

# wwPDB X-ray Structure Validation Summary Report (i)

#### Dec 9, 2023 - 05:46 pm GMT

PDB ID	:	1H83
Title	:	STRUCTURE OF POLYAMINE OXIDASE IN COMPLEX WITH 1,8-
		DIAMINOOCTANE
Authors	:	Binda, C.; Coda, A.; Angelini, R.; Federico, R.; Ascenzi, P.; Mattevi, A.
Deposited on	:	2001-01-24
Resolution	:	1.90  Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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:

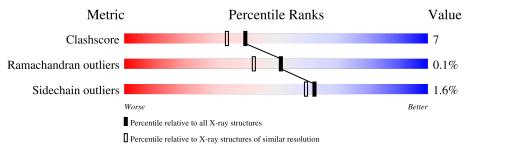
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as $541$ be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
$\mathrm{EDS}$	:	NOT EXECUTED
buster-report	:	1.1.7(2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber $(2001)$
Ideal geometry (DNA, RNA)		
Validation Pipeline (wwPDB-VP)	:	2.36

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $X\text{-}RAY \, DIFFRACTION$ 

The reported resolution of this entry is 1.90 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760(1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)

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	А	472	80%	16%	•••
1	В	472	84%	13%	•••
1	С	472	85%	12%	•••
2	D	2	100%		
2	Е	2	100%		
3	F	5	40% 60%		



# 2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 12183 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.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	Δ	459	Total	С	Ν	0	$\mathbf{S}$	65	0	0
	A		3684	2353	621	696	14	05	0	0
1	В	462	Total	С	Ν	0	S	66	0	0
	D	402	3715	2374	627	700	14			
1	C	462	Total	С	Ν	0	S	58	0	0
			3715	2374	627	700	14			

• Molecule 1 is a protein called POLYAMINE OXIDASE.

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



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
2	D	2	$\begin{bmatrix} Total & C & N & O \\ 28 & 16 & 2 & 10 \end{bmatrix}$	0	0	0
2	E	2	Total         C         N         O           28         16         2         10	0	0	0

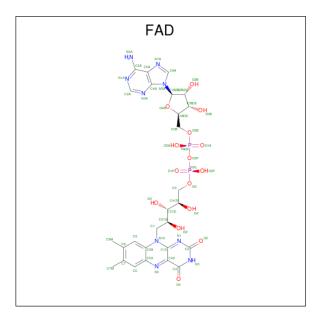
• Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-alpha-D-mannopyran ose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-3)]2-ac etamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	F	5	Total 60 3	C N 34 2	O 24	0	0	0

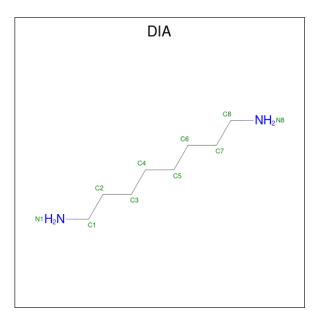


• Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
4	Δ	1	Total	С	Ν	Ο	Р	0	0
4	4 A	1	53	27	9	15	2	0	0
4	В	1	Total	С	Ν	Ο	Р	0	0
4	4 D	1	53	27	9	15	2	0	
4	С	1	Total	С	Ν	Ο	Р	0	0
4	U	1	53	27	9	15	2		0

• Molecule 5 is OCTANE 1,8-DIAMINE (three-letter code: DIA) (formula:  $C_8H_{20}N_2$ ).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	1	Total         C         N           20         16         4	0	1
5	В	1	Total         C         N           20         16         4	0	1
5	С	1	Total         C         N           20         16         4	0	1

• Molecule 6 is water.

