



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:54 AM UTC

PDB ID : 2IID / pdb_00002iid
Title : Structure of L-amino acid oxidase from Calloselasma rhodostoma in complex with L-phenylalanine
Authors : Moustafa, I.M.; Foster, S.; Lyubimov, A.Y.; Vrielink, A.
Deposited on : 2006-09-27
Resolution : 1.80 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

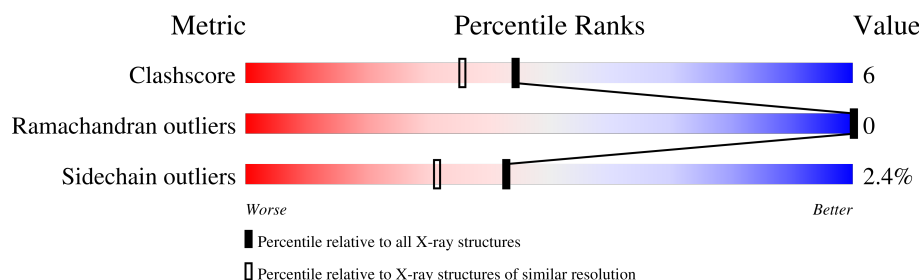
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)

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	498	86% 11% .
1	B	498	85% 11% ..
1	C	498	87% 10% ...
1	D	498	87% 10% ...
2	E	2	100%
2	F	2	50% 50%
2	G	2	100%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 18055 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 L-amino-acid oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	483	Total	C	N	O	S	0	4	0
			3892	2479	666	735	12			
1	B	483	Total	C	N	O	S	0	4	0
			3892	2479	666	735	12			
1	C	483	Total	C	N	O	S	0	4	0
			3892	2479	666	735	12			
1	D	483	Total	C	N	O	S	0	4	0
			3892	2479	666	735	12			

- Molecule 2 is an oligosaccharide called alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose.



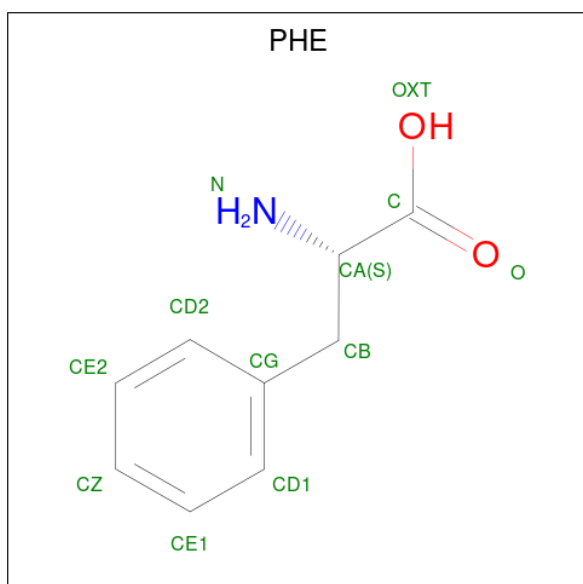
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	2	Total	C	N	O	0	0	0
			24	14	1	9			
2	F	2	Total	C	N	O	0	0	0
			24	14	1	9			
2	G	2	Total	C	N	O	0	0	0
			24	14	1	9			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



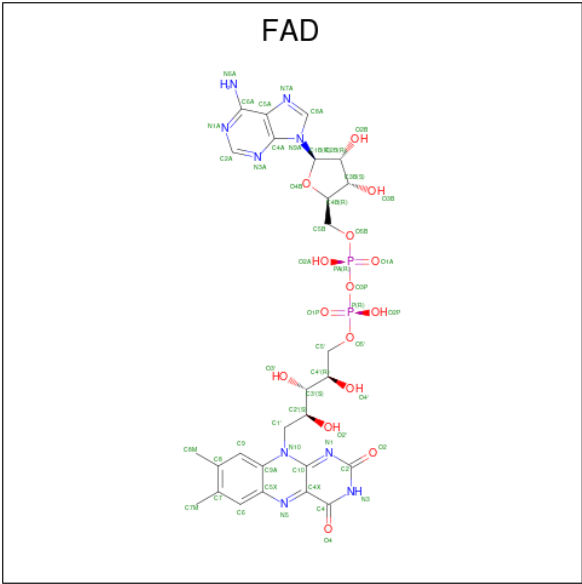
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is PHENYLALANINE (CCD ID: PHE) (formula: $C_9H_{11}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			12	9	1	2		
4	B	1	Total	C	N	O	0	0
			12	9	1	2		
4	C	1	Total	C	N	O	0	0
			12	9	1	2		
4	D	1	Total	C	N	O	0	0
			12	9	1	2		

- Molecule 5 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
5	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
5	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
5	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	538	Total	O	0	0
			538	538		
6	B	541	Total	O	0	0
			541	541		

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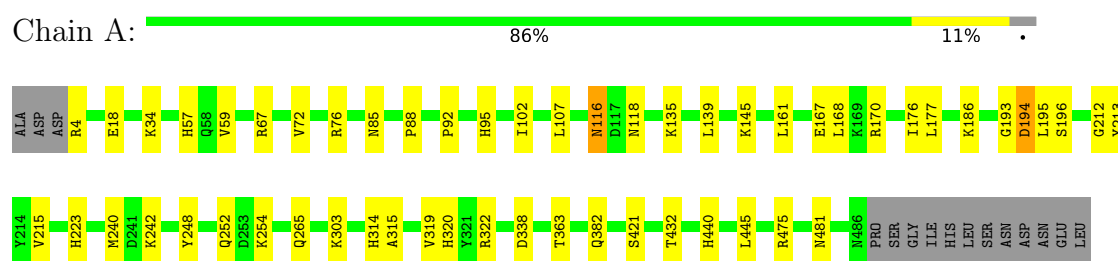
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	528	Total 528	O 528	0	0
6	D	478	Total 478	O 478	0	0

3 Residue-property plots

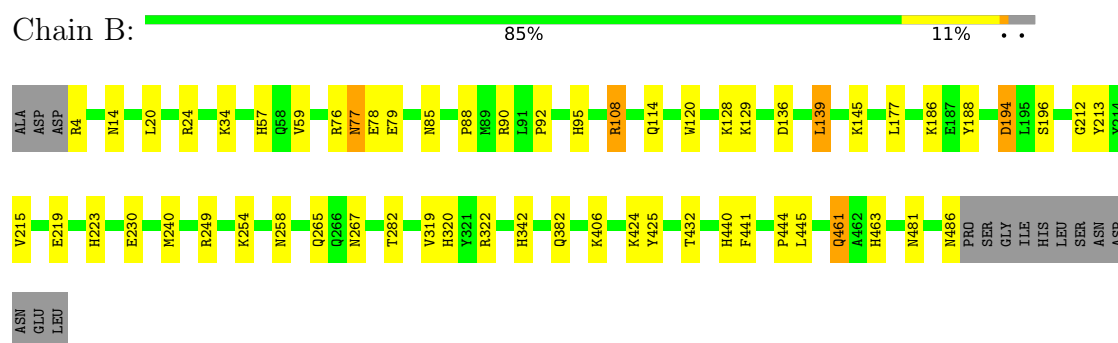
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

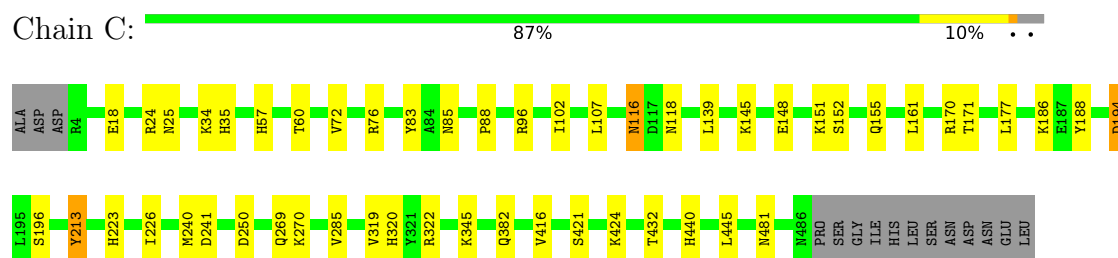
• Molecule 1: L-amino-acid oxidase



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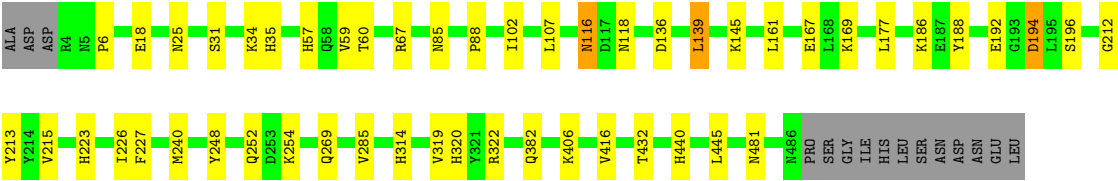


