



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 08:46 AM UTC

PDB ID : 1CQE / pdb_00001cqe
Title : PROSTAGLANDIN H2 SYNTHASE-1 COMPLEX WITH FLURBIPROFEN
Authors : Picot, D.; Loll, P.J.; Mulichak, A.M.; Garavito, R.M.
Deposited on : 1999-06-15
Resolution : 3.10 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

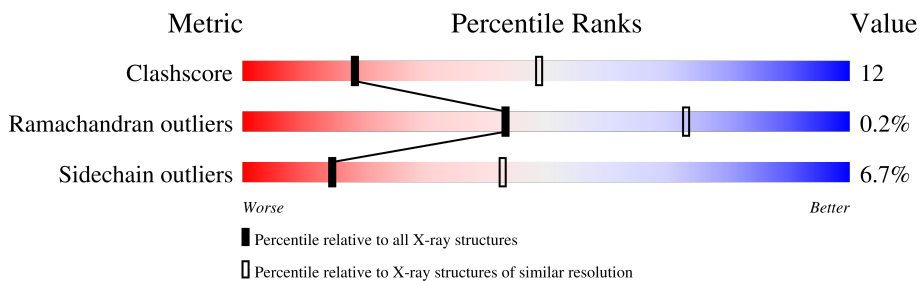
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)

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	580	
1	B	580	
2	C	2	
2	D	2	
2	E	2	
2	F	2	
2	G	2	
2	H	2	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 9413 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 PROTEIN (PROSTAGLANDIN H2 SYNTHASE-1).

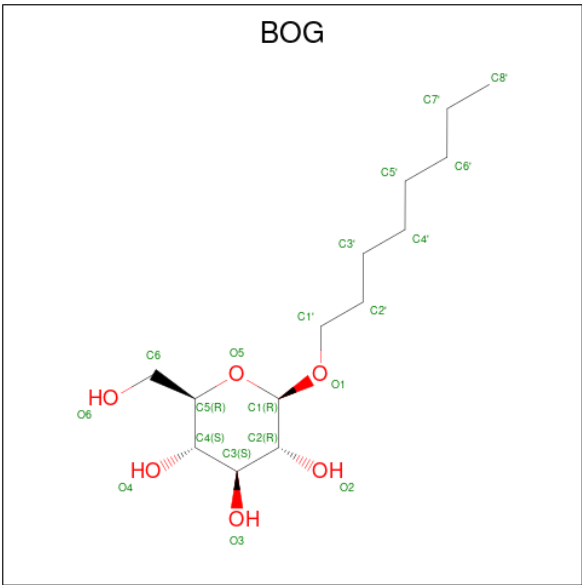
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	552	Total	C	N	O	S	0	0	0
			4458	2891	750	789	28			
1	B	553	Total	C	N	O	S	0	0	0
			4465	2896	751	790	28			

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



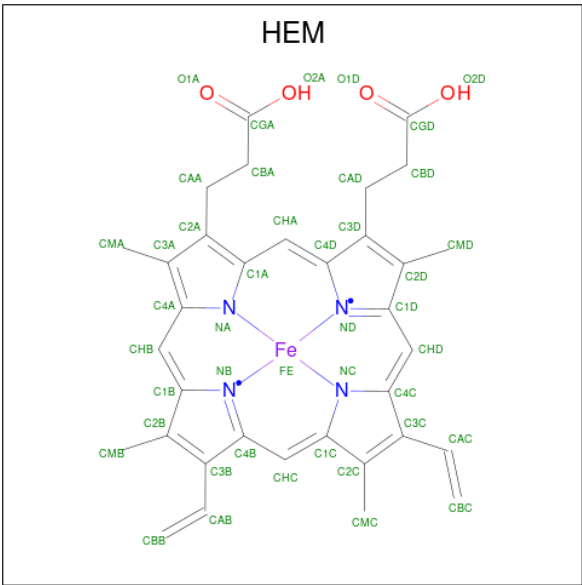
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	D	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	H	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is octyl beta-D-glucopyranoside (CCD ID: BOG) (formula: C₁₄H₂₈O₆).



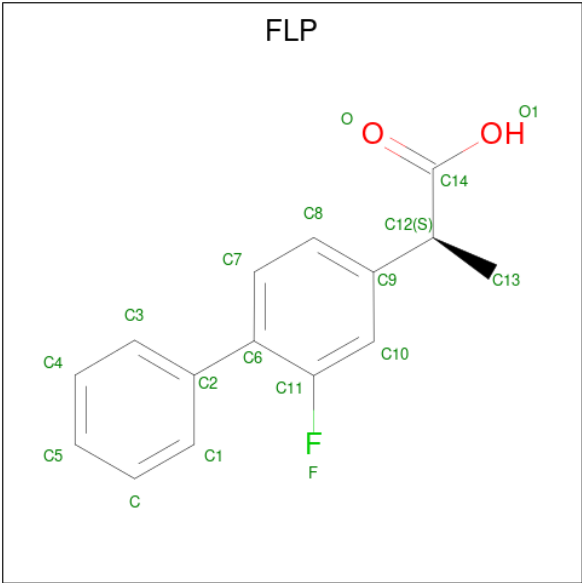
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			20	14	6		
3	A	1	Total	C	O	0	0
			10	9	1		
3	A	1	Total	C	O	0	0
			20	14	6		
3	B	1	Total	C	O	0	0
			10	9	1		
3	B	1	Total	C	O	0	0
			10	9	1		

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: C₃₄H₃₂FeN₄O₄).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 5 is FLURBIPROFEN (CCD ID: FLP) (formula: C₁₅H₁₃FO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	F	O	0	0
			18	15	1	2		
5	B	1	Total	C	F	O	0	0
			18	15	1	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	72	Total 72	O 72	0	0
6	B	58	Total 58	O 58	0	0

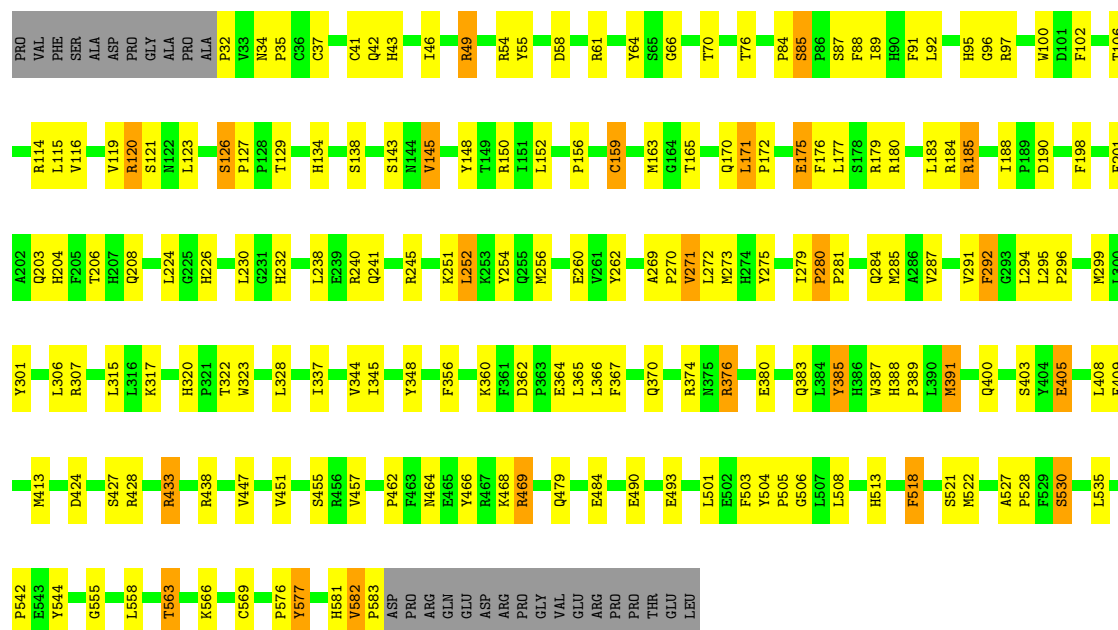
3 Residue-property plots [i](#)

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.

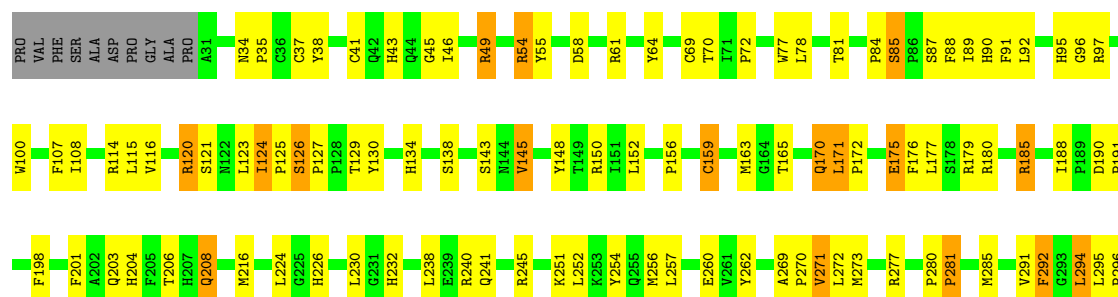
• Molecule 1: PROTEIN (PROSTAGLANDIN H2 SYNTHASE-1)

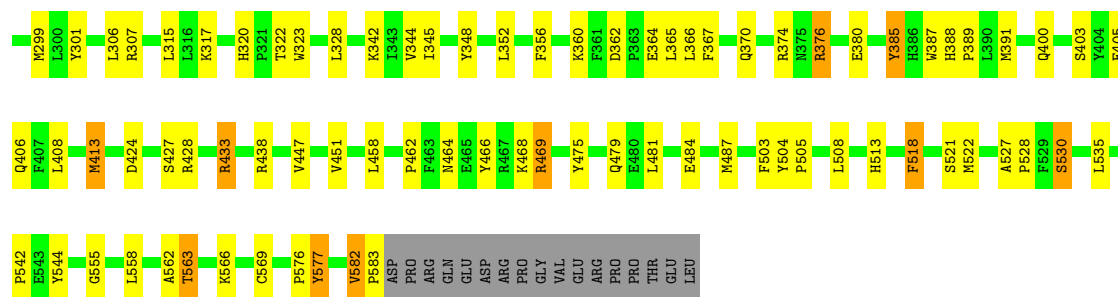
Chain A: 



• Molecule 1: PROTEIN (PROSTAGLANDIN H2 SYNTHASE-1)

Chain B: 





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

Chain C: 100%



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

Chain D: 50% 50%



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

Chain E: 50% 50%



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

Chain F: 100%



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

Chain G: 100%



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

Chain H:

100%

3AG1
3AG2

4 Data and refinement statistics

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

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	99.40Å 210.30Å 233.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.10	Depositor
% Data completeness (in resolution range)	94.3 (20.00-3.10)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.186 , 0.226	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9413	wwPDB-VP
Average B, all atoms (Å ²)	27.0	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: NAG, FLP, BOG, HEM

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	0.76	0/4596	1.14	36/6245 (0.6%)
1	B	0.79	0/4604	1.12	32/6257 (0.5%)
All	All	0.77	0/9200	1.13	68/12502 (0.5%)

There are no bond length outliers.

