

wwPDB X-ray Structure Validation Summary Report (i)

Jul 4, 2022 – 04:31 pm BST

PDB ID : 7P0K

Title: Crystal structure of Autotaxin (ENPP2) with 18F-labeled positron emission

tomography ligand

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Deposited on : 2021-06-29

Resolution : 2.20 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 as541be (2020)

Xtriage (Phenix) : 1.13

EDS : 2.29

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0267$

CCP4 : 7.1.010 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

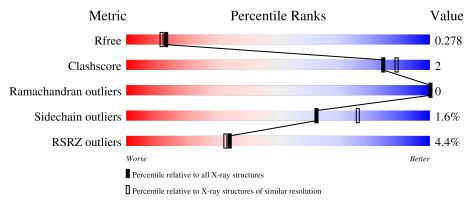
Validation Pipeline (wwPDB-VP) : 2.29

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 2.20 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	Similar resolution
Metric	$(\# ext{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
R_{free}	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain					
1	AAA	770	4%	92%	7%			
2	AbA	7	29%	71%				

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	MAN	AbA	6	X	-	-	X
2	MAN	AbA	7	X	-	-	-
8	4I0	AAA	910	X	-	-	-



2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 12835 atoms, of which 6144 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 Isoform 2 of Ectonucleotide pyrophosphatase/phosphodiester ase family member 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
1	AAA	770	Total 12252	C 3955	H 6018	N 1074	O 1156	S 49	202	2	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AAA	?	-	ASN	deletion	UNP Q64610
AAA	?	-	ASN	deletion	UNP Q64610
AAA	?	-	SER	deletion	UNP Q64610
AAA	?	-	LYS	deletion	UNP Q64610
AAA	410	ALA	ASN	engineered mutation	UNP Q64610
AAA	?	-	LEU	deletion	UNP Q64610
AAA	?	-	ASP	deletion	UNP Q64610
AAA	?	-	VAL	deletion	UNP Q64610
AAA	?	-	TYR	deletion	UNP Q64610
AAA	?	-	LYS	deletion	UNP Q64610
AAA	?	-	LYS	deletion	UNP Q64610
AAA	?	-	PRO	deletion	UNP Q64610
AAA	?	-	SER	deletion	UNP Q64610
AAA	?	-	GLY	deletion	UNP Q64610
AAA	?	-	LYS	deletion	UNP Q64610
AAA	?	-	GLY	deletion	UNP Q64610
AAA	?	-	ASP	deletion	UNP Q64610
AAA	?	-	LYS	deletion	UNP Q64610
AAA	?	-	VAL	deletion	UNP Q64610
AAA	?	-	GLU	deletion	UNP Q64610
AAA	?	-	PRO	deletion	UNP Q64610
AAA	?	-	LYS	deletion	UNP Q64610
AAA	?	-	ASN	deletion	UNP Q64610
AAA	?	-	LYS	deletion	UNP Q64610
AAA	?	-	LEU	deletion	UNP Q64610
AAA	?	-	GLU	deletion	UNP Q64610

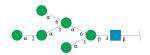
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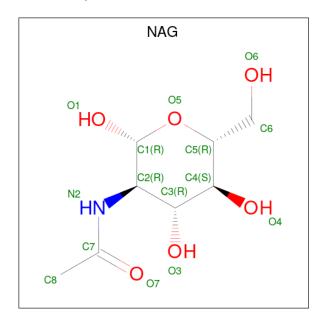
Chain	Residue	Modelled	Actual	${f Comment}$	Reference
AAA	?	-	GLU	deletion	UNP Q64610
AAA	?	-	LEU	deletion	UNP Q64610
AAA	?	-	ASN	deletion	UNP Q64610
AAA	?	-	LYS	deletion	UNP Q64610
AAA	?	-	ARG	deletion	UNP Q64610
AAA	?	-	LEU	deletion	UNP Q64610
AAA	?	-	HIS	deletion	UNP Q64610
AAA	?	-	THR	deletion	UNP Q64610
AAA	?	-	LYS	deletion	UNP Q64610
AAA	591	THR	ARG	conflict	UNP Q64610

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



Mol	Chain	Residues	${f Atoms}$			ZeroOcc	AltConf	Trace		
2	AbA	7	Total 160	C 44	H 80	N 1	O 35	22	0	0

