A:183:MET:HE3	1:A:286:ILE:HG13	1.93	0.50
1:B:119:LEU:O	1:B:123:ILE:HG13	2.11	0.50
1:B:315:ASN:O	1:B:318:VAL:HG13	2.11	0.50
1:B:156:GLY:O	1:B:159:PHE:HB3	2.12	0.50
1:A:197:GLU:O	1:A:197:GLU:CG	2.60	0.49
1:B:6:GLY:C	1:B:8:ASN:H	2.19	0.49
1:A:133:ILE:O	1:A:134:GLU:C	2.55	0.49
1:A:220:PHE:O	1:A:223:TYR:HB3	2.13	0.49
1:A:305:ILE:HA	5:A:1500:BNG:H8'2	1.94	0.49
1:B:262:LEU:HB3	1:B:266:LEU:HD22	1.95	0.49
1:A:54:ILE:HG12	1:B:49:MET:CE	2.31	0.49
1:A:126:TRP:NE1	1:A:163:MET:HB3	2.28	0.49
1:A:64:GLN:NE2	1:A:320:THR:OG1	2.45	0.49
1:A:189:ILE:HB	1:A:191:TYR:CE1	2.48	0.49
1:B:59:LEU:HD12	1:B:77:LEU:CD1	2.42	0.49
1:A:68:LEU:HG	1:A:337:VAL:HG21	1.93	0.49
1:B:42:ALA:O	1:B:45:PHE:HB3	2.13	0.49
1:B:107:PRO:O	1:B:111:ASN:ND2	2.46	0.49
1:B:87:VAL:HA	1:B:91:PHE:HB2	1.94	0.49
1:B:75:ILE:CG1	1:B:131:LEU:HD21	2.42	0.48
1:B:87:VAL:O	1:B:91:PHE:HB2	2.11	0.48
1:B:136:TYR:HB2	1:B:222:CYS:SG	2.52	0.48
1:B:195:HIS:HB3	1:B:200:ASN:ND2	2.28	0.48
1:A:308:MET:CE	1:B:99:LEU:HD21	2.44	0.48
1:B:180:PRO:HA	1:B:186:SER:O	2.13	0.48
1:B:244:GLN:HB2	1:B:245:LYS:CD	2.42	0.48
1:A:46:LEU:HD13	10:A:1400:HTO:H71	1.96	0.48
1:A:139:VAL:O	1:A:229:THR:HG21	2.13	0.48
1:B:59:LEU:O	1:B:63:VAL:HG12	2.13	0.48
1:A:252:ARG:O	1:A:256:ILE:HG12	2.13	0.48
1:A:17:THR:OG1	1:A:19:VAL:HG12	2.14	0.48
1:A:76:LEU:HD22	1:A:306:TYR:CG	2.48	0.48
1:A:32:ALA:HB1	1:A:36:GLN:OE1	2.14	0.48
1:A:55:ASN:ND2	1:A:83:ASP:HB3	2.29	0.48
1:A:75:ILE:HG13	1:A:131:LEU:CD2	2.43	0.48
1:A:230:VAL:CA	1:A:248:LYS:HE3	2.44	0.48
1:B:287:PHE:HB2	5:B:1506:BNG:H4'1	1.96	0.48
1:B:195:HIS:HB3	1:B:200:ASN:HD21	1.78	0.47
1:A:9:PHE:C	1:A:179:ILE:HD11	2.40	0.47
1:B:50:LEU:HD23	1:B:54:ILE:CD1	2.40	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:VAL:HG12	1:A:229:THR:HB	1.96	0.47
1:A:311:LYS:NZ	1:A:330:ASP:H	2.12	0.47
1:B:77:LEU:O	1:B:81:VAL:HG23	2.14	0.47
1:A:107:PRO:O	1:A:111:ASN:ND2	2.47	0.47
1:A:271:VAL:HG21	1:A:291:PRO:HG3	1.96	0.47
1:A:328:LEU:HB2	1:B:96:TYR:CD2	2.39	0.47
1:B:68:LEU:CA	1:B:73:ASN:HD22	2.27	0.47
1:A:134:GLU:HA	1:A:148:PHE:HE2	1.75	0.47
1:A:167:CYS:HB2	1:A:211:HIS:CG	2.50	0.47
1:A:199:ASN:N	1:A:199:ASN:ND2	2.55	0.47
1:B:70:THR:O	1:B:71:PRO:C	2.56	0.47
1:B:91:PHE:O	1:B:94:THR:HG23	2.15	0.47
1:B:183:MET:HG3	1:B:289:THR:OG1	2.14	0.47
1:A:66:LYS:O	1:A:69:ARG:HD3	2.15	0.47
3:D:2:NAG:H3	3:D:2:NAG:H83	1.97	0.47
1:A:170:PRO:HB2	1:A:171:PRO:HD3	1.96	0.47
1:B:72:LEU:HD11	1:B:253:MET:CE	2.45	0.47
1:B:187:CYS:O	9:B:1296:RET:H12	2.15	0.47
1:B:244:GLN:C	1:B:245:LYS:HD2	2.40	0.47
1:B:287:PHE:HB2	5:B:1506:BNG:H3'2	1.97	0.47
1:A:52:PHE:HB3	1:A:53:PRO:CD	2.46	0.46
1:A:45:PHE:O	1:A:49:MET:HB2	2.15	0.46
1:A:230:VAL:HB	1:A:248:LYS:HE3	1.97	0.46
1:A:252:ARG:HH21	5:A:1501:BNG:H62	1.78	0.46
1:B:65:HIS:HB2	1:B:68:LEU:HD12	1.97	0.46
1:B:125:LEU:HB2	1:B:261:PHE:CZ	2.50	0.46
1:A:148:PHE:HD1	1:A:152:HIS:CD2	2.34	0.46
1:A:194:PRO:O	1:A:195:HIS:C	2.58	0.46
1:A:22:SER:C	1:A:24:PHE:H	2.23	0.46
1:A:42:ALA:O	1:A:45:PHE:HB3	2.15	0.46
1:A:156:GLY:O	1:A:160:THR:HG23	2.15	0.46
1:A:327:PRO:HA	1:A:328:LEU:HD22	1.96	0.46
1:B:167:CYS:HB2	1:B:211:HIS:CG	2.50	0.46
1:B:301:TYR:HE1	11:B:2008:HOH:O	1.98	0.46
1:A:75:ILE:HG21	1:A:131:LEU:CD2	2.36	0.46
1:A:114:GLY:HA2	1:A:187:CYS:O	2.16	0.46
1:A:18:GLY:HA2	2:C:1:NAG:C1	2.46	0.46
1:A:180:PRO:HA	1:A:186:SER:O	2.15	0.46
5:A:1500:BNG:C3'	1:B:35:TRP:HB2	2.46	0.46
1:B:298:SER:HA	1:B:301:TYR:CD2	2.49	0.46
1:A:72:LEU:HD22	1:A:250:VAL:HG13	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:ILE:CG1	1:A:131:LEU:HD21	2.46	0.45
1:A:152:HIS:HD1	1:A:152:HIS:N	2.14	0.45
1:B:9:PHE:CA	1:B:179:ILE:HD11	2.40	0.45
1:A:167:CYS:HB2	1:A:211:HIS:CE1	2.51	0.45
1:A:263:ILE:O	1:A:294:PHE:HE2	1.99	0.45
1:B:44:MET:HE2	1:B:94:THR:CG2	2.46	0.45
1:B:194:PRO:O	1:B:195:HIS:C	2.60	0.45
1:A:139:VAL:HG11	1:A:230:VAL:HG12	1.97	0.45
1:B:128:LEU:HD22	1:B:258:VAL:CG2	2.44	0.45
1:A:68:LEU:CB	1:A:73:ASN:HD22	2.28	0.45
1:A:230:VAL:HG22	1:A:230:VAL:O	2.17	0.45
1:A:19:VAL:HG22	1:A:19:VAL:O	2.16	0.45
1:A:70:THR:O	1:A:71:PRO:C	2.59	0.45
1:A:179:ILE:HG13	1:A:180:PRO:HD2	1.99	0.45
5:A:1500:BNG:H1'1	1:B:35:TRP:HA	1.97	0.45
1:B:41:ALA:O	1:B:95:LEU:HD13	2.17	0.45
1:B:163:MET:O	1:B:166:ALA:HB3	2.16	0.45
1:A:87:VAL:O	1:A:91:PHE:HB2	2.17	0.45
1:A:230:VAL:O	1:A:230:VAL:HG13	2.16	0.45
1:B:75:ILE:HG22	1:B:79:LEU:HD21	1.99	0.45
1:B:322:CYS:HA	8:B:1322:PLM:O2	2.15	0.45
1:B:57:LEU:HD23	1:B:57:LEU:O	2.17	0.44
1:B:183:MET:HE3	1:B:286:ILE:HG13	1.98	0.44
1:B:266:LEU:N	1:B:267:PRO:CD	2.79	0.44
1:B:167:CYS:HB2	1:B:211:HIS:CE1	2.53	0.44
2:C:1:NAG:H61	2:C:2:NAG:N2	2.33	0.44
1:A:128:LEU:HB3	1:A:219:ILE:HD13	1.98	0.44
1:A:298:SER:HA	1:A:301:TYR:CD2	2.53	0.44
1:B:213:ILE:C	1:B:217:ILE:HD13	2.42	0.44
1:A:126:TRP:CZ2	1:A:163:MET:HE3	2.51	0.44
1:B:9:PHE:C	1:B:179:ILE:HD11	2.42	0.44
1:B:64:GLN:O	1:B:64:GLN:HG2	2.17	0.44
1:B:91:PHE:HA	1:B:94:THR:HG23	1.96	0.44
1:B:214:ILE:HB	1:B:215:PRO:CD	2.44	0.44
1:A:87:VAL:HA	1:A:91:PHE:HB2	2.00	0.44
1:A:266:LEU:N	1:A:267:PRO:CD	2.79	0.44
4:E:1:NAG:H62	4:E:2:NAG:H82	2.00	0.44
1:A:6:GLY:C	1:A:8:ASN:H	2.25	0.44
1:A:291:PRO:O	1:A:295:ALA:HB2	2.17	0.44
1:B:195:HIS:HB3	1:B:198:THR:HG22	2.00	0.44
1:B:253:MET:HA	1:B:309:MET:CE	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:52:PHE:HB3	1:B:53:PRO:CD	2.48	0.44
1:A:67:LYS:O	1:A:69:ARG:N	2.51	0.44
1:A:126:TRP:CZ3	1:A:215:PRO:HG3	2.53	0.44
1:B:50:LEU:HD21	1:B:54:ILE:CD1	2.47	0.44
1:B:209:VAL:O	1:B:214:ILE:HG13	2.18	0.43
1:A:57:LEU:C	1:A:57:LEU:CD2	2.91	0.43
1:A:50:LEU:CD2	1:A:54:ILE:HD12	2.47	0.43
1:A:68:LEU:CA	1:A:73:ASN:HD22	2.31	0.43
1:A:163:MET:O	1:A:166:ALA:HB3	2.18	0.43
1:B:244:GLN:HB2	1:B:245:LYS:CE	2.49	0.43
1:A:11:VAL:C	1:A:13:PHE:H	2.27	0.43
1:A:203:PHE:O	1:A:206:TYR:HB3	2.19	0.43
1:A:334:SER:N	10:A:1404:HTO:H2	2.34	0.43
1:A:347:PRO:HG3	5:A:1504:BNG:O3	2.19	0.43
8:A:1322:PLM:H72	1:B:88:PHE:HE1	1.84	0.43
1:A:195:HIS:HB3	1:A:198:THR:HG22	2.01	0.42
1:B:136:TYR:CB	1:B:226:LEU:HD11	2.48	0.42
1:B:284:GLY:HA3	5:B:1506:BNG:H6'1	2.00	0.42
1:A:121:GLY:O	1:A:124:ALA:HB3	2.18	0.42
1:A:169:ALA:N	1:A:170:PRO:CD	2.81	0.42
1:B:39:MET:HA	1:B:39:MET:CE	2.44	0.42
1:A:253:MET:HA	1:A:309:MET:CE	2.49	0.42
1:A:302:ASN:CB	1:A:303:PRO:HD3	2.50	0.42
1:B:20:VAL:HA	1:B:30:TYR:CZ	2.55	0.42
1:B:57:LEU:C	1:B:57:LEU:CD2	2.93	0.42
1:A:167:CYS:HB2	1:A:211:HIS:CD2	2.55	0.42
1:A:176:SER:OG	1:A:189:ILE:HG23	2.19	0.42
1:A:97:THR:HG21	1:A:185:CYS:HA	2.02	0.42
1:B:57:LEU:HD23	1:B:61:VAL:HG23	2.02	0.42
1:B:79:LEU:HG	1:B:302:ASN:ND2	2.33	0.42
1:B:121:GLY:O	1:B:124:ALA:HB3	2.20	0.42
1:A:64:GLN:O	1:A:64:GLN:HG2	2.19	0.42
1:A:77:LEU:O	1:A:81:VAL:HG23	2.20	0.42
1:B:156:GLY:O	1:B:160:THR:HG23	2.20	0.42
1:B:285:PRO:CD	5:B:1506:BNG:H8'1	2.47	0.42
1:B:126:TRP:CE2	1:B:163:MET:HB3	2.55	0.41
2:C:1:NAG:H62	2:C:2:NAG:C8	2.44	0.41
1:A:4:THR:HG23	2:C:1:NAG:H83	2.01	0.41
1:A:66:LYS:HA	1:A:69:ARG:NH1	2.35	0.41
1:A:152:HIS:ND1	1:A:152:HIS:N	2.67	0.41
1:B:76:LEU:HA	1:B:79:LEU:HD23	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:TYR:O	1:B:140:CYS:HB2	2.20	0.41
1:A:97:THR:CG2	1:A:185:CYS:HA	2.51	0.41
1:A:230:VAL:CB	1:A:248:LYS:HE3	2.51	0.41
1:A:318:VAL:HG11	1:A:327:PRO:HA	2.02	0.41
1:A:342:THR:C	1:A:344:GLN:H	2.27	0.41
1:B:22:SER:C	1:B:24:PHE:H	2.28	0.41
1:B:40:LEU:HD22	1:B:293:PHE:CE2	2.55	0.41
1:B:169:ALA:N	1:B:170:PRO:CD	2.83	0.41
1:A:115:PHE:CD1	1:A:172:LEU:HD11	2.55	0.41
1:B:176:SER:OG	1:B:189:ILE:HG23	2.20	0.41
1:B:253:MET:O	1:B:254:VAL:C	2.64	0.41
1:B:316:CYS:O	1:B:319:THR:HB	2.21	0.41
1:A:221:PHE:O	1:A:224:GLY:N	2.53	0.41
1:B:28:GLN:C	1:B:30:TYR:H	2.29	0.41
1:B:135:ARG:HG3	1:B:226:LEU:HD22	2.03	0.41
1:A:131:LEU:O	1:A:132:ALA:C	2.63	0.41
1:B:264:CYS:HA	1:B:294:PHE:CE2	2.56	0.41
1:B:301:TYR:HA	10:B:1406:HTO:H11	2.03	0.41
1:A:91:PHE:O	1:A:94:THR:HG23	2.21	0.41
1:A:93:THR:HG21	1:A:109:GLY:CA	2.51	0.41
1:A:135:ARG:HH12	1:A:247:GLU:HG3	1.85	0.41
1:A:264:CYS:HA	1:A:294:PHE:CE2	2.55	0.41
1:A:307:ILE:HD13	1:A:317:MET:SD	2.60	0.41
1:A:308:MET:HE3	1:B:99:LEU:HD21	2.01	0.41
1:B:263:ILE:O	1:B:294:PHE:HE2	2.03	0.41
1:B:296:LYS:C	1:B:298:SER:N	2.79	0.41
1:B:315:ASN:HD22	1:B:315:ASN:HA	1.63	0.41
1:B:11:VAL:C	1:B:13:PHE:H	2.28	0.40
1:B:68:LEU:HA	1:B:73:ASN:ND2	2.36	0.40
1:B:197:GLU:O	1:B:197:GLU:HG3	2.21	0.40
1:A:158:ALA:O	1:A:162:VAL:HG23	2.21	0.40
1:A:286:ILE:HB	5:A:1505:BNG:H3'1	2.03	0.40
1:A:76:LEU:HD13	1:A:306:TYR:CE1	2.56	0.40
1:A:213:ILE:C	1:A:217:ILE:HD13	2.47	0.40
1:B:129:VAL:HG13	1:B:218:VAL:HG11	2.03	0.40
1:B:150:GLU:HA	1:B:153:ALA:HB3	2.03	0.40
1:A:65:HIS:HB2	1:A:68:LEU:CD1	2.52	0.40
1:A:109:GLY:O	1:A:113:GLU:HB2	2.21	0.40
1:A:233:ALA:O	1:A:234:ALA:HB2	2.22	0.40
1:B:167:CYS:HB2	1:B:211:HIS:CD2	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:GLU:OE2	7:B:958:ZN:ZN[2_654]	1.56	0.64

