



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:19 PM UTC

PDB ID : 1E8V / pdb_00001e8v
Title : Structure of the multifunctional paramyxovirus hemagglutinin-neuraminidase
Authors : Crennell, S.; Takimoto, T.; Portner, A.; Taylor, G.
Deposited on : 2000-10-01
Resolution : 2.00 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

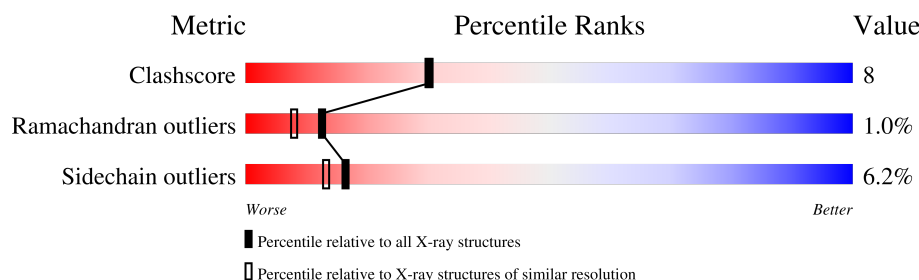
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)

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	454	
1	B	454	
2	C	2	

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
2	NAG	C	1	X	-	-	-
5	NAG	A	1572	X	-	-	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7245 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 HEMAGGLUTININ-NEURAMINIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	447	Total	C	N	O	S	0	0	0
			3449	2174	588	668	19			
1	B	446	Total	C	N	O	S	0	0	0
			3445	2172	587	667	19			

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

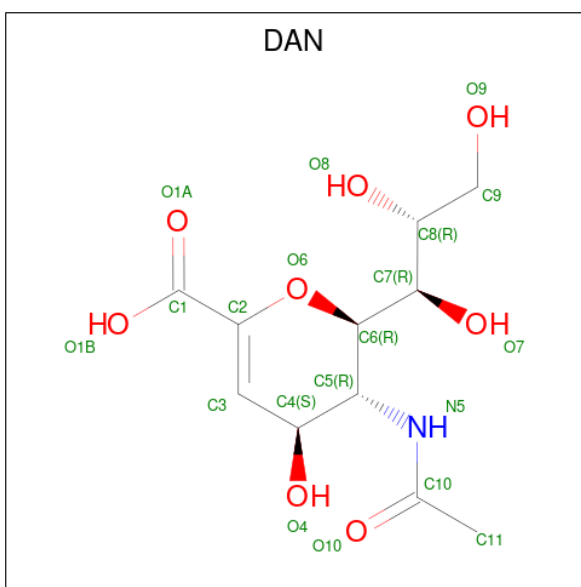


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

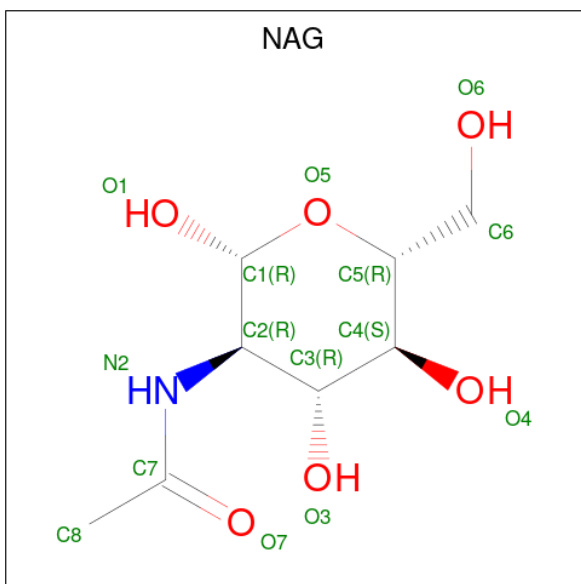
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	B	1	Total	Ca	0	0
			1	1		

- Molecule 4 is 2-DEOXY-2,3-DEHYDRO-N-ACETYL-NEURAMINIC ACID (CCD ID: DAN) (formula: C₁₁H₁₇NO₈).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			20	11	1	8		
4	B	1	Total	C	N	O	0	0
			20	11	1	8		

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is water.