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	280	PRO	CA-C-N	9.62	129.24	119.24
1	A	280	PRO	C-N-CA	9.62	129.24	119.24
1	A	190	ASP	CA-C-N	8.90	128.55	119.56
1	A	190	ASP	C-N-CA	8.90	128.55	119.56
1	B	148	TYR	N-CA-C	-8.30	98.19	110.48
1	A	148	TYR	N-CA-C	-7.93	97.71	110.32
1	B	269	ALA	CA-C-N	7.85	129.66	119.84
1	B	269	ALA	C-N-CA	7.85	129.66	119.84
1	A	569	CYS	N-CA-C	7.81	120.69	111.71
1	A	85	SER	CA-C-N	7.65	128.04	119.32
1	A	85	SER	C-N-CA	7.65	128.04	119.32
1	B	582	VAL	N-CA-C	-7.62	101.94	109.02
1	A	513	HIS	N-CA-C	-7.61	99.24	110.20
1	B	85	SER	CA-C-N	7.53	127.91	119.32
1	B	85	SER	C-N-CA	7.53	127.91	119.32
1	B	513	HIS	N-CA-C	-7.40	99.54	110.20
1	B	569	CYS	N-CA-C	7.40	120.22	111.71
1	A	582	VAL	N-CA-C	-7.14	102.38	109.02
1	A	32	PRO	N-CA-CB	6.47	110.12	103.00
1	A	129	THR	N-CA-C	6.30	118.87	108.48
1	B	41	CYS	N-CA-C	6.07	117.89	109.15
1	A	269	ALA	CA-C-N	6.05	127.41	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	269	ALA	C-N-CA	6.05	127.41	119.84
1	B	129	THR	N-CA-C	6.03	118.44	108.48
1	A	66	GLY	CA-C-N	5.99	125.87	119.28
1	A	66	GLY	C-N-CA	5.99	125.87	119.28
1	B	577	TYR	N-CA-C	-5.96	100.83	109.59
1	B	345	ILE	N-CA-C	5.94	116.11	110.53
1	A	43	HIS	N-CA-C	5.93	119.58	111.39
1	B	43	HIS	N-CA-C	5.92	119.56	111.39
1	B	530	SER	N-CA-C	5.91	117.52	111.14
1	A	530	SER	N-CA-C	5.88	117.49	111.14
1	A	345	ILE	N-CA-C	5.87	116.04	110.53
1	B	190	ASP	CA-C-N	5.82	125.50	119.56
1	B	190	ASP	C-N-CA	5.82	125.50	119.56
1	A	292	PHE	N-CA-C	5.80	119.83	112.87
1	A	577	TYR	N-CA-C	-5.79	101.08	109.59
1	B	121	SER	N-CA-C	5.74	117.54	111.28
1	A	121	SER	N-CA-C	5.64	117.43	111.28
1	B	159	CYS	CA-C-N	5.59	126.83	119.84
1	B	159	CYS	C-N-CA	5.59	126.83	119.84
1	B	281	PRO	N-CA-C	-5.52	106.34	113.57
1	B	208	GLN	N-CA-C	-5.50	106.69	113.41
1	B	521	SER	N-CA-C	5.49	117.26	111.28
1	A	171	LEU	CA-C-N	5.45	125.46	119.90
1	A	171	LEU	C-N-CA	5.45	125.46	119.90
1	B	503	PHE	N-CA-C	5.43	117.64	111.02
1	B	171	LEU	CA-C-N	5.39	125.40	119.90
1	B	171	LEU	C-N-CA	5.39	125.40	119.90
1	A	370	GLN	N-CA-C	-5.36	99.78	108.41
1	A	521	SER	N-CA-C	5.36	117.12	111.28
1	A	143	SER	N-CA-C	5.36	119.94	113.41
1	A	41	CYS	N-CA-C	5.35	116.97	108.67
1	B	143	SER	N-CA-C	5.31	119.89	113.41
1	B	292	PHE	N-CA-C	5.31	119.24	112.87
1	A	238	LEU	N-CA-C	5.30	117.47	111.11
1	A	503	PHE	N-CA-C	5.29	117.48	111.02
1	A	457	VAL	N-CA-C	-5.19	105.65	110.53
1	B	124	ILE	CA-C-N	5.19	125.11	119.76
1	B	124	ILE	C-N-CA	5.19	125.11	119.76
1	B	238	LEU	N-CA-C	5.10	117.22	111.11
1	A	337	ILE	N-CA-C	-5.09	105.42	110.62
1	A	159	CYS	CA-C-N	5.07	126.18	119.84
1	A	159	CYS	C-N-CA	5.07	126.18	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	370	GLN	N-CA-C	-5.07	99.53	108.20
1	B	413	MET	N-CA-C	5.03	116.76	111.28
1	A	391	MET	CA-C-N	5.01	126.34	120.98
1	A	391	MET	C-N-CA	5.01	126.34	120.98

