

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

#### Aug 21, 2023 – 08:55 PM EDT

PDB ID	:	2PGZ
Title	:	Crystal structure of Cocaine bound to an ACh-Binding Protein
Authors	:	Hansen, S.B.; Taylor, P.
Deposited on		
Resolution	:	1.76 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:

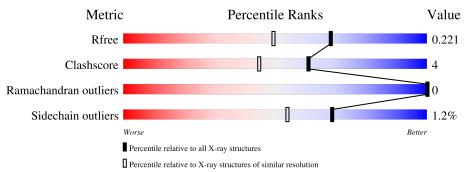
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
$\mathrm{EDS}$	:	2.35
buster-report	:	1.1.7(2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber $(2001)$
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

# 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 1.76 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R <sub>free</sub>	130704	2340(1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)

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%

Mol	Chain	Length	Quality of chain	
1	А	230	87%	5% 8%
1	В	230	83%	8% 8%
1	С	230	82%	9% • 8%
1	D	230	86%	7% 7%
1	Е	230	81%	10% 9%
2	F	5	20% 80%	



#### 2 PGZ

# 2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 9900 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	211	Total	С	Ν	0	S	0	6	0
	A	211	1721	1085	285	343	8	0	0	0
1	В	211	Total	С	Ν	0	S	0	2	0
	D	211	1695	1072	276	339	8	0		0
1	С	211	Total	С	Ν	0	S	0	7	0
	U	211	1728	1091	286	341	10	0		
1	D	214	Total	С	Ν	0	S	0	5	0
	D	214	1742	1100	285	348	9	0	5	0
1	Е	209	Total	С	Ν	0	S	0	9	0
	Ľ	209	1718	1086	282	341	9	0	9	0

• Molecule 1 is a protein called Soluble acetylcholine receptor.

There are 55 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	-8	ASP	-	cloning artifact	UNP Q8WSF8
А	-7	TYR	-	cloning artifact	UNP Q8WSF8
А	-6	LYS	-	cloning artifact	UNP Q8WSF8
А	-5	ASP	-	cloning artifact	UNP Q8WSF8
А	-4	ASP	-	cloning artifact	UNP Q8WSF8
А	-3	ASP	-	cloning artifact	UNP Q8WSF8
А	-2	ASP	-	cloning artifact	UNP Q8WSF8
А	-1	LYS	-	cloning artifact	UNP Q8WSF8
А	0	LEU	-	cloning artifact	UNP Q8WSF8
А	220	SER	-	cloning artifact	UNP Q8WSF8
А	221	ARG	-	cloning artifact	UNP Q8WSF8
В	-8	ASP	-	cloning artifact	UNP Q8WSF8
В	-7	TYR	-	cloning artifact	UNP Q8WSF8
В	-6	LYS	-	cloning artifact	UNP Q8WSF8
В	-5	ASP	-	cloning artifact	UNP Q8WSF8
В	-4	ASP	-	cloning artifact	UNP Q8WSF8
В	-3	ASP	-	cloning artifact	UNP Q8WSF8
В	-2	ASP	-	cloning artifact	UNP Q8WSF8
В	-1	LYS	-	cloning artifact	UNP Q8WSF8