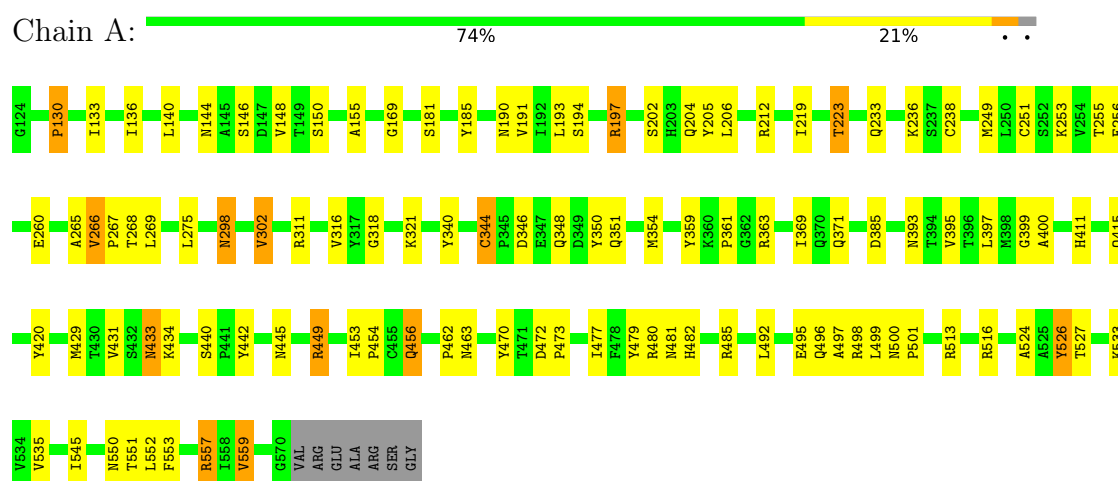
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	129	Total	O	0	0
			129	129		
6	B	110	Total	O	0	0
			110	110		

3 Residue-property plots

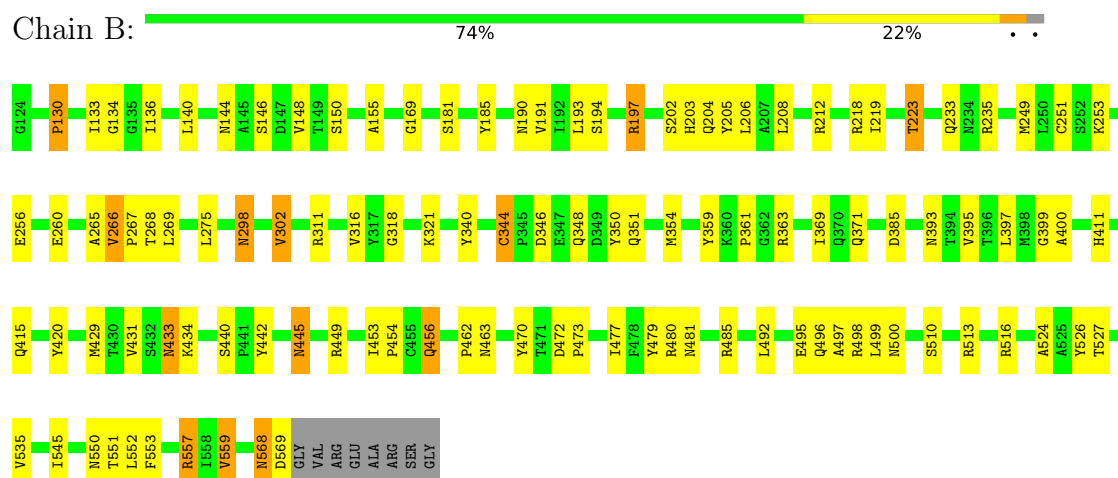
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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



• Molecule 1: HEMAGGLUTININ-NEURAMINIDASE



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





4 Data and refinement statistics

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

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	137.46Å 137.46Å 116.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 2.00	Depositor
% Data completeness (in resolution range)	97.0 (6.00-2.00)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.232 , 0.267	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7245	wwPDB-VP
Average B, all atoms (Å ²)	44.0	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: CA, DAN, NAG

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	0.46	0/3534	1.02	22/4807 (0.5%)
1	B	0.45	0/3530	1.01	22/4802 (0.5%)
All	All	0.45	0/7064	1.01	44/9609 (0.5%)

There are no bond length outliers.

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	449	ARG	NE-CZ-NH2	-8.29	111.73	119.20
1	A	449	ARG	CD-NE-CZ	7.51	134.91	124.40
1	A	193	LEU	N-CA-C	7.34	118.93	111.07
1	B	193	LEU	N-CA-C	7.26	118.83	111.07
1	B	445	ASN	N-CA-C	6.79	120.44	111.75
1	A	449	ARG	NE-CZ-NH1	6.58	128.08	121.50
1	A	440	SER	CA-C-N	6.51	126.49	119.78
1	A	440	SER	C-N-CA	6.51	126.49	119.78
1	B	140	LEU	N-CA-C	6.40	118.34	111.36
1	A	140	LEU	N-CA-C	6.30	118.23	111.36
1	B	440	SER	CA-C-N	6.06	126.02	119.78
1	B	440	SER	C-N-CA	6.06	126.02	119.78
1	A	472	ASP	N-CA-C	6.03	118.44	110.36
1	A	559	VAL	N-CA-C	6.01	114.45	107.89
1	B	559	VAL	N-CA-C	5.97	114.40	107.89
1	B	194	SER	N-CA-C	5.94	119.71	112.23
1	B	449	ARG	NE-CZ-NH2	-5.86	113.92	119.20
1	A	146	SER	N-CA-C	5.82	117.80	110.24
1	B	495	GLU	N-CA-C	5.77	117.38	111.14
1	B	155	ALA	N-CA-C	-5.77	102.36	110.50
1	B	449	ARG	CD-NE-CZ	5.75	132.45	124.40
1	B	442	TYR	N-CA-C	-5.75	100.30	109.96

