

Full wwPDB X-ray Structure Validation Report (i)

Aug 10, 2020 – 02:23 AM BST

PDB ID : 5WKE

Title: Crystal Structure of Human CD1b in Complex with PS

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Deposited on : 2017-07-25

Resolution : 1.69 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org
A user guide is available at

https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The 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

EDS : 2.13.1 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)

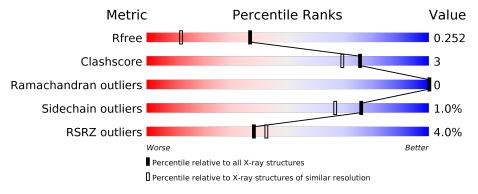
Validation Pipeline (wwPDB-VP) : 2.13.1

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.69 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
R_{free}	130704	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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 on 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 ch	ain		
1	A	300	4%	86%		6%	7%
2	В	99	4%	94%			6%
3	С	3	33%		67%		
4	D	7	57%		43%	1	



2 Entry composition (i)

There are 11 unique types of molecules in this entry. The entry contains 3644 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 T-cell surface glycoprotein CD1b.

Mol	Chain	Residues		$\mathbf{A}\mathbf{t}$	oms			ZeroOcc	AltConf	Trace	
1	A	279	Total 2241	C 1435	N 381	O 415	S 10	3	10	0	

There are 23 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	279	GLY	-	expression tag	UNP P29016
A	280	SER	-	expression tag	UNP P29016
A	281	GLY	_	expression tag	UNP P29016
A	282	LEU	-	expression tag	UNP P29016
A	283	ASN	_	expression tag	UNP P29016
A	284	ASP	_	expression tag	UNP P29016
A	285	ILE	-	expression tag	UNP P29016
A	286	PHE	_	expression tag	UNP P29016
A	287	GLU	_	expression tag	UNP P29016
A	288	ALA	_	expression tag	UNP P29016
A	289	GLN	_	expression tag	UNP P29016
A	290	LYS	_	expression tag	UNP P29016
A	291	ILE	_	expression tag	UNP P29016
A	292	GLU	_	expression tag	UNP P29016
A	293	TRP	_	expression tag	UNP P29016
A	294	HIS	_	expression tag	UNP P29016
A	295	GLU	_	expression tag	UNP P29016
A	296	HIS	_	expression tag	UNP P29016
A	297	HIS	-	expression tag	UNP P29016
A	298	HIS	-	expression tag	UNP P29016
A	299	HIS	-	expression tag	UNP P29016
A	300	HIS	-	expression tag	UNP P29016
A	301	HIS	-	expression tag	UNP P29016

• Molecule 2 is a protein called Beta-2-microglobulin.



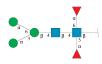
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
2	В	99	Total 831	C 529	N 141	O 159	S 2	0	2	0

• Molecule 3 is an oligosaccharide called alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-be ta-D-glucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	A	\ton	ıs		ZeroOcc	AltConf	Trace
3	С	3	Total 38	C 22	N 2	O 14	4	0	0

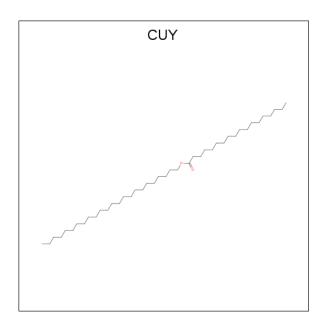
• Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)][alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
4	D	7	Total 81	C N 46 2	O 33	0	0	0

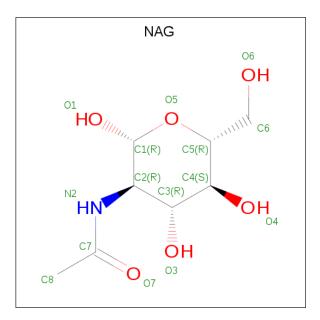
• Molecule 5 is tetracosyl octadecanoate (three-letter code: CUY) (formula: C₄₂H₈₄O₂).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	$\begin{array}{c} { m Total} \\ 35 \end{array}$	C 33	O 2	0	0

• Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	A	ton	ns		ZeroOcc	AltConf
6	A	1	Total	C 8	N 1	O 5	0	0

• Molecule 7 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).