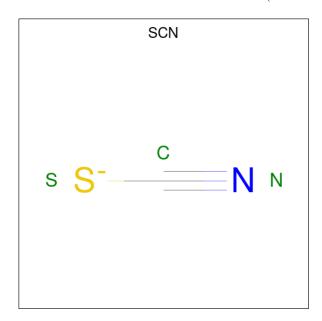
• Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	Λ Λ Λ	1	Total	С	Н	N	О	ე	0
3	AAA	1	28	8	14	1	5	2	U

• Molecule 4 is THIOCYANATE ION (three-letter code: SCN) (formula: CNS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	AAA	1	Total 3		N 1		0	0
4	AAA	1	Total 3		N 1		0	0
4	AAA	1	Total 3	C 1	N 1	S 1	0	0

• Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	AAA	2	Total Zn 2 2	0	0

• Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

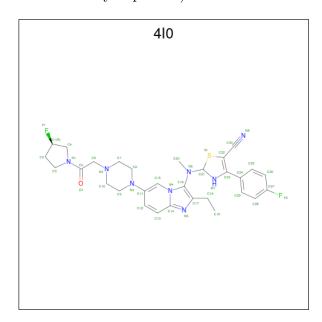
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	AAA	1	Total Ca 1 1	0	0

• Molecule 7 is IODIDE ION (three-letter code: IOD) (formula: I).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	AAA	5	Total I 5 5	0	0

• Molecule 8 is 2-[[2-ethyl-6-[4-[2-[(3 $\{R\})-3$ -fluoranylpyrrolidin-1-yl]-2-oxidanylidene-ethyl]pi perazin-1-yl]imidazo[1,2-a]pyridin-3-yl]-methyl-amino]-4-(4-fluorophenyl)-2,3-dihydro-1,3-th iazole-5-carbonitrile (three-letter code: 4I0) (formula: $C_{30}H_{34}F_2N_8OS$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf			
0	A A A	1	Total	С	F	Н	N	О	S	0	0
0	AAA	1	74	30	2	32	8	1	1	0	0

• Molecule 9 is water.

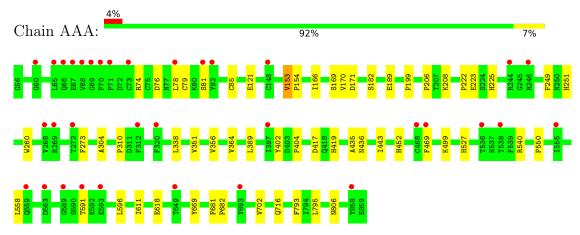
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	AAA	304	Total O 304 304	0	0



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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). 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.

• Molecule 1: Isoform 2 of Ectonucleotide pyrophosphatase/phosphodiesterase family member 2



 $\bullet \ \, Molecule \ 2: \ alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose$

Chain AbA: 29% 71%





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	61.57Å 87.15Å 145.36Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.57 - 2.20	Depositor
resolution (A)	47.52 - 2.20	EDS
% Data completeness	97.9 (47.57-2.20)	Depositor
(in resolution range)	97.9 (47.52-2.20)	EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.44 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0267, REFMAC 5.8.0267	Depositor
P. P.	0.212 , 0.271	Depositor
R, R_{free}	0.222 , 0.278	DCC
R_{free} test set	1926 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.888	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	(Not available), (Not available)	EDS
L-test for twinning ²	$< L > = 0.51, < L^2> = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12835	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.39% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities 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: MAN, SCN, NAG, CA, 4I0, ZN, IOD, BMA

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).

1/4		Chain	Bond	lengths	Bond angles	
101	OI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	L	AAA	0.64	0/6416	0.74	0/8699

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	AAA	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	AAA	436	ASN	Peptide

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	AAA	6234	6018	5985	29	0
2	AbA	80	80	67	0	0

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Continued	trom	mromonie	maaa
-	110116	DICULUUS	Duuc
	J	1	1

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	AAA	14	14	12	0	0
4	AAA	9	0	0	0	0
5	AAA	2	0	0	0	0
6	AAA	1	0	0	0	0
7	AAA	5	0	0	1	0
8	AAA	42	32	0	0	0
9	AAA	304	0	0	4	0
All	All	6691	6144	6064	29	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 2.