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	4458	0	4330	104	0
1	B	4465	0	4340	114	0
2	C	28	0	25	1	0
2	D	28	0	25	0	0
2	E	28	0	25	0	0
2	F	28	0	25	3	0
2	G	28	0	25	0	0
2	H	28	0	25	0	0
3	A	50	0	73	3	0
3	B	20	0	34	1	0
4	A	43	0	30	5	0
4	B	43	0	30	4	0
5	A	18	0	12	1	0
5	B	18	0	12	0	0
6	A	72	0	0	6	0
6	B	58	0	0	4	0
All	All	9413	0	9011	220	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (220) 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:B:185:ARG:HH11	1:B:438:ARG:HH11	1.19	0.89
1:B:185:ARG:NH1	1:B:438:ARG:HH11	1.74	0.85
1:B:84:PRO:HG2	1:B:89:ILE:HD11	1.59	0.82
1:B:185:ARG:NH1	1:B:438:ARG:NH1	2.29	0.81
1:A:84:PRO:HG2	1:A:89:ILE:HD11	1.63	0.80
1:A:150:ARG:HD3	1:A:152:LEU:O	1.82	0.80
1:B:150:ARG:HD3	1:B:152:LEU:O	1.83	0.78
1:B:176:PHE:CZ	1:B:180:ARG:HD2	2.18	0.78
1:A:116:VAL:O	1:A:120:ARG:HB2	1.86	0.76
1:B:116:VAL:O	1:B:120:ARG:HB2	1.87	0.74
1:B:391:MET:HG3	4:B:601:HEM:HAB	1.69	0.74
1:A:88:PHE:CE2	1:A:92:LEU:HD11	2.23	0.73
1:A:391:MET:HG3	4:A:601:HEM:HAB	1.72	0.72
1:A:163:MET:HE1	6:A:878:HOH:O	1.89	0.71
1:B:88:PHE:CE2	1:B:92:LEU:HD11	2.27	0.70
1:B:277:ARG:HG2	1:B:277:ARG:HH11	1.57	0.70
1:A:577:TYR:CE2	1:A:583:PRO:HD3	2.28	0.69
1:B:563:THR:HG22	1:B:566:LYS:H	1.58	0.68
3:A:1702:BOG:C1	6:A:1124:HOH:O	2.40	0.68
1:A:563:THR:HG22	1:A:566:LYS:H	1.58	0.67
1:B:577:TYR:CE2	1:B:583:PRO:HD3	2.28	0.67
2:F:1:NAG:H3	2:F:2:NAG:HN2	1.60	0.66
1:B:156:PRO:HB2	1:B:159:CYS:SG	2.36	0.65
1:A:123:LEU:O	1:A:469:ARG:NH2	2.31	0.64
1:B:226:HIS:CE1	1:B:376:ARG:HD2	2.34	0.63
1:B:84:PRO:CG	1:B:89:ILE:HD11	2.30	0.62
1:B:208:GLN:NE2	1:B:230:LEU:H	1.97	0.62
1:A:175:GLU:O	1:A:179:ARG:HG3	1.99	0.62
1:A:208:GLN:NE2	1:A:230:LEU:H	1.96	0.62
1:A:226:HIS:CE1	1:A:376:ARG:HD2	2.35	0.61
1:A:364:GLU:HG2	1:A:367:PHE:CE2	2.35	0.61
1:B:364:GLU:HG2	1:B:367:PHE:CE2	2.36	0.61
1:B:172:PRO:HB2	1:B:177:LEU:HD22	1.83	0.61
1:A:156:PRO:HB2	1:A:159:CYS:SG	2.41	0.60
1:B:366:LEU:HD12	1:B:535:LEU:HD12	1.83	0.60
1:A:49:ARG:NH1	6:A:842:HOH:O	2.34	0.60
1:A:294:LEU:HD22	1:A:409:PHE:CE1	2.36	0.60
1:B:145:VAL:HG12	1:B:224:LEU:HD22	1.84	0.59
1:A:88:PHE:CZ	1:A:92:LEU:HD21	2.37	0.59
1:A:145:VAL:HG12	1:A:224:LEU:HD22	1.84	0.59
1:B:240:ARG:HG3	1:B:271:VAL:CG2	2.33	0.59
1:B:294:LEU:O	1:B:294:LEU:HG	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:PRO:CG	1:A:89:ILE:HD11	2.32	0.59
1:A:275:TYR:CE2	1:A:284:GLN:HB3	2.38	0.59
1:B:77:TRP:CE2	1:B:81:THR:HG21	2.38	0.58
1:B:123:LEU:O	1:B:469:ARG:NH2	2.37	0.58
1:B:91:PHE:CE2	1:B:95:HIS:CD2	2.93	0.57
1:B:91:PHE:CZ	1:B:95:HIS:HD2	2.23	0.57
1:B:185:ARG:NH1	1:B:438:ARG:HG2	2.19	0.57
1:B:175:GLU:O	1:B:179:ARG:HG3	2.05	0.56
1:B:241:GLN:HE21	1:B:245:ARG:HH11	1.54	0.56
1:A:163:MET:HB3	1:A:462:PRO:HG3	1.88	0.55
1:B:438:ARG:HE	1:B:487:MET:HE2	1.71	0.55
1:A:504:TYR:HB3	1:A:505:PRO:HD3	1.89	0.55
1:B:315:LEU:HD12	1:B:558:LEU:HD11	1.87	0.54
1:A:360:LYS:HE2	1:A:362:ASP:HB2	1.89	0.54
1:A:366:LEU:HD12	1:A:535:LEU:HD12	1.88	0.54
1:A:241:GLN:HE21	1:A:245:ARG:HH11	1.55	0.54
1:B:91:PHE:CE2	1:B:95:HIS:HD2	2.26	0.54
1:A:172:PRO:HB2	1:A:177:LEU:HD22	1.89	0.54
1:A:102:PHE:O	1:A:106:THR:HG23	2.07	0.53
1:A:403:SER:OG	1:A:405:GLU:HG2	2.07	0.53
1:B:88:PHE:CZ	1:B:92:LEU:HD21	2.44	0.53
1:A:287:VAL:HA	6:A:833:HOH:O	2.08	0.53
1:B:45:GLY:HA3	1:B:69:CYS:SG	2.49	0.53
1:B:518:PHE:CD2	1:B:522:MET:HG2	2.44	0.53
1:A:240:ARG:HD3	6:A:884:HOH:O	2.09	0.52
1:A:315:LEU:HD12	1:A:558:LEU:HD11	1.90	0.52
1:A:91:PHE:CE2	1:A:95:HIS:CD2	2.97	0.52
1:A:279:ILE:O	1:A:284:GLN:NE2	2.43	0.52
1:B:388:HIS:N	1:B:389:PRO:HD2	2.24	0.52
1:A:91:PHE:CZ	1:A:95:HIS:HD2	2.27	0.52
1:B:100:TRP:CD1	1:B:356:PHE:HB2	2.44	0.52
1:B:176:PHE:CE1	1:B:180:ARG:HD2	2.45	0.52
3:A:1702:BOG:C1	6:A:885:HOH:O	2.57	0.51
1:B:380:GLU:HG2	1:B:466:TYR:CE1	2.45	0.51
1:A:518:PHE:CD2	1:A:522:MET:HG2	2.45	0.51
1:B:61:ARG:NH2	6:B:1005:HOH:O	2.43	0.51
1:B:582:VAL:HG22	1:B:583:PRO:HD2	1.93	0.50
1:B:360:LYS:HE2	1:B:362:ASP:HB2	1.92	0.50
1:B:527:ALA:HB3	1:B:528:PRO:HD3	1.92	0.50
1:B:464:ASN:O	1:B:468:LYS:HG3	2.11	0.50
1:B:504:TYR:CZ	1:B:508:LEU:HD11	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:ASN:HB3	1:B:37:CYS:SG	2.52	0.50
1:B:504:TYR:HB3	1:B:505:PRO:HD3	1.93	0.50
1:B:163:MET:HB3	1:B:462:PRO:HG3	1.93	0.50
1:B:403:SER:OG	1:B:406:GLN:HG3	2.12	0.50
1:A:114:ARG:HD3	1:A:365:LEU:O	2.11	0.50
1:A:176:PHE:CZ	1:A:180:ARG:HG3	2.47	0.50
1:B:387:TRP:HB2	4:B:601:HEM:HAC	1.94	0.50
1:B:424:ASP:HB2	1:B:576:PRO:HB2	1.94	0.50
1:A:388:HIS:N	1:A:389:PRO:HD2	2.27	0.49
1:A:504:TYR:CZ	1:A:508:LEU:HD11	2.47	0.49
1:A:126:SER:HA	1:A:127:PRO:C	2.37	0.49
1:A:582:VAL:HG22	1:A:583:PRO:HD2	1.93	0.49
1:B:317:LYS:HD3	1:B:328:LEU:HD11	1.95	0.49
1:A:91:PHE:CE2	1:A:95:HIS:HD2	2.30	0.49
1:A:380:GLU:HG2	1:A:466:TYR:CE1	2.48	0.49
1:A:527:ALA:HB3	1:A:528:PRO:HD3	1.95	0.49
1:A:555:GLY:HA2	1:A:558:LEU:HD12	1.95	0.48
1:A:294:LEU:HD22	1:A:409:PHE:HE1	1.77	0.48
1:A:100:TRP:CD1	1:A:356:PHE:HB2	2.48	0.48
1:A:61:ARG:NH1	1:B:542:PRO:O	2.47	0.48
1:A:179:ARG:HH11	1:A:179:ARG:HB3	1.78	0.48
3:A:1701:BOG:H1'2	3:A:1701:BOG:O2	2.13	0.48
2:F:1:NAG:H61	2:F:2:NAG:C1	2.43	0.48
1:B:447:VAL:O	1:B:451:VAL:HG23	2.13	0.48
1:A:180:ARG:HD3	1:A:490:GLU:OE2	2.14	0.48
1:B:85:SER:O	1:B:89:ILE:HG12	2.14	0.48
1:A:34:ASN:HB3	1:A:37:CYS:SG	2.53	0.48
1:A:179:ARG:O	1:A:185:ARG:NH2	2.47	0.48
1:A:391:MET:CG	4:A:601:HEM:HAB	2.42	0.48
1:B:188:ILE:HD13	1:B:433:ARG:HH21	1.79	0.48
1:A:294:LEU:HD22	1:A:409:PHE:CD1	2.49	0.47
1:A:184:ARG:HA	1:A:438:ARG:O	2.14	0.47
1:A:427:SER:HB3	1:A:577:TYR:CD2	2.50	0.47
1:B:134:HIS:HD2	1:B:138:SER:OG	1.98	0.47
1:A:134:HIS:HD2	1:A:138:SER:OG	1.96	0.47
1:A:387:TRP:HB2	4:A:601:HEM:HAC	1.97	0.47
1:A:424:ASP:HB2	1:A:576:PRO:HB2	1.97	0.47
1:A:542:PRO:O	1:B:61:ARG:NH1	2.47	0.47
1:B:49:ARG:NH1	6:B:942:HOH:O	2.46	0.47
1:A:256:MET:HA	1:A:260:GLU:O	2.15	0.47
1:A:262:TYR:HB3	1:A:285:MET:CE	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:464:ASN:O	1:A:468:LYS:HG3	2.14	0.46
1:B:77:TRP:CE3	1:B:78:LEU:HD23	2.50	0.46
1:A:317:LYS:HD3	1:A:328:LEU:HD11	1.96	0.46
1:B:126:SER:HA	1:B:127:PRO:C	2.39	0.46
1:A:85:SER:O	1:A:89:ILE:HG12	2.16	0.46
1:B:114:ARG:HD3	1:B:365:LEU:O	2.15	0.46
1:B:277:ARG:HG2	1:B:277:ARG:NH1	2.27	0.46
1:B:163:MET:HE3	1:B:171:LEU:HD11	1.97	0.46
1:A:127:PRO:HG2	1:B:544:TYR:CE1	2.51	0.46
1:B:555:GLY:HA2	1:B:558:LEU:HD12	1.97	0.46
1:A:240:ARG:HG3	1:A:271:VAL:CG2	2.47	0.45
1:A:367:PHE:CD1	1:A:542:PRO:HG3	2.50	0.45
1:A:203:GLN:HA	4:A:601:HEM:HBC2	1.98	0.45
1:A:320:HIS:HB3	1:A:323:TRP:CD1	2.51	0.45
1:B:91:PHE:CZ	1:B:95:HIS:CD2	3.03	0.45
1:B:198:PHE:CD1	1:B:198:PHE:C	2.94	0.45
1:B:64:TYR:CE2	1:B:72:PRO:HB3	2.52	0.45
1:B:115:LEU:HD23	3:B:2802:BOG:H4'2	1.98	0.45
1:B:367:PHE:CD1	1:B:542:PRO:HG3	2.51	0.45
1:A:163:MET:HE3	1:A:171:LEU:HD11	1.98	0.45
1:A:204:HIS:CD2	1:A:292:PHE:CE2	3.05	0.45
1:B:35:PRO:HB2	1:B:55:TYR:CD2	2.52	0.45
1:A:344:VAL:HA	1:A:348:TYR:HB3	1.99	0.45
1:B:180:ARG:HH11	1:B:180:ARG:HB3	1.82	0.45
1:A:208:GLN:HE21	1:A:230:LEU:H	1.65	0.45
1:A:42:GLN:HG3	1:A:70:THR:HG22	1.99	0.44
1:A:198:PHE:CD1	1:A:198:PHE:C	2.95	0.44
1:A:254:TYR:HB3	1:A:306:LEU:HD21	1.99	0.44
1:B:203:GLN:HA	4:B:601:HEM:HBC2	1.99	0.44
1:B:256:MET:HA	1:B:260:GLU:O	2.17	0.44
1:B:391:MET:CG	4:B:601:HEM:HAB	2.41	0.44
1:B:475:TYR:CE2	1:B:481:LEU:HD12	2.52	0.44
1:B:198:PHE:HZ	1:B:352:LEU:HD13	1.82	0.44
1:B:413:MET:HE3	1:B:413:MET:HB2	1.80	0.44
1:A:273:MET:HE2	1:A:285:MET:O	2.17	0.44
1:B:273:MET:HE2	1:B:285:MET:O	2.18	0.44
1:A:447:VAL:O	1:A:451:VAL:HG23	2.18	0.44
1:B:179:ARG:HB3	1:B:179:ARG:HH11	1.83	0.44
1:B:427:SER:HB3	1:B:577:TYR:CD2	2.52	0.44
1:B:38:TYR:CZ	2:F:2:NAG:H61	2.52	0.43
1:A:391:MET:SD	4:A:601:HEM:HAB	2.58	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:GLY:O	1:A:97:ARG:C	2.61	0.43
1:A:292:PHE:O	1:A:299:MET:HE2	2.18	0.43
1:A:413:MET:HE3	1:A:413:MET:HB2	1.81	0.43
1:B:150:ARG:NH2	1:B:458:LEU:O	2.48	0.43
1:A:183:LEU:HD12	1:A:184:ARG:H	1.84	0.43
1:B:206:THR:HG21	1:B:385:TYR:CE1	2.54	0.43
1:B:208:GLN:HE21	1:B:230:LEU:H	1.64	0.43
1:A:280:PRO:HA	1:A:281:PRO:HD3	1.78	0.43
1:A:295:LEU:CD2	1:A:408:LEU:HD23	2.49	0.43
1:A:581:HIS:ND1	1:A:582:VAL:O	2.49	0.43
1:B:344:VAL:HA	1:B:348:TYR:HB3	2.00	0.43
1:A:88:PHE:O	1:A:92:LEU:HG	2.18	0.42
1:A:35:PRO:HB2	1:A:55:TYR:CD2	2.55	0.42
1:A:252:LEU:HD12	1:A:252:LEU:HA	1.91	0.42
1:A:544:TYR:CE1	1:B:127:PRO:HG2	2.54	0.42
1:B:280:PRO:HA	1:B:281:PRO:HD3	1.87	0.42
1:A:46:ILE:HB	1:A:58:ASP:HB3	2.01	0.42
1:B:90:HIS:CE1	6:B:1207:HOH:O	2.72	0.42
1:B:180:ARG:HB3	1:B:180:ARG:NH1	2.34	0.42
1:B:204:HIS:CD2	1:B:292:PHE:CE2	3.08	0.42
1:B:46:ILE:HB	1:B:58:ASP:HB3	2.02	0.42
1:B:54:ARG:HB2	1:B:54:ARG:HH11	1.85	0.42
1:B:320:HIS:HB3	1:B:323:TRP:CD1	2.55	0.42
1:A:501:LEU:HD21	1:A:506:GLY:HA2	2.02	0.42
2:C:1:NAG:H61	2:C:2:NAG:H2	2.01	0.42
1:B:254:TYR:HB3	1:B:306:LEU:HD21	2.01	0.42
1:A:115:LEU:O	1:A:119:VAL:HB	2.20	0.41
1:A:240:ARG:HG3	1:A:271:VAL:HG22	2.02	0.41
1:A:64:TYR:CE1	1:A:76:THR:HG21	2.55	0.41
1:B:124:ILE:HA	1:B:125:PRO:HD2	1.96	0.41
1:A:91:PHE:CZ	1:A:95:HIS:CD2	3.08	0.41
1:A:183:LEU:HD12	1:A:184:ARG:N	2.36	0.41
1:A:424:ASP:O	1:A:428:ARG:HG3	2.20	0.41
1:B:201:PHE:HD2	1:B:301:TYR:CE1	2.39	0.41
1:B:96:GLY:O	1:B:97:ARG:C	2.64	0.41
1:B:107:PHE:CD1	1:B:107:PHE:C	2.98	0.41
1:B:257:LEU:HB2	1:B:262:TYR:CD2	2.56	0.41
1:A:188:ILE:HD13	1:A:433:ARG:HH21	1.86	0.41
1:B:240:ARG:HG3	1:B:271:VAL:HG22	2.02	0.41
1:B:295:LEU:HD21	1:B:408:LEU:CD2	2.51	0.41
1:A:385:TYR:OH	5:A:1650:FLP:H	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:ARG:HG3	1:B:271:VAL:HG21	2.02	0.41
1:B:292:PHE:O	1:B:299:MET:HE2	2.20	0.41
1:A:201:PHE:HD2	1:A:301:TYR:CE1	2.39	0.40
1:A:206:THR:HG21	1:A:385:TYR:CE1	2.56	0.40
1:A:383:GLN:OE1	1:A:455:SER:HB2	2.21	0.40
1:B:108:ILE:O	1:B:108:ILE:HG22	2.21	0.40
1:B:216:MET:HE1	6:B:980:HOH:O	2.20	0.40
1:B:130:TYR:HB2	1:B:150:ARG:HG3	2.03	0.40
1:B:226:HIS:ND1	1:B:376:ARG:HD2	2.36	0.40
1:B:342:LYS:HD3	1:B:562:ALA:HB3	2.02	0.40
1:B:191:PRO:HD2	1:B:433:ARG:HG3	2.04	0.40
1:B:424:ASP:O	1:B:428:ARG:HG3	2.21	0.40
1:B:170:GLN:HE21	1:B:170:GLN:HB2	1.75	0.40
1:B:438:ARG:HH21	1:B:487:MET:HG3	1.85	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	550/580 (95%)	523 (95%)	26 (5%)	1 (0%)	43	73
1	B	551/580 (95%)	524 (95%)	26 (5%)	1 (0%)	43	73
All	All	1101/1160 (95%)	1047 (95%)	52 (5%)	2 (0%)	43	73