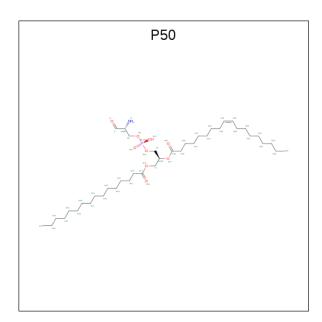
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C O 4 2 2	0	0
7	A	1	Total C O 4 2 2	0	0
7	В	1	Total C O 4 2 2	0	0

• Molecule 8 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	В	2	Total Cl 2 2	0	0
8	A	5	Total Cl 5 5	0	0

• Molecule 9 is (19S,22S,25R)-25-amino-22-hydroxy-16,22,26-trioxo-17,21,23-trioxa-22lam bda 5 -phosphahexacosan-19-yl (9Z)-octadec-9-enoate (three-letter code: P50) (formula: $C_{40}H_{76}NO_9P$).





Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
0	Α	1	Total	С	N	О	Р	0	0
9	A	1	51	40	1	9	1	U	0

• Molecule 10 is SODIUM ION (three-letter code: NA) (formula: Na).

\mathbf{Mol}	Chain	Residues	${f Atoms}$	ZeroOcc	AltConf
10	В	3	Total Na 3 3	0	0

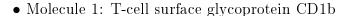
• Molecule 11 is water.

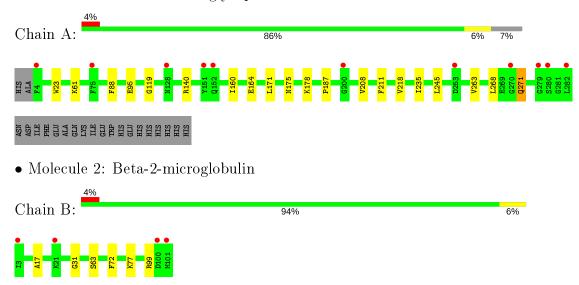
Mol	Chain	Residues	${f Atoms}$	ZeroOcc	AltConf
11	A	238	Total O 238 238	0	0
11	В	93	Total O 93 93	0	0



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 3: alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C: 33% 67%

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





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	57.25Å 78.55Å 84.79Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.64 - 1.69	Depositor
resolution (A)	35.64 - 1.69	EDS
% Data completeness	100.0 (35.64-1.69)	Depositor
(in resolution range)	100.0 (35.64-1.69)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.84 (at 1.69Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
D D.	0.198 , 0.239	Depositor
R, R_{free}	0.209 , 0.252	DCC
R_{free} test set	2155 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	18.2	Xtriage
Anisotropy	0.182	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.34 , 48.0	EDS
L-test for twinning ²	$ < L > = 0.48, < L^2> = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3644	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.62% 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: BMA, NAG, CL, NA, EDO, CUY, FUC, P50, 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	Bond	lengths	Bond angles		
MIOI	Chain	RMSZ	# Z >5	RMSZ	# Z > 5	
1	A	0.49	0/2303	0.63	0/3130	
2	В	0.48	0/854	0.64	0/1158	
All	All	0.49	0/3157	0.63	0/4288	

There are no bond length outliers.

There are no bond angle outliers.

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)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
1	A	2241	0	2133	17	0
2	В	831	0	779	2	0
3	С	38	0	34	1	0
4	D	81	0	70	1	0
5	A	35	0	0	0	0
6	A	14	0	13	1	0
7	A	8	0	12	4	0
7	В	4	0	6	0	0
8	A	5	0	0	1	0
8	В	2	0	0	0	0
9	A	51	0	0	0	0

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Mol	Chain	Non-H	H(model)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
10	В	3	0	0	0	0
11	A	238	0	0	0	0
11	В	93	0	0	0	0
All	All	3644	0	3047	20	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 3.

