Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 04:05 PM EDT

PDB ID	:	1E8V
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 (i) 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 (1)) were used in the production of this report:

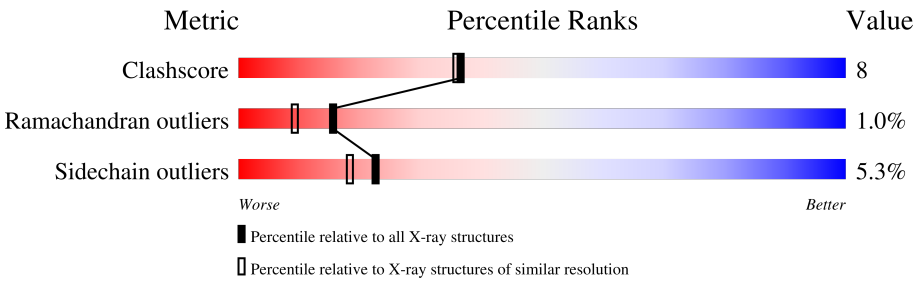
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

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	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	454	79% 18% • •
1	B	454	79% 17% • •
2	C	2	50% 50%

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
			Total	C	N	O	S			
1	A	447	3449	2174	588	668	19	0	0	0
1	B	446	3445	2172	587	667	19	0	0	0

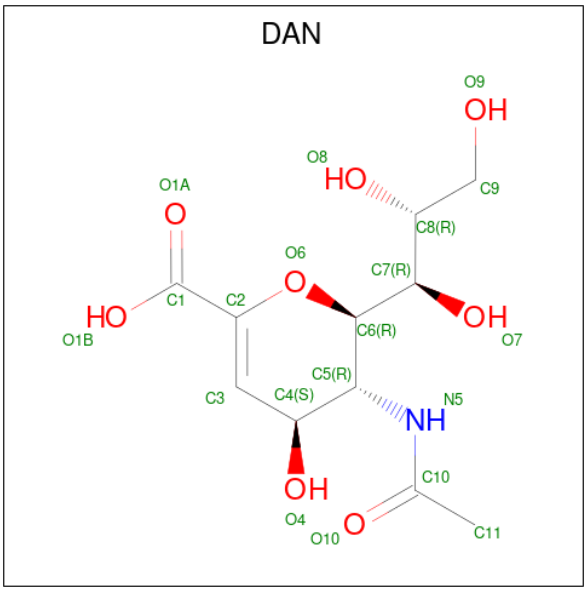
- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

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

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

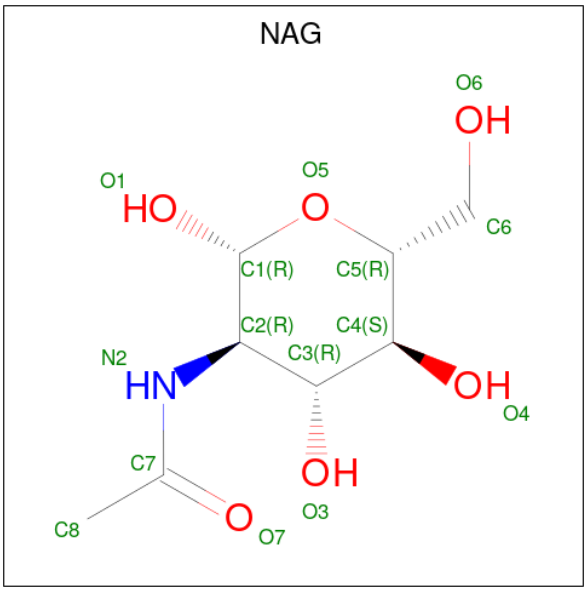
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
3	A	1	1	1	0	0
3	B	1	1	1	0	0

- Molecule 4 is 2-DEOXY-2,3-DEHYDRO-N-ACETYL-NEURAMINIC ACID (three-letter code: DAN) (formula: $C_{11}H_{17}NO_8$).



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

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



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

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

- Molecule 6 is water.