• Molecule 1: L-amino-acid oxidase



• Molecule 1: L-amino-acid oxidase





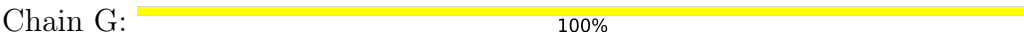
- Molecule 2: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 2: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 2: alpha-L-fucopyranose-(1-6)-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 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.76Å 154.00Å 103.18Å 90.00° 109.52° 90.00°	Depositor
Resolution (Å)	50.00 – 1.80	Depositor
% Data completeness (in resolution range)	99.5 (50.00-1.80)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.172 , 0.210	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	18055	wwPDB-VP
Average B, all atoms (Å ²)	19.0	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: NAG, FUC, FAD

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	0.75	0/3984	0.84	1/5391 (0.0%)
1	B	0.76	0/3984	0.83	0/5391
1	C	0.75	0/3984	0.82	0/5391
1	D	0.75	0/3984	0.83	0/5391
All	All	0.75	0/15936	0.83	1/21564 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	195	LEU	N-CA-C	7.13	120.34	110.35

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	3892	0	3808	53	0
1	B	3892	0	3808	56	0
1	C	3892	0	3808	46	0
1	D	3892	0	3808	44	0
2	E	24	0	22	0	0
2	F	24	0	22	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	G	24	0	22	0	0
3	A	14	0	13	0	0
3	B	28	0	26	0	0
3	C	14	0	13	1	0
3	D	14	0	13	0	0
4	A	12	0	8	1	0
4	B	12	0	8	0	0
4	C	12	0	8	1	0
4	D	12	0	8	0	0
5	A	53	0	31	2	0
5	B	53	0	31	1	0
5	C	53	0	31	3	0
5	D	53	0	31	1	0
6	A	538	0	0	18	0
6	B	541	0	0	14	0
6	C	528	0	0	15	0
6	D	478	0	0	14	0
All	All	18055	0	15519	190	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 6.