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	270	PRO
1	B	270	PRO

5.3.2 Protein sidechains ⓘ

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	481/510 (94%)	449 (93%)	32 (7%)	15	43
1	B	482/510 (94%)	449 (93%)	33 (7%)	14	42
All	All	963/1020 (94%)	898 (93%)	65 (7%)	15	43

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	ARG
1	A	54	ARG
1	A	87	SER
1	A	120	ARG
1	A	126	SER
1	A	145	VAL
1	A	165	THR
1	A	170	GLN
1	A	175	GLU
1	A	185	ARG
1	A	232	HIS
1	A	251	LYS
1	A	252	LEU
1	A	271	VAL
1	A	272	LEU
1	A	291	VAL
1	A	296	PRO
1	A	307	ARG
1	A	322	THR
1	A	374	ARG
1	A	376	ARG
1	A	385	TYR
1	A	400	GLN
1	A	405	GLU
1	A	433	ARG
1	A	469	ARG
1	A	479	GLN

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Mol	Chain	Res	Type
1	A	484	GLU
1	A	493	GLU
1	A	518	PHE
1	A	530	SER
1	A	563	THR
1	B	49	ARG
1	B	54	ARG
1	B	70	THR
1	B	87	SER
1	B	120	ARG
1	B	126	SER
1	B	145	VAL
1	B	165	THR
1	B	170	GLN
1	B	175	GLU
1	B	185	ARG
1	B	232	HIS
1	B	251	LYS
1	B	252	LEU
1	B	271	VAL
1	B	272	LEU
1	B	291	VAL
1	B	294	LEU
1	B	296	PRO
1	B	307	ARG
1	B	322	THR
1	B	374	ARG
1	B	376	ARG
1	B	385	TYR
1	B	400	GLN
1	B	405	GLU
1	B	433	ARG
1	B	469	ARG
1	B	479	GLN
1	B	484	GLU
1	B	518	PHE
1	B	530	SER
1	B	563	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	GLN
1	A	90	HIS
1	A	95	HIS
1	A	134	HIS
1	A	170	GLN
1	A	208	GLN
1	A	232	HIS
1	A	241	GLN
1	A	243	GLN
1	A	358	GLN
1	A	400	GLN
1	A	479	GLN
1	A	513	HIS
1	B	56	GLN
1	B	95	HIS
1	B	134	HIS
1	B	170	GLN
1	B	208	GLN
1	B	232	HIS
1	B	241	GLN
1	B	243	GLN
1	B	282	GLN
1	B	375	ASN
1	B	479	GLN
1	B	513	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

12 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	1,2	14,14,15	0.79	0	17,19,21	1.49	2 (11%)
2	NAG	C	2	2	14,14,15	0.91	1 (7%)	17,19,21	0.96	1 (5%)
2	NAG	D	1	1,2	14,14,15	0.49	0	17,19,21	0.95	2 (11%)
2	NAG	D	2	2	14,14,15	0.61	0	17,19,21	0.89	0
2	NAG	E	1	1,2	14,14,15	0.64	0	17,19,21	0.96	2 (11%)
2	NAG	E	2	2	14,14,15	0.71	0	17,19,21	0.74	0
2	NAG	F	1	1,2	14,14,15	0.74	0	17,19,21	1.12	1 (5%)
2	NAG	F	2	2	14,14,15	0.99	1 (7%)	17,19,21	1.51	4 (23%)
2	NAG	G	1	1,2	14,14,15	0.61	0	17,19,21	1.13	2 (11%)
2	NAG	G	2	2	14,14,15	0.76	0	17,19,21	1.02	2 (11%)
2	NAG	H	1	1,2	14,14,15	0.57	0	17,19,21	0.67	0
2	NAG	H	2	2	14,14,15	0.47	0	17,19,21	0.79	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	1,2	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	NAG	D	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	0/6/23/26	0/1/1/1
2	NAG	E	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	NAG	F	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	F	2	2	-	3/6/23/26	0/1/1/1
2	NAG	G	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
2	NAG	H	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	2/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	NAG	C1-C2	2.43	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	2	NAG	C3-C2	2.38	1.57	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	C4-C3-C2	3.80	116.58	111.02
2	F	1	NAG	C2-N2-C7	-3.58	118.10	122.90
2	F	2	NAG	C1-C2-N2	-3.28	105.27	110.43
2	C	1	NAG	O4-C4-C3	-3.24	102.75	110.38
2	F	2	NAG	C3-C4-C5	2.78	115.28	110.23
2	F	2	NAG	C4-C3-C2	2.72	115.00	111.02
2	G	1	NAG	C2-N2-C7	-2.63	119.37	122.90
2	D	1	NAG	C4-C3-C2	-2.41	107.48	111.02
2	C	2	NAG	C4-C3-C2	2.32	114.41	111.02
2	F	2	NAG	C2-N2-C7	2.22	125.87	122.90
2	G	2	NAG	C1-C2-N2	-2.20	106.96	110.43
2	G	1	NAG	C4-C3-C2	-2.10	107.93	111.02
2	E	1	NAG	C4-C3-C2	-2.09	107.95	111.02
2	E	1	NAG	C2-N2-C7	-2.04	120.17	122.90
2	G	2	NAG	C2-N2-C7	-2.03	120.18	122.90
2	D	1	NAG	C2-N2-C7	-2.02	120.19	122.90

There are no chirality outliers.

All (12) torsion outliers are listed below:

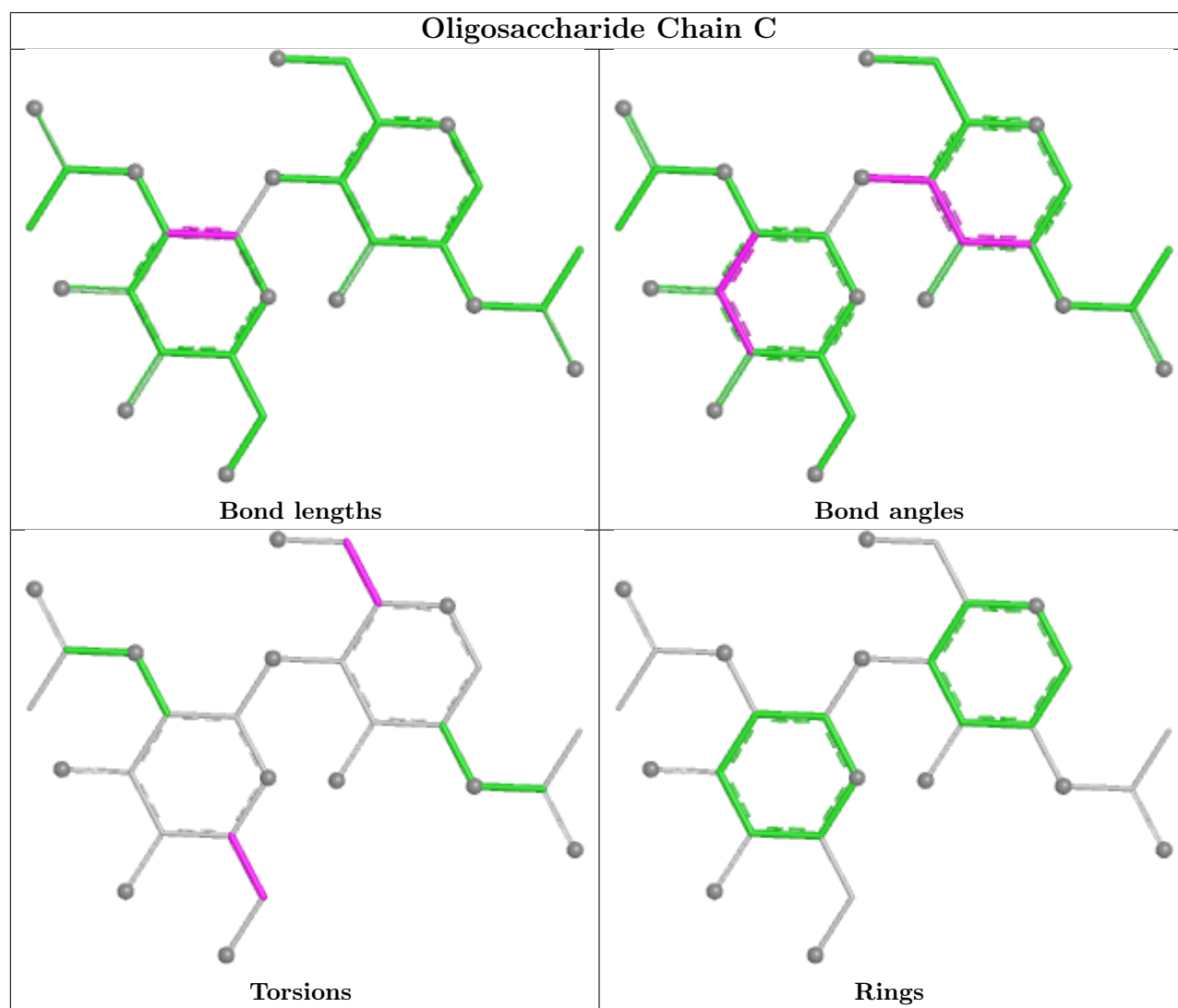
Mol	Chain	Res	Type	Atoms
2	F	2	NAG	C3-C2-N2-C7
2	H	2	NAG	O5-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6
2	H	2	NAG	C4-C5-C6-O6
2	E	2	NAG	O5-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6
2	F	2	NAG	C4-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
2	F	1	NAG	C4-C5-C6-O6
2	F	2	NAG	O5-C5-C6-O6