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

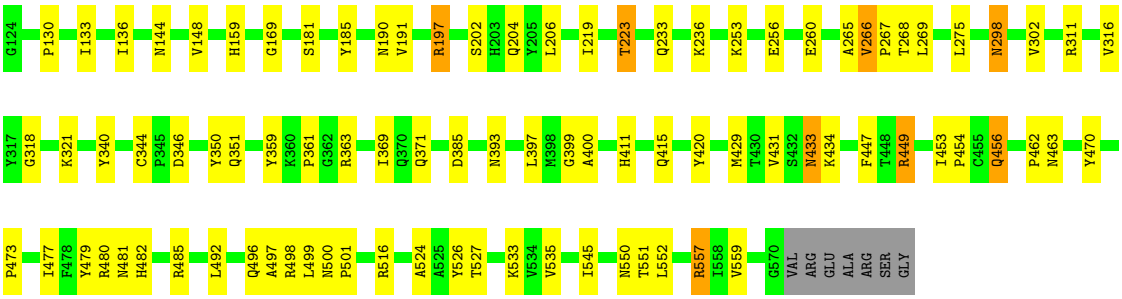
3 Residue-property plots (i)

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.

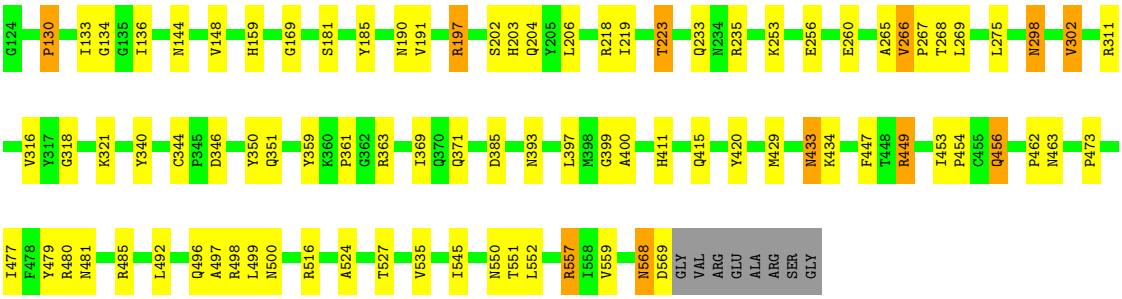
• Molecule 1: HEMAGGLUTININ-NEURAMINIDASE

Chain A:



• Molecule 1: HEMAGGLUTININ-NEURAMINIDASE

Chain B:



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

Chain C:

4 Data and refinement statistics (i)

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 (i)

5.1 Standard geometry (i)

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	0.39	0/3534	0.77	3/4807 (0.1%)
1	B	0.39	0/3530	0.74	3/4802 (0.1%)
All	All	0.39	0/7064	0.75	6/9609 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	449	ARG	NE-CZ-NH2	-17.13	111.73	120.30
1	A	449	ARG	NE-CZ-NH1	15.55	128.08	120.30
1	B	449	ARG	NE-CZ-NH2	-12.75	113.92	120.30
1	B	449	ARG	NE-CZ-NH1	12.12	126.36	120.30
1	A	449	ARG	CD-NE-CZ	8.08	134.91	123.60
1	B	449	ARG	CD-NE-CZ	6.32	132.45	123.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3449	0	3342	53	0
1	B	3445	0	3339	54	0
2	C	28	0	25	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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
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	109	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 (109) 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: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:B:481:ASN:CG	2:C:1:NAG:HN2	2.06	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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: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: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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:501:PRO:HD2	6:A:2106:HOH:O	2.16	0.45
1:A:456:GLN:NE2	1:A:456:GLN:H	2.15	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: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:A:185:TYR:HH	1:A:223:THR:HG22	1.86	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: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
1:B:346:ASP:HB3	1:B:350:TYR:HD2	1.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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%)	17 11
1	B	444/454 (98%)	408 (92%)	31 (7%)	5 (1%)	14 8
All	All	889/908 (98%)	818 (92%)	62 (7%)	9 (1%)	15 9

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 (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	387/392 (99%)	367 (95%)	20 (5%)	23	19
1	B	387/392 (99%)	366 (95%)	21 (5%)	22	18
All	All	774/784 (99%)	733 (95%)	41 (5%)	22	18