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	335/349 (96%)	275 (82%)	50 (15%)	10 (3%)	3	12
1	B	296/349 (85%)	256 (86%)	38 (13%)	2 (1%)	18	47
All	All	631/698 (90%)	531 (84%)	88 (14%)	12 (2%)	6	22

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	233	ALA
1	A	328	LEU
1	A	341	GLU
1	A	195	HIS
1	A	232	GLU
1	A	324	GLY
1	A	337	VAL
1	A	323	CYS
1	A	347	PRO
1	B	195	HIS
1	A	338	SER
1	B	300	VAL

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	289/296 (98%)	262 (91%)	27 (9%)	8	27
1	B	257/296 (87%)	234 (91%)	23 (9%)	9	29
All	All	546/592 (92%)	496 (91%)	50 (9%)	8	27

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	39	MET
1	A	50	LEU
1	A	59	LEU
1	A	70	THR
1	A	76	LEU
1	A	79	LEU
1	A	94	THR
1	A	143	MET
1	A	147	ARG
1	A	151	ASN
1	A	152	HIS
1	A	199	ASN
1	A	226	LEU
1	A	232	GLU
1	A	247	GLU
1	A	248	LYS
1	A	252	ARG
1	A	264	CYS
1	A	266	LEU
1	A	286	ILE
1	A	302	ASN
1	A	328	LEU
1	A	330	ASP
1	A	337	VAL
1	A	341	GLU
1	A	344	GLN
1	B	8	ASN
1	B	27	PRO
1	B	39	MET
1	B	50	LEU
1	B	59	LEU
1	B	66	LYS

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Mol	Chain	Res	Type
1	B	69	ARG
1	B	70	THR
1	B	76	LEU
1	B	79	LEU
1	B	94	THR
1	B	96	TYR
1	B	152	HIS
1	B	177	ARG
1	B	199	ASN
1	B	249	GLU
1	B	252	ARG
1	B	264	CYS
1	B	266	LEU
1	B	286	ILE
1	B	302	ASN
1	B	311	LYS
1	B	318	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	64	GLN
1	A	111	ASN
1	A	145	ASN
1	A	199	ASN
1	A	302	ASN
1	A	312	GLN
1	A	315	ASN
1	B	8	ASN
1	B	64	GLN
1	B	111	ASN
1	B	195	HIS
1	B	199	ASN
1	B	302	ASN
1	B	312	GLN
1	B	315	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

11 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	C	1	2,1	14,14,15	0.66	0	17,19,21	0.72	0
2	NAG	C	2	2	14,14,15	0.66	0	17,19,21	0.97	0
2	MAN	C	3	2	11,11,12	0.69	0	15,15,17	0.69	0
3	NAG	D	1	3,1	14,14,15	0.55	0	17,19,21	0.81	1 (5%)
3	NAG	D	2	3	14,14,15	0.83	0	17,19,21	0.88	0
4	NAG	E	1	4,1	14,14,15	0.46	0	17,19,21	0.84	0
4	NAG	E	2	4	14,14,15	0.51	0	17,19,21	0.87	1 (5%)
4	BMA	E	3	4	11,11,12	0.74	0	15,15,17	0.53	0
4	BMA	E	4	4	11,11,12	0.68	0	15,15,17	0.63	0
3	NAG	F	1	3,1	14,14,15	0.66	0	17,19,21	0.75	0
3	NAG	F	2	3	14,14,15	0.81	1 (7%)	17,19,21	0.66	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
2	NAG	C	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	C	2	2	-	4/6/23/26	0/1/1/1
2	MAN	C	3	2	-	2/2/19/22	0/1/1/1
3	NAG	D	1	3,1	-	4/6/23/26	0/1/1/1
3	NAG	D	2	3	-	5/6/23/26	0/1/1/1
4	NAG	E	1	4,1	-	4/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	E	2	4	-	4/6/23/26	0/1/1/1
4	BMA	E	3	4	-	2/2/19/22	0/1/1/1
4	BMA	E	4	4	-	2/2/19/22	0/1/1/1
3	NAG	F	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	4/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	2	NAG	C1-C2	2.08	1.55	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	2	NAG	C2-N2-C7	-2.58	119.45	122.90
3	D	1	NAG	C2-N2-C7	-2.11	120.07	122.90