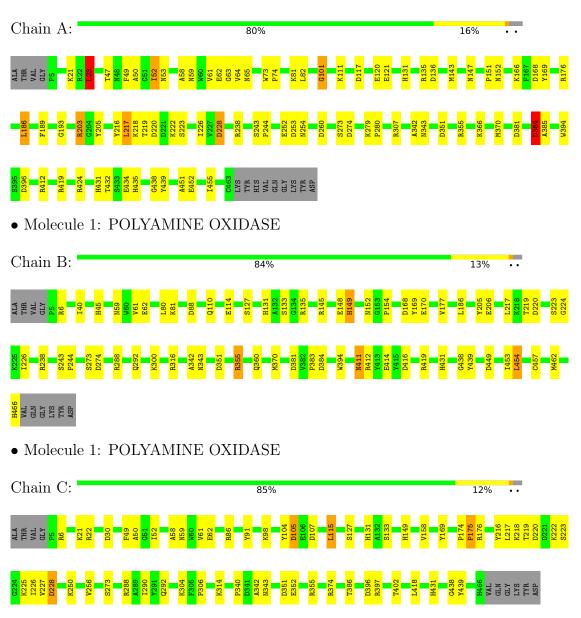
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	223	Total O 223 223	0	0
6	В	242	Total O 242 242	0	0
6	С	269	Total O 269 269	0	0



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

• Molecule 1: POLYAMINE OXIDASE



Chain D:	100%		
NAG2 NAG2			
• Molecule 2: 2-acetamido-2-de opyranose	eoxy-beta-D-glucopyran	ose-(1-4)-2-acetamide	o-2-deoxy-beta-D-gluc
Chain E:	100%		
NAG1 NAG2			
• Melecule 2. elube D. menuer	(1.6) alpha D	(1, 4)	2 acetomide 2 decem

 $\label{eq:mannopyranose-(1-6)-alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-D-fucopyranose-(1-4)]2-acetam$ 

Chain F:	40%	60%
NAG1 NAC2 MAN3 MAN4 FCA5		



## 4 Data and refinement statistics (i)

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

Property	Value	Source	
Space group	P 65 2 2	Depositor	
Cell constants	184.69Å 184.69Å 282.22Å	Depositor	
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $120.00^{\circ}$	Depositor	
Resolution (Å)	20.00 - 1.90	Depositor	
% Data completeness	97.9 (20.00-1.90)	Depositor	
(in resolution range)	51.5 (20.00 1.50)	Depositor	
$R_{merge}$	0.08	Depositor	
$R_{sym}$	(Not available)	Depositor	
Refinement program	TNT 5D	Depositor	
$R, R_{free}$	0.196 , $0.237$	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	12183	wwPDB-VP	
Average B, all atoms $(Å^2)$	15.0	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: FCA, DIA, FAD, MAN, 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	ol Chain Bond lengths		Bond angles		
MOI	Unam	RMSZ	# Z  > 5	RMSZ	# Z  > 5
1	А	0.55	0/3775	1.21	15/5116~(0.3%)
1	В	0.55	0/3808	1.21	15/5160~(0.3%)
1	С	0.56	0/3808	1.22	17/5160~(0.3%)
All	All	0.56	0/11391	1.21	47/15436~(0.3%)

There are no bond length outliers.

The worst 5 of 47 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	В	355	ARG	NE-CZ-NH2	-10.02	115.29	120.30
1	А	355	ARG	NE-CZ-NH2	-8.79	115.91	120.30
1	С	176	ARG	NE-CZ-NH2	-8.47	116.06	120.30
1	А	424	ARG	NE-CZ-NH1	8.23	124.42	120.30
1	С	176	ARG	NE-CZ-NH1	8.12	124.36	120.30

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	А	3684	0	3585	56	0
1	В	3715	0	3614	47	0
1	С	3715	0	3614	41	0

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Conti	Continuea from previous page						
Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes	
2	D	28	0	26	5	0	
2	Е	28	0	26	4	0	
3	F	60	0	53	6	0	
4	А	53	0	31	2	0	
4	В	53	0	31	1	0	
4	С	53	0	31	4	0	
5	А	20	0	40	5	0	
5	В	20	0	40	4	0	
5	С	20	0	40	5	0	
6	А	223	0	0	7	1	
6	В	242	0	0	4	0	
6	С	269	0	0	3	2	
All	All	12183	0	11131	158	2	

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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 7.