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Chain	Residue	Modelled	Actual	Comment	Reference
В	0	LEU	-	cloning artifact	UNP Q8WSF8
В	220	SER	-	cloning artifact	UNP Q8WSF8
В	221	ARG	-	cloning artifact	UNP Q8WSF8
С	-8	ASP	-	cloning artifact	UNP Q8WSF8
С	-7	TYR	-	cloning artifact	UNP Q8WSF8
С	-6	LYS	-	cloning artifact	UNP Q8WSF8
С	-5	ASP	-	cloning artifact	UNP Q8WSF8
С	-4	ASP	-	cloning artifact	UNP Q8WSF8
С	-3	ASP	-	cloning artifact	UNP Q8WSF8
С	-2	ASP	-	cloning artifact	UNP Q8WSF8
С	-1	LYS	-	cloning artifact	UNP Q8WSF8
С	0	LEU	-	cloning artifact	UNP Q8WSF8
С	220	SER	-	cloning artifact	UNP Q8WSF8
С	221	ARG	-	cloning artifact	UNP Q8WSF8
D	-8	ASP	-	cloning artifact	UNP Q8WSF8
D	-7	TYR	-	cloning artifact	UNP Q8WSF8
D	-6	LYS	-	cloning artifact	UNP Q8WSF8
D	-5	ASP	-	cloning artifact	UNP Q8WSF8
D	-4	ASP	-	cloning artifact	UNP Q8WSF8
D	-3	ASP	-	cloning artifact	UNP Q8WSF8
D	-2	ASP	-	cloning artifact	UNP Q8WSF8
D	-1	LYS	-	cloning artifact	UNP Q8WSF8
D	0	LEU	-	cloning artifact	UNP Q8WSF8
D	220	SER	-	cloning artifact	UNP Q8WSF8
D	221	ARG	-	cloning artifact	UNP Q8WSF8
E	-8	ASP	-	cloning artifact	UNP Q8WSF8
E	-7	TYR	-	cloning artifact	UNP Q8WSF8
E	-6	LYS	-	cloning artifact	UNP Q8WSF8
E	-5	ASP	-	cloning artifact	UNP Q8WSF8
E	-4	ASP	-	cloning artifact	UNP Q8WSF8
E	-3	ASP	-	cloning artifact	UNP Q8WSF8
Е	-2	ASP	-	cloning artifact	UNP Q8WSF8
Е	-1	LYS	-	cloning artifact	UNP Q8WSF8
Е	0	LEU	-	cloning artifact	UNP Q8WSF8
Е	220	SER	-	cloning artifact	UNP Q8WSF8
Е	221	ARG	_	cloning artifact	UNP Q8WSF8

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

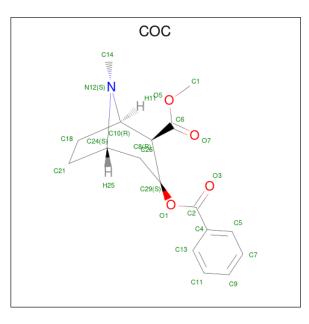






Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
2	F	5	Total 61	C 34	-	O 25	0	0	0

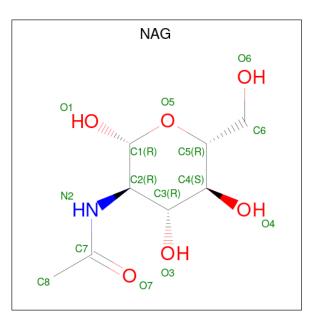
• Molecule 3 is COCAINE (three-letter code: COC) (formula:  $C_{17}H_{21}NO_4$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	А	1	Total 22			0 4	0	0
3	D	1	Total 22	C 17	N 1	0 4	0	0

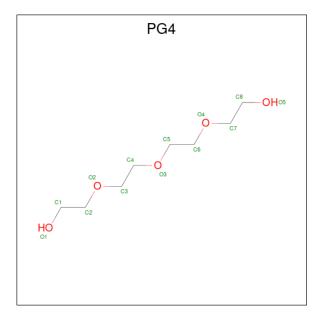
• Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).





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

• Molecule 5 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula:  $C_8H_{18}O_5$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	В	1	Total         C         O           13         8         5	0	0
5	С	1	Total         C         O           10         6         4	0	0
5	Е	1	Total         C         O           13         8         5	0	0



• Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	251	Total O 251 251	0	6
6	В	229	Total         O           229         229	0	2
6	С	230	Total         O           230         230	0	8
6	D	223	Total         O           223         223	0	2
6	Е	208	Total         O           208         208	0	6

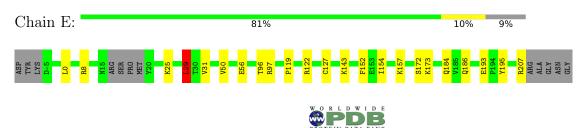


# 3 Residue-property plots (i)

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.