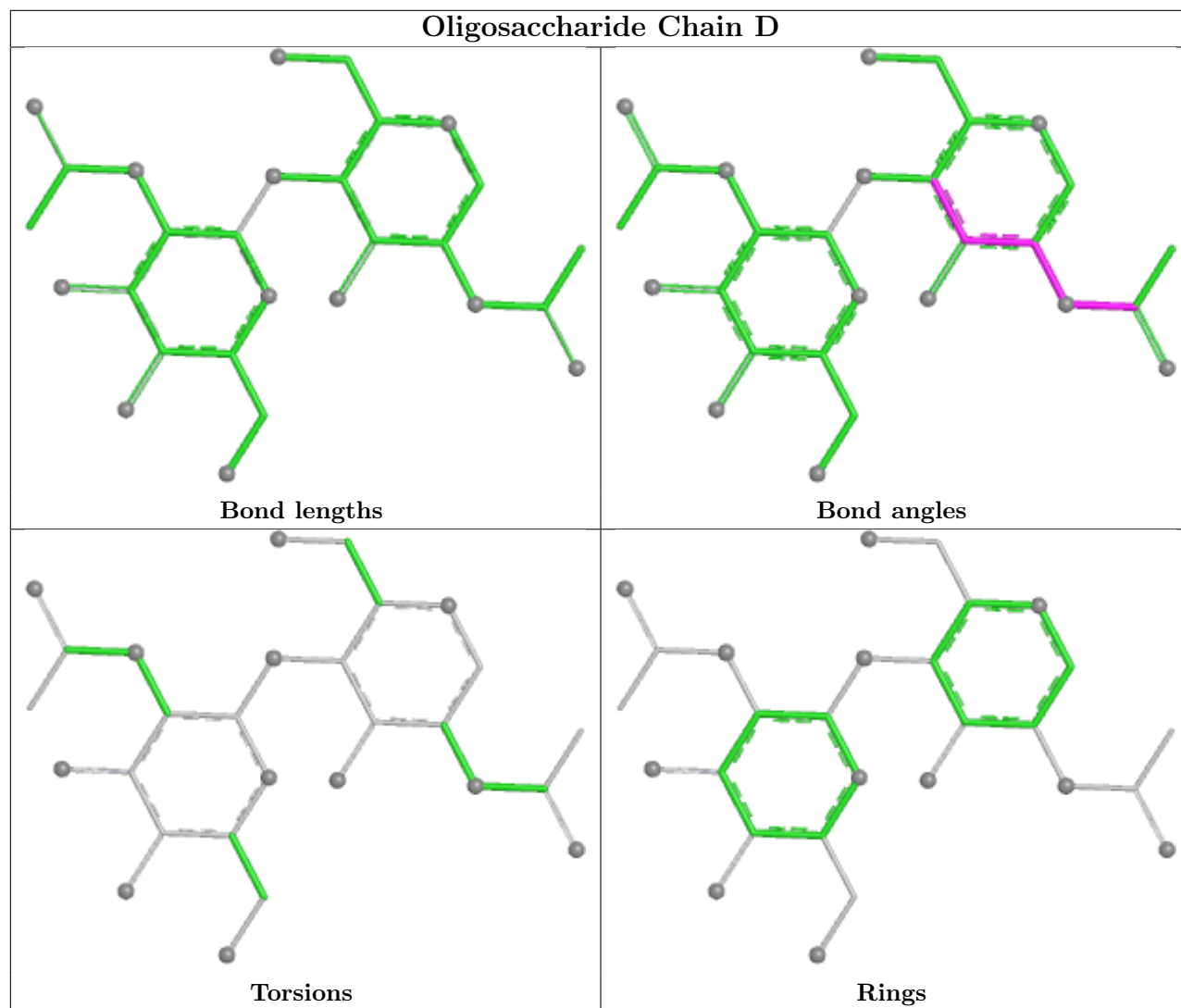
There are no ring outliers.

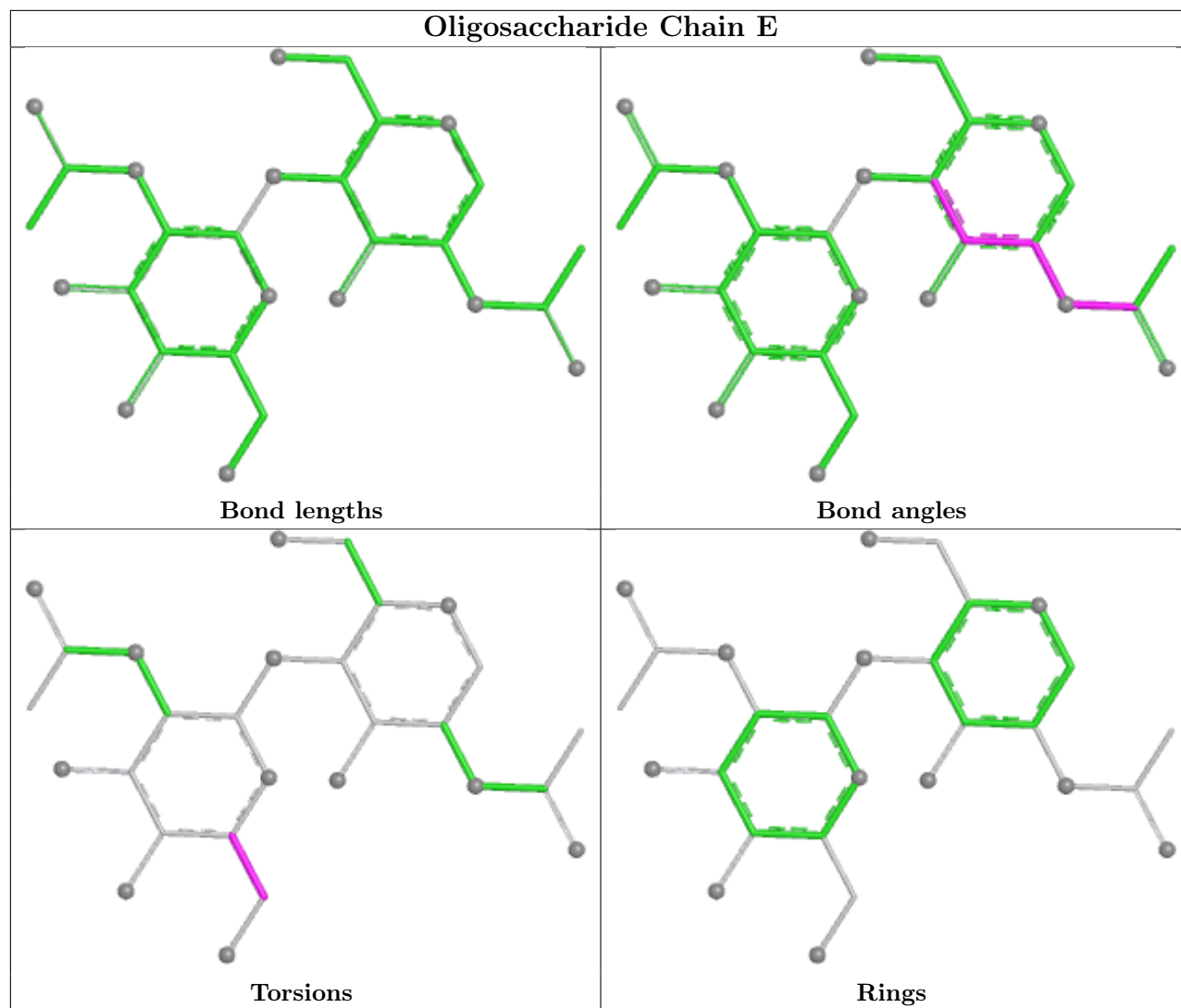
4 monomers are involved in 4 short contacts:

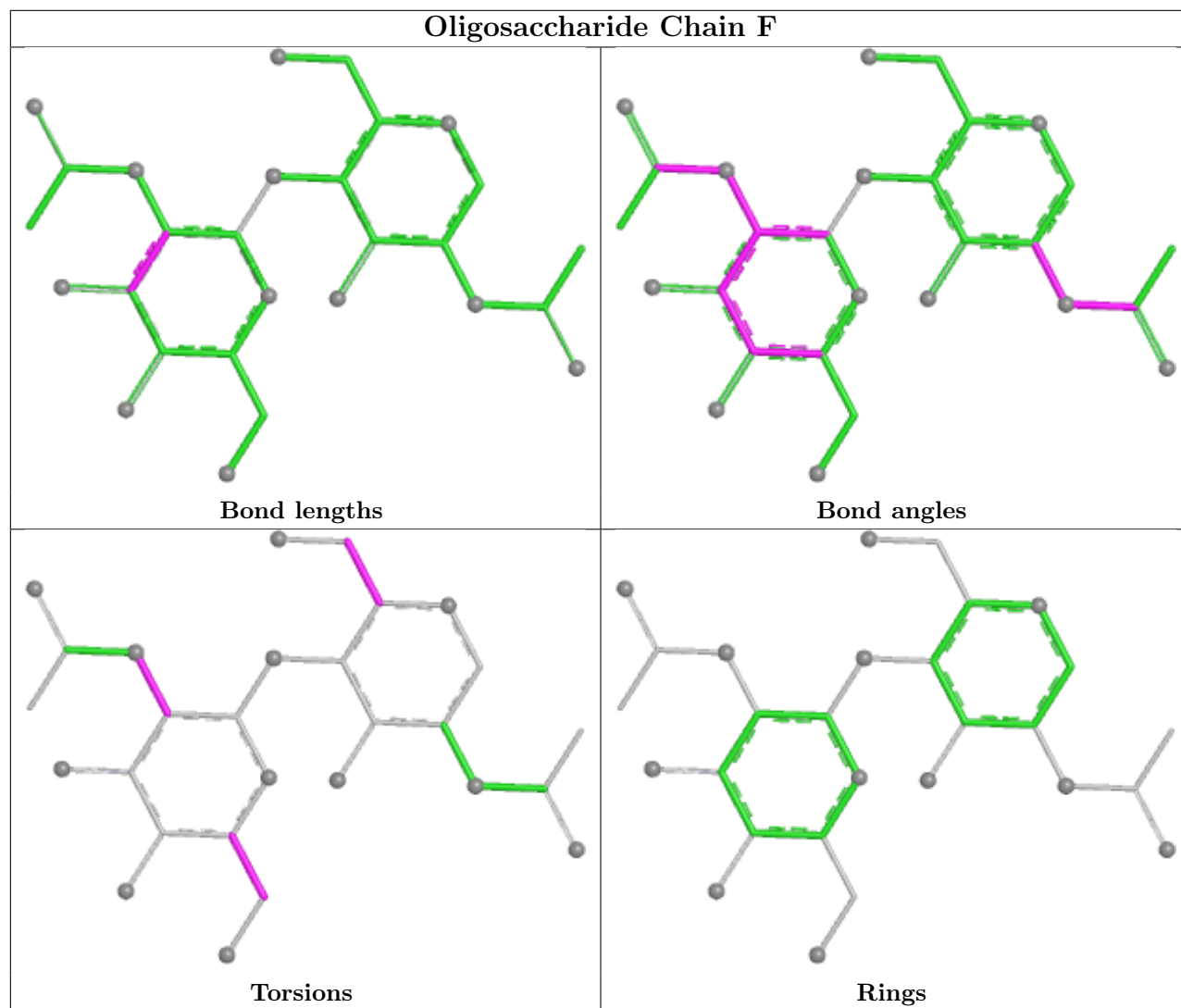
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	NAG	1	0
2	F	1	NAG	2	0
2	C	2	NAG	1	0
2	F	2	NAG	3	0