The worst 5 of 158 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)
5:C:590[A]:DIA:HC82	6:C:2227:HOH:O	1.55	1.06
1:A:273:SER:O	1:A:274:ASP:HB2	1.72	0.88
3:F:1:NAG:C4	3:F:2:NAG:C1	2.54	0.86
2:E:1:NAG:C4	2:E:2:NAG:C1	2.55	0.84
1:C:91:TYR:OH	1:C:314:LYS:HE2	1.80	0.81

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:C:2094:HOH:O	6:C:2094:HOH:O[10_665]	0.73	1.47
6:A:2131:HOH:O	6:C:2058:HOH:O[6_654]	1.30	0.90

## 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	Perce	ntiles
1	А	457/472~(97%)	438 (96%)	18 (4%)	1 (0%)	47	38
1	В	460/472~(98%)	444 (96%)	16 (4%)	0	100	100
1	С	460/472~(98%)	442 (96%)	18 (4%)	0	100	100
All	All	1377/1416~(97%)	1324 (96%)	52 (4%)	1 (0%)	51	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	101	GLY

#### 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	А	394/404~(98%)	387~(98%)	7~(2%)	59 55
1	В	397/404~(98%)	390~(98%)	7 (2%)	59 55
1	С	397/404~(98%)	392~(99%)	5 (1%)	69 68
All	All	1188/1212 (98%)	1169 (98%)	19 (2%)	62 60

 $5~{\rm of}~19$  residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
1	С	59	ASN
1	С	149	HIS
1	С	175	PRO
1	С	127	SER
1	В	127	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 19 such sidechains are listed below:



Mol	Chain	Res	Type
1	В	466	HIS
1	С	360	GLN
1	С	431	HIS
1	С	152	ASN
1	В	152	ASN

#### 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)

9 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	Turne	Chain	Res	Link	Bo	ond leng	ths	В	ond ang	les
10101	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z >2
2	NAG	D	1	1,2	14,14,15	0.94	1 (7%)	17,19,21	1.44	4 (23%)
2	NAG	D	2	2	14,14,15	0.88	1 (7%)	17,19,21	1.77	4 (23%)
2	NAG	Е	1	1,2	14,14,15	1.21	2 (14%)	17,19,21	1.58	3 (17%)
2	NAG	Е	2	2	14,14,15	0.95	1 (7%)	17,19,21	1.33	2 (11%)
3	NAG	F	1	3,1	14,14,15	1.06	2 (14%)	17,19,21	2.20	7 (41%)
3	NAG	F	2	3	14,14,15	0.96	1 (7%)	17,19,21	1.53	4 (23%)
3	MAN	F	3	3	11,11,12	0.72	0	$15,\!15,\!17$	<mark>3.03</mark>	5 (33%)
3	MAN	F	4	3	11,11,12	0.55	0	15, 15, 17	1.59	3 (20%)
3	FCA	F	5	3	10,10,11	1.21	1 (10%)	14,14,16	2.12	6 (42%)

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



Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	D	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	D	2	2	-	2/6/23/26	0/1/1/1
2	NAG	Е	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	Е	2	2	-	2/6/23/26	0/1/1/1
3	NAG	F	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	MAN	F	3	3	-	0/2/19/22	0/1/1/1
3	MAN	F	4	3	-	2/2/19/22	0/1/1/1
3	FCA	F	5	3	-	-	0/1/1/1

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms		Observed(Å)	Ideal(Å)
3	F	2	NAG	O5-C1	-2.93	1.39	1.43
2	Е	1	NAG	C1-C2	-2.86	1.48	1.52
3	F	5	FCA	C2-C3	-2.69	1.48	1.52
2	Е	1	NAG	O5-C1	-2.57	1.39	1.43
2	Е	2	NAG	O5-C1	-2.48	1.39	1.43

The worst 5 of 38 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	F	3	MAN	O2-C2-C3	10.34	130.85	110.14
2	D	2	NAG	C2-N2-C7	-5.10	115.64	122.90
3	F	1	NAG	C1-C2-N2	-4.66	102.53	110.49
3	F	5	FCA	O5-C1-C2	-4.40	103.98	110.77
3	F	4	MAN	C1-C2-C3	-4.32	104.35	109.67

There are no chirality outliers.