All (190) 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:322[A]:ARG:HD3	6:B:976:HOH:O	1.25	1.32
1:A:34:LYS:H	1:A:57:HIS:HD2	1.12	0.94
1:B:34:LYS:H	1:B:57:HIS:HD2	1.16	0.93
1:C:34:LYS:H	1:C:57:HIS:HD2	1.17	0.88
1:B:90:ARG:HH11	1:B:114:GLN:HE22	1.24	0.85
1:B:145:LYS:HG3	1:B:196:SER:HB3	1.59	0.84
1:A:254:LYS:HE3	6:A:849:HOH:O	1.78	0.82
1:B:219:GLU:OE2	1:B:463:HIS:HD2	1.62	0.81
1:D:145:LYS:HG3	1:D:196:SER:HB3	1.63	0.81
1:A:116:ASN:HD22	1:A:118:ASN:H	1.30	0.80
1:C:322[A]:ARG:HD3	6:C:778:HOH:O	1.80	0.79
1:D:34:LYS:H	1:D:57:HIS:HD2	1.31	0.79
1:C:145:LYS:HG3	1:C:196:SER:HB3	1.64	0.78
1:B:319:VAL:HG21	1:B:445:LEU:HD11	1.67	0.76
1:B:108:ARG:HH11	1:B:108:ARG:HG3	1.48	0.76
1:C:223[A]:HIS:CE1	6:C:995:HOH:O	2.39	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:59:VAL:HG22	1:B:254:LYS:HD3	1.69	0.75
1:D:116:ASN:HD22	1:D:118:ASN:H	1.33	0.75
1:D:322[A]:ARG:HD3	6:D:994:HOH:O	1.87	0.74
1:C:116:ASN:HD22	1:C:118:ASN:H	1.36	0.74
1:B:282:THR:O	1:B:282:THR:HG23	1.88	0.73
1:A:315:ALA:O	1:A:319:VAL:HG23	1.89	0.73
1:B:186:LYS:NZ	1:D:440:HIS:HD2	1.87	0.73
1:A:145:LYS:HG3	1:A:196:SER:HB3	1.71	0.72
1:D:319:VAL:HG21	1:D:445:LEU:HD11	1.70	0.72
1:B:14:ASN:H	1:B:461:GLN:NE2	1.87	0.71
1:B:108:ARG:HD3	6:B:917:HOH:O	1.89	0.71
1:A:440:HIS:HD2	1:C:186:LYS:NZ	1.89	0.70
1:B:85:ASN:HD21	1:B:240:MET:H	1.39	0.70
1:D:85:ASN:HD21	1:D:240:MET:H	1.38	0.70
1:C:25:ASN:ND2	6:C:702:HOH:O	2.24	0.69
1:A:475:ARG:HG2	6:A:994:HOH:O	1.91	0.69
1:B:57:HIS:HE1	1:B:481:ASN:OD1	1.77	0.67
1:D:57:HIS:HE1	1:D:481:ASN:OD1	1.77	0.67
1:D:322[A]:ARG:HG2	6:D:994:HOH:O	1.94	0.66
1:A:34:LYS:N	1:A:57:HIS:HD2	1.90	0.65
1:A:319:VAL:HG21	1:A:445:LEU:HD11	1.78	0.65
1:C:85:ASN:HD21	1:C:240:MET:H	1.42	0.65
1:C:319:VAL:HG21	1:C:445:LEU:HD11	1.77	0.65
1:D:59:VAL:HG22	1:D:254:LYS:HD3	1.79	0.64
1:B:14:ASN:H	1:B:461:GLN:HE21	1.45	0.64
1:A:76:ARG:NH1	6:A:1054:HOH:O	2.30	0.63
1:C:18:GLU:HG3	6:C:975:HOH:O	1.98	0.63
1:B:440:HIS:HD2	1:D:186:LYS:NZ	1.97	0.62
1:A:85:ASN:HD21	1:A:240:MET:H	1.47	0.62
1:A:248:TYR:CZ	1:A:252:GLN:HG3	2.36	0.61
1:B:34:LYS:H	1:B:57:HIS:CD2	2.08	0.61
1:A:116:ASN:ND2	1:A:118:ASN:H	1.96	0.61
4:A:526:PHE:HA	5:A:527:FAD:N5	2.15	0.60
1:D:116:ASN:ND2	1:D:118:ASN:H	1.99	0.60
1:A:322[A]:ARG:HD2	6:A:878:HOH:O	2.02	0.60
4:C:526:PHE:HA	5:C:527:FAD:N5	2.17	0.59
1:A:88:PRO:HA	5:A:527:FAD:C4X	2.32	0.59
1:C:57:HIS:HE1	1:C:481:ASN:OD1	1.85	0.58
1:B:85:ASN:ND2	1:B:240:MET:H	2.01	0.58
1:D:136:ASP:HB3	1:D:139:LEU:HD22	1.86	0.58
1:A:382:GLN:NE2	1:C:382:GLN:NE2	2.52	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:ARG:NH1	1:B:114:GLN:HE22	1.98	0.58
1:D:88:PRO:HA	5:D:527:FAD:C4X	2.34	0.57
1:A:194:ASP:C	1:A:194:ASP:OD1	2.47	0.57
1:C:76:ARG:NH2	1:C:241:ASP:OD1	2.28	0.57
1:D:85:ASN:ND2	1:D:240:MET:H	2.01	0.57
1:B:219:GLU:OE2	1:B:463:HIS:CD2	2.51	0.56
3:C:522:NAG:O7	6:C:676:HOH:O	2.17	0.56
1:A:34:LYS:H	1:A:57:HIS:CD2	2.05	0.56
1:D:223[A]:HIS:CE1	1:D:227:PHE:CD1	2.94	0.56
1:B:77:ASN:C	1:B:77:ASN:HD22	2.14	0.55
1:D:34:LYS:H	1:D:57:HIS:CD2	2.19	0.55
1:B:76:ARG:NH1	6:B:1054:HOH:O	2.36	0.55
1:C:85:ASN:ND2	1:C:240:MET:H	2.04	0.55
1:B:186:LYS:HZ1	1:D:440:HIS:HD2	1.52	0.55
1:B:223[B]:HIS:CD2	6:B:1044:HOH:O	2.58	0.54
1:A:85:ASN:ND2	1:A:240:MET:H	2.05	0.54
1:D:406:LYS:HE2	6:D:995:HOH:O	2.07	0.54
1:A:314:HIS:HD2	6:C:688:HOH:O	1.90	0.54
1:D:102:ILE:HG23	1:D:107:LEU:HB2	1.89	0.54
1:D:25:ASN:ND2	6:D:915:HOH:O	2.40	0.54
1:B:463:HIS:HE1	6:B:540:HOH:O	1.89	0.54
1:C:345:LYS:NZ	6:C:961:HOH:O	2.39	0.53
1:B:322[A]:ARG:NE	1:B:432:THR:HG21	2.23	0.53
1:C:35:HIS:HE1	1:C:60:THR:OG1	1.92	0.53
1:D:322[A]:ARG:CG	6:D:994:HOH:O	2.54	0.52
1:A:363:THR:HG22	6:A:872:HOH:O	2.09	0.52
1:A:475:ARG:NH1	6:A:634:HOH:O	2.42	0.52
1:B:319:VAL:CG2	1:B:445:LEU:HD11	2.39	0.52
1:D:186:LYS:NZ	6:D:600:HOH:O	2.42	0.52
1:C:116:ASN:ND2	1:C:118:ASN:H	2.07	0.52
1:B:59:VAL:CG2	1:B:254:LYS:HD3	2.39	0.51
1:A:303:LYS:HE3	6:A:988:HOH:O	2.11	0.51
1:B:78:GLU:HG2	6:B:1026:HOH:O	2.10	0.51
1:A:440:HIS:HD2	1:C:186:LYS:HZ1	1.54	0.51
1:B:258:ASN:OD1	6:B:1041:HOH:O	2.20	0.50
6:B:635:HOH:O	1:D:314:HIS:HD2	1.94	0.50
1:A:475:ARG:NH2	6:A:763:HOH:O	2.45	0.50
1:C:226:ILE:HD11	6:C:934:HOH:O	2.10	0.50
1:A:57:HIS:HE1	1:A:481:ASN:OD1	1.94	0.50
1:C:24:ARG:HD2	6:C:702:HOH:O	2.10	0.50
1:D:226:ILE:HD11	6:D:880:HOH:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:THR:O	1:B:282:THR:CG2	2.59	0.49
1:C:194:ASP:OD1	1:C:194:ASP:C	2.55	0.49
1:A:76:ARG:HG2	1:A:76:ARG:HH11	1.78	0.48
1:D:31:SER:HA	6:D:977:HOH:O	2.12	0.48
1:B:322[A]:ARG:CD	6:B:976:HOH:O	2.11	0.48
1:A:186:LYS:NZ	1:C:440:HIS:HD2	2.11	0.48
1:B:249:ARG:HD2	6:B:883:HOH:O	2.13	0.48
1:D:35:HIS:HE1	1:D:60:THR:OG1	1.96	0.48
1:C:116:ASN:HD21	1:C:118:ASN:HD22	1.60	0.48
1:B:76:ARG:HH11	1:B:76:ARG:HG2	1.77	0.47
1:B:92:PRO:HG2	1:B:95:HIS:CG	2.49	0.47
1:D:18:GLU:HG3	6:D:804:HOH:O	2.14	0.47
1:A:102:ILE:HG23	1:A:107:LEU:HB2	1.95	0.47
1:B:382:GLN:NE2	1:D:382:GLN:HE22	2.12	0.47
1:A:72:VAL:HG22	1:A:88:PRO:HG2	1.96	0.47
1:A:135:LYS:HG3	6:A:778:HOH:O	2.14	0.47
1:C:226:ILE:HD12	6:C:942:HOH:O	2.14	0.47
1:C:34:LYS:H	1:C:57:HIS:CD2	2.10	0.47
1:C:72:VAL:HG22	1:C:88:PRO:HG2	1.97	0.47
1:A:382:GLN:CD	1:C:382:GLN:NE2	2.73	0.47
1:B:440:HIS:HD2	1:D:186:LYS:HZ1	1.62	0.47
1:A:382:GLN:NE2	1:C:382:GLN:CD	2.72	0.47
1:D:188:TYR:CD2	1:D:188:TYR:C	2.92	0.47
1:A:475:ARG:CG	6:A:994:HOH:O	2.57	0.46
1:B:342:HIS:HD2	6:B:686:HOH:O	1.98	0.46
1:C:88:PRO:HA	5:C:527:FAD:C4X	2.45	0.46
1:C:424:LYS:HD2	6:C:891:HOH:O	2.14	0.46
1:B:88:PRO:HA	5:B:525:FAD:C4X	2.46	0.46
1:C:102:ILE:HG23	1:C:107:LEU:HB2	1.97	0.46
1:D:194:ASP:C	1:D:194:ASP:OD1	2.58	0.46
1:B:186:LYS:NZ	6:B:1024:HOH:O	2.49	0.46
1:D:248:TYR:CZ	1:D:252:GLN:HG3	2.51	0.46
1:A:116:ASN:HD22	1:A:116:ASN:C	2.23	0.46
1:A:319:VAL:CG2	1:A:445:LEU:HD11	2.45	0.46
1:A:18:GLU:HG3	6:A:700:HOH:O	2.14	0.46
1:B:136:ASP:HB3	1:B:139:LEU:HD22	1.97	0.46
1:D:322[A]:ARG:CD	6:D:994:HOH:O	2.53	0.45
1:A:76:ARG:NH1	1:A:76:ARG:HG2	2.31	0.45
1:D:320:HIS:HB3	1:D:432:THR:O	2.16	0.45
1:C:170:ARG:NH1	2:F:2:FUC:H62	2.31	0.45
1:D:192:GLU:OE2	6:D:824:HOH:O	2.21	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:319:VAL:CG2	1:D:445:LEU:HD11	2.43	0.45
1:D:212:GLY:O	1:D:215:VAL:HG22	2.16	0.45
1:B:120:TRP:CH2	1:B:129:LYS:HE2	2.52	0.44
1:B:382:GLN:CD	1:D:382:GLN:NE2	2.75	0.44
1:D:169:LYS:HG3	6:D:649:HOH:O	2.17	0.44
1:C:322[A]:ARG:NE	1:C:432:THR:HG21	2.32	0.44
1:A:167:GLU:CD	1:A:170:ARG:HH21	2.26	0.44
1:B:424:LYS:HE3	1:B:425:TYR:CE2	2.52	0.44
1:B:194:ASP:HA	6:B:932:HOH:O	2.17	0.44
1:B:265:GLN:HE21	1:B:267:ASN:HD21	1.66	0.43
1:C:96:ARG:HD2	6:C:924:HOH:O	2.18	0.43
1:A:92:PRO:HG2	1:A:95:HIS:CG	2.53	0.43
1:C:83:TYR:C	1:C:83:TYR:CD1	2.97	0.43
1:B:128:LYS:HE3	1:B:139:LEU:HG	2.00	0.43
1:A:440:HIS:HD2	1:C:186:LYS:HZ3	1.66	0.43
1:D:223[A]:HIS:NE2	1:D:227:PHE:CE1	2.87	0.43
1:A:67:ARG:NH2	6:A:598:HOH:O	2.51	0.43
1:B:212:GLY:O	1:B:215:VAL:HG22	2.18	0.43
1:B:20:LEU:O	1:B:24:ARG:HG2	2.19	0.43
1:A:223[B]:HIS:CD2	6:A:907:HOH:O	2.72	0.42
1:B:461:GLN:HG2	6:B:734:HOH:O	2.19	0.42
1:A:193:GLY:O	1:A:194:ASP:HB3	2.18	0.42
1:C:171:THR:HG22	2:F:2:FUC:H5	2.02	0.42
1:B:441:PHE:C	1:B:444:PRO:HD2	2.45	0.42
1:B:320:HIS:HB3	1:B:432:THR:O	2.20	0.42
1:C:320:HIS:HB3	1:C:432:THR:O	2.20	0.41
1:A:4:ARG:N	6:A:953:HOH:O	2.53	0.41
1:B:79:GLU:HG2	6:C:718:HOH:O	2.19	0.41
1:C:213:TYR:HD1	1:C:213:TYR:HA	1.73	0.41
1:D:67:ARG:NH2	6:D:604:HOH:O	2.49	0.41
1:C:269:GLN:O	1:C:270:LYS:HD3	2.20	0.41
1:B:188:TYR:CD2	1:B:188:TYR:C	2.98	0.41
1:C:34:LYS:N	1:C:57:HIS:HD2	1.99	0.41
1:B:76:ARG:NH1	1:B:76:ARG:HG2	2.34	0.41
1:B:92:PRO:HB3	1:B:230:GLU:HG2	2.03	0.41
1:A:322[A]:ARG:CD	6:A:878:HOH:O	2.66	0.41
1:C:188:TYR:CD2	1:C:188:TYR:C	2.98	0.41
5:C:527:FAD:H8A	6:C:534:HOH:O	2.19	0.41
1:D:116:ASN:HD22	1:D:118:ASN:N	2.10	0.41
1:A:168:LEU:HB2	1:A:176:ILE:HD13	2.03	0.41
1:A:265:GLN:HG2	6:A:1064:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:338:ASP:HB2	6:A:1012:HOH:O	2.21	0.41
1:C:223[A]:HIS:HE1	6:C:995:HOH:O	1.93	0.41
1:C:152:SER:OG	1:C:155:GLN:HG3	2.20	0.40
1:A:212:GLY:O	1:A:215:VAL:HG22	2.22	0.40
1:A:303:LYS:NZ	6:A:932:HOH:O	2.54	0.40
1:A:382:GLN:NE2	1:A:421[B]:SER:OG	2.55	0.40
1:A:320:HIS:HB3	1:A:432:THR:O	2.21	0.40
1:C:382:GLN:NE2	1:C:421[B]:SER:OG	2.54	0.40
1:D:6:PRO:HG2	6:D:951:HOH:O	2.21	0.40
1:C:148:GLU:HA	1:C:151:LYS:HD2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	485/498 (97%)	474 (98%)	11 (2%)	0	100	100
1	B	485/498 (97%)	474 (98%)	11 (2%)	0	100	100
1	C	485/498 (97%)	471 (97%)	14 (3%)	0	100	100
1	D	485/498 (97%)	474 (98%)	11 (2%)	0	100	100
All	All	1940/1992 (97%)	1893 (98%)	47 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	418/427 (98%)	410 (98%)	8 (2%)	50	41
1	B	418/427 (98%)	408 (98%)	10 (2%)	43	31
1	C	418/427 (98%)	409 (98%)	9 (2%)	45	34
1	D	418/427 (98%)	408 (98%)	10 (2%)	43	31
All	All	1672/1708 (98%)	1635 (98%)	37 (2%)	43	34