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	442	TYR	N-CA-C	-5.74	100.31	109.96
1	A	155	ALA	N-CA-C	-5.71	102.45	110.50
1	A	513	ARG	N-CA-C	-5.69	102.47	110.50
1	B	146	SER	N-CA-C	5.63	117.56	110.24
1	B	513	ARG	N-CA-C	-5.52	102.71	110.50
1	A	553	PHE	N-CA-C	5.46	117.56	107.73
1	B	553	PHE	N-CA-C	5.40	117.44	107.73
1	A	495	GLU	N-CA-C	5.39	116.96	111.14
1	A	150	SER	N-CA-C	-5.34	106.27	112.89
1	B	472	ASP	N-CA-C	5.26	117.41	110.36
1	B	150	SER	N-CA-C	-5.25	106.38	112.89
1	B	205	TYR	N-CA-C	-5.25	101.16	109.76
1	A	526	TYR	N-CA-C	5.24	117.05	110.24
1	A	205	TYR	N-CA-C	-5.20	101.24	109.76
1	A	194	SER	N-CA-C	5.19	119.69	112.90
1	B	526	TYR	N-CA-C	5.18	116.98	110.24
1	B	344	CYS	CA-C-N	5.16	124.66	119.19
1	B	344	CYS	C-N-CA	5.16	124.66	119.19
1	B	510	SER	N-CA-C	5.09	117.02	109.69
1	A	238	CYS	N-CA-C	5.06	117.77	110.28
1	A	344	CYS	CA-C-N	5.05	124.54	119.19
1	A	344	CYS	C-N-CA	5.05	124.54	119.19

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	3449	0	3342	55	0
1	B	3445	0	3339	59	0
2	C	28	0	25	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	20	0	16	0	0
4	B	20	0	16	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	28	0	26	2	0
5	B	14	0	13	0	0
6	A	129	0	0	5	0
6	B	110	0	0	4	0
All	All	7245	0	6777	116	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 8.