There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	C	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O7-C7-N2-C2
3	D	1	NAG	C8-C7-N2-C2
3	D	1	NAG	O7-C7-N2-C2
3	D	2	NAG	C3-C2-N2-C7
3	D	2	NAG	C8-C7-N2-C2
3	D	2	NAG	O7-C7-N2-C2
3	F	1	NAG	C8-C7-N2-C2
3	F	1	NAG	O7-C7-N2-C2
3	F	2	NAG	C8-C7-N2-C2
3	F	2	NAG	O7-C7-N2-C2
4	E	1	NAG	C8-C7-N2-C2
4	E	1	NAG	O7-C7-N2-C2
2	C	2	NAG	O5-C5-C6-O6
4	E	1	NAG	O5-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
4	E	2	NAG	O5-C5-C6-O6

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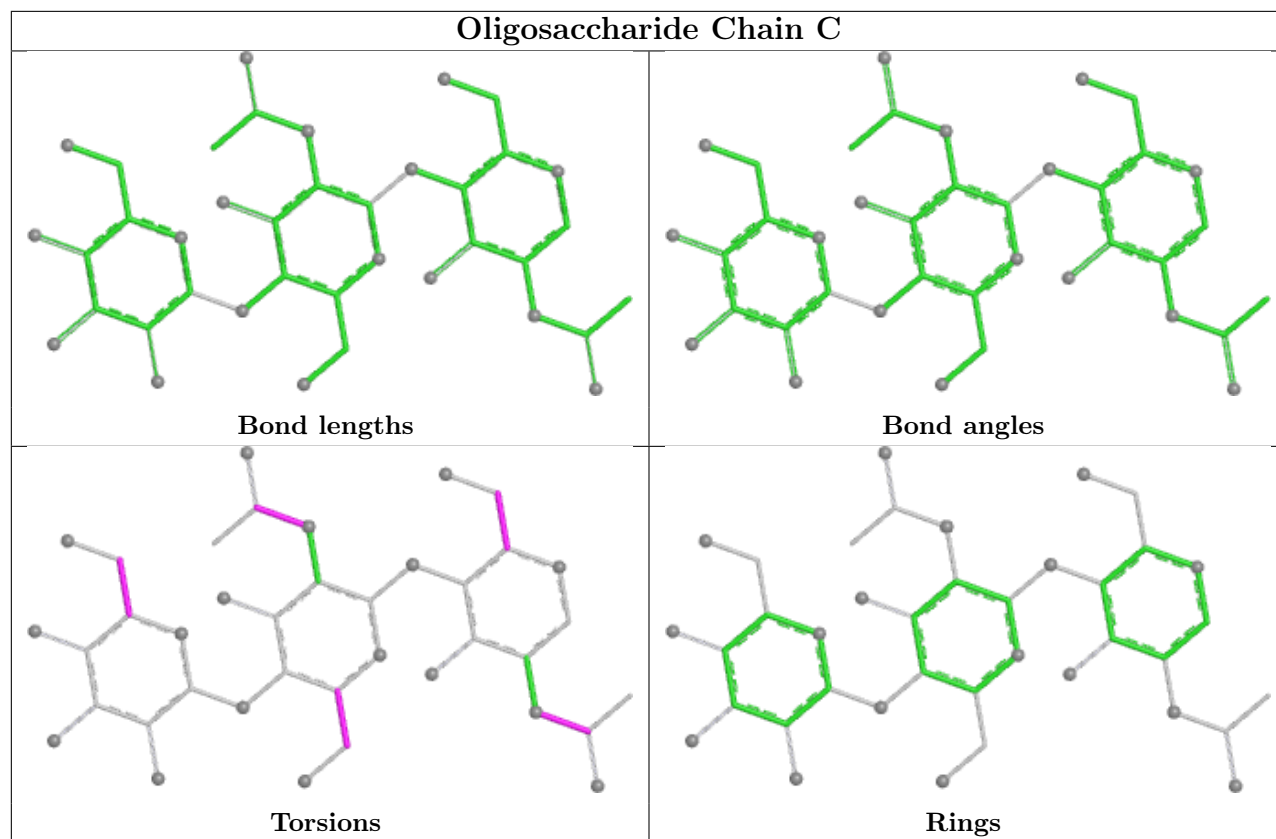
Mol	Chain	Res	Type	Atoms
4	E	1	NAG	C4-C5-C6-O6
3	D	2	NAG	O5-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
4	E	2	NAG	C4-C5-C6-O6
2	C	3	MAN	C4-C5-C6-O6
4	E	4	BMA	O5-C5-C6-O6
3	D	2	NAG	C4-C5-C6-O6
4	E	3	BMA	C4-C5-C6-O6
2	C	3	MAN	O5-C5-C6-O6
4	E	2	NAG	C8-C7-N2-C2
3	F	2	NAG	O5-C5-C6-O6
4	E	3	BMA	O5-C5-C6-O6
4	E	4	BMA	C4-C5-C6-O6
3	D	1	NAG	C4-C5-C6-O6
4	E	2	NAG	O7-C7-N2-C2
3	D	1	NAG	O5-C5-C6-O6
3	F	2	NAG	C4-C5-C6-O6

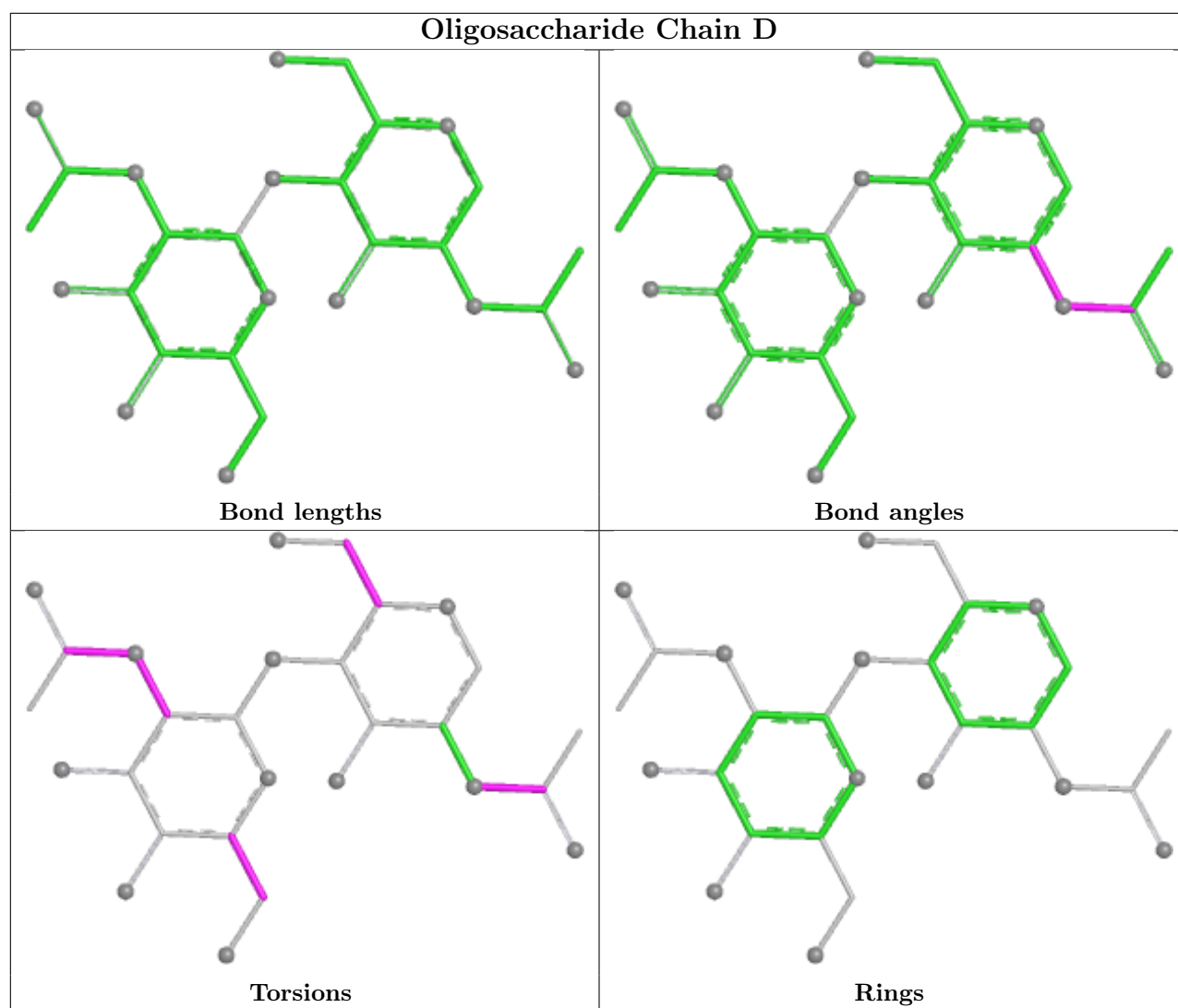
There are no ring outliers.