- Chain A: 87% 5% 8% • Molecule 1: Soluble acetylcholine receptor Chain B: 83% 8% 8% SER • Molecule 1: Soluble acetylcholine receptor Chain C: 82% 9% 8% ASP SER • Molecule 1: Soluble acetylcholine receptor Chain D: 86% 7% 7% ALA ASN ASN
- Molecule 1: Soluble acetylcholine receptor

• Molecule 1: Soluble acetylcholine receptor



#### PHE PHE ARG ASN LEU PHE ASP SER ARG

 • Molecule 2: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)] beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F: 20% 80%

NAG1 NAG2 BMA3 MAN4 MAN5 MAN5



# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	86.82Å 115.59Å 130.88Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 - 1.76	Depositor
Resolution (A)	47.62 - 1.76	EDS
% Data completeness	99.5 (50.00-1.76)	Depositor
(in resolution range)	99.5 (47.62 - 1.76)	EDS
R <sub>merge</sub>	0.06	Depositor
R <sub>sym</sub>	0.05	Depositor
$< I/\sigma(I) > 1$	$2.37 (at 1.76 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
D D.	0.180 , $0.210$	Depositor
$R, R_{free}$	0.191 , $0.221$	DCC
$R_{free}$ test set	1314 reflections $(1.01\%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	25.8	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.32 , $36.8$	EDS
L-test for twinning <sup>2</sup>	$ \langle L  \rangle = 0.51, \langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	9900	wwPDB-VP
Average B, all atoms $(Å^2)$	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.27% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

# 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: BMA, PG4, NAG, COC, MAN

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	Bo	Bond lengths		ond angles
	Unam	RMSZ	RMSZ $\# Z  > 5$		# Z  > 5
1	А	0.53	0/1777	0.61	0/2418
1	В	0.54	0/1740	0.63	0/2369
1	С	0.49	0/1787	0.64	0/2428
1	D	0.61	1/1796~(0.1%)	0.66	0/2443
1	Е	0.65	3/1787~(0.2%)	0.76	3/2430~(0.1%)
All	All	0.57	4/8887~(0.0%)	0.66	3/12088~(0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	Ε	8	ARG	NE-CZ	11.91	1.48	1.33
1	Е	8	ARG	CG-CD	8.34	1.72	1.51
1	D	-8	ASP	N-CA	6.38	1.59	1.46
1	Е	8	ARG	CZ-NH2	5.43	1.40	1.33

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	Ε	8	ARG	NE-CZ-NH1	-16.74	111.93	120.30
1	Е	8	ARG	NE-CZ-NH2	10.15	125.37	120.30
1	Е	29	LEU	CA-CB-CG	5.31	127.52	115.30

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



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	1721	0	1659	9	0
1	В	1695	0	1615	19	0
1	С	1728	0	1675	21	0
1	D	1742	0	1675	15	1
1	Ε	1718	0	1661	16	1
2	F	61	0	52	0	0
3	А	22	0	21	1	0
3	D	22	0	21	0	0
4	В	14	0	13	0	0
5	В	13	0	18	3	0
5	С	10	0	13	0	0
5	Ε	13	0	18	2	0
6	А	251	0	0	3	0
6	В	229	0	0	8	0
6	С	230	0	0	7	0
6	D	223	0	0	2	0
6	Ε	208	0	0	5	0
All	All	9900	0	8441	77	1

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.

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 4.

The worst 5 of 77 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:C:97[B]:ARG:HG3	6:C:633[B]:HOH:O	1.45	1.12
1:A:97[B]:ARG:NE	6:A:650[B]:HOH:O	1.65	1.00
5:B:403:PG4:H62	6:B:611:HOH:O	1.78	0.83
1:A:97[B]:ARG:CZ	6:A:650[B]:HOH:O	2.16	0.79
1:C:207[A]:ARG:HD3	6:C:635[A]:HOH:O	1.82	0.78

All (1) 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)
1:D:-7:TYR:OH	1:E:173:LYS:O[2_554]	2.12	0.08



### 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	212/230~(92%)	210 (99%)	2(1%)	0	100	100
1	В	208/230~(90%)	206 (99%)	2(1%)	0	100	100
1	С	213/230~(93%)	208 (98%)	5(2%)	0	100	100
1	D	214/230~(93%)	212 (99%)	2(1%)	0	100	100
1	Е	213/230~(93%)	210 (99%)	3 (1%)	0	100	100
All	All	1060/1150~(92%)	1046 (99%)	14 (1%)	0	100	100

There are no Ramachandran outliers to report.