5 of 14 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	Ε	2	NAG	C4-C5-C6-O6
2	D	1	NAG	C4-C5-C6-O6
2	D	2	NAG	C4-C5-C6-O6
2	Е	1	NAG	C4-C5-C6-O6
3	F	2	NAG	C4-C5-C6-O6

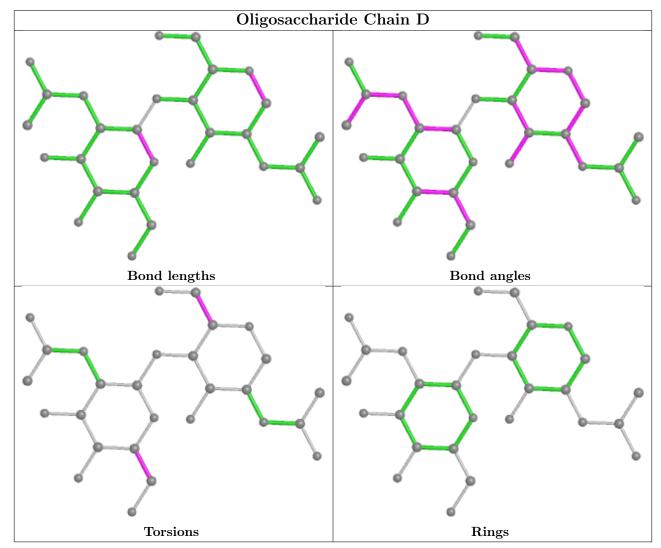


There are no ring outliers.

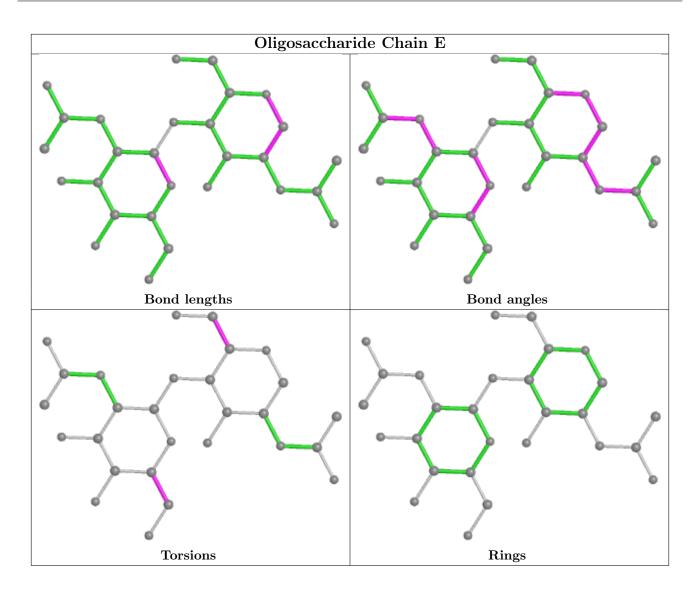
7 monomers are involved in 15 sl	nort contacts:
----------------------------------	----------------