The worst 5 of 29 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{aligned} & & & & & & & & & & & & & & & & & & &$	Clash overlap (Å)
		distance (A)	overlap (A)
1:AAA:223:GLU:OE1	1:AAA:435:ALA:O	2.11	0.68
1:AAA:222:PRO:HA	1:AAA:225:HIS:CE1	2.32	0.64
1:AAA:669:TYR:O	9:AAA:1001:HOH:O	2.16	0.59
1:AAA:169:SER:HA	1:AAA:356:VAL:O	2.05	0.57
1:AAA:81:SER:HB3	1:AAA:249:PHE:CZ	2.41	0.56

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	Perce	ntiles
1	AAA	762/770~(99%)	731 (96%)	31 (4%)	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	Analysed	Analysed Rotameric		Percentiles	
1	AAA	699/697 (100%)	688 (98%)	11 (2%)	62 76	

5 of 11 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	558	LEU
1	AAA	591	THR
1	AAA	806	ASN
1	AAA	702	VAL
1	AAA	251	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

7 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	Tuno	Chain	Res	Link	Вс	ond leng	ths	В	ond ang	les
MIOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	AbA	1	2	14,14,15	0.50	0	17,19,21	1.35	2 (11%)
2	BMA	AbA	2	2	11,11,12	0.38	0	15,15,17	1.43	3 (20%)
2	MAN	AbA	3	2	11,11,12	0.42	0	15,15,17	1.39	2 (13%)
2	MAN	AbA	4	2	11,11,12	0.65	0	15,15,17	0.86	0
2	MAN	AbA	5	2	11,11,12	0.46	0	15,15,17	0.88	0
2	MAN	AbA	6	2	11,11,12	0.63	0	15,15,17	1.44	3 (20%)
2	MAN	AbA	7	2	11,11,12	0.39	0	15,15,17	1.27	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	AbA	1	2	-	0/6/23/26	0/1/1/1
2	BMA	AbA	2	2	-	0/2/19/22	0/1/1/1
2	MAN	AbA	3	2	-	1/2/19/22	0/1/1/1
2	MAN	AbA	4	2	-	1/2/19/22	0/1/1/1
2	MAN	AbA	5	2	-	0/2/19/22	0/1/1/1
2	MAN	AbA	6	2	1/1/5/5	1/2/19/22	0/1/1/1
2	MAN	AbA	7	2	1/1/5/5	0/2/19/22	0/1/1/1

There are no bond length outliers.

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
2	AbA	3	MAN	C1-O5-C5	4.05	117.67	112.19
2	AbA	1	NAG	C1-O5-C5	3.62	117.10	112.19
2	AbA	6	MAN	C1-O5-C5	3.45	116.87	112.19
2	AbA	2	BMA	O5-C5-C6	3.21	112.24	107.20
2	AbA	7	MAN	C1-O5-C5	2.66	115.80	112.19

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	AbA	6	MAN	C1
2	AbA	7	MAN	C1

All (3) torsion outliers are listed below:

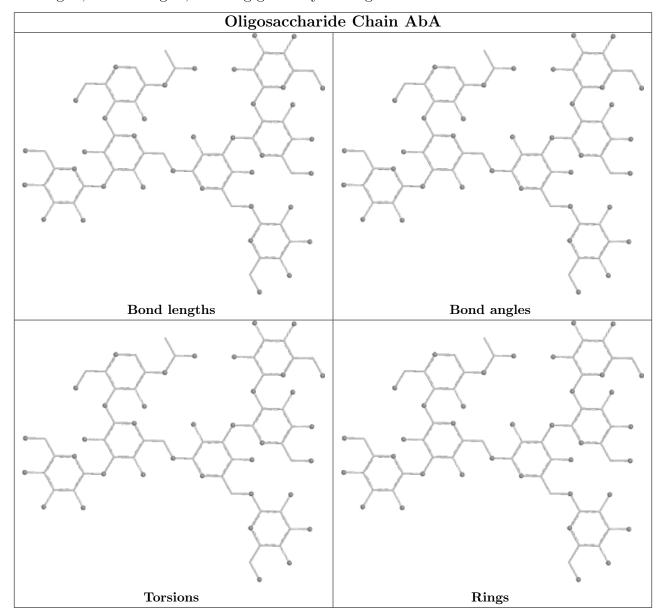


Mol	Chain	Res	Type	Atoms
2	AbA	6	MAN	O5-C5-C6-O6
2	AbA	3	MAN	O5-C5-C6-O6
2	AbA	4	MAN	C4-C5-C6-O6

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)

Of 13 ligands modelled in this entry, 8 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	True	Chain	Des	Link	В	ond leng	$_{ m gths}$	Bond angles			
MIOI	Type	Chain	Res	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2	
4	SCN	AAA	913	-	1,2,2	1.15	0	0,1,1	-	-	
3	NAG	AAA	901	1	14,14,15	0.40	0	17,19,21	1.06	1 (5%)	
4	SCN	AAA	903	-	1,2,2	0.69	0	0,1,1	-	-	
8	4I0	AAA	910	-	40,47,47	1.02	2 (5%)	45,68,68	1.40	5 (11%)	
4	SCN	AAA	902	-	1,2,2	2.26	1 (100%)	0,1,1	-	-	

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.