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	VAL
1	A	116	ASN
1	A	139	LEU
1	A	161	LEU
1	A	177	LEU
1	A	194	ASP
1	A	213	TYR
1	A	242	LYS
1	B	4	ARG
1	B	77	ASN
1	B	108	ARG
1	B	139	LEU
1	B	177	LEU
1	B	194	ASP
1	B	213	TYR
1	B	406	LYS
1	B	461	GLN
1	B	486	ASN
1	C	116	ASN
1	C	139	LEU
1	C	161	LEU
1	C	177	LEU
1	C	194	ASP
1	C	213	TYR
1	C	250	ASP
1	C	285	VAL
1	C	416	VAL
1	D	116	ASN
1	D	139	LEU
1	D	161	LEU

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Mol	Chain	Res	Type
1	D	167	GLU
1	D	177	LEU
1	D	194	ASP
1	D	213	TYR
1	D	269	GLN
1	D	285	VAL
1	D	416	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	HIS
1	A	57	HIS
1	A	85	ASN
1	A	110	ASN
1	A	116	ASN
1	A	125	ASN
1	A	252	GLN
1	A	269	GLN
1	A	314	HIS
1	A	382	GLN
1	A	439	GLN
1	A	440	HIS
1	B	57	HIS
1	B	77	ASN
1	B	85	ASN
1	B	110	ASN
1	B	114	GLN
1	B	118	ASN
1	B	178	ASN
1	B	258	ASN
1	B	267	ASN
1	B	342	HIS
1	B	440	HIS
1	B	461	GLN
1	B	463	HIS
1	C	35	HIS
1	C	57	HIS
1	C	85	ASN
1	C	116	ASN
1	C	208	ASN
1	C	252	GLN

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Mol	Chain	Res	Type
1	C	269	GLN
1	C	314	HIS
1	C	382	GLN
1	C	440	HIS
1	D	35	HIS
1	D	57	HIS
1	D	58	GLN
1	D	85	ASN
1	D	110	ASN
1	D	116	ASN
1	D	118	ASN
1	D	252	GLN
1	D	258	ASN
1	D	265	GLN
1	D	269	GLN
1	D	314	HIS
1	D	382	GLN
1	D	440	HIS

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 ⓘ

6 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	E	1	2,1	14,14,15	0.56	0	17,19,21	1.17	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FUC	E	2	2	10,10,11	0.78	0	14,14,16	1.70	4 (28%)
2	NAG	F	1	2,1	14,14,15	0.49	0	17,19,21	0.70	0
2	FUC	F	2	2	10,10,11	0.67	0	14,14,16	0.55	0
2	NAG	G	1	2,1	14,14,15	0.56	0	17,19,21	1.10	1 (5%)
2	FUC	G	2	2	10,10,11	0.72	0	14,14,16	1.84	4 (28%)

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	E	1	2,1	-	2/6/23/26	0/1/1/1
2	FUC	E	2	2	-	-	0/1/1/1
2	NAG	F	1	2,1	-	2/6/23/26	0/1/1/1
2	FUC	F	2	2	-	-	0/1/1/1
2	NAG	G	1	2,1	-	0/6/23/26	0/1/1/1
2	FUC	G	2	2	-	-	0/1/1/1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	2	FUC	O5-C5-C4	3.74	116.28	109.55
2	E	2	FUC	O5-C5-C4	3.44	115.74	109.55
2	G	2	FUC	C3-C4-C5	3.32	114.86	109.81
2	G	2	FUC	C1-O5-C5	2.98	119.99	112.97
2	E	2	FUC	C3-C4-C5	2.95	114.30	109.81
2	E	2	FUC	C1-O5-C5	2.88	119.76	112.97
2	G	1	NAG	C2-N2-C7	2.55	126.32	122.90
2	E	1	NAG	O5-C1-C2	-2.46	107.48	111.29
2	G	2	FUC	O2-C2-C3	-2.34	105.30	110.15
2	E	2	FUC	C1-C2-C3	-2.18	106.47	109.64

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	1	NAG	C8-C7-N2-C2
2	F	1	NAG	O7-C7-N2-C2