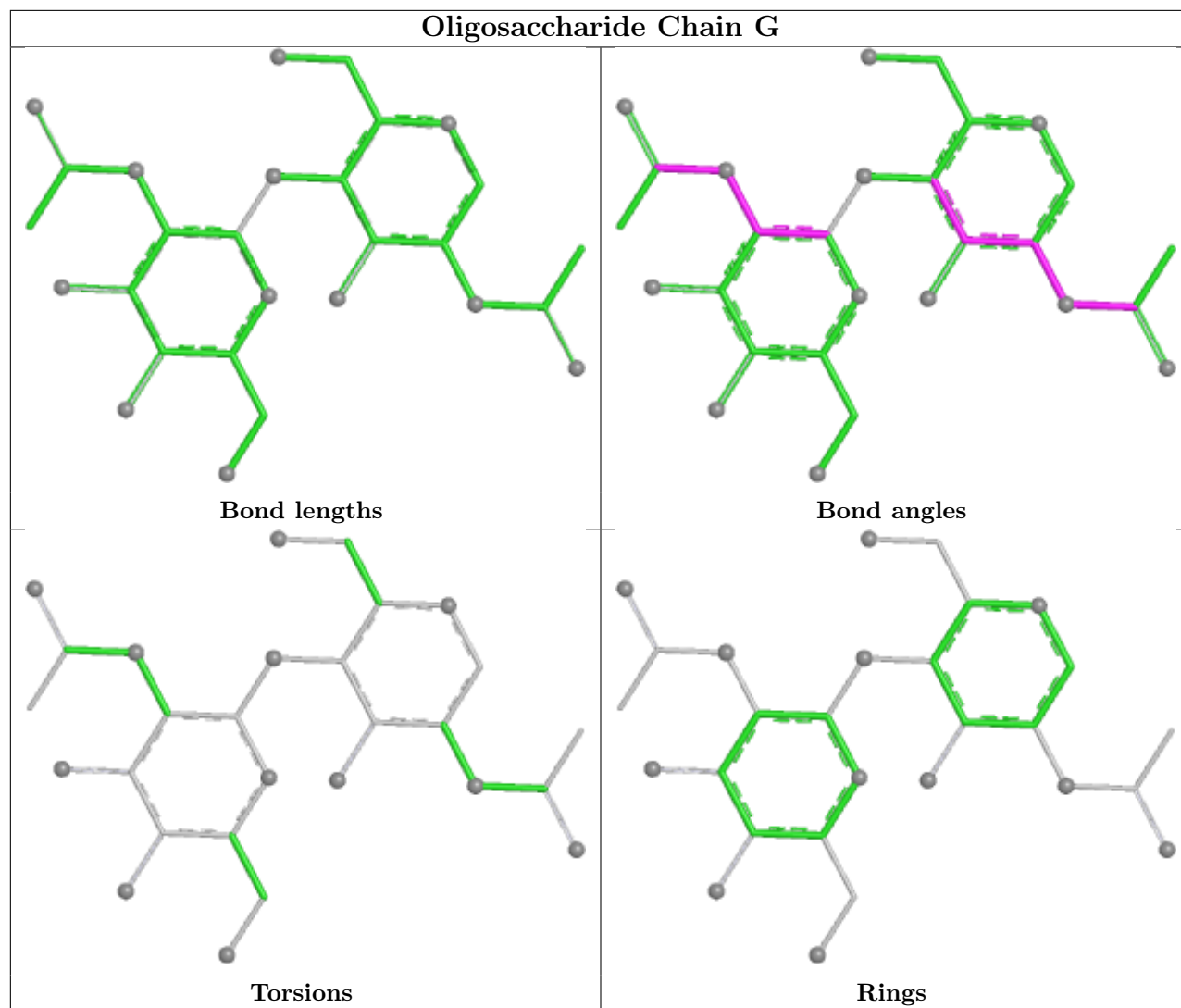
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

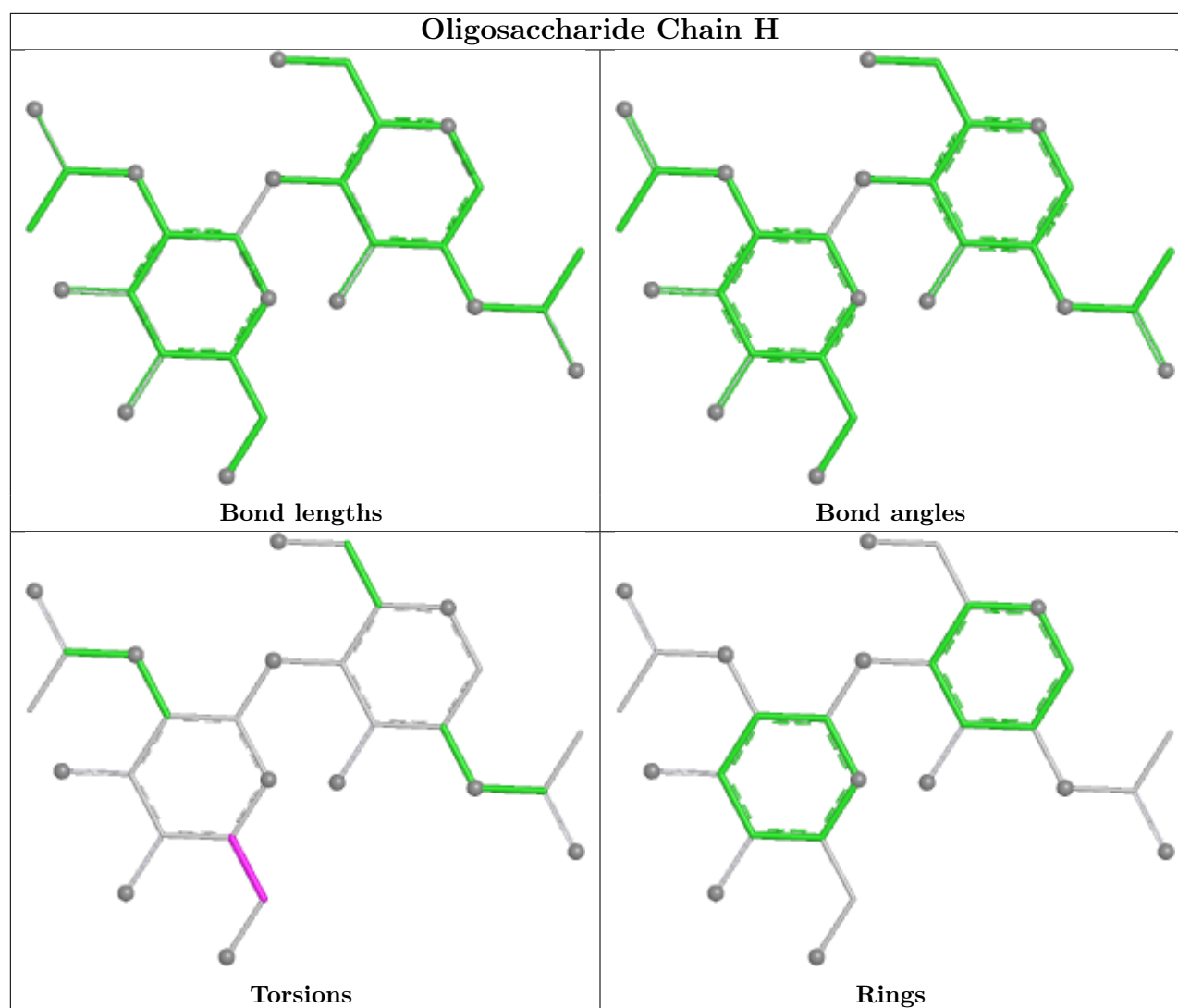












5.6 Ligand geometry [i](#)

9 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	$\# Z > 2$
5	FLP	A	1650	-	19,19,19	1.58	4 (21%)	26,26,26	1.18	2 (7%)
3	BOG	B	2802	-	9,9,20	1.46	2 (22%)	8,8,25	0.80	0
3	BOG	A	1703	-	20,20,20	0.51	0	25,25,25	0.92	1 (4%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BOG	A	1702	-	9,9,20	1.25	2 (22%)	8,8,25	0.97	1 (12%)
5	FLP	B	2650	-	19,19,19	1.87	6 (31%)	26,26,26	1.00	1 (3%)
3	BOG	A	1701	-	20,20,20	0.54	0	25,25,25	1.02	3 (12%)
3	BOG	B	2803	-	9,9,20	0.95	1 (11%)	8,8,25	0.78	0
4	HEM	B	601	1	50,50,50	1.37	7 (14%)	67,82,82	1.41	10 (14%)
4	HEM	A	601	1	50,50,50	1.39	7 (14%)	67,82,82	1.42	7 (10%)

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
5	FLP	A	1650	-	-	0/12/12/12	0/2/2/2
3	BOG	B	2802	-	-	1/7/7/31	-
3	BOG	A	1703	-	-	6/11/31/31	0/1/1/1
3	BOG	A	1702	-	-	2/7/7/31	-
5	FLP	B	2650	-	-	0/12/12/12	0/2/2/2
3	BOG	A	1701	-	-	2/11/31/31	0/1/1/1
3	BOG	B	2803	-	-	0/7/7/31	-
4	HEM	B	601	1	-	5/14/54/54	-
4	HEM	A	601	1	-	5/14/54/54	-