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	159	HIS
1	A	197	ARG
1	A	219	ILE
1	A	223	THR
1	A	266	VAL
1	A	269	LEU
1	A	298	ASN
1	A	316	VAL
1	A	340	TYR
1	A	397	LEU
1	A	431	VAL
1	A	433	ASN
1	A	447	PHE
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	159	HIS
1	B	197	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	397	LEU
1	B	433	ASN

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Mol	Chain	Res	Type
1	B	447	PHE
1	B	449	ARG
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 (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	204	GLN
1	A	298	ASN
1	A	351	GLN
1	A	371	GLN
1	A	456	GLN
1	A	463	ASN
1	A	496	GLN
1	B	204	GLN
1	B	298	ASN
1	B	351	GLN
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	2,1	14,14,15	1.58	3 (21%)	17,19,21	0.74	0
2	NAG	C	2	2	14,14,15	0.52	0	17,19,21	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
2	NAG	C	1	2,1	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	-3.99	1.44	1.53
2	C	1	NAG	C4-C3	-2.75	1.45	1.52
2	C	1	NAG	C1-C2	2.25	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:

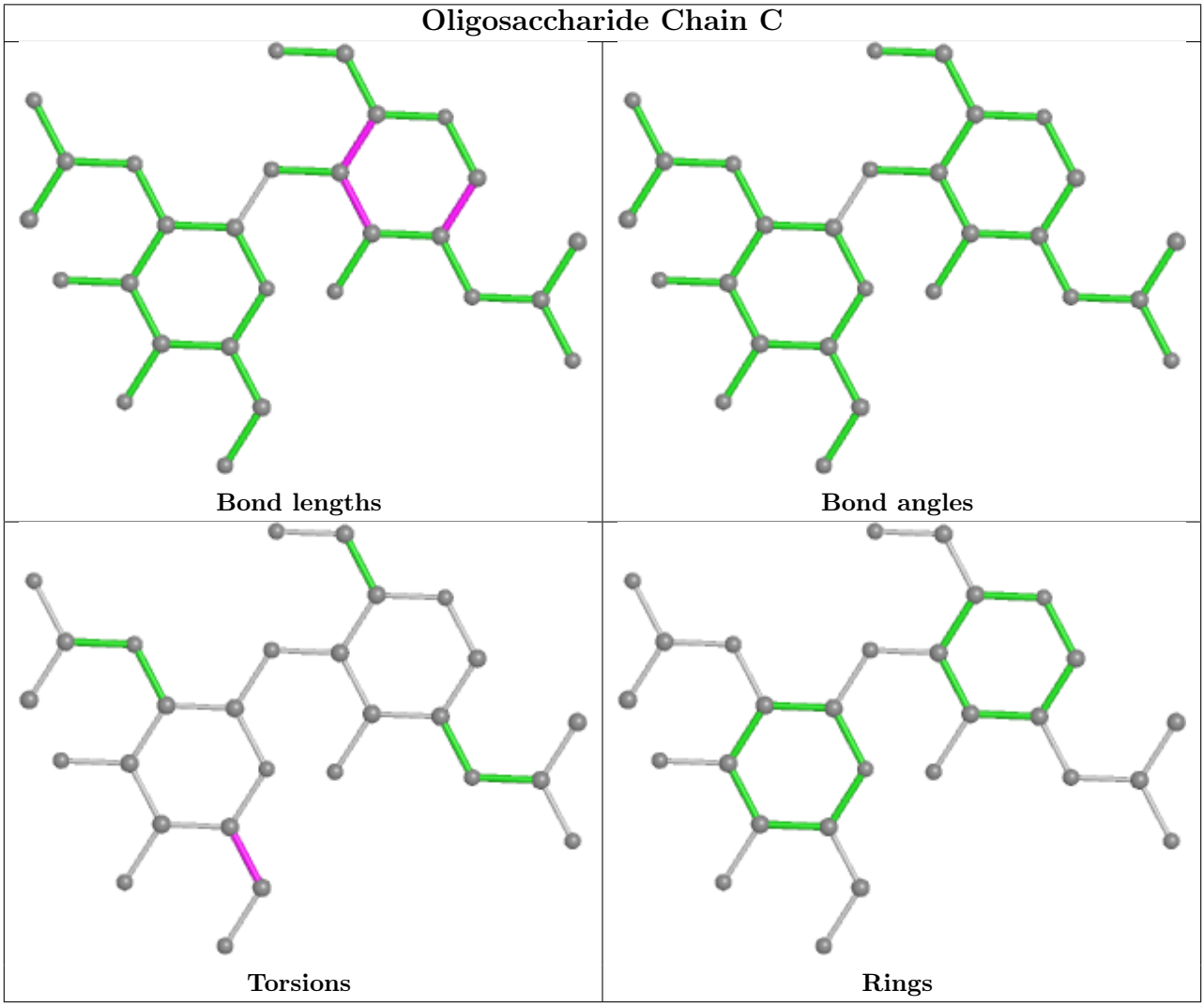
Mol	Chain	Res	Type	Atoms
2	C	2	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	NAG	1	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	$\#\|Z\| > 2$
4	DAN	B	1570	-	20,20,20	2.00	4 (20%)	23,28,28	2.12	4 (17%)
5	NAG	A	1573	1	14,14,15	0.84	0	17,19,21	0.60	0
4	DAN	A	1571	-	20,20,20	2.52	7 (35%)	23,28,28	1.63	4 (17%)
5	NAG	B	841	1	14,14,15	1.38	1 (7%)	17,19,21	0.73	0
5	NAG	A	1572	1	14,14,15	1.46	2 (14%)	17,19,21	0.88	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
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
4	DAN	A	1571	-	-	0/18/34/34	0/1/1/1
5	NAG	B	841	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1572	1	1/1/5/7	0/6/23/26	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1571	DAN	C3-C2	6.09	1.42	1.33
4	A	1571	DAN	C6-C5	5.68	1.62	1.53
4	B	1570	DAN	C3-C2	5.54	1.41	1.33
4	A	1571	DAN	C4-C3	4.47	1.56	1.50
5	B	841	NAG	C3-C2	4.36	1.61	1.52
4	B	1570	DAN	C4-C3	4.24	1.55	1.50
5	A	1572	NAG	C3-C2	3.81	1.60	1.52
5	A	1572	NAG	C1-C2	3.08	1.56	1.52
4	A	1571	DAN	C2-C1	-3.07	1.41	1.48
4	B	1570	DAN	C2-C1	-2.95	1.41	1.48
4	A	1571	DAN	C4-C5	2.76	1.56	1.53
4	A	1571	DAN	C5-N5	2.74	1.50	1.45
4	B	1570	DAN	C7-C6	-2.09	1.50	1.53
4	A	1571	DAN	O4-C4	-2.05	1.39	1.43