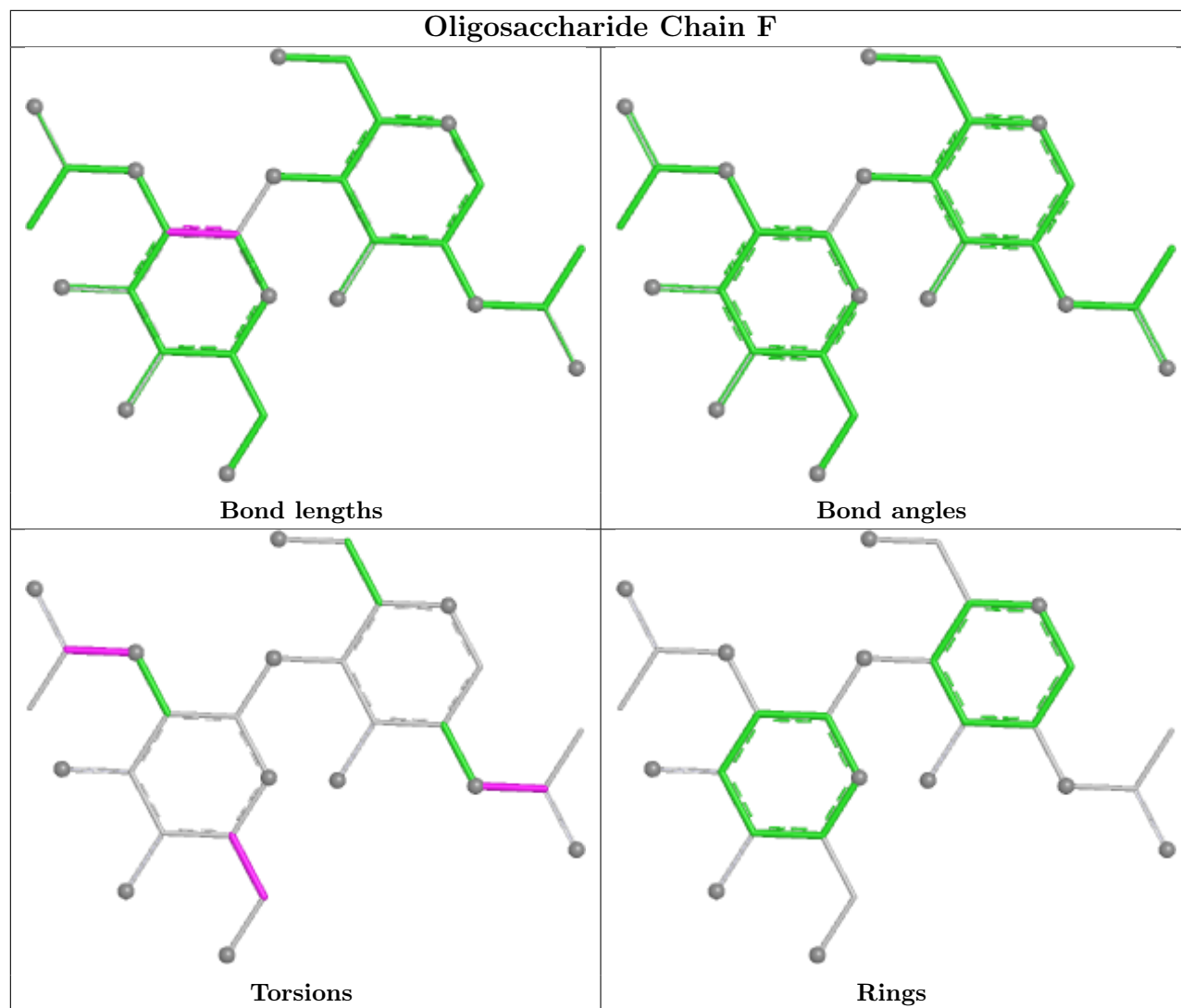
7 monomers are involved in 10 short contacts:

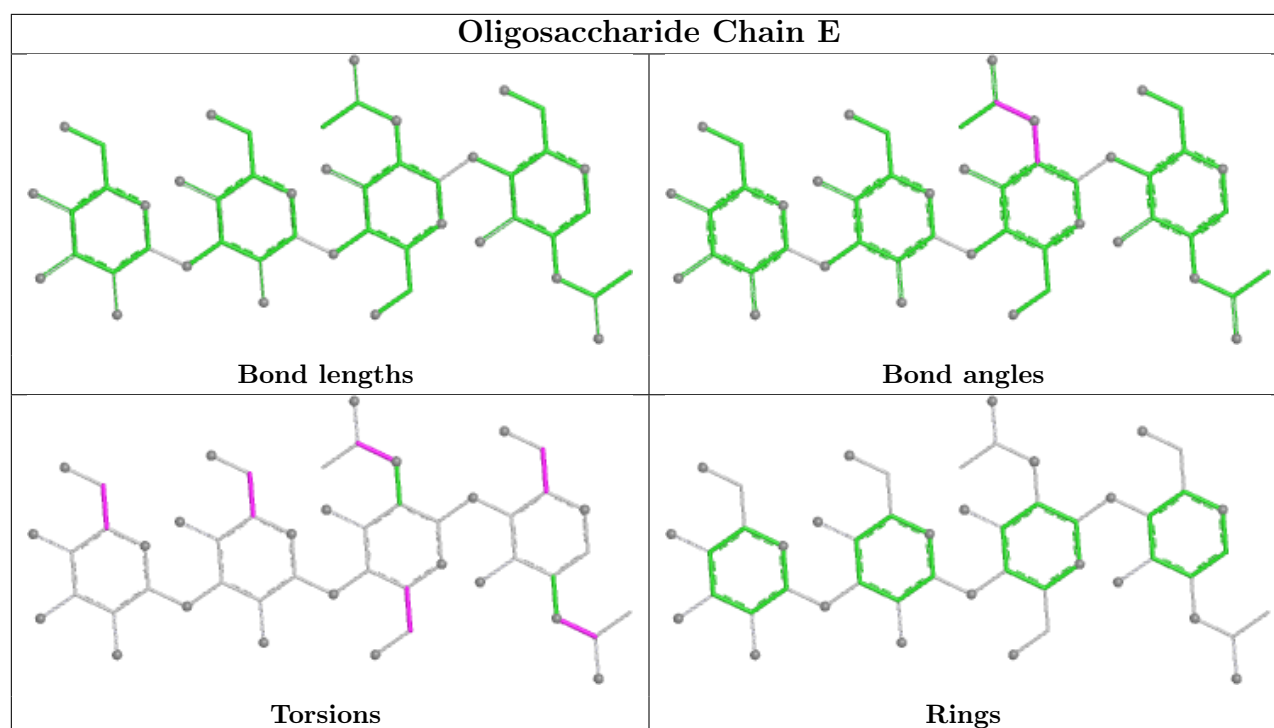
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	NAG	3	0
3	F	1	NAG	1	0
3	D	2	NAG	2	0
4	E	2	NAG	1	0
2	C	1	NAG	5	0
3	D	1	NAG	1	0
4	E	1	NAG	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.









5.6 Ligand geometry [i](#)

Of 31 ligands modelled in this entry, 13 are monoatomic - leaving 18 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	$\# Z > 2$
8	PLM	A	1322	1	15,16,17	0.21	0	14,15,17	0.59	0
5	BNG	A	1500	-	21,21,21	1.29	3 (14%)	26,26,26	0.60	0
5	BNG	A	1505	-	21,21,21	1.40	4 (19%)	26,26,26	0.66	0
5	BNG	B	1502	-	21,21,21	1.43	3 (14%)	26,26,26	0.64	0
10	HTO	A	1404	-	9,9,9	1.72	2 (22%)	10,10,10	0.70	0
10	HTO	A	1400	-	9,9,9	1.71	2 (22%)	10,10,10	0.66	0
8	PLM	A	1323	1	15,16,17	0.35	0	14,15,17	0.50	0
5	BNG	A	1504	-	21,21,21	1.31	3 (14%)	26,26,26	0.60	0
9	RET	A	1296	1	20,20,21	2.00	6 (30%)	27,27,28	1.95	10 (37%)
5	BNG	A	1501	-	21,21,21	1.26	2 (9%)	26,26,26	0.65	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	BNG	A	1503	-	21,21,21	1.23	3 (14%)	26,26,26	0.60	0
8	PLM	B	1322	1	15,16,17	0.34	0	14,15,17	0.52	0
10	HTO	A	1403	-	9,9,9	1.58	1 (11%)	10,10,10	0.50	0
10	HTO	A	1405	-	9,9,9	1.37	1 (11%)	10,10,10	0.45	0
10	HTO	B	1401	-	9,9,9	1.61	1 (11%)	10,10,10	0.81	0
9	RET	B	1296	1	20,20,21	1.95	6 (30%)	27,27,28	2.00	9 (33%)
10	HTO	B	1406	-	9,9,9	1.71	2 (22%)	10,10,10	0.78	0
5	BNG	B	1506	-	21,21,21	1.37	3 (14%)	26,26,26	0.64	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
8	PLM	A	1322	1	-	10/14/14/15	-
5	BNG	A	1500	-	-	8/12/32/32	0/1/1/1
5	BNG	A	1505	-	-	6/12/32/32	0/1/1/1
5	BNG	B	1502	-	-	7/12/32/32	0/1/1/1
10	HTO	A	1404	-	-	4/10/10/10	-
10	HTO	A	1400	-	-	3/10/10/10	-
8	PLM	A	1323	1	-	8/14/14/15	-
5	BNG	A	1504	-	-	8/12/32/32	0/1/1/1
9	RET	A	1296	1	-	7/13/30/31	0/1/1/1
5	BNG	A	1501	-	-	6/12/32/32	0/1/1/1
5	BNG	A	1503	-	-	7/12/32/32	0/1/1/1
8	PLM	B	1322	1	-	10/14/14/15	-
10	HTO	A	1403	-	-	3/10/10/10	-
10	HTO	A	1405	-	-	3/10/10/10	-
10	HTO	B	1401	-	-	2/10/10/10	-
9	RET	B	1296	1	-	5/13/30/31	0/1/1/1
10	HTO	B	1406	-	-	3/10/10/10	-
5	BNG	B	1506	-	-	10/12/32/32	0/1/1/1