All (116) 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:223:THR:HG21	6:A:2032:HOH:O	1.69	0.91
1:B:223:THR:HG21	6:B:2034:HOH:O	1.89	0.72
1:A:197:ARG:HD3	1:A:197:ARG:H	1.61	0.66
1:B:197:ARG:HD3	1:B:197:ARG:H	1.63	0.64
1:B:481:ASN:CG	2:C:1:NAG:HN2	2.06	0.64
1:A:453:ILE:HG13	1:A:496:GLN:NE2	2.12	0.63
1:B:453:ILE:HG13	1:B:496:GLN:NE2	2.14	0.61
6:A:2047:HOH:O	1:B:223:THR:HG23	1.99	0.61
1:B:185:TYR:OH	1:B:223:THR:HG22	2.01	0.60
1:B:568:ASN:O	1:B:569:ASP:HB2	2.02	0.60
1:B:371:GLN:HE22	1:B:400:ALA:H	1.48	0.59
1:A:453:ILE:HA	1:A:456:GLN:NE2	2.18	0.58
1:A:371:GLN:HE22	1:A:400:ALA:H	1.49	0.58
1:B:190:ASN:HD22	1:B:204:GLN:NE2	2.00	0.58
1:A:550:ASN:HB2	1:A:557:ARG:HG2	1.86	0.58
1:A:185:TYR:OH	1:A:223:THR:HG22	2.03	0.58
1:A:190:ASN:HD22	1:A:204:GLN:NE2	2.02	0.57
5:A:1573:NAG:O7	5:A:1573:NAG:C3	2.52	0.57
1:B:453:ILE:HA	1:B:456:GLN:NE2	2.19	0.56
1:A:233:GLN:NE2	1:A:253:LYS:HD3	2.20	0.56
1:A:371:GLN:NE2	1:A:400:ALA:H	2.04	0.55
1:B:479:TYR:CD1	1:B:485:ARG:HD2	2.41	0.55
1:A:266:VAL:HG12	1:A:267:PRO:HD2	1.90	0.54
1:B:371:GLN:NE2	1:B:400:ALA:H	2.04	0.54
1:A:344:CYS:SG	1:A:351:GLN:NE2	2.81	0.54
1:B:134:GLY:HA2	6:B:2003:HOH:O	2.08	0.54
1:A:223:THR:HG23	6:B:2044:HOH:O	2.07	0.53
1:A:479:TYR:CD1	1:A:485:ARG:HD2	2.43	0.53
1:A:454:PRO:HA	1:A:456:GLN:HE22	1.72	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:169:GLY:HA2	1:B:552:LEU:O	2.08	0.53
1:B:266:VAL:HG12	1:B:267:PRO:HD2	1.89	0.53
1:B:233:GLN:NE2	1:B:253:LYS:HD3	2.24	0.53
1:A:169:GLY:HA2	1:A:552:LEU:O	2.09	0.53
1:B:454:PRO:HA	1:B:456:GLN:HE22	1.73	0.53
1:B:371:GLN:HE21	1:B:399:GLY:HA2	1.75	0.52
1:B:550:ASN:HB2	1:B:557:ARG:HG2	1.91	0.52
1:A:191:VAL:O	1:A:202:SER:HB2	2.11	0.51
1:B:361:PRO:HD2	1:B:369:ILE:HD13	1.93	0.50
1:B:249:MET:HE3	1:B:251:CYS:SG	2.52	0.50
1:B:191:VAL:O	1:B:202:SER:HB2	2.11	0.50
1:B:344:CYS:SG	1:B:351:GLN:NE2	2.84	0.50
1:B:369:ILE:N	1:B:369:ILE:HD12	2.27	0.49
1:B:354:MET:HA	1:B:354:MET:HE2	1.95	0.49
1:A:361:PRO:HD2	1:A:369:ILE:HD13	1.95	0.49
1:B:420:TYR:CG	1:B:462:PRO:HA	2.48	0.49
1:B:498:ARG:HB3	1:B:524:ALA:O	2.13	0.49
1:A:371:GLN:HE21	1:A:399:GLY:HA2	1.78	0.48
1:A:354:MET:HE2	1:A:354:MET:HA	1.95	0.48
1:B:527:THR:HA	1:B:545:ILE:O	2.13	0.48
1:A:420:TYR:CG	1:A:462:PRO:HA	2.48	0.48
1:A:144:ASN:OD1	1:A:480:ARG:HG3	2.13	0.48
1:A:369:ILE:HD12	1:A:369:ILE:N	2.29	0.48
1:A:411:HIS:HB2	1:A:429:MET:O	2.14	0.47
1:B:268:THR:H	1:B:298:ASN:HD21	1.62	0.47
1:B:479:TYR:CE2	1:B:485:ARG:HG2	2.49	0.47
1:B:136:ILE:HG21	1:B:181:SER:O	2.15	0.47
1:A:206:LEU:HD13	1:A:275:LEU:HB2	1.96	0.47
1:A:268:THR:H	1:A:298:ASN:HD21	1.62	0.46
1:B:144:ASN:OD1	1:B:480:ARG:HG3	2.15	0.46
1:B:148:VAL:HG13	1:B:477:ILE:HD12	1.98	0.46
1:A:479:TYR:CE2	1:A:485:ARG:HG2	2.50	0.46
1:A:527:THR:HA	1:A:545:ILE:O	2.15	0.46
1:A:533:LYS:HE2	6:A:2115:HOH:O	2.15	0.46
1:A:318:GLY:O	1:A:369:ILE:HA	2.16	0.46
1:B:456:GLN:NE2	1:B:456:GLN:H	2.14	0.46
1:B:497:ALA:HB3	1:B:499:LEU:HD12	1.97	0.46
1:B:206:LEU:HD13	1:B:275:LEU:HB2	1.97	0.45
1:A:498:ARG:HB3	1:A:524:ALA:O	2.16	0.45
1:B:318:GLY:O	1:B:369:ILE:HA	2.16	0.45
1:A:501:PRO:HD2	6:A:2106:HOH:O	2.16	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:456:GLN:NE2	1:A:456:GLN:H	2.15	0.45
1:A:249:MET:HE3	1:A:251:CYS:SG	2.56	0.45
1:A:497:ALA:HB3	1:A:499:LEU:HD12	1.97	0.45
1:B:551:THR:O	1:B:552:LEU:HB2	2.16	0.45
6:A:2009:HOH:O	1:B:203:HIS:HE1	2.01	0.44
1:B:371:GLN:NE2	1:B:415:GLN:HE21	2.15	0.44
1:A:516:ARG:HG2	1:A:516:ARG:HH11	1.82	0.44
1:A:256:GLU:H	1:A:256:GLU:HG2	1.69	0.43
1:B:218:ARG:NH2	6:B:2041:HOH:O	2.51	0.43
1:B:433:ASN:HB3	1:B:434:LYS:H	1.46	0.43
1:A:371:GLN:NE2	1:A:415:GLN:HE21	2.15	0.43
1:B:208:LEU:HD23	1:B:208:LEU:C	2.44	0.43
1:A:202:SER:OG	1:A:236:LYS:HE3	2.18	0.43
1:B:516:ARG:HG2	1:B:516:ARG:HH11	1.83	0.43
1:A:197:ARG:H	1:A:197:ARG:CD	2.23	0.43
1:B:256:GLU:HB2	1:B:260:GLU:HB2	2.01	0.43
1:A:136:ILE:HG21	1:A:181:SER:O	2.20	0.42
1:A:268:THR:H	1:A:298:ASN:ND2	2.18	0.42
1:A:148:VAL:HG13	1:A:477:ILE:HD12	2.02	0.42
1:A:346:ASP:HB3	1:A:350:TYR:HD2	1.85	0.42
1:A:433:ASN:HB3	1:A:434:LYS:H	1.46	0.42
1:B:411:HIS:HB2	1:B:429:MET:O	2.20	0.42
1:B:130:PRO:HA	1:B:133:ILE:HG12	2.01	0.42
1:B:311:ARG:NH2	1:B:385:ASP:O	2.52	0.42
1:B:190:ASN:HD22	1:B:204:GLN:HE21	1.67	0.42
1:A:311:ARG:NH2	1:A:385:ASP:O	2.52	0.41
1:A:551:THR:O	1:A:552:LEU:HB2	2.20	0.41
1:B:268:THR:H	1:B:298:ASN:ND2	2.19	0.41
1:A:130:PRO:HA	1:A:133:ILE:HG12	2.01	0.41
1:B:235:ARG:HH11	1:B:235:ARG:HG3	1.86	0.41
1:B:256:GLU:H	1:B:256:GLU:HG2	1.72	0.41
1:A:256:GLU:HB2	1:A:260:GLU:HB2	2.02	0.41
1:A:265:ALA:O	1:A:321:LYS:HE2	2.21	0.41
5:A:1573:NAG:O7	5:A:1573:NAG:H3	2.20	0.41
1:B:359:TYR:CE2	1:B:393:ASN:HB3	2.56	0.41
1:B:420:TYR:CD1	1:B:462:PRO:HA	2.56	0.41
1:A:359:TYR:CE2	1:A:393:ASN:HB3	2.56	0.41
1:B:453:ILE:HA	1:B:456:GLN:HE21	1.85	0.40
1:B:481:ASN:CG	2:C:1:NAG:N2	2.75	0.40
1:B:265:ALA:O	1:B:321:LYS:HE2	2.21	0.40
1:A:190:ASN:HD22	1:A:204:GLN:HE21	1.66	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:470:TYR:CD1	1:B:470:TYR:C	2.99	0.40
1:A:470:TYR:CD1	1:A:470:TYR:C	3.00	0.40
1:B:346:ASP:HB3	1:B:350:TYR:HD2	1.85	0.40
1:A:470:TYR:O	1:A:526:TYR:HA	2.22	0.40
1:A:481:ASN:O	1:A:482:HIS:HB2	2.22	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	445/454 (98%)	410 (92%)	31 (7%)	4 (1%)	14	9
1	B	444/454 (98%)	408 (92%)	31 (7%)	5 (1%)	11	7
All	All	889/908 (98%)	818 (92%)	62 (7%)	9 (1%)	12	8