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	1570	DAN	O6-C2-C1	6.16	124.39	112.06
4	B	1570	DAN	O6-C2-C3	-5.70	117.97	124.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	1571	DAN	O6-C2-C1	3.68	119.43	112.06
4	A	1571	DAN	O4-C4-C3	3.52	117.25	109.31
4	A	1571	DAN	C3-C2-C1	-2.87	117.46	123.65
4	B	1570	DAN	C3-C2-C1	-2.79	117.63	123.65
4	A	1571	DAN	O6-C2-C3	-2.49	121.70	124.61
4	B	1570	DAN	C8-C7-C6	-2.13	109.00	113.03

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

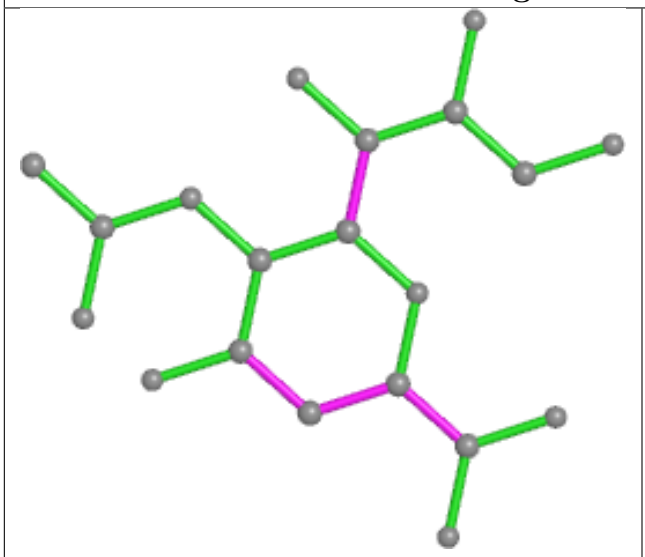
There are no ring outliers.