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	B	1296	RET	C14-C13	5.46	1.37	1.33
9	A	1296	RET	C14-C13	4.48	1.36	1.33
10	B	1401	HTO	C3-C2	4.02	1.62	1.53
10	A	1404	HTO	C3-C2	3.97	1.62	1.53
5	B	1502	BNG	O1-C1	3.97	1.46	1.40
10	A	1400	HTO	C3-C2	3.96	1.62	1.53
10	A	1403	HTO	C3-C2	3.82	1.62	1.53
10	B	1406	HTO	C3-C2	3.76	1.62	1.53
5	A	1505	BNG	O1-C1	3.71	1.46	1.40
10	A	1405	HTO	C3-C2	3.39	1.61	1.53
9	A	1296	RET	C2-C1	3.32	1.61	1.54
5	B	1506	BNG	O1-C1	3.31	1.45	1.40
9	A	1296	RET	C18-C5	3.29	1.56	1.50
5	B	1502	BNG	O5-C1	3.29	1.50	1.41
5	B	1506	BNG	O5-C1	3.25	1.50	1.41
5	A	1503	BNG	O5-C1	3.20	1.50	1.41
5	A	1500	BNG	O5-C1	3.15	1.49	1.41
5	A	1505	BNG	O5-C1	3.13	1.49	1.41
5	A	1504	BNG	O5-C1	3.08	1.49	1.41
5	A	1501	BNG	O1-C1	3.08	1.45	1.40
5	A	1501	BNG	O5-C1	3.02	1.49	1.41
5	A	1500	BNG	O1-C1	2.94	1.45	1.40
5	A	1504	BNG	O1-C1	2.90	1.45	1.40
5	A	1503	BNG	O1-C1	2.80	1.44	1.40
9	A	1296	RET	C1-C6	-2.76	1.50	1.53
9	B	1296	RET	C17-C1	2.72	1.58	1.53
10	B	1406	HTO	C4-C3	2.71	1.57	1.52
9	A	1296	RET	C17-C1	2.54	1.58	1.53
5	A	1504	BNG	C4-C5	2.41	1.58	1.53
10	A	1400	HTO	C4-C3	2.40	1.56	1.52
5	B	1506	BNG	C4-C5	2.37	1.58	1.53
9	A	1296	RET	C5-C6	2.35	1.38	1.34
9	B	1296	RET	C18-C5	2.31	1.54	1.50
9	B	1296	RET	C7-C6	2.29	1.52	1.45
5	A	1500	BNG	C4-C5	2.26	1.57	1.53
9	B	1296	RET	C12-C13	2.23	1.50	1.46
10	A	1404	HTO	C4-C3	2.19	1.56	1.52
5	A	1503	BNG	C4-C5	2.15	1.57	1.53
9	B	1296	RET	C16-C1	2.09	1.57	1.53
5	A	1505	BNG	C4-C5	2.04	1.57	1.53
5	B	1502	BNG	C4-C5	2.01	1.57	1.53
5	A	1505	BNG	C3-C2	2.01	1.57	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	1296	RET	C11-C12-C13	4.81	139.56	126.36
9	B	1296	RET	C11-C12-C13	4.76	139.42	126.36
9	A	1296	RET	C17-C1-C6	3.95	116.44	110.24
9	B	1296	RET	C2-C1-C6	3.84	116.02	110.44
9	B	1296	RET	C19-C9-C8	-3.45	112.82	118.09
9	A	1296	RET	C2-C1-C6	3.40	115.38	110.44
9	B	1296	RET	C8-C7-C6	-3.10	118.71	127.00
9	B	1296	RET	C7-C8-C9	3.09	130.81	126.23
9	A	1296	RET	C7-C8-C9	2.97	130.62	126.23
9	B	1296	RET	C8-C9-C10	2.89	123.56	119.01
9	A	1296	RET	C19-C9-C8	-2.88	113.69	118.09
9	B	1296	RET	C10-C11-C12	2.85	131.47	123.20
9	B	1296	RET	C17-C1-C6	2.72	114.51	110.24
9	A	1296	RET	C8-C9-C10	2.53	122.99	119.01
9	A	1296	RET	C10-C11-C12	2.35	130.01	123.20
9	A	1296	RET	C17-C1-C2	-2.21	100.46	108.95
9	A	1296	RET	C3-C4-C5	-2.15	110.22	114.06
9	A	1296	RET	C8-C7-C6	-2.04	121.56	127.00
9	B	1296	RET	C18-C5-C6	2.00	126.67	124.48

There are no chirality outliers.

All (110) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1506	BNG	C2'-C1'-O1-C1
8	A	1323	PLM	C1-C2-C3-C4
8	B	1322	PLM	O2-C1-C2-C3
9	A	1296	RET	C20-C13-C14-C15
9	B	1296	RET	C11-C12-C13-C14
9	B	1296	RET	C20-C13-C14-C15
5	B	1506	BNG	O5-C5-C6-O6
5	A	1504	BNG	O5-C5-C6-O6
5	A	1500	BNG	O5-C5-C6-O6
5	B	1506	BNG	C4-C5-C6-O6
5	A	1504	BNG	C4-C5-C6-O6
9	A	1296	RET	C7-C8-C9-C19
9	B	1296	RET	C11-C12-C13-C20
5	A	1503	BNG	O1-C1'-C2'-C3'
10	A	1403	HTO	C2-C3-C4-C5
9	A	1296	RET	C11-C12-C13-C20
9	A	1296	RET	C7-C8-C9-C10
5	A	1503	BNG	C2'-C3'-C4'-C5'
5	A	1501	BNG	C1'-C2'-C3'-C4'

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Mol	Chain	Res	Type	Atoms
5	B	1506	BNG	C5'-C6'-C7'-C8'
8	A	1322	PLM	C9-CA-CB-CC
5	A	1504	BNG	C3'-C4'-C5'-C6'
5	A	1505	BNG	C2'-C3'-C4'-C5'
8	A	1323	PLM	C4-C5-C6-C7
8	A	1322	PLM	C4-C5-C6-C7
5	A	1501	BNG	C4'-C5'-C6'-C7'
5	A	1500	BNG	C1'-C2'-C3'-C4'
5	B	1502	BNG	C1'-C2'-C3'-C4'
8	A	1323	PLM	CB-CC-CD-CE
5	A	1503	BNG	C1'-C2'-C3'-C4'
5	A	1505	BNG	C5'-C6'-C7'-C8'
8	B	1322	PLM	CB-CC-CD-CE
8	A	1323	PLM	C7-C8-C9-CA
5	A	1500	BNG	C4-C5-C6-O6
5	B	1502	BNG	C5'-C6'-C7'-C8'
5	A	1504	BNG	C2'-C3'-C4'-C5'
5	A	1504	BNG	C1'-C2'-C3'-C4'
5	A	1501	BNG	C2'-C3'-C4'-C5'
5	B	1506	BNG	C1'-C2'-C3'-C4'
5	A	1500	BNG	C2'-C3'-C4'-C5'
5	A	1503	BNG	C5'-C6'-C7'-C8'
5	A	1504	BNG	C5'-C6'-C7'-C8'
5	A	1504	BNG	C4'-C5'-C6'-C7'
10	A	1403	HTO	O3-C3-C4-C5
8	A	1322	PLM	C6-C7-C8-C9
8	B	1322	PLM	C6-C7-C8-C9
5	A	1503	BNG	C4'-C5'-C6'-C7'
5	A	1501	BNG	C5'-C6'-C7'-C8'
5	B	1506	BNG	C2'-C3'-C4'-C5'
5	A	1500	BNG	C5'-C6'-C7'-C8'
5	B	1506	BNG	O1-C1'-C2'-C3'
8	B	1322	PLM	C8-C9-CA-CB
8	B	1322	PLM	C3-C4-C5-C6
5	A	1500	BNG	C3'-C4'-C5'-C6'
8	A	1323	PLM	CA-CB-CC-CD
8	A	1323	PLM	C2-C3-C4-C5
8	B	1322	PLM	C4-C5-C6-C7
5	B	1506	BNG	C3'-C4'-C5'-C6'
5	B	1506	BNG	C4'-C5'-C6'-C7'
8	A	1322	PLM	C2-C3-C4-C5
8	A	1322	PLM	C3-C4-C5-C6

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Mol	Chain	Res	Type	Atoms
5	B	1502	BNG	O1-C1'-C2'-C3'
5	A	1500	BNG	C4'-C5'-C6'-C7'
9	A	1296	RET	C12-C13-C14-C15
9	B	1296	RET	C12-C13-C14-C15
5	A	1500	BNG	C6'-C7'-C8'-C9'
5	A	1503	BNG	C6'-C7'-C8'-C9'
5	B	1506	BNG	C6'-C7'-C8'-C9'
5	A	1505	BNG	C4'-C5'-C6'-C7'
5	A	1504	BNG	C6'-C7'-C8'-C9'
5	A	1505	BNG	C6'-C7'-C8'-C9'
5	A	1501	BNG	C6'-C7'-C8'-C9'
8	B	1322	PLM	C5-C6-C7-C8
5	B	1502	BNG	C6'-C7'-C8'-C9'
10	B	1406	HTO	C4-C5-C6-C7
8	A	1323	PLM	C3-C4-C5-C6
10	A	1404	HTO	O3-C3-C4-C5
5	A	1503	BNG	C3'-C4'-C5'-C6'
8	A	1322	PLM	CA-CB-CC-CD
8	A	1322	PLM	CC-CD-CE-CF
8	A	1323	PLM	CD-CE-CF-CG
5	B	1502	BNG	C2'-C3'-C4'-C5'
5	B	1502	BNG	O5-C5-C6-O6
5	A	1505	BNG	C3'-C4'-C5'-C6'
10	A	1403	HTO	C1-C2-C3-O3
10	A	1404	HTO	C1-C2-C3-O3
10	A	1405	HTO	C1-C2-C3-O3
10	B	1401	HTO	C1-C2-C3-O3
8	B	1322	PLM	C1-C2-C3-C4
10	A	1400	HTO	C3-C4-C5-C6
10	A	1404	HTO	C2-C3-C4-C5
8	A	1322	PLM	C5-C6-C7-C8
5	B	1502	BNG	C3'-C4'-C5'-C6'
9	A	1296	RET	C10-C11-C12-C13
9	B	1296	RET	C10-C11-C12-C13
10	A	1400	HTO	C4-C5-C6-C7
9	A	1296	RET	C11-C12-C13-C14
8	B	1322	PLM	C2-C3-C4-C5
8	B	1322	PLM	CC-CD-CE-CF
10	B	1406	HTO	C3-C4-C5-C6
10	A	1404	HTO	C4-C5-C6-C7
5	A	1501	BNG	C3'-C4'-C5'-C6'
10	A	1400	HTO	C1-C2-C3-O3