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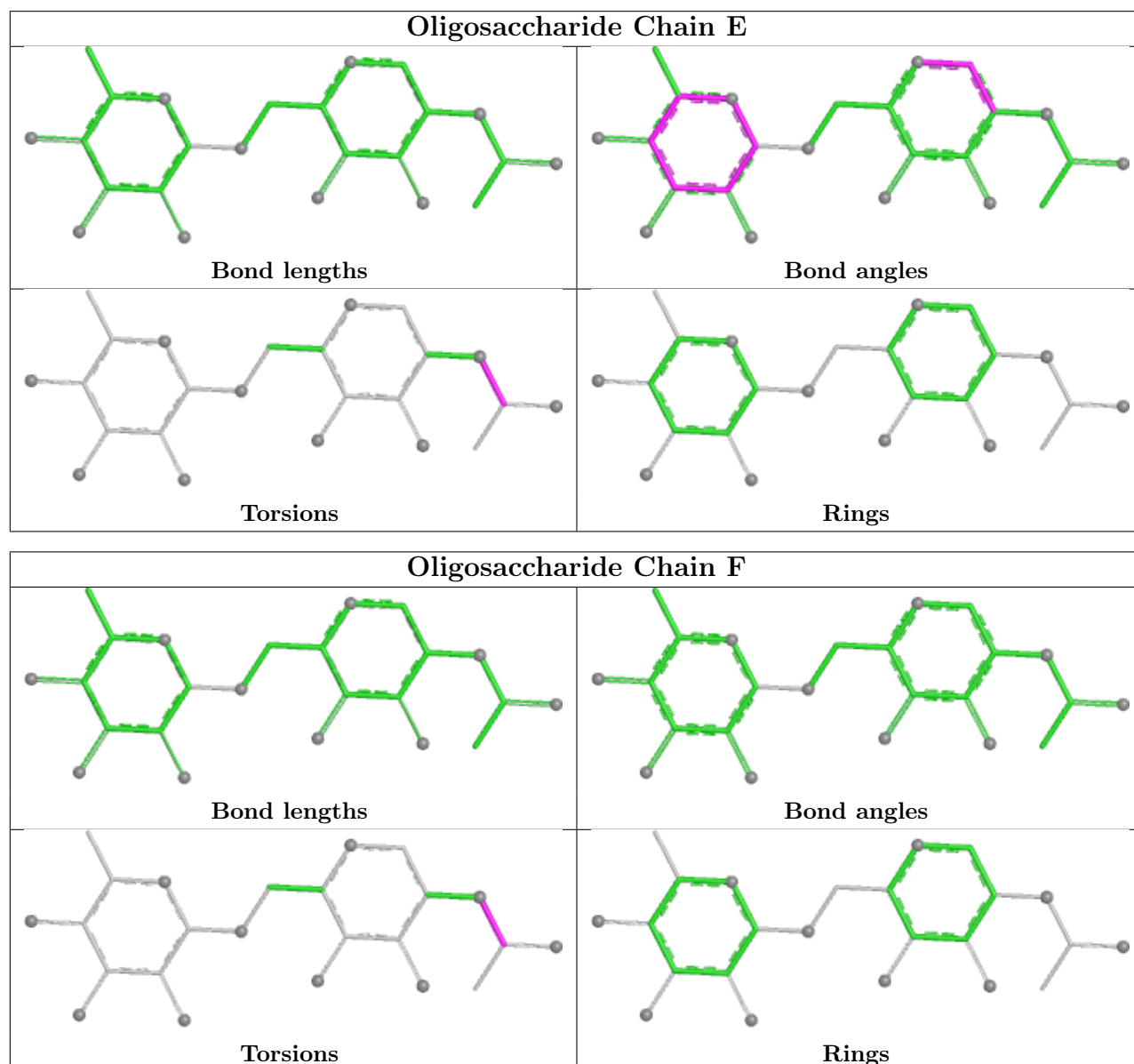
Mol	Chain	Res	Type	Atoms
2	E	1	NAG	C8-C7-N2-C2
2	E	1	NAG	O7-C7-N2-C2

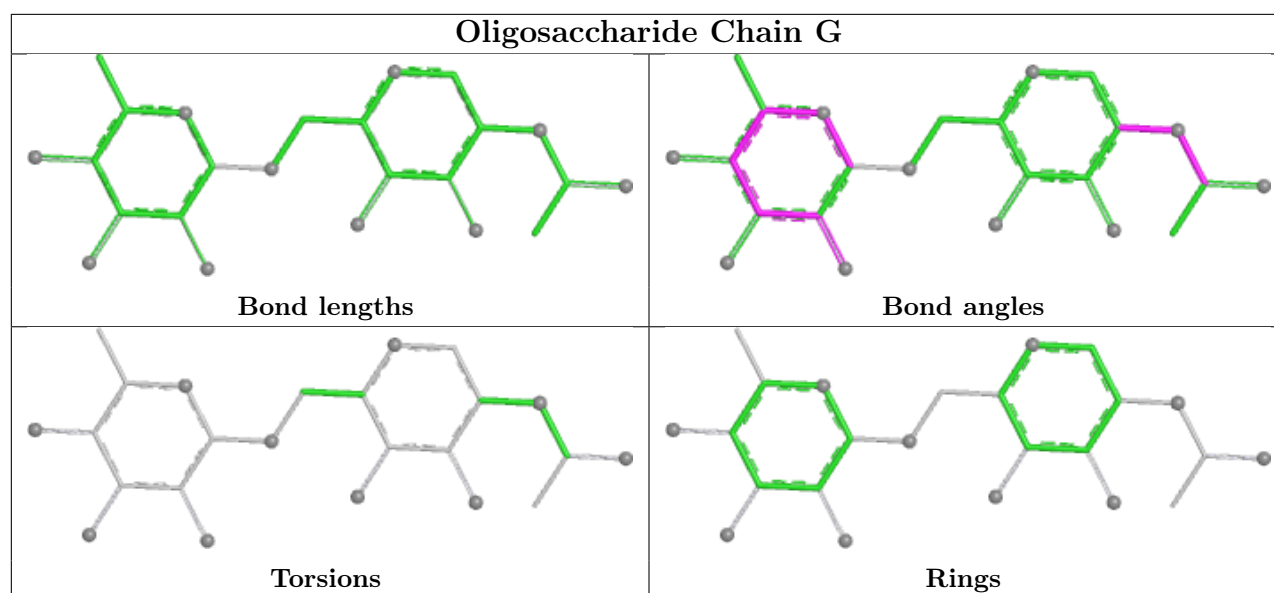
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	2	FUC	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](#)

13 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	FAD	A	527	-	58,58,58	1.28	7 (12%)	85,89,89	1.59	17 (20%)
5	FAD	B	525	-	58,58,58	1.24	5 (8%)	85,89,89	1.59	16 (18%)
4	PHE	C	526	-	11,12,12	0.80	0	11,15,15	1.58	2 (18%)
4	PHE	A	526	-	11,12,12	0.81	1 (9%)	11,15,15	1.45	3 (27%)
4	PHE	B	524	-	11,12,12	1.02	1 (9%)	11,15,15	2.50	3 (27%)
3	NAG	C	522	1	14,14,15	0.62	0	17,19,21	1.73	5 (29%)
3	NAG	B	523	1	14,14,15	0.51	0	17,19,21	1.14	1 (5%)
5	FAD	C	527	-	58,58,58	1.21	5 (8%)	85,89,89	1.53	17 (20%)
5	FAD	D	527	-	58,58,58	1.29	5 (8%)	85,89,89	1.59	17 (20%)
3	NAG	A	522	1	14,14,15	0.46	0	17,19,21	1.24	2 (11%)
4	PHE	D	526	-	11,12,12	0.89	1 (9%)	11,15,15	1.66	3 (27%)
3	NAG	D	522	1	14,14,15	0.47	0	17,19,21	1.33	2 (11%)
3	NAG	B	522	1	14,14,15	0.54	0	17,19,21	0.89	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	FAD	A	527	-	-	5/34/50/50	0/6/6/6
5	FAD	B	525	-	-	6/34/50/50	0/6/6/6
4	PHE	C	526	-	-	3/8/8/8	0/1/1/1
4	PHE	A	526	-	-	5/8/8/8	0/1/1/1
4	PHE	B	524	-	-	5/8/8/8	0/1/1/1
3	NAG	C	522	1	-	1/6/23/26	0/1/1/1
3	NAG	B	523	1	-	2/6/23/26	0/1/1/1
5	FAD	C	527	-	-	5/34/50/50	0/6/6/6
5	FAD	D	527	-	-	4/34/50/50	0/6/6/6
3	NAG	A	522	1	-	0/6/23/26	0/1/1/1
4	PHE	D	526	-	-	4/8/8/8	0/1/1/1
3	NAG	D	522	1	-	2/6/23/26	0/1/1/1
3	NAG	B	522	1	-	0/6/23/26	0/1/1/1

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	525	FAD	C4X-N5	4.84	1.41	1.30
5	D	527	FAD	C4X-N5	4.46	1.40	1.30
5	C	527	FAD	C4X-N5	4.26	1.40	1.30
5	A	527	FAD	C4X-N5	4.15	1.39	1.30
5	D	527	FAD	C10-N1	3.73	1.40	1.33
5	D	527	FAD	P-O3P	3.72	1.63	1.59
5	A	527	FAD	P-O3P	3.55	1.63	1.59
5	C	527	FAD	P-O3P	3.41	1.63	1.59
5	C	527	FAD	C10-N1	3.33	1.39	1.33
5	B	525	FAD	P-O3P	3.31	1.63	1.59
5	A	527	FAD	C10-N1	3.03	1.39	1.33
5	A	527	FAD	C2A-N3A	3.01	1.39	1.33
5	B	525	FAD	C10-N1	2.91	1.39	1.33
5	B	525	FAD	C2A-N1A	2.88	1.39	1.33
5	A	527	FAD	C1'-C2'	2.63	1.56	1.52
5	A	527	FAD	C8A-N7A	2.59	1.36	1.31
5	D	527	FAD	C2A-N3A	2.55	1.38	1.33
5	B	525	FAD	C8A-N7A	2.49	1.36	1.31
5	A	527	FAD	C2A-N1A	2.36	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	527	FAD	C8A-N7A	2.34	1.36	1.31
5	D	527	FAD	C1'-C2'	2.24	1.55	1.52
5	C	527	FAD	C2A-N3A	2.15	1.37	1.33
4	D	526	PHE	OXT-C	-2.08	1.24	1.30
4	A	526	PHE	OXT-C	-2.07	1.24	1.30
4	B	524	PHE	OXT-C	-2.05	1.24	1.30