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	363	ARG
1	B	363	ARG
1	B	568	ASN
1	A	302	VAL
1	A	473	PRO
1	B	473	PRO
1	B	302	VAL
1	A	500	ASN
1	B	500	ASN

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	387/392 (99%)	362 (94%)	25 (6%)	15	12
1	B	387/392 (99%)	364 (94%)	23 (6%)	18	14
All	All	774/784 (99%)	726 (94%)	48 (6%)	16	13

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	130	PRO
1	A	197	ARG
1	A	212	ARG
1	A	219	ILE
1	A	223	THR
1	A	255	THR
1	A	266	VAL
1	A	269	LEU
1	A	298	ASN
1	A	302	VAL
1	A	316	VAL
1	A	340	TYR
1	A	348	GLN
1	A	395	VAL
1	A	397	LEU
1	A	431	VAL
1	A	433	ASN
1	A	445	ASN
1	A	449	ARG
1	A	456	GLN
1	A	463	ASN
1	A	492	LEU
1	A	535	VAL
1	A	557	ARG
1	A	559	VAL
1	B	130	PRO
1	B	197	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	212	ARG
1	B	219	ILE
1	B	223	THR
1	B	266	VAL
1	B	269	LEU
1	B	298	ASN
1	B	302	VAL
1	B	316	VAL
1	B	340	TYR
1	B	348	GLN
1	B	395	VAL
1	B	397	LEU
1	B	431	VAL
1	B	433	ASN
1	B	445	ASN
1	B	456	GLN
1	B	463	ASN
1	B	492	LEU
1	B	535	VAL
1	B	557	ARG
1	B	559	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	157	GLN
1	A	204	GLN
1	A	233	GLN
1	A	282	HIS
1	A	298	ASN
1	A	351	GLN
1	A	371	GLN
1	A	456	GLN
1	A	463	ASN
1	A	482	HIS
1	A	496	GLN
1	B	157	GLN
1	B	204	GLN
1	B	233	GLN
1	B	282	HIS
1	B	298	ASN
1	B	351	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	371	GLN
1	B	456	GLN
1	B	463	ASN
1	B	496	GLN

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](#)

2 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	1.58	3 (21%)	17,19,21	0.77	0
2	NAG	C	2	2	14,14,15	0.51	0	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	1,2	1/1/5/7	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	1/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1	NAG	C4-C5	-3.93	1.44	1.53
2	C	1	NAG	C4-C3	-2.70	1.45	1.52
2	C	1	NAG	C1-C2	2.47	1.55	1.52

There are no bond angle outliers.