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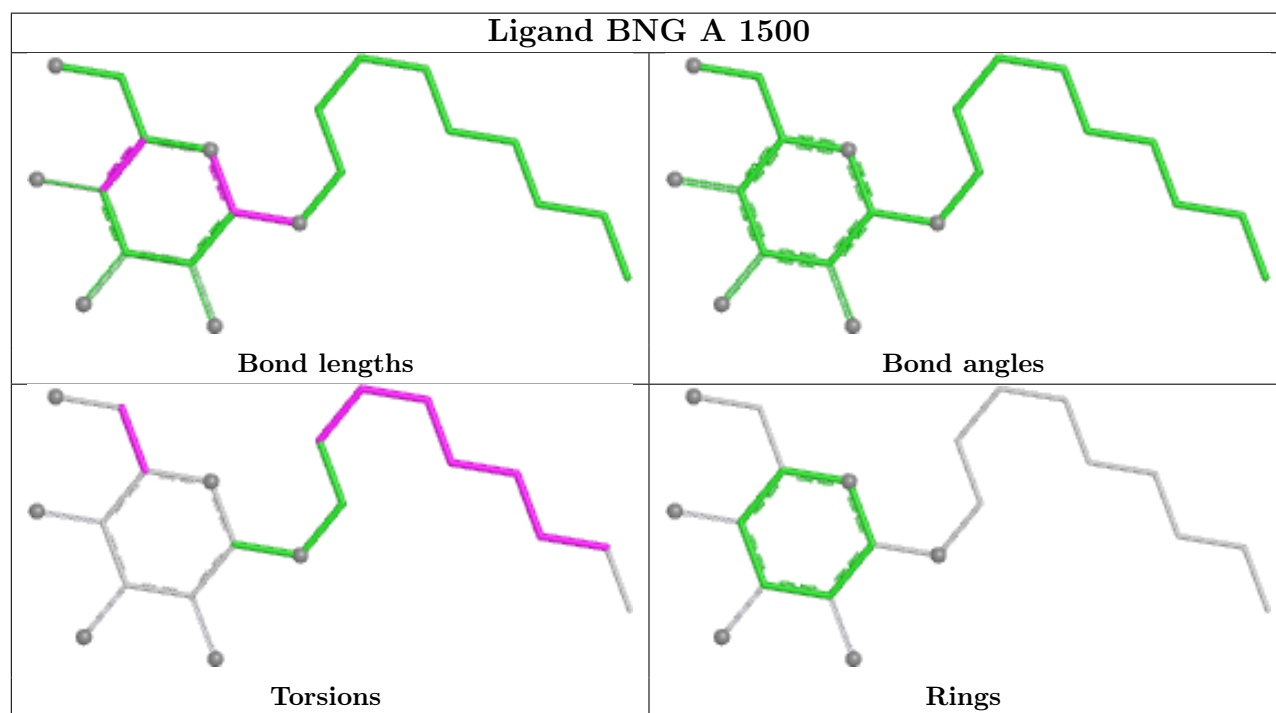
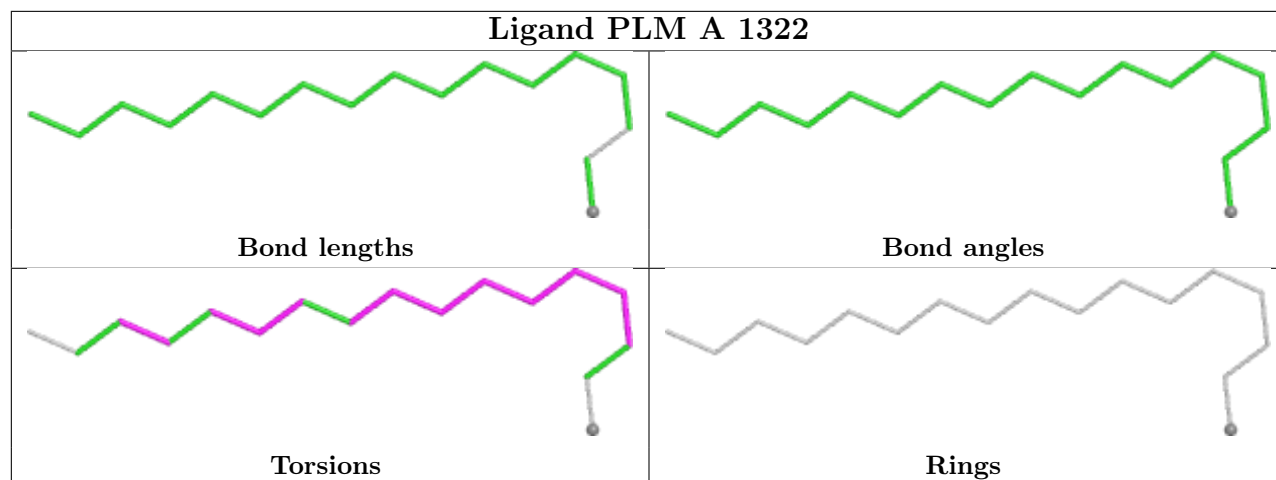
Mol	Chain	Res	Type	Atoms
10	B	1406	HTO	C1-C2-C3-O3
8	A	1322	PLM	C1-C2-C3-C4
10	A	1405	HTO	C4-C5-C6-C7
10	A	1405	HTO	C3-C4-C5-C6
10	B	1401	HTO	C3-C4-C5-C6
8	A	1322	PLM	C7-C8-C9-CA
5	A	1505	BNG	O1-C1'-C2'-C3'

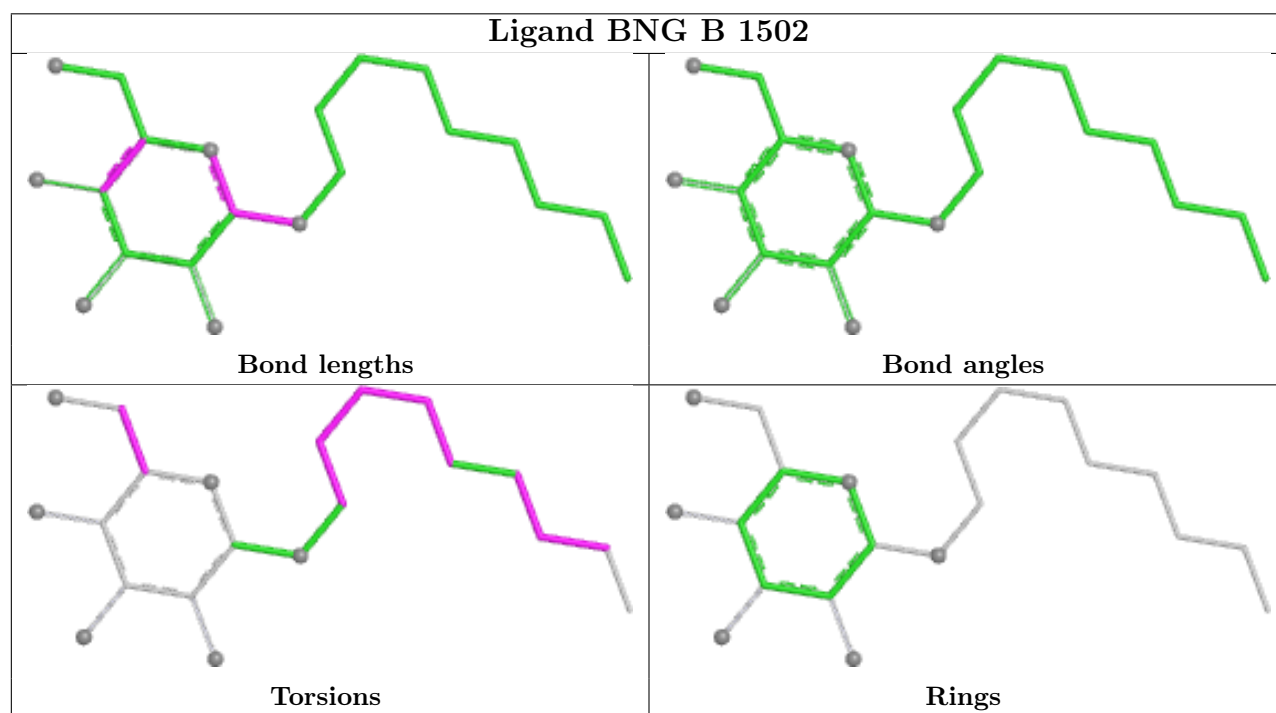
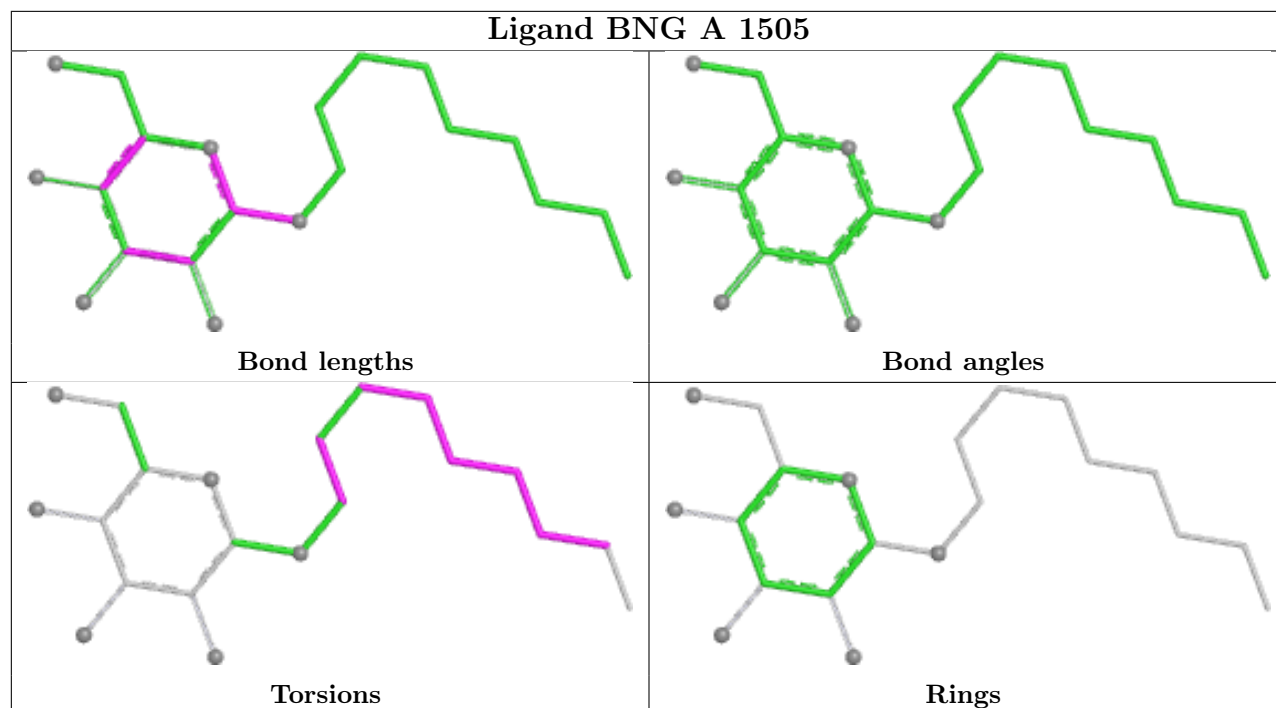
There are no ring outliers.