\mathbf{Mol}	\mathbf{Type}	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	AAA	901	1	-	0/6/23/26	0/1/1/1
8	4I0	AAA	910	-	1/1/9/10	4/19/59/59	0/6/6/6

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\textup{\AA})$	$\operatorname{Ideal}(ext{\AA})$
8	AAA	910	4I0	C21-N7	-5.21	1.30	1.44
4	AAA	902	SCN	C-N	2.26	1.23	1.15
8	AAA	910	4I0	C17-C16	-2.06	1.38	1.42

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
8	AAA	910	4I0	C22-C30-N8	-5.58	168.34	177.71
8	AAA	910	4I0	C24-C23-N7	-3.80	116.81	121.39
8	AAA	910	4I0	C15-C11-N3	-2.69	118.81	121.89
8	AAA	910	4I0	C9-N3-C8	2.61	117.27	111.52
8	AAA	910	4I0	C22-C23-N7	2.51	113.23	107.92

All (1) chirality outliers are listed below:



Mol	Chain	Res	Type	Atom	
8	AAA	910	4I0	C21	

All (4) torsion outliers are listed below:

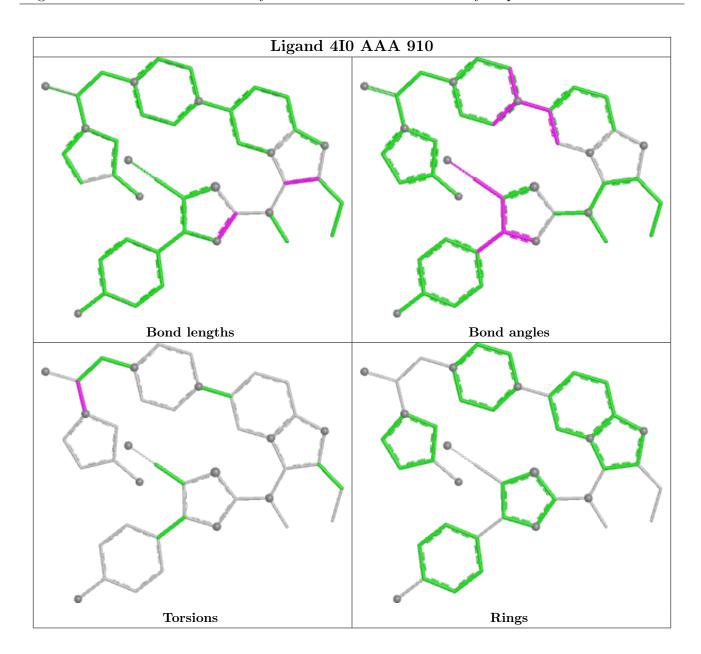
Mol	Chain	Res	Type	Atoms
8	AAA	910	4I0	C6-C5-N1-C3
8	AAA	910	4I0	O1-C5-N1-C3
8	AAA	910	4I0	C6-C5-N1-C4
8	AAA	910	4I0	O1-C5-N1-C4

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 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)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	AAA	4

All chain breaks are listed below:



Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	AAA	569:ASP	С	589:GLY	N	30.60
1	AAA	457:PRO	С	468:CYS	N	13.75
1	AAA	397:ILE	С	402:TYR	N	8.68
1	AAA	553:PRO	С	555:ILE	N	3.27



6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ>2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	$\langle { m RSRZ} \rangle$	# RSRZ > 2	$OWAB(Å^2)$	Q<0.9
1	AAA	770/770 (100%)	0.42	34 (4%) 34 32	21, 34, 60, 96	0

The worst 5 of 34 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	82	TYR	7.6
1	AAA	71	PRO	7.0
1	AAA	70	PRO	6.6
1	AAA	65	LEU	5.7
1	AAA	468	CYS	5.3

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

There are no non-standard protein/DNA/RNA residues in this entry.