All (88) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	525	FAD	N3A-C2A-N1A	-6.92	118.10	128.58
5	D	527	FAD	N3A-C2A-N1A	-5.97	119.55	128.58
4	B	524	PHE	OXT-C-O	5.53	136.63	124.08
5	A	527	FAD	N3A-C2A-N1A	-5.44	120.35	128.58
4	B	524	PHE	CB-CA-N	-5.38	90.80	111.40
5	C	527	FAD	N3A-C2A-N1A	-4.44	121.86	128.58
5	C	527	FAD	C5A-C4A-N3A	-4.23	120.89	126.72
5	A	527	FAD	N9A-C8A-N7A	-4.18	108.01	113.94
4	C	526	PHE	CB-CA-N	-4.02	96.00	111.40
4	D	526	PHE	OXT-C-O	3.97	133.08	124.08
5	D	527	FAD	N9A-C8A-N7A	-3.80	108.55	113.94
5	C	527	FAD	N9A-C8A-N7A	-3.71	108.68	113.94
5	B	525	FAD	C2A-N3A-C4A	3.69	120.85	111.83
5	B	525	FAD	N9A-C8A-N7A	-3.68	108.72	113.94
3	C	522	NAG	O5-C1-C2	-3.54	105.82	111.29
5	A	527	FAD	C5A-N7A-C8A	3.44	108.86	103.45
5	D	527	FAD	C5A-C4A-N3A	-3.31	122.15	126.72
5	D	527	FAD	O3P-P-O1P	-3.24	100.97	110.70
5	C	527	FAD	C10-C4X-N5	-3.21	118.25	124.81
5	D	527	FAD	O3P-PA-O1A	-3.17	101.15	110.70
3	C	522	NAG	C2-N2-C7	3.13	127.09	122.90
5	D	527	FAD	C5A-N7A-C8A	3.13	108.36	103.45
5	B	525	FAD	C5A-C4A-N3A	-3.09	122.46	126.72
5	C	527	FAD	C9A-C5X-N5	-3.08	119.19	122.45
5	D	527	FAD	O2A-PA-O3P	3.06	115.56	107.27
4	C	526	PHE	OXT-C-O	3.04	130.99	124.08
5	A	527	FAD	C5A-C4A-N3A	-3.01	122.57	126.72
5	C	527	FAD	C5A-N7A-C8A	2.99	108.15	103.45
4	D	526	PHE	CG-CB-CA	2.99	120.24	114.13
5	C	527	FAD	C2A-N3A-C4A	2.98	119.11	111.83
3	C	522	NAG	C1-C2-N2	2.98	115.13	110.43
3	A	522	NAG	O5-C5-C6	2.97	113.45	107.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	527	FAD	C2A-N3A-C4A	2.95	119.04	111.83
4	A	526	PHE	OXT-C-O	2.95	130.77	124.08
3	D	522	NAG	C1-O5-C5	2.92	116.10	112.19
5	A	527	FAD	C4A-C5A-N7A	-2.90	107.27	110.58
5	A	527	FAD	C4X-C10-N10	2.88	120.60	116.48
4	B	524	PHE	CG-CB-CA	2.87	120.01	114.13
5	B	525	FAD	C10-C4X-N5	-2.86	118.97	124.81
5	B	525	FAD	O3P-PA-O1A	-2.76	102.41	110.70
5	A	527	FAD	C2A-N3A-C4A	2.74	118.52	111.83
5	B	525	FAD	C5A-N7A-C8A	2.74	107.76	103.45
4	A	526	PHE	CG-CB-CA	2.73	119.71	114.13
5	C	527	FAD	O3P-P-O1P	-2.70	102.58	110.70
5	B	525	FAD	C4-C4X-N5	2.68	121.91	118.21
5	D	527	FAD	C4X-C10-N10	2.63	120.25	116.48
5	A	527	FAD	C9A-C5X-N5	-2.62	119.67	122.45
5	A	527	FAD	O3P-PA-O1A	-2.58	102.95	110.70
4	A	526	PHE	CB-CA-N	-2.56	101.61	111.40
5	D	527	FAD	C10-C4X-N5	-2.56	119.59	124.81
3	B	523	NAG	C1-O5-C5	2.53	115.58	112.19
5	B	525	FAD	C4X-C10-N10	2.50	120.06	116.48
5	A	527	FAD	C10-C4X-N5	-2.49	119.72	124.81
5	B	525	FAD	O4-C4-C4X	-2.48	119.99	126.53
5	C	527	FAD	C4-C4X-N5	2.43	121.57	118.21
5	C	527	FAD	N3A-C4A-N9A	2.41	131.27	127.17
5	C	527	FAD	C4A-C5A-N7A	-2.41	107.83	110.58
5	D	527	FAD	C5X-C9A-N10	2.41	120.14	117.97
5	A	527	FAD	C4A-N9A-C8A	2.40	108.26	105.74
5	A	527	FAD	O3P-P-O1P	-2.37	103.58	110.70
5	D	527	FAD	C4X-C4-N3	2.34	119.22	113.25
5	B	525	FAD	O3P-P-O1P	-2.34	103.67	110.70
5	A	527	FAD	C4X-C4-N3	2.32	119.17	113.25
5	A	527	FAD	C4-N3-C2	-2.32	121.53	125.64
5	D	527	FAD	C9A-C5X-N5	-2.32	120.00	122.45
3	C	522	NAG	C4-C3-C2	-2.31	107.63	111.02
3	A	522	NAG	C2-N2-C7	2.30	125.98	122.90
5	D	527	FAD	C2A-N1A-C6A	2.30	122.50	118.73
5	C	527	FAD	O3'-C3'-C2'	2.28	114.11	108.93
5	C	527	FAD	O4-C4-C4X	-2.26	120.56	126.53
5	B	525	FAD	C4X-C4-N3	2.26	119.00	113.25
5	C	527	FAD	O3P-PA-O1A	-2.25	103.92	110.70
5	B	525	FAD	C4A-C5A-N7A	-2.25	108.01	110.58
3	C	522	NAG	O3-C3-C2	2.25	114.07	109.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	527	FAD	C5X-C9A-N10	2.24	119.99	117.97
5	A	527	FAD	O4-C4-C4X	-2.24	120.62	126.53
5	D	527	FAD	O4-C4-C4X	-2.23	120.65	126.53
4	D	526	PHE	CB-CA-N	-2.19	103.03	111.40
5	D	527	FAD	C4A-C5A-N7A	-2.18	108.09	110.58
5	C	527	FAD	O2A-PA-O3P	2.18	113.15	107.27
5	C	527	FAD	C4X-C10-N10	2.15	119.56	116.48
5	A	527	FAD	C5'-C4'-C3'	-2.14	108.18	112.22
5	C	527	FAD	C2B-C3B-C4B	2.14	106.74	102.61
3	D	522	NAG	C3-C4-C5	2.10	114.04	110.23
5	D	527	FAD	C9-C9A-N10	-2.07	119.07	121.85
5	B	525	FAD	C5'-C4'-C3'	-2.06	108.33	112.22
5	B	525	FAD	C9A-C5X-N5	-2.04	120.29	122.45
5	B	525	FAD	C2A-N1A-C6A	2.01	122.02	118.73

There are no chirality outliers.