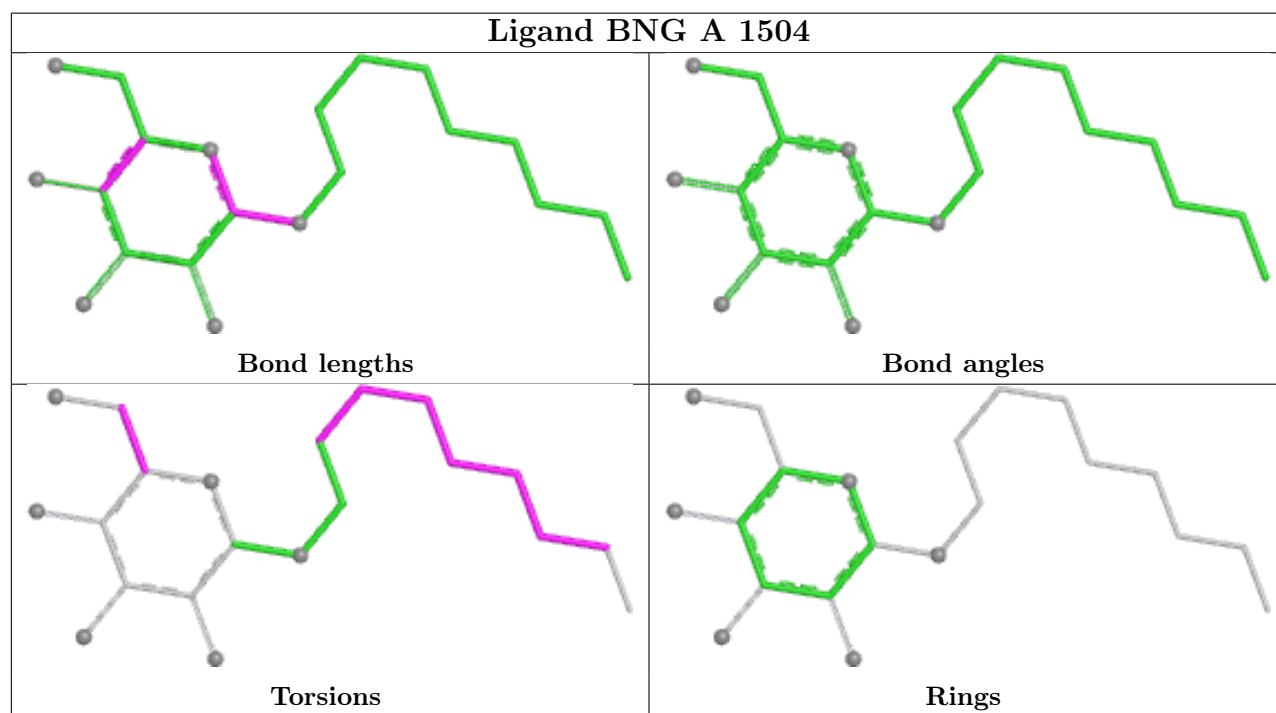
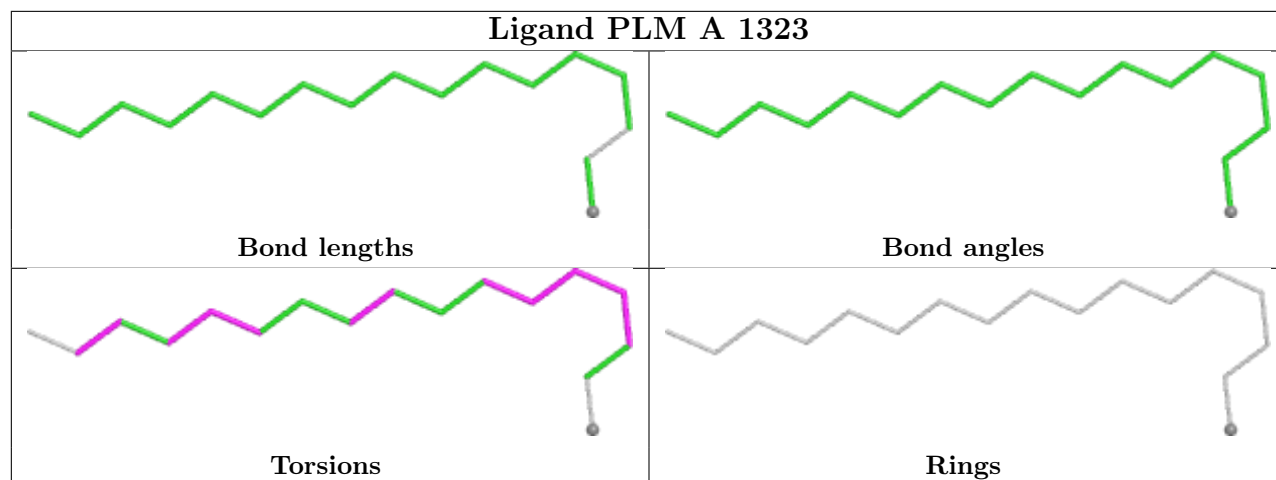
15 monomers are involved in 29 short contacts:

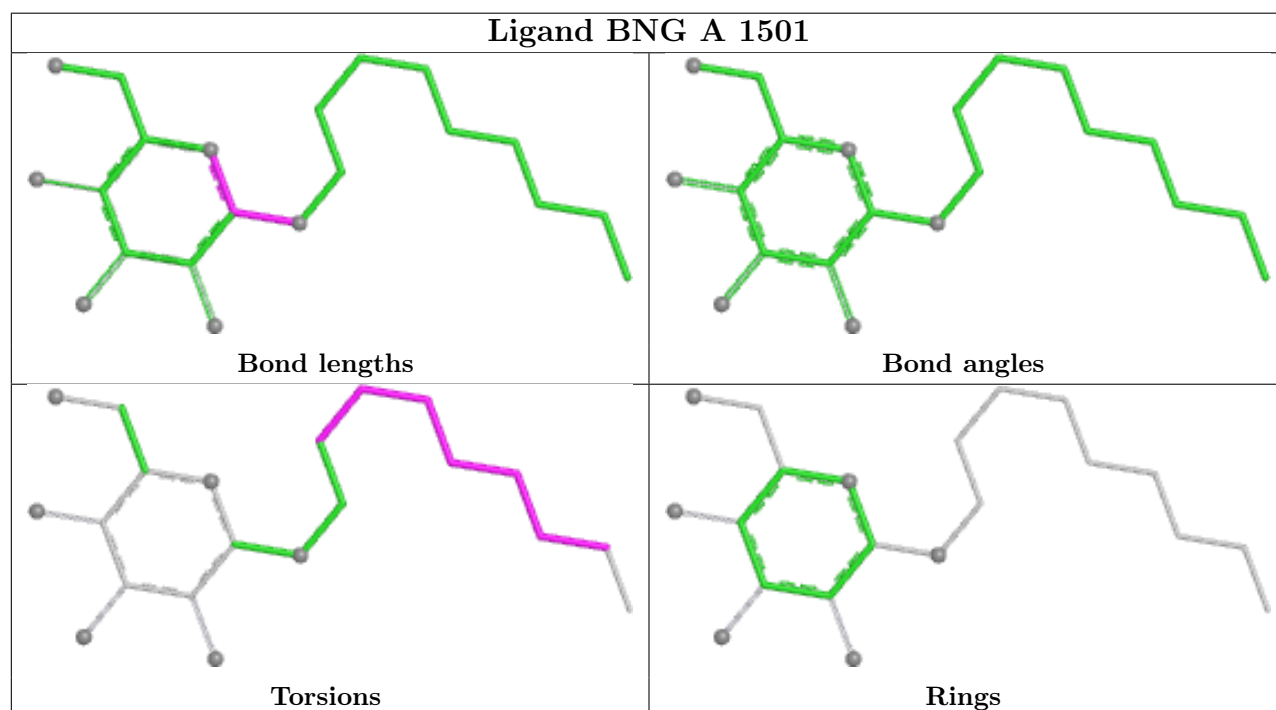
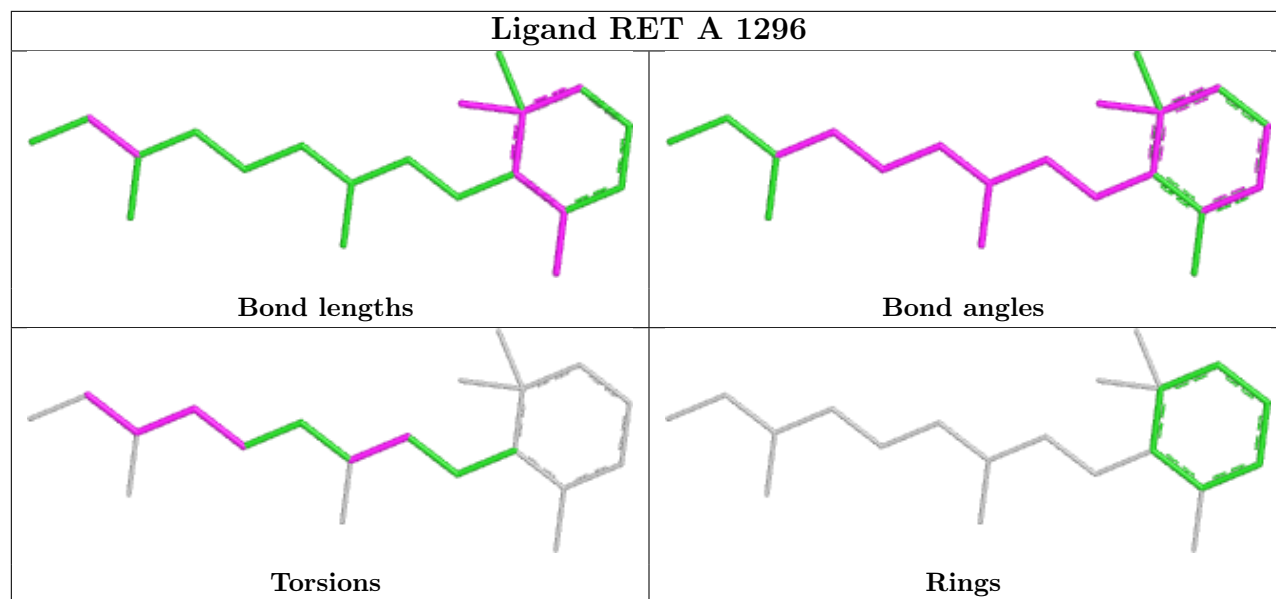
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	1322	PLM	1	0
5	A	1500	BNG	4	0
5	A	1505	BNG	1	0
10	A	1404	HTO	1	0
10	A	1400	HTO	1	0
5	A	1504	BNG	1	0
9	A	1296	RET	1	0
5	A	1501	BNG	2	0
5	A	1503	BNG	1	0
8	B	1322	PLM	1	0
10	A	1403	HTO	2	0
10	B	1401	HTO	1	0
9	B	1296	RET	1	0
10	B	1406	HTO	1	0
5	B	1506	BNG	10	0