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	1	NAG	C1

All (1) torsion outliers are listed below:

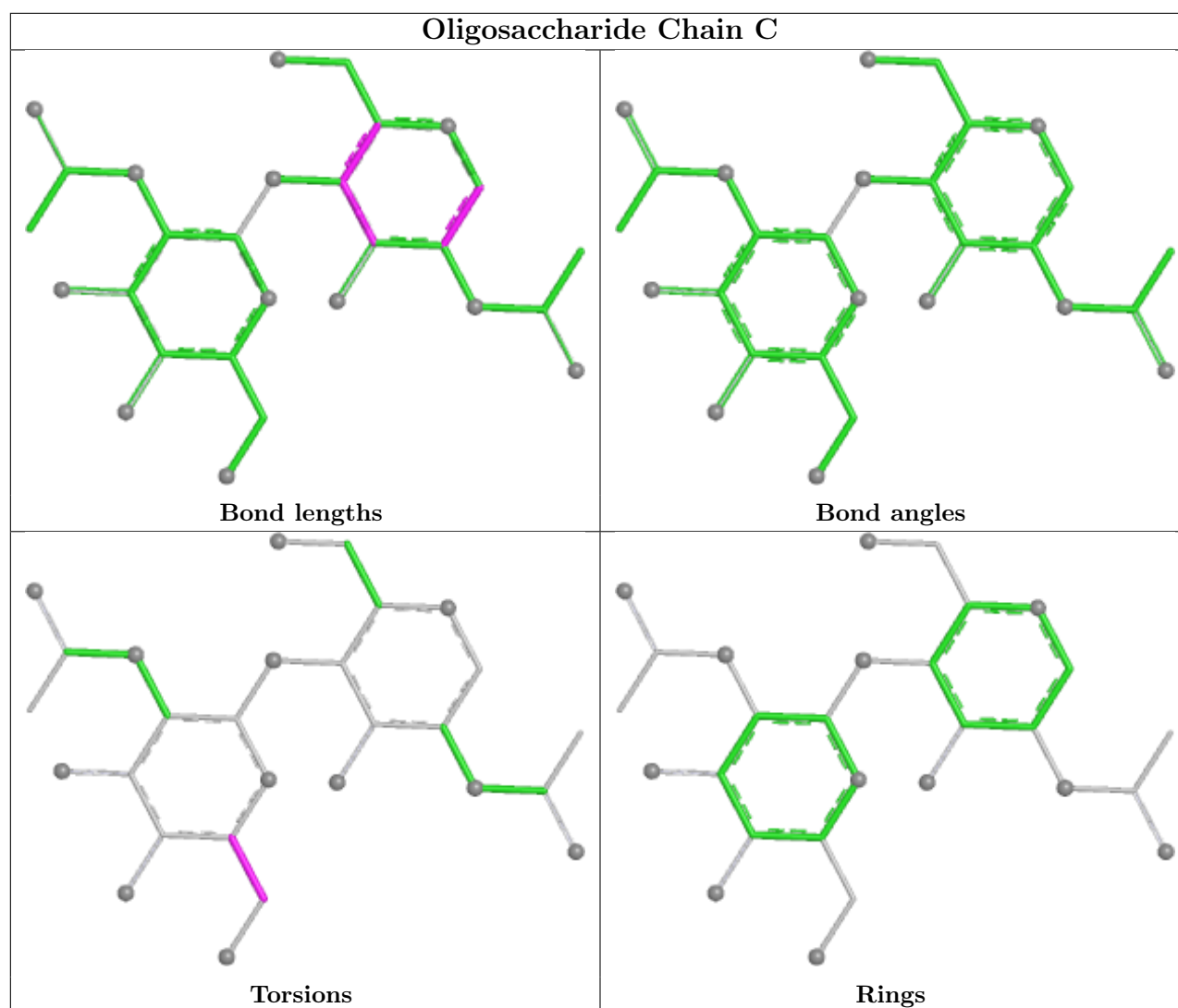
Mol	Chain	Res	Type	Atoms
2	C	2	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	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 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	$\# Z > 2$
5	NAG	B	841	1	14,14,15	1.42	2 (14%)	17,19,21	0.75	0
4	DAN	A	1571	-	20,20,20	2.51	7 (35%)	24,28,28	1.73	5 (20%)
5	NAG	A	1572	1	14,14,15	1.52	2 (14%)	17,19,21	0.91	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	DAN	B	1570	-	20,20,20	1.97	5 (25%)	24,28,28	2.28	5 (20%)
5	NAG	A	1573	1	14,14,15	0.86	0	17,19,21	0.61	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
5	NAG	B	841	1	-	0/6/23/26	0/1/1/1
4	DAN	A	1571	-	-	0/18/34/34	0/1/1/1
5	NAG	A	1572	1	1/1/5/7	0/6/23/26	0/1/1/1
4	DAN	B	1570	-	-	0/18/34/34	0/1/1/1
5	NAG	A	1573	1	-	3/6/23/26	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1571	DAN	C6-C5	5.90	1.62	1.53
4	A	1571	DAN	C3-C2	5.36	1.42	1.33
4	B	1570	DAN	C3-C2	4.88	1.41	1.33
4	A	1571	DAN	C4-C3	4.55	1.56	1.50
5	B	841	NAG	C3-C2	4.42	1.61	1.52
4	B	1570	DAN	C4-C3	4.31	1.55	1.50
5	A	1572	NAG	C3-C2	3.86	1.60	1.52
4	A	1571	DAN	C2-C1	-3.38	1.41	1.48
5	A	1572	NAG	C1-C2	3.38	1.56	1.52
4	B	1570	DAN	C2-C1	-3.25	1.41	1.48
4	A	1571	DAN	C5-N5	2.81	1.50	1.45
4	A	1571	DAN	C4-C5	2.81	1.56	1.53
5	B	841	NAG	C1-C2	2.16	1.55	1.52
4	A	1571	DAN	O4-C4	-2.16	1.39	1.43
4	B	1570	DAN	C7-C6	-2.12	1.50	1.52
4	B	1570	DAN	O1B-C1	-2.03	1.25	1.30