#### 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	hain Analysed Rotameric Outlier		Outliers	Percentiles
1	А	198/208~(95%)	196~(99%)	2(1%)	76 63
1	В	193/208~(93%)	190~(98%)	3~(2%)	62 45
1	С	199/208~(96%)	194 (98%)	5(2%)	47 25
1	D	200/208~(96%)	199 (100%)	1 (0%)	88 83
1	Ε	199/208~(96%)	196~(98%)	3~(2%)	65 49
All	All	989/1040~(95%)	975~(99%)	14 (1%)	69 52

5 of 14 residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	С	186	GLN
1	С	207[A]	ARG
1	Ε	186	GLN
1	Е	0	LEU
1	Е	29	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 14 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	15	ASN
1	D	74	ASN
1	Е	186	GLN
1	Е	15	ASN
1	Е	184	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)

5 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	Bo	ond leng	$\mathbf{ths}$	В	ond ang	les
	туре	Unam	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z  > 2
2	NAG	F	1	2,1	$14,\!14,\!15$	0.45	0	17,19,21	1.25	1 (5%)
2	NAG	F	2	2	14,14,15	0.33	0	17,19,21	1.03	1 (5%)
2	BMA	F	3	2	11,11,12	0.40	0	$15,\!15,\!17$	0.69	0
2	MAN	F	4	2	$11,\!11,\!12$	0.47	0	$15,\!15,\!17$	1.22	1 (6%)



ſ	Mol	Type	Chain	Dog	Link	Bond lengths			Bond angles		
	WIOI			Res		Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z >2
	2	MAN	F	5	2	11,11,12	0.53	0	$15,\!15,\!17$	0.99	2 (13%)

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	F	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	F	2	2	-	0/6/23/26	0/1/1/1
2	BMA	F	3	2	-	0/2/19/22	0/1/1/1
2	MAN	F	4	2	-	0/2/19/22	0/1/1/1
2	MAN	F	5	2	-	1/2/19/22	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	F	4	MAN	C1-O5-C5	3.99	117.60	112.19
2	F	1	NAG	C2-N2-C7	3.99	128.58	122.90
2	F	2	NAG	C1-C2-N2	-3.20	105.02	110.49
2	F	5	MAN	C1-O5-C5	2.41	115.46	112.19
2	F	5	MAN	C1-C2-C3	2.38	112.59	109.67

There are no chirality outliers.

All (3) torsion outliers are listed below:

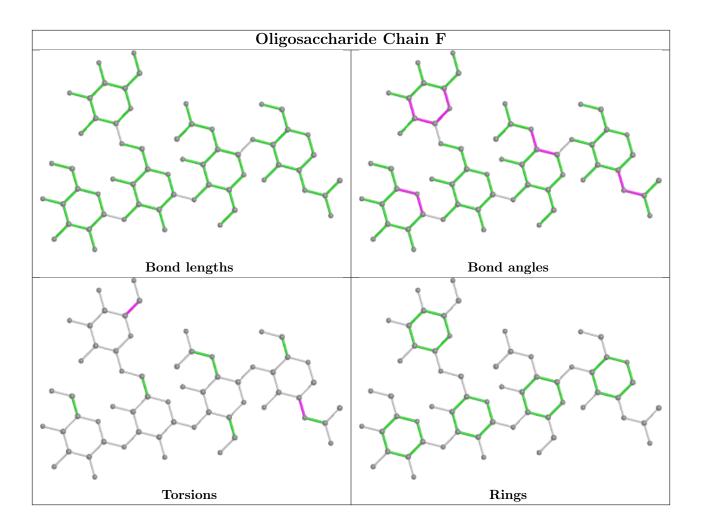
Mol	Chain	Res	Type	Atoms
2	F	1	NAG	C1-C2-N2-C7
2	F	5	MAN	C4-C5-C6-O6
2	F	1	NAG	C3-C2-N2-C7

There are no ring outliers.