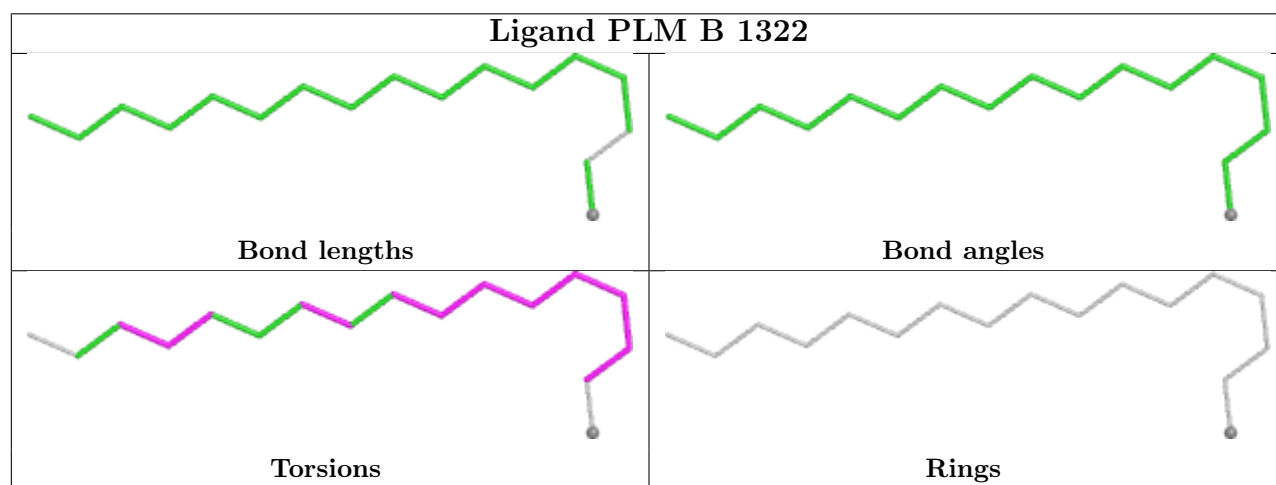
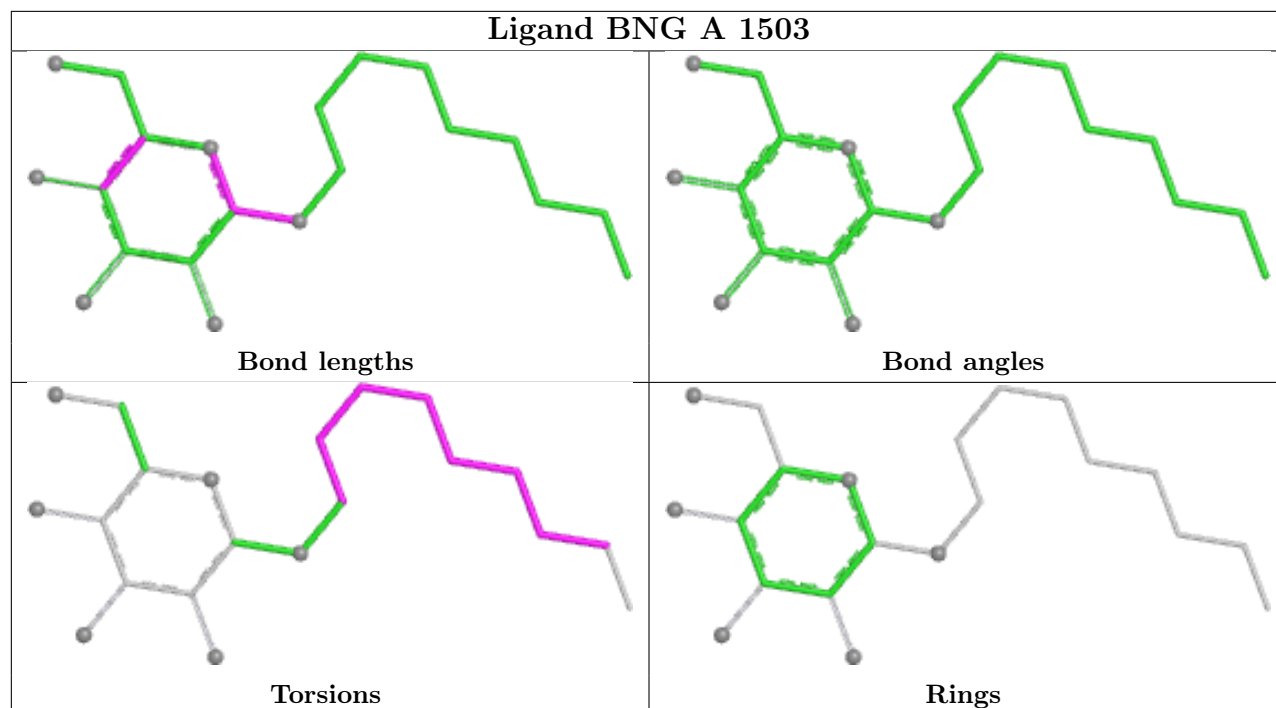
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

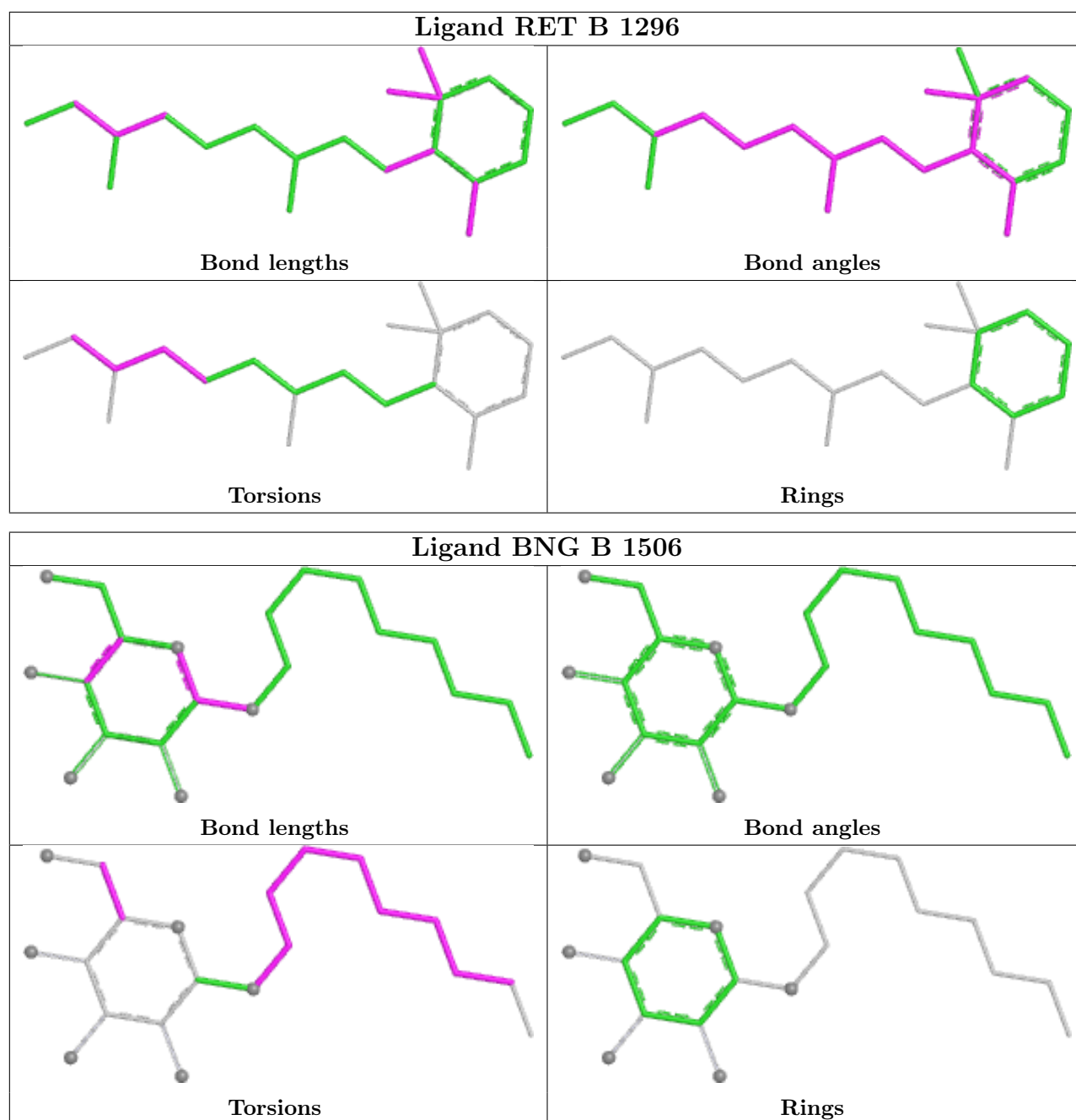












5.7 Other polymers [i](#)

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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.