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	2650	FLP	C6-C2	-4.97	1.40	1.49
4	B	601	HEM	CAC-C3C	-3.95	1.36	1.47
4	A	601	HEM	CAC-C3C	-3.84	1.37	1.47
4	A	601	HEM	CAB-C3B	-3.73	1.37	1.47
5	A	1650	FLP	C6-C2	-3.56	1.43	1.49
4	B	601	HEM	CAB-C3B	-3.24	1.38	1.47
3	B	2802	BOG	O1-C1	3.19	1.57	1.42
4	A	601	HEM	C1B-C2B	2.97	1.50	1.44
4	B	601	HEM	C1B-C2B	2.69	1.50	1.44
3	B	2802	BOG	O1-C1'	2.67	1.55	1.40
3	A	1702	BOG	O1-C1'	2.65	1.55	1.40
5	A	1650	FLP	C6-C11	2.56	1.43	1.39
5	B	2650	FLP	C6-C11	2.52	1.43	1.39
4	A	601	HEM	CBB-CAB	2.52	1.42	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	601	HEM	C2A-C3A	-2.44	1.32	1.38
5	B	2650	FLP	C8-C7	2.42	1.42	1.38
4	B	601	HEM	CBB-CAB	2.41	1.42	1.30
5	A	1650	FLP	C4-C3	-2.40	1.34	1.38
5	B	2650	FLP	C4-C3	-2.37	1.34	1.38
3	B	2803	BOG	O1-C1'	2.28	1.53	1.40
3	A	1702	BOG	O1-C1	2.27	1.53	1.42
4	A	601	HEM	C2A-C3A	-2.27	1.33	1.38
4	B	601	HEM	CBC-CAC	2.22	1.41	1.30
5	B	2650	FLP	C-C1	2.20	1.42	1.38
5	B	2650	FLP	C10-C9	-2.13	1.35	1.39
5	A	1650	FLP	C8-C7	2.12	1.42	1.38
4	B	601	HEM	FE-ND	2.05	2.01	1.94
4	A	601	HEM	C4B-NB	2.01	1.42	1.38
4	A	601	HEM	C1C-C2C	-2.01	1.41	1.45

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	601	HEM	C3B-C4B-NB	5.16	113.17	109.47
4	B	601	HEM	C3B-C4B-NB	4.52	112.71	109.47
4	A	601	HEM	CBA-CAA-C2A	3.78	122.98	112.53
4	B	601	HEM	CBA-CAA-C2A	3.51	122.24	112.53
4	B	601	HEM	CAA-C2A-C3A	-3.25	119.78	127.07
4	B	601	HEM	CAA-C2A-C1A	3.08	130.95	124.94
3	A	1703	BOG	C1'-O1-C1	2.99	118.78	113.68
4	A	601	HEM	CAA-C2A-C3A	-2.96	120.43	127.07
5	A	1650	FLP	F-C11-C6	2.90	123.59	119.00
4	A	601	HEM	CAA-C2A-C1A	2.80	130.40	124.94
4	B	601	HEM	C1B-NB-C4B	-2.49	102.26	105.21
3	A	1702	BOG	C1-O1-C1'	2.47	127.79	112.90
5	B	2650	FLP	F-C11-C6	2.44	122.87	119.00
4	B	601	HEM	C4D-ND-C1D	-2.40	102.37	105.21
4	A	601	HEM	C1D-C2D-C3D	-2.38	104.48	106.98
4	A	601	HEM	C4B-C3B-C2B	-2.35	105.12	107.28
4	A	601	HEM	C1B-NB-C4B	-2.34	102.44	105.21
4	B	601	HEM	C2B-C1B-NB	2.33	112.51	109.84
3	A	1701	BOG	C4-C3-C2	-2.29	106.81	110.83
4	B	601	HEM	C1D-C2D-C3D	-2.23	104.63	106.98
5	A	1650	FLP	C5-C-C1	-2.12	117.62	120.24
4	B	601	HEM	C2D-C1D-ND	2.11	112.34	109.90
3	A	1701	BOG	C6-C5-C4	-2.08	107.92	113.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1701	BOG	C8'-C7'-C6'	-2.08	99.33	113.36
4	B	601	HEM	CHB-C1B-NB	-2.03	121.86	124.37

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1701	BOG	O5-C1-O1-C1'
3	A	1703	BOG	C2-C1-O1-C1'
3	A	1703	BOG	O5-C1-O1-C1'
4	A	601	HEM	C2C-C3C-CAC-CBC
4	A	601	HEM	C4C-C3C-CAC-CBC
4	B	601	HEM	C2C-C3C-CAC-CBC
4	B	601	HEM	C4C-C3C-CAC-CBC
3	A	1701	BOG	C2-C1-O1-C1'
3	A	1703	BOG	O1-C1'-C2'-C3'
3	B	2802	BOG	C2'-C3'-C4'-C5'
3	A	1702	BOG	C2'-C3'-C4'-C5'
3	A	1703	BOG	O5-C5-C6-O6
3	A	1703	BOG	C2'-C1'-O1-C1
4	B	601	HEM	C3A-C2A-CAA-CBA
3	A	1702	BOG	C1'-C2'-C3'-C4'
4	A	601	HEM	C3A-C2A-CAA-CBA
3	A	1703	BOG	C5'-C6'-C7'-C8'
4	B	601	HEM	CAA-CBA-CGA-O2A
4	A	601	HEM	CAA-CBA-CGA-O2A
4	A	601	HEM	CAA-CBA-CGA-O1A
4	B	601	HEM	CAA-CBA-CGA-O1A

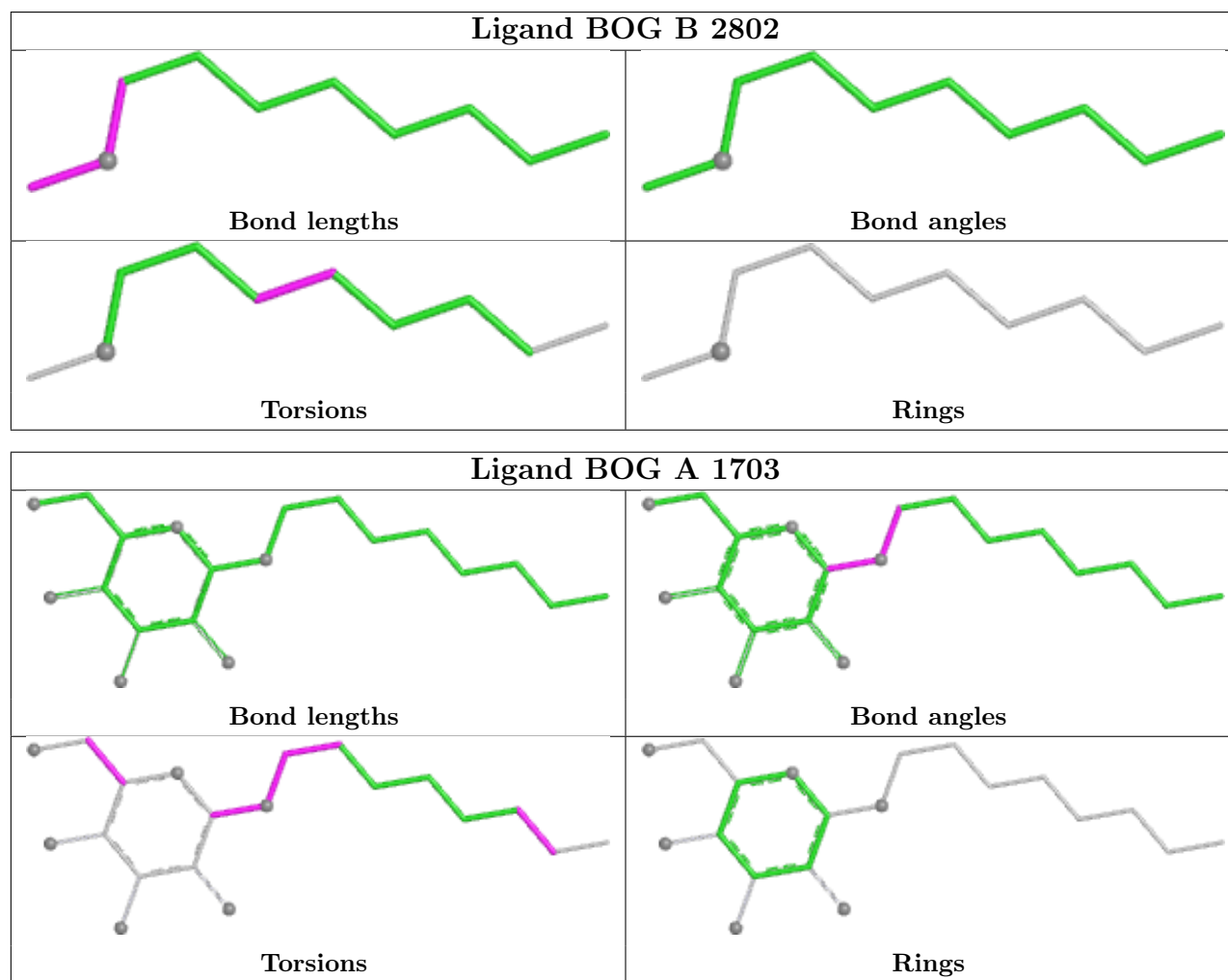
There are no ring outliers.

6 monomers are involved in 14 short contacts:

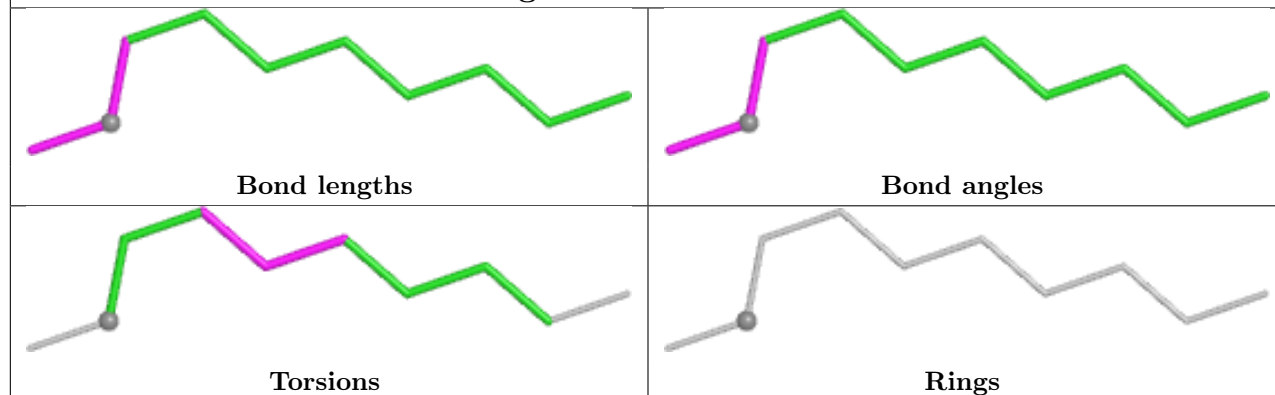
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1650	FLP	1	0
3	B	2802	BOG	1	0
3	A	1702	BOG	2	0
3	A	1701	BOG	1	0
4	B	601	HEM	4	0
4	A	601	HEM	5	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