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1570	DAN	O6-C2-C1	6.21	124.39	112.02
4	B	1570	DAN	O6-C2-C3	-5.95	117.97	124.62
4	B	1570	DAN	C4-C5-N5	-4.05	106.59	111.13
4	A	1571	DAN	O6-C2-C1	3.72	119.43	112.02

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1571	DAN	O4-C4-C3	3.53	117.25	109.27
4	A	1571	DAN	C4-C5-N5	-2.92	107.86	111.13
4	A	1571	DAN	C3-C2-C1	-2.82	117.46	123.56
4	B	1570	DAN	C3-C2-C1	-2.74	117.63	123.56
4	A	1571	DAN	O6-C2-C3	-2.61	121.70	124.62
4	B	1570	DAN	C8-C7-C6	-2.16	109.00	113.05

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	A	1572	NAG	C1

All (3) torsion outliers are listed below:

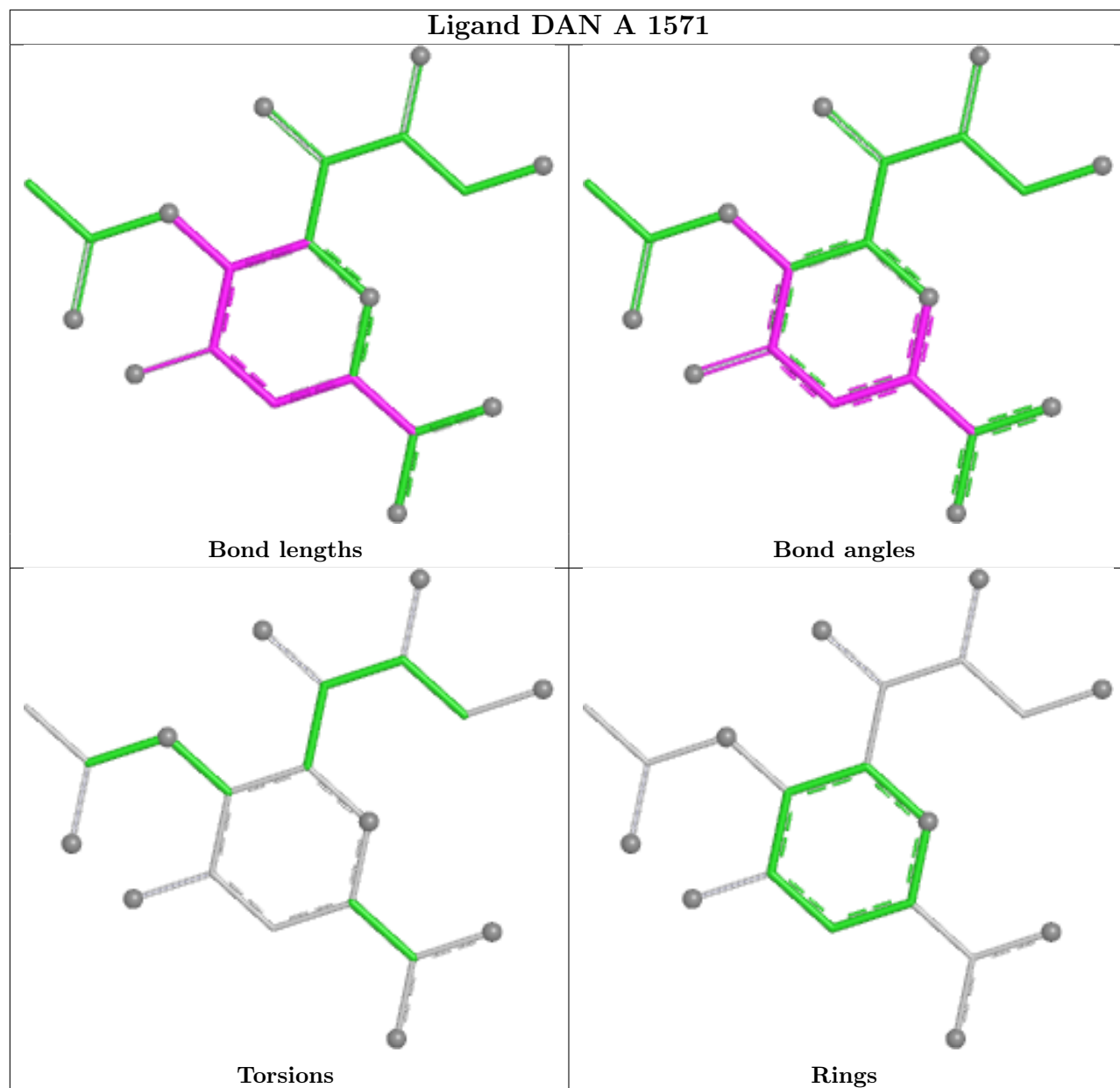
Mol	Chain	Res	Type	Atoms
5	A	1573	NAG	C3-C2-N2-C7
5	A	1573	NAG	O5-C5-C6-O6
5	A	1573	NAG	C4-C5-C6-O6