No monomer is involved in short contacts.

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)

6 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	Mol Type Chain Res		Link	Bo	ond leng	ths	Bond angles			
IVIOI	туре	Chain	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z  > 2
5	PG4	В	403	-	12,12,12	1.54	3 (25%)	11,11,11	0.64	0
5	PG4	С	407	-	9,9,12	1.18	1 (11%)	8,8,11	0.82	0
3	COC	А	401	-	24,24,24	1.62	3 (12%)	32,34,34	2.95	8 (25%)
4	NAG	В	402	1	14,14,15	<mark>3.39</mark>	4 (28%)	17,19,21	3.04	5 (29%)
3	COC	D	401	-	24,24,24	1.76	3 (12%)	32,34,34	2.55	12 (37%)



Mol Type		Chain	Res	Link	Bond lengths			Bond angles		
IVIOI	туре	Ullalli	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
5	PG4	Е	403	-	12,12,12	0.52	0	11,11,11	0.27	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	PG4	В	403	-	-	7/10/10/10	-
5	PG4	С	407	-	-	3/7/7/10	-
3	COC	А	401	-	-	0/14/39/39	0/4/3/3
4	NAG	В	402	1	-	2/6/23/26	0/1/1/1
3	COC	D	401	-	-	1/14/39/39	0/4/3/3
5	PG4	Ε	403	-	_	8/10/10/10	_

The worst 5 of 14 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
4	В	402	NAG	C7-N2	8.95	1.65	1.34
4	В	402	NAG	O7-C7	7.06	1.39	1.23
3	D	401	COC	O5-C6	5.40	1.46	1.33
3	D	401	COC	O1-C2	5.31	1.45	1.34
3	А	401	COC	O1-C2	4.60	1.44	1.34

The worst 5 of 25 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	А	401	COC	C18-C10-N12	-12.54	92.26	105.18
4	В	402	NAG	C2-N2-C7	-9.43	109.47	122.90
3	D	401	COC	C18-C10-N12	-8.16	96.77	105.18
4	В	402	NAG	C8-C7-N2	-6.42	105.23	116.10
3	D	401	COC	O5-C6-C8	5.73	119.38	111.03

There are no chirality outliers.

5 of 21 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	С	407	PG4	O4-C7-C8-O5
4	В	402	NAG	O7-C7-N2-C2
5	Е	403	PG4	O2-C3-C4-O3
5	С	407	PG4	O2-C3-C4-O3

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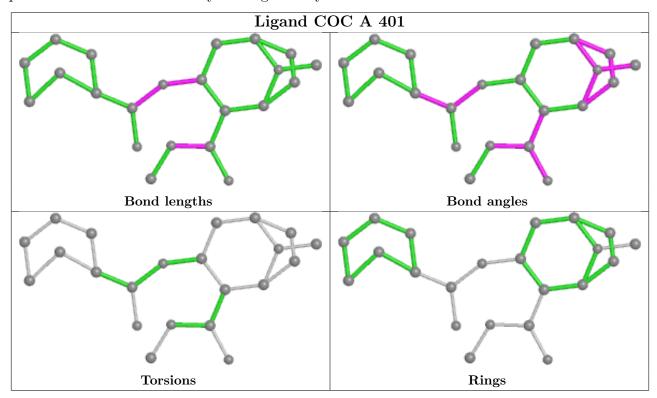
Mol	Chain	Res	Type	Atoms
4	В	402	NAG	O5-C5-C6-O6

There are no ring outliers.