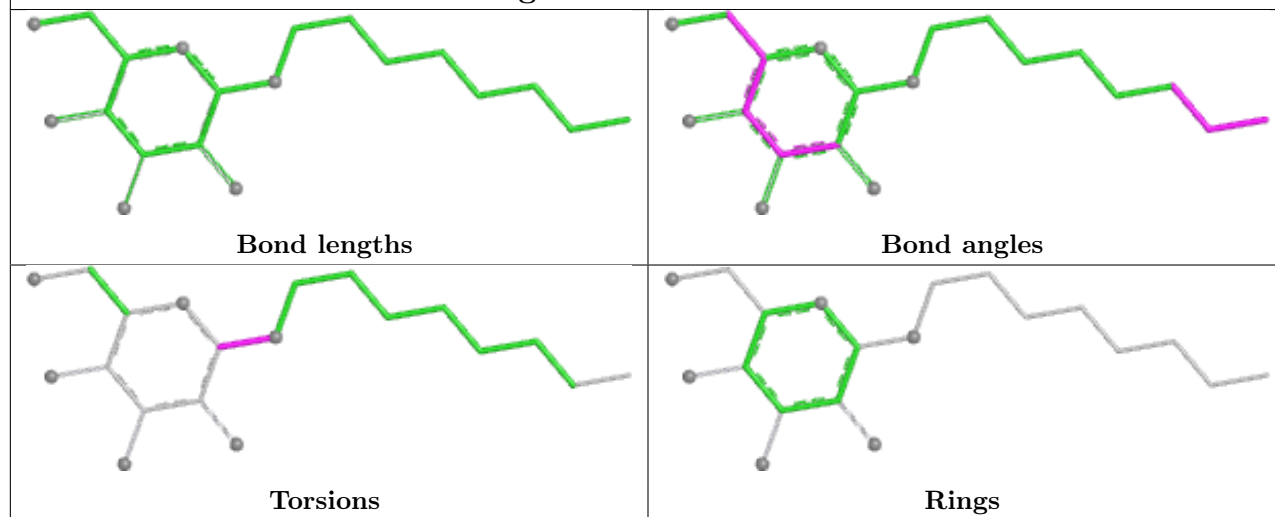
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 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.



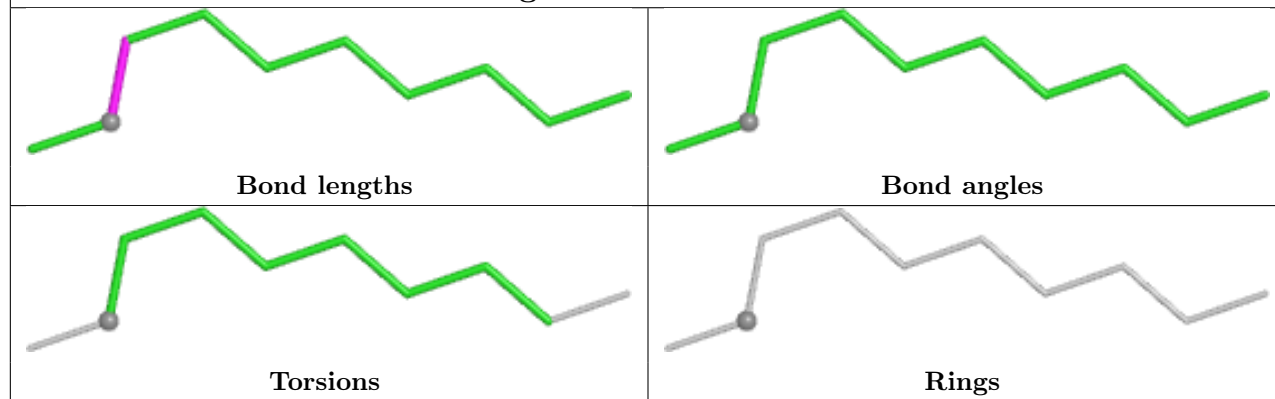
Ligand BOG A 1702

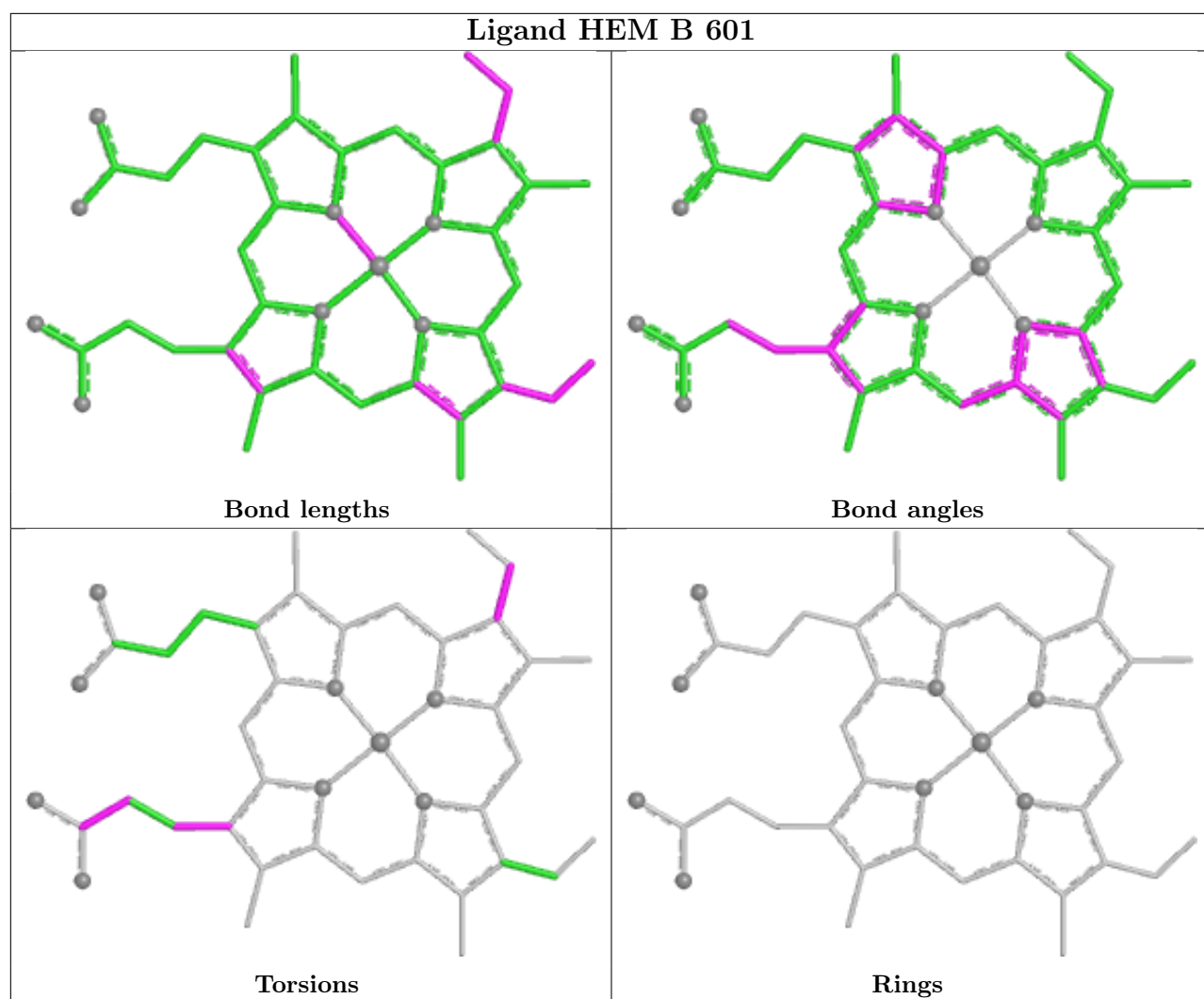


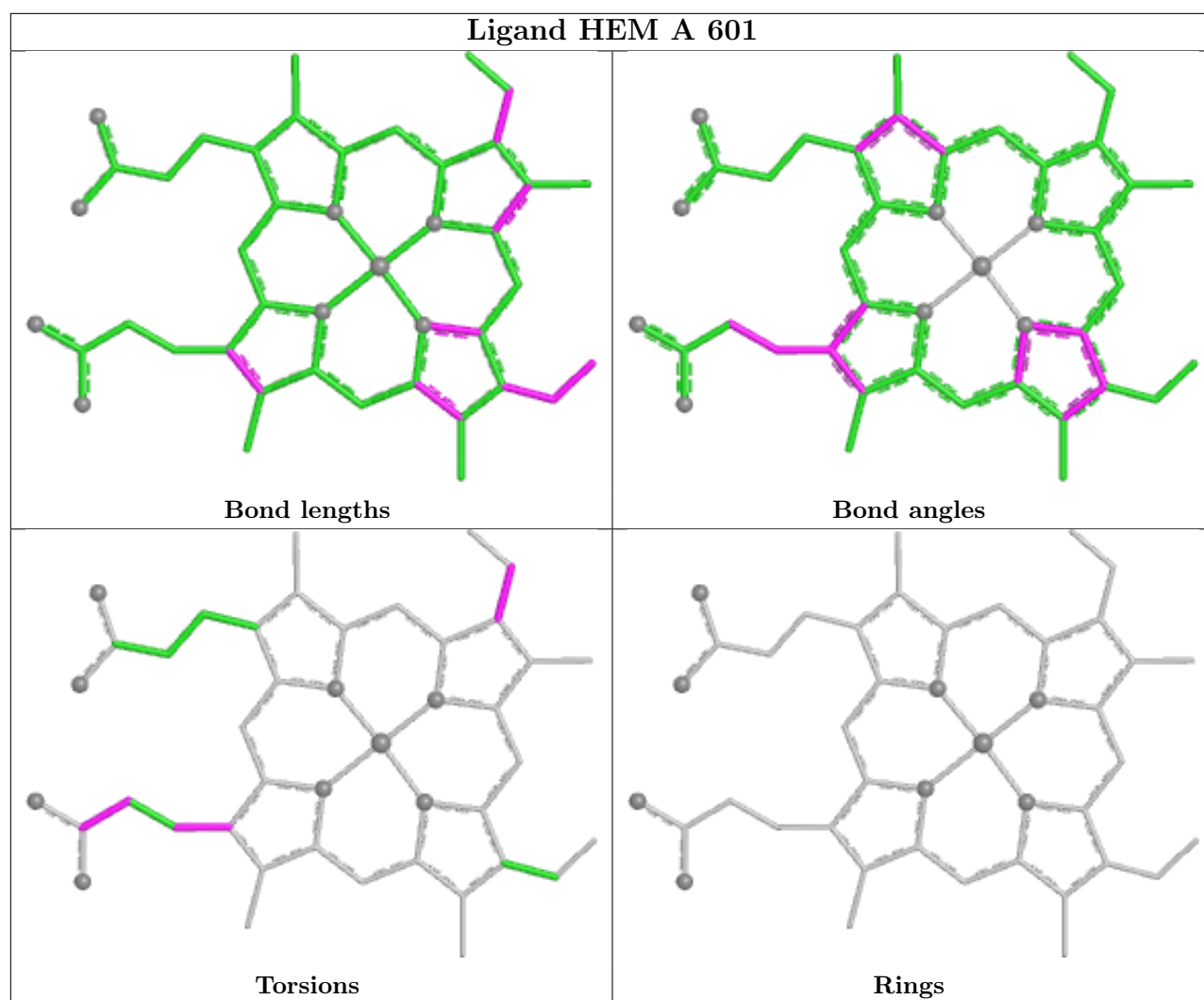
Ligand BOG A 1701



Ligand BOG B 2803







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.