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

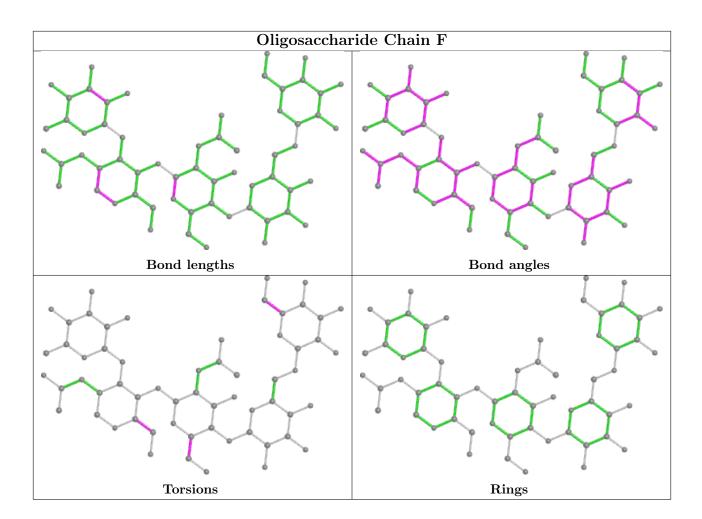
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











## 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).

Mal	Mol Type C		Res	Link	Bo	ond leng	ths	E	ond ang	gles
NIOI	туре	Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
4	FAD	А	579	-	$53,\!58,\!58$	0.87	2 (3%)	68,89,89	1.47	11 (16%)
5	DIA	А	590[B]	-	$9,\!9,\!9$	0.51	0	8,8,8	0.30	0
4	FAD	С	579	-	$53,\!58,\!58$	1.05	1 (1%)	68,89,89	1.29	8 (11%)
5	DIA	С	590[B]	-	$9,\!9,\!9$	0.50	0	8,8,8	0.35	0
4	FAD	В	579	-	$53,\!58,\!58$	0.80	1 (1%)	68,89,89	1.26	7 (10%)
5	DIA	В	590[A]	-	$9,\!9,\!9$	0.48	0	8,8,8	0.44	0



Mol	Type	Chain	Res	Link	Bo	Bond lengths			Bond angles		
IVIOI	туре	Unam	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z  > 2	
5	DIA	А	590[A]	-	$9,\!9,\!9$	0.47	0	8,8,8	0.28	0	
5	DIA	С	590[A]	-	$9,\!9,\!9$	0.49	0	8,8,8	0.42	0	
5	DIA	В	590[B]	-	$9,\!9,\!9$	0.51	0	8,8,8	0.35	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	FAD	А	579	-	-	2/30/50/50	0/6/6/6
5	DIA	А	590[B]	-	-	3/7/7/7	-
4	FAD	С	579	-	-	1/30/50/50	0/6/6/6
5	DIA	С	590[B]	-	-	4/7/7/7	-
4	FAD	В	579	-	-	5/30/50/50	0/6/6/6
5	DIA	В	590[A]	-	-	5/7/7/7	-
5	DIA	А	590[A]	-	-	4/7/7/7	-
5	DIA	С	590[A]	-	-	6/7/7/7	-
5	DIA	В	590[B]	-	-	3/7/7/7	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
4	С	579	FAD	C4X-N5	3.49	1.37	1.30
4	В	579	FAD	C4X-N5	2.55	1.35	1.30
4	А	579	FAD	C4X-N5	2.50	1.35	1.30
4	А	579	FAD	C5X-N5	-2.10	1.35	1.39

The worst 5 of 26 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
4	А	579	FAD	C9A-C5X-N5	-5.36	116.60	122.43
4	С	579	FAD	O4B-C1B-C2B	-4.21	100.77	106.93
4	В	579	FAD	O4B-C1B-C2B	-3.94	101.17	106.93
4	А	579	FAD	O4'-C4'-C3'	3.12	116.68	109.10
4	А	579	FAD	C6-C5X-C9A	3.07	123.28	118.94

There are no chirality outliers.