1 monomer is involved in 2 short contacts:

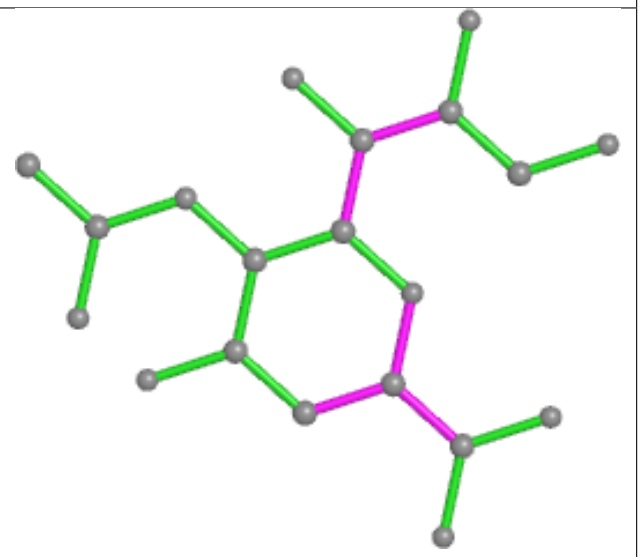
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1573	NAG	2	0

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 then 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.

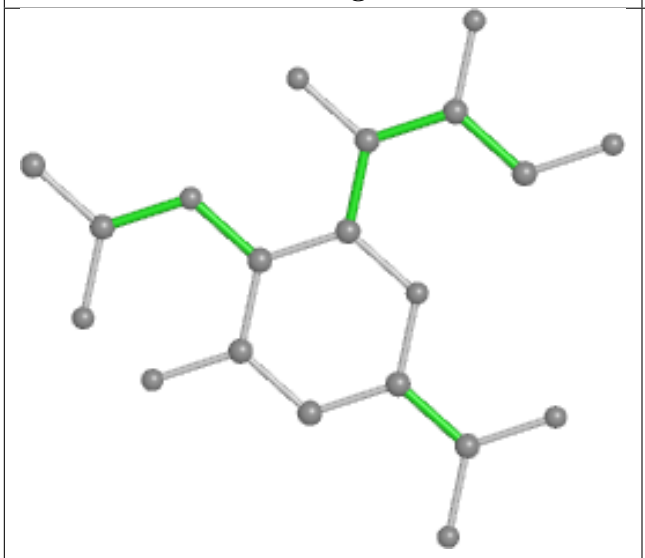
Ligand DAN B 1570



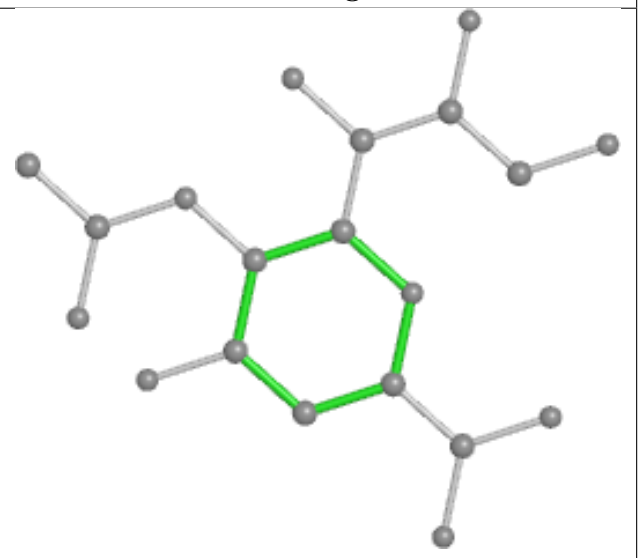
Bond lengths



Bond angles

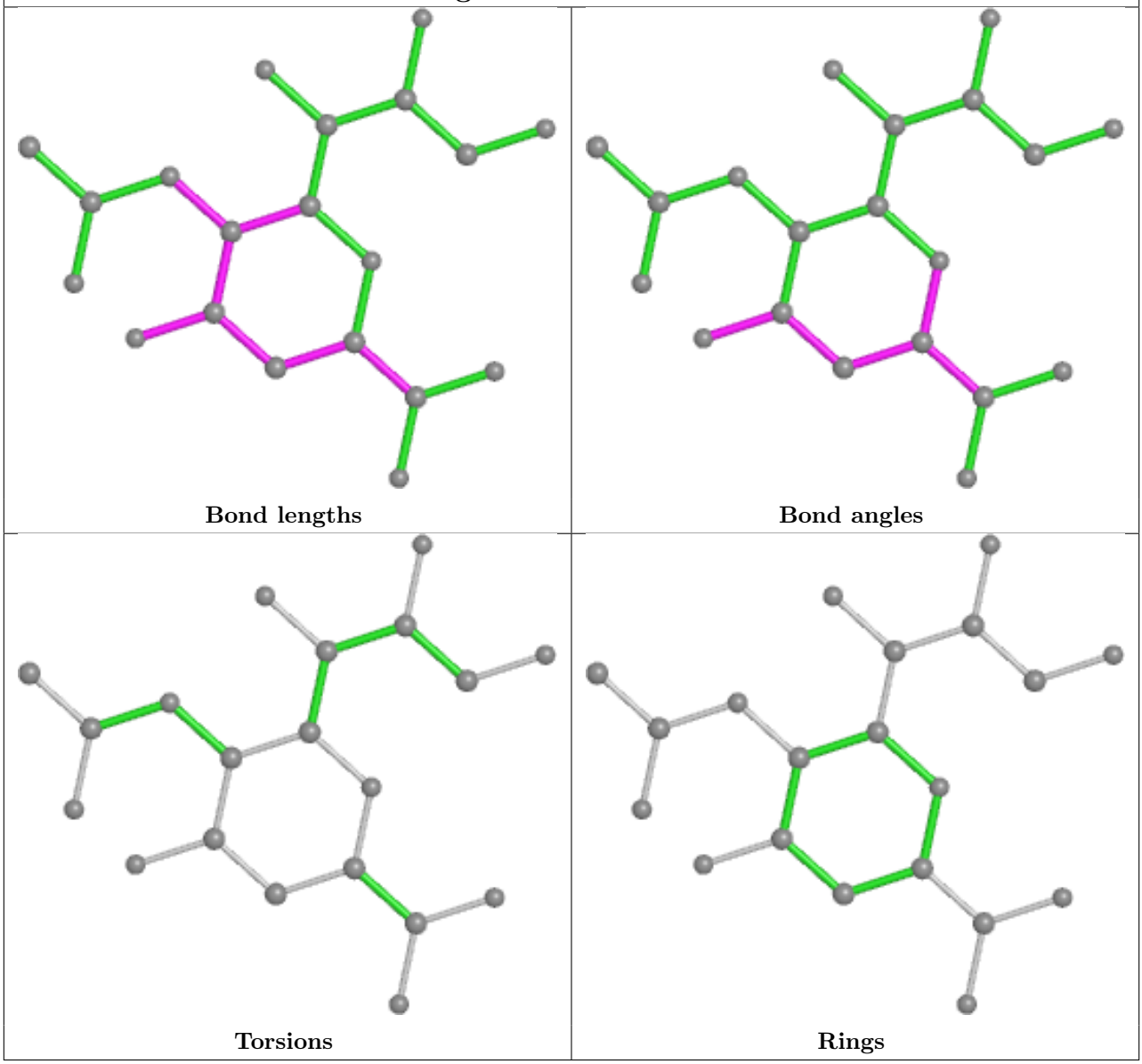


Torsions



Rings

Ligand DAN A 1571



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.