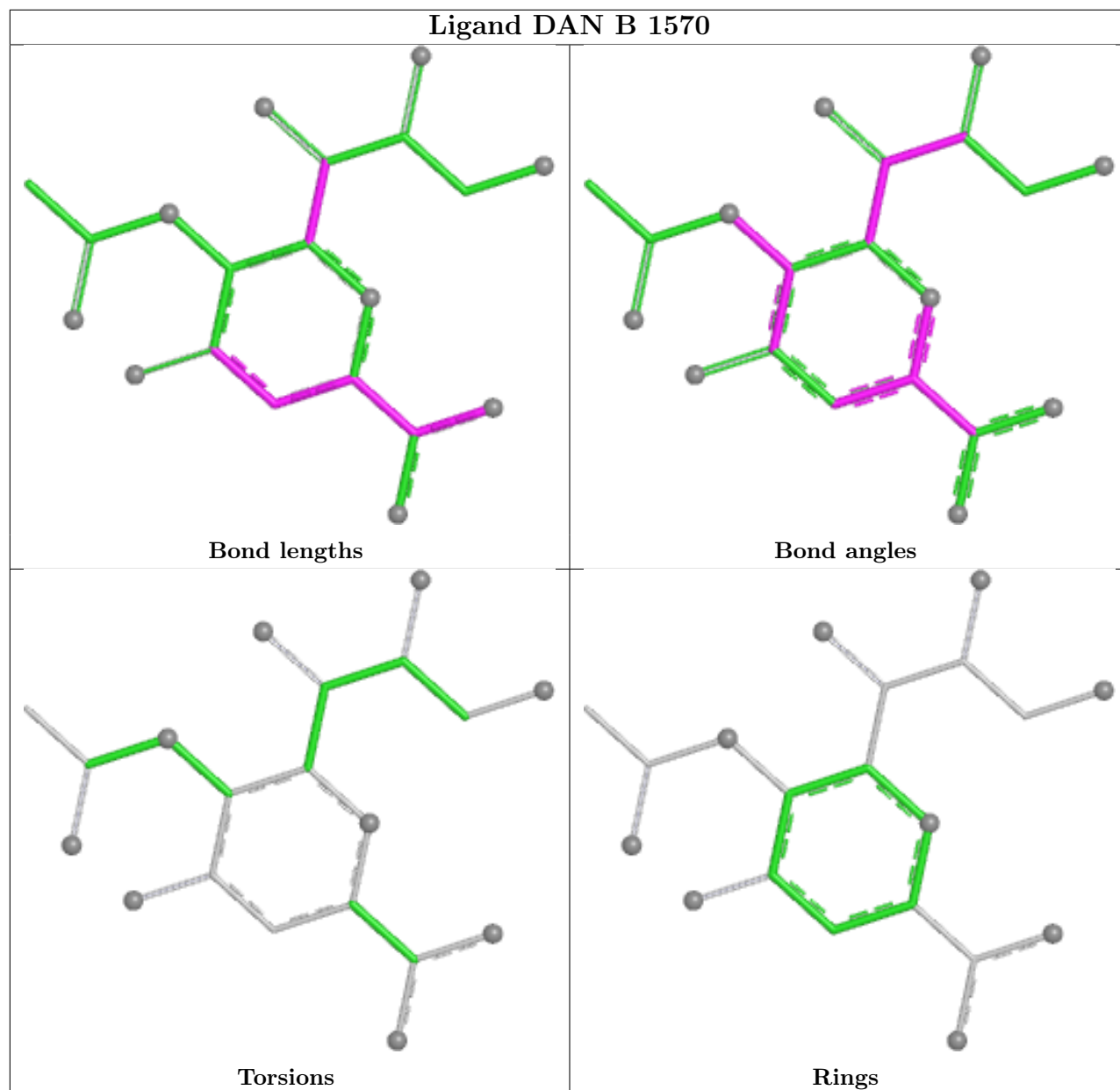
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1573	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 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.