5 of 33 torsion outliers are listed below:



Mol	Chain	Res	Type	Atoms
4	А	579	FAD	C5B-O5B-PA-O1A
5	А	590[B]	DIA	C2-C3-C4-C5
5	В	590[A]	DIA	C3-C4-C5-C6
5	В	590[B]	DIA	C2-C3-C4-C5
5	В	590[A]	DIA	C2-C3-C4-C5

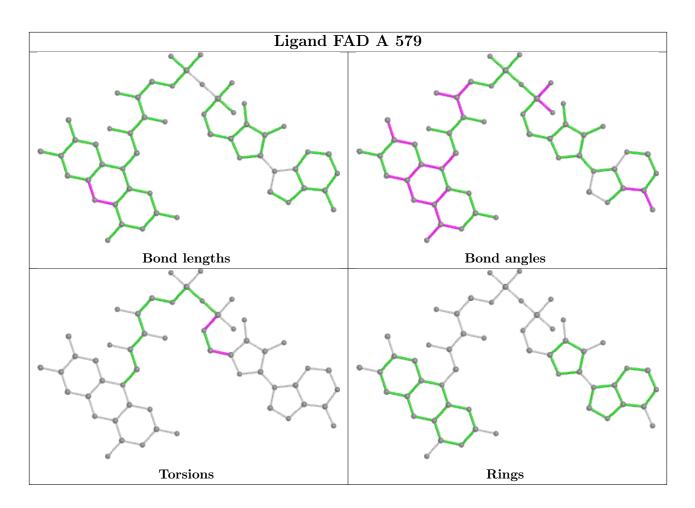
There are no ring outliers.

9 monomers are involved in 21 short contacts:

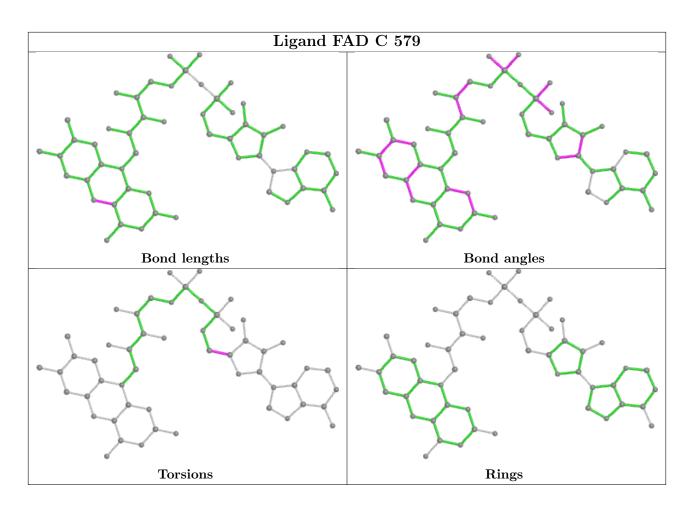
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	А	579	FAD	2	0
5	А	590[B]	DIA	1	0
4	С	579	FAD	4	0
5	С	590[B]	DIA	2	0
4	В	579	FAD	1	0
5	В	590[A]	DIA	3	0
5	А	590[A]	DIA	4	0
5	С	590[A]	DIA	3	0
5	В	590[B]	DIA	1	0

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 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.



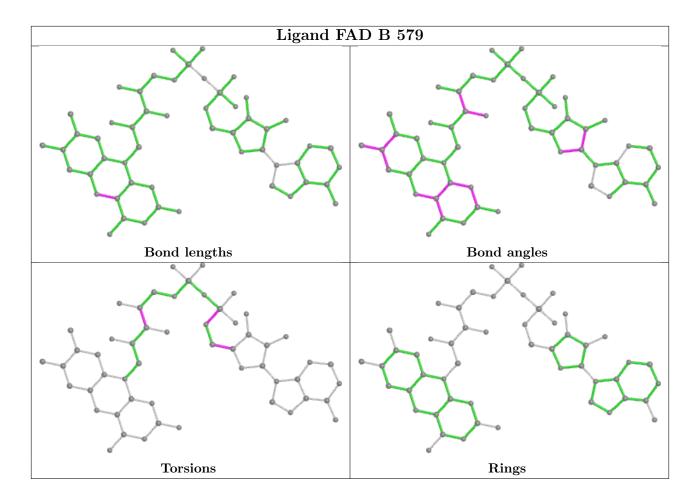












## 5.7 Other polymers (i)

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.