All (42) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	522	NAG	C3-C2-N2-C7
5	C	527	FAD	PA-O3P-P-O5'
5	D	527	FAD	PA-O3P-P-O5'
3	D	522	NAG	O5-C5-C6-O6
5	C	527	FAD	C2'-C3'-C4'-O4'
5	A	527	FAD	C2'-C3'-C4'-C5'
5	C	527	FAD	C2'-C3'-C4'-C5'
3	D	522	NAG	C4-C5-C6-O6
5	A	527	FAD	C2'-C3'-C4'-O4'
5	B	525	FAD	C2'-C3'-C4'-O4'
5	D	527	FAD	C2'-C3'-C4'-O4'
5	B	525	FAD	O3'-C3'-C4'-C5'
3	B	523	NAG	O5-C5-C6-O6
5	B	525	FAD	C2'-C3'-C4'-C5'
4	B	524	PHE	CA-CB-CG-CD1
4	B	524	PHE	CA-CB-CG-CD2
5	A	527	FAD	PA-O3P-P-O5'
5	B	525	FAD	PA-O3P-P-O5'
5	C	527	FAD	O3'-C3'-C4'-O4'
4	A	526	PHE	OXT-C-CA-CB
4	B	524	PHE	N-CA-CB-CG
4	A	526	PHE	O-C-CA-CB
5	A	527	FAD	O3'-C3'-C4'-C5'

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Mol	Chain	Res	Type	Atoms
3	B	523	NAG	C4-C5-C6-O6
5	D	527	FAD	C2'-C3'-C4'-C5'
4	C	526	PHE	OXT-C-CA-CB
4	C	526	PHE	O-C-CA-CB
5	C	527	FAD	O3'-C3'-C4'-C5'
5	B	525	FAD	O3'-C3'-C4'-O4'
4	A	526	PHE	CA-CB-CG-CD1
4	B	524	PHE	OXT-C-CA-CB
4	D	526	PHE	OXT-C-CA-CB
4	A	526	PHE	CA-CB-CG-CD2
4	B	524	PHE	O-C-CA-CB
4	D	526	PHE	O-C-CA-CB
4	A	526	PHE	OXT-C-CA-N
5	A	527	FAD	O4B-C4B-C5B-O5B
5	D	527	FAD	O3'-C3'-C4'-O4'
4	C	526	PHE	CA-CB-CG-CD1
4	D	526	PHE	O-C-CA-N
5	B	525	FAD	O4B-C4B-C5B-O5B
4	D	526	PHE	OXT-C-CA-N

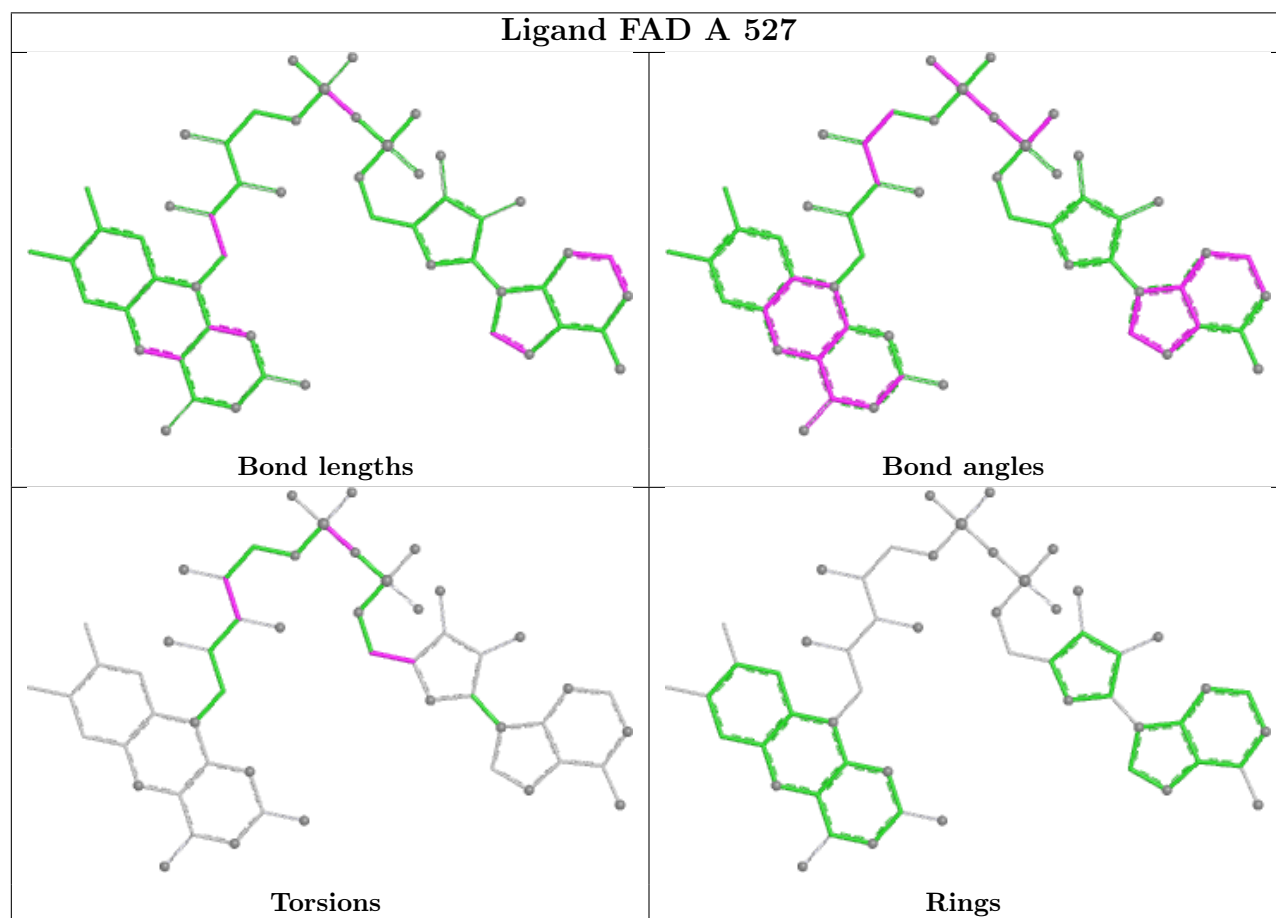
There are no ring outliers.

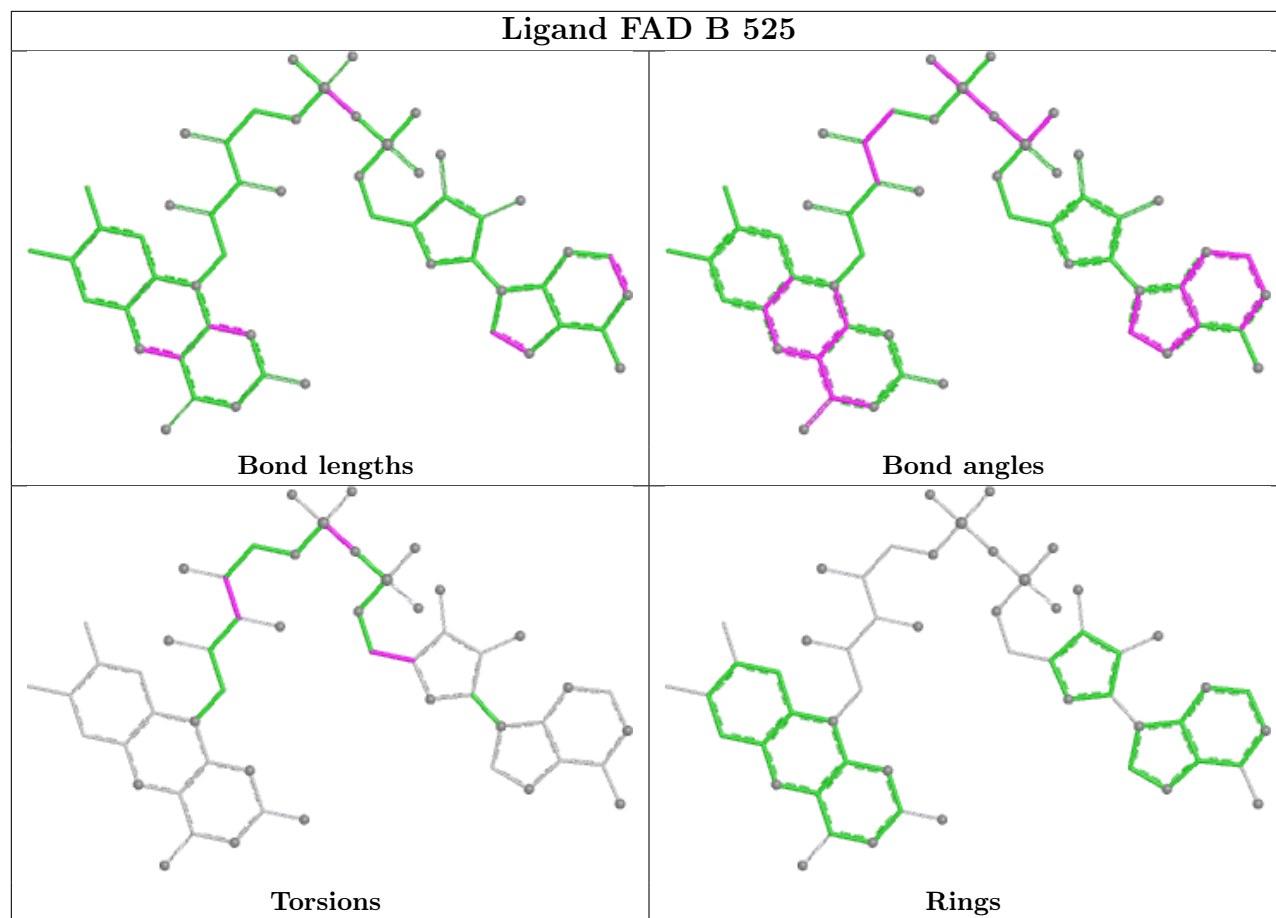
7 monomers are involved in 8 short contacts:

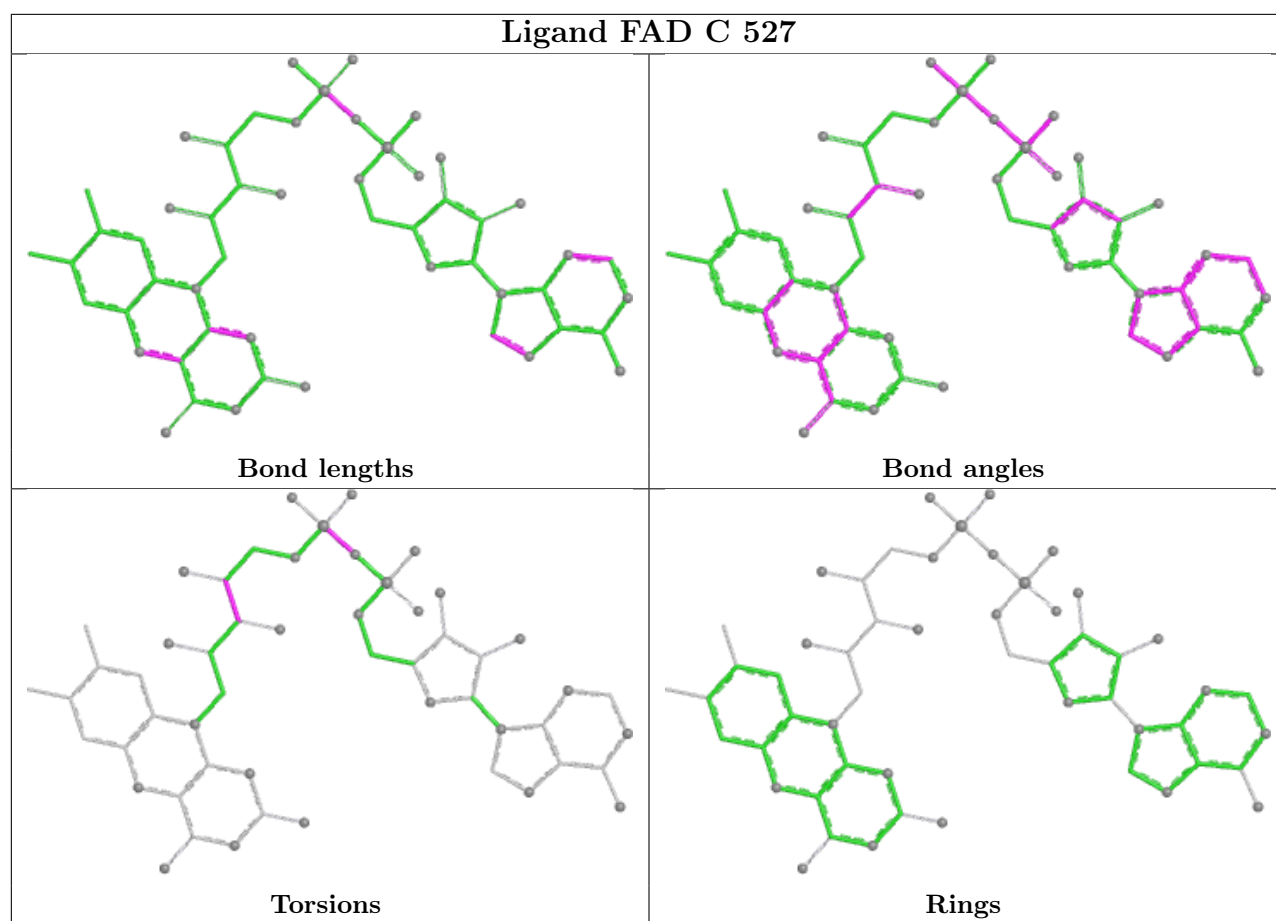
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	527	FAD	2	0
5	B	525	FAD	1	0
4	C	526	PHE	1	0
4	A	526	PHE	1	0
3	C	522	NAG	1	0
5	C	527	FAD	3	0
5	D	527	FAD	1	0

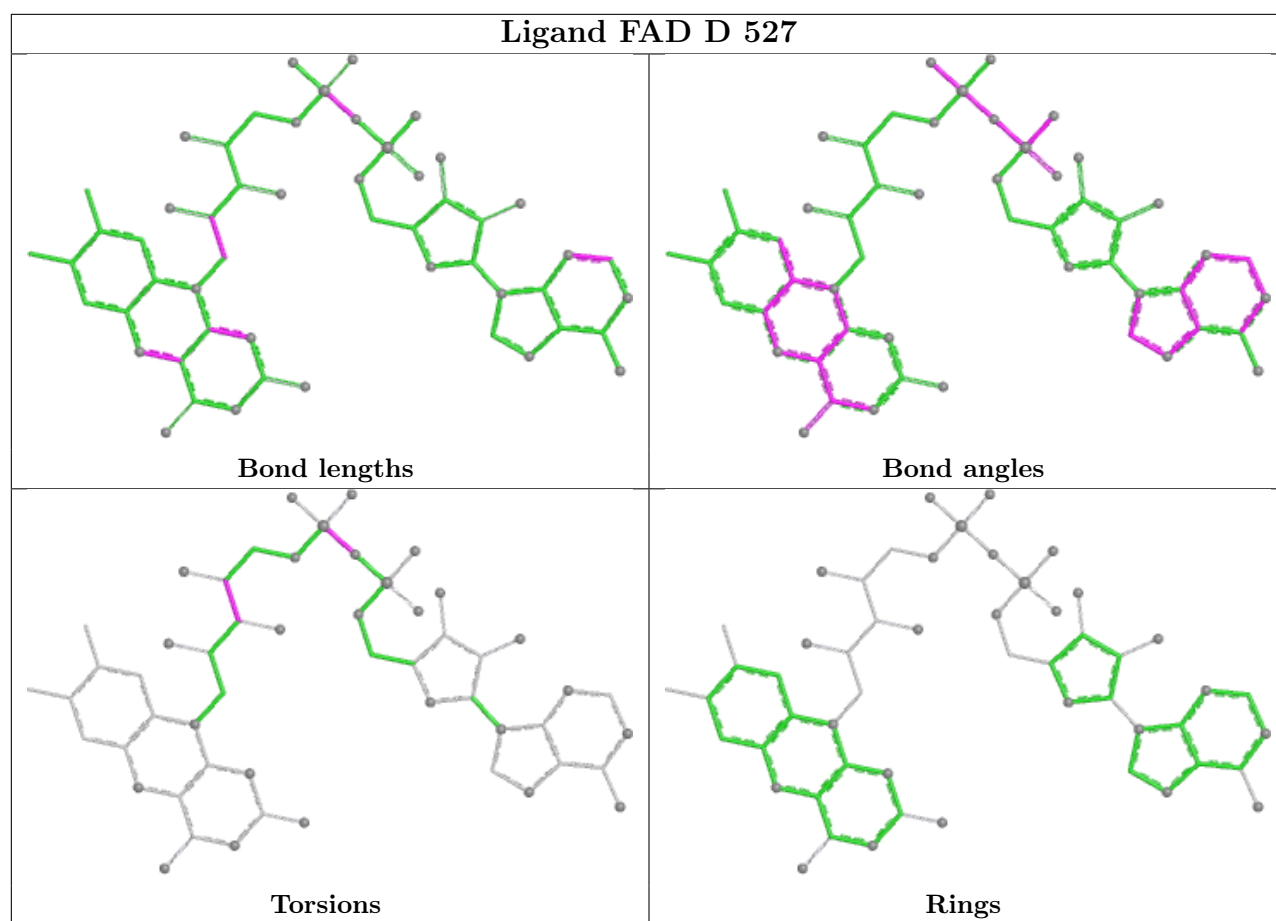
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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