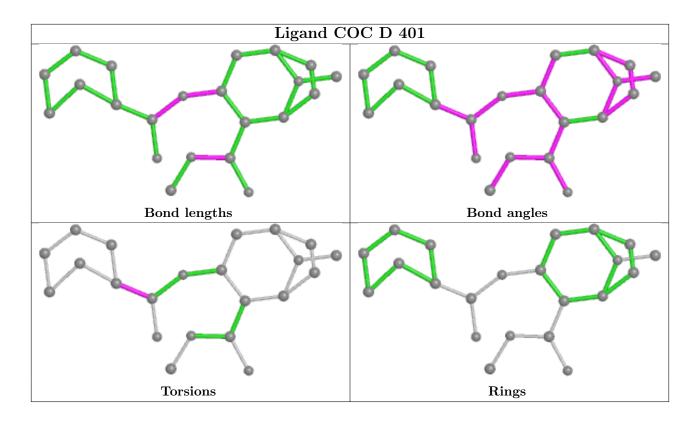
3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	В	403	PG4	3	0
3	А	401	COC	1	0
5	Е	403	PG4	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.







### 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)

Unable to reproduce the depositors R factor - this section is therefore empty.

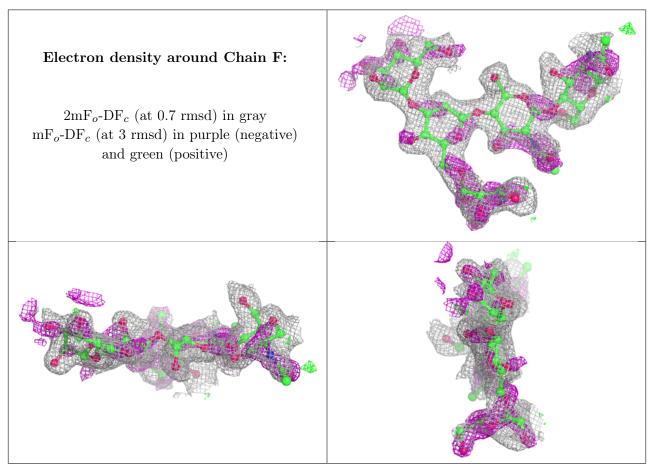
## 6.2 Non-standard residues in protein, DNA, RNA chains (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

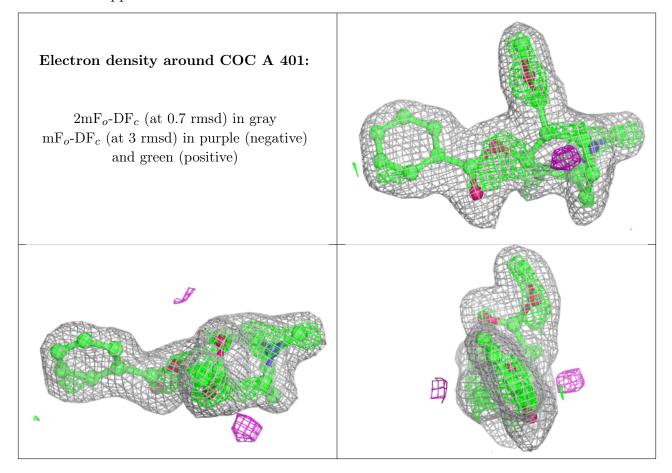


# 6.4 Ligands (i)

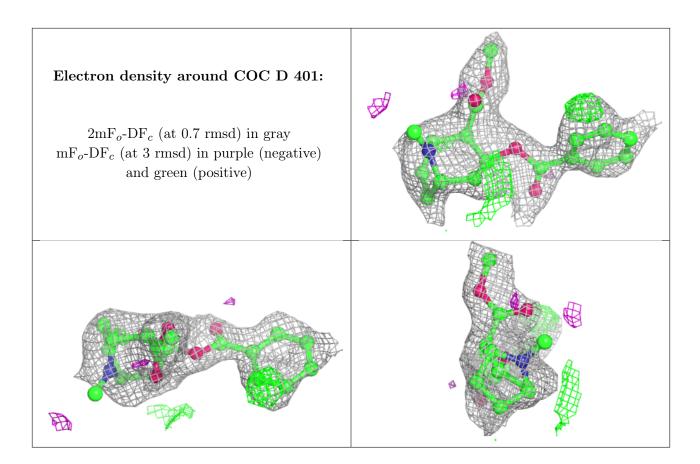
Unable to reproduce the depositors R factor - this section is therefore empty.



The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.







### 6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