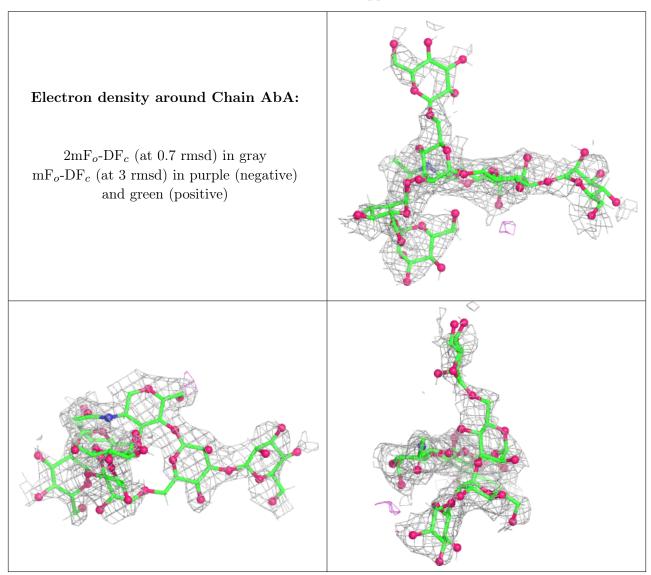
6.3 Carbohydrates (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
2	MAN	AbA	7	11/12	0.61	0.30	30,83,87,91	4
2	MAN	AbA	6	11/12	0.64	0.51	30,82,88,91	4
2	BMA	AbA	2	11/12	0.73	0.27	30,63,67,67	3
2	MAN	AbA	5	11/12	0.78	0.23	30,59,62,65	4
2	MAN	AbA	4	11/12	0.81	0.26	30,59,62,64	3
2	MAN	AbA	3	11/12	0.83	0.21	30,64,72,79	2
2	NAG	AbA	1	14/15	0.87	0.12	30,39,44,47	2



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)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	\mathbf{Type}	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
4	SCN	AAA	902	3/3	0.58	0.24	47,47,62,63	0
4	SCN	AAA	903	3/3	0.67	0.18	50,50,51,60	0
4	SCN	AAA	913	3/3	0.71	0.21	52,52,56,70	0
8	4I0	AAA	910	42/42	0.90	0.14	26,32,50,50	0

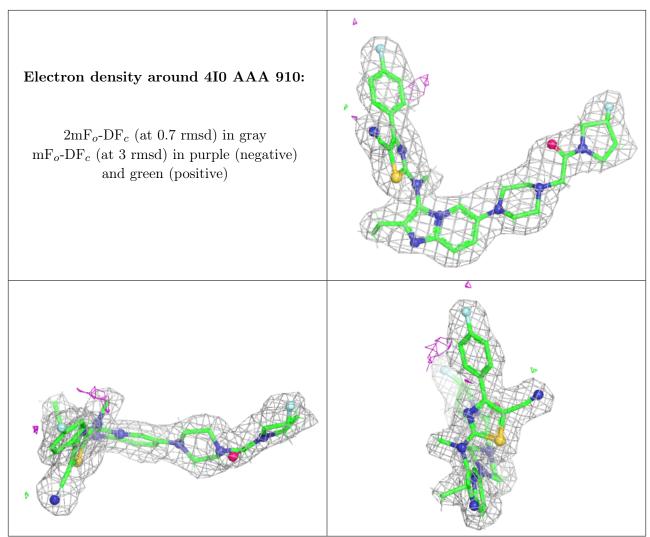
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
3	NAG	AAA	901	14/15	0.93	0.14	23,26,30,31	2
7	IOD	AAA	909	1/1	0.96	0.05	80,80,80,80	0
5	ZN	AAA	905	1/1	0.96	0.07	61,61,61,61	0
7	IOD	AAA	908	1/1	0.97	0.09	83,83,83,83	0
7	IOD	AAA	911	1/1	0.98	0.04	60,60,60,60	0
5	ZN	AAA	904	1/1	0.98	0.11	31,31,31,31	0
7	IOD	AAA	912	1/1	0.99	0.07	64,64,64,64	0
6	CA	AAA	906	1/1	0.99	0.20	29,29,29,29	0
7	IOD	AAA	907	1/1	1.00	0.12	37,37,37,37	0

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)

There are no such residues in this entry.