All (20) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{array}{ll} ext{Interatomic} \ ext{distance} & (ext{Å}) \end{array}$	Clash overlap (Å)
1:A:61:LYS:HB2	7:A:414:EDO:H22	1.32	1.05
1:A:268:LEU:O	1:A:271:GLN:HG3	1.66	0.94
1:A:208:VAL:HG21	1:A:218[B]:VAL:HG11	1.55	0.89
1:A:61:LYS:HB2	7:A:414:EDO:C2	2.08	0.83
1:A:61:LYS:CB	7:A:414:EDO:H22	2.12	0.78
1:A:268:LEU:O	1:A:271:GLN:CG	2.41	0.69
1:A:95:GLU:HB2	7:A:413:EDO:H12	1.84	0.60
1:A:23:TRP:CD1	6:A:402:NAG:H62	2.36	0.60
1:A:218[B]:VAL:HG12	1:A:263:VAL:HG22	1.83	0.59
1:A:187:PRO:HB3	1:A:211:PHE:HB3	1.87	0.57
1:A:119:GLY:HA2	8:A:419:CL:CL	2.43	0.55
1:A:235[A]:ILE:HG13	1:A:245:LEU:HD13	1.89	0.55
3:C:1:NAG:H61	3:C:3:NAG:HN2	1.79	0.48
1:A:160:ILE:HA	1:A:164:GLU:HG3	1.95	0.48
2:B:17:ALA:HB3	2:B:99:ARG:HG3	1.94	0.48
2:B:31:GLY:HA2	2:B:63[A]:SER:HB3	1.98	0.45
1:A:171:LEU:HD12	4:D:2:NAG:H82	1.99	0.44
1:A:88:PHE:HD2	1:A:140:ARG:HG3	1.84	0.43
1:A:175:ASN:O	1:A:178:LYS:HG2	2.20	0.41
1:A:208:VAL:HG21	1:A:218[B]:VAL:CG1	2.39	0.41

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	${f ntiles}$
1	A	287/300 (96%)	284 (99%)	3 (1%)	0	100	100
2	В	99/99 (100%)	99 (100%)	0	0	100	100
All	All	386/399 (97%)	383 (99%)	3 (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	Analysed	Rotameric	Rotameric Outliers	
1	A	$234/247 \ (95\%)$	233 (100%)	1 (0%)	91 87
2	В	93/94 (99%)	91 (98%)	2 (2%)	52 34
All	All	327/341 (96%)	324 (99%)	3 (1%)	76 70

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	271	GLN
2	В	72	PHE
2	В	77	LYS

Some 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)

10 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	Trno	Chain	Res	Link	Вс	ond leng	ths	Bond angles		
MIOI	Type	Chain	nes	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	С	1	1,3	14,14,15	0.48	0	17,19,21	1.28	1 (5%)
3	FUC	С	2	3	10,10,11	0.60	0	14,14,16	1.26	1 (7%)
3	NAG	С	3	3	14,14,15	0.34	0	17,19,21	1.18	1 (5%)
4	NAG	D	1	1,4	14,14,15	0.43	0	17,19,21	0.92	0
4	NAG	D	2	4	14,14,15	0.31	0	17,19,21	0.61	0
4	BMA	D	3	4	11,11,12	0.34	0	15,15,17	0.48	0
4	MAN	D	4	4	11,11,12	0.38	0	15,15,17	0.98	1 (6%)
4	MAN	D	5	4	11,11,12	0.27	0	15,15,17	0.67	1 (6%)
4	FUC	D	6	4	10,10,11	0.34	0	14,14,16	0.75	0
4	FUC	D	7	4	10,10,11	0.41	0	14,14,16	0.66	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
3	NAG	С	1	1,3	-	2/6/23/26	0/1/1/1
3	FUC	С	2	3	-	-	0/1/1/1
3	NAG	С	3	3	-	4/6/23/26	0/1/1/1
4	NAG	D	1	1,4	-	1/6/23/26	0/1/1/1
4	NAG	D	2	4	-	0/6/23/26	0/1/1/1
4	BMA	D	3	4	-	0/2/19/22	0/1/1/1
4	MAN	D	4	4	-	0/2/19/22	0/1/1/1
4	MAN	D	5	4	-	0/2/19/22	0/1/1/1
4	FUC	D	6	4	-	-	0/1/1/1
4	FUC	D	7	4	-	-	0/1/1/1

There are no bond length outliers.



All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
3	С	3	NAG	C1-C2-N2	-4.28	103.18	110.49
3	С	2	FUC	C1-C2-C3	3.94	114.51	109.67
3	С	1	NAG	O3-C3-C4	3.70	118.89	110.35
4	D	4	MAN	C1-O5-C5	3.46	116.89	112.19
4	D	5	MAN	C1-O5-C5	2.23	115.21	112.19

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	С	3	NAG	C8-C7-N2-C2
3	С	3	NAG	O7-C7-N2-C2
3	С	3	NAG	O5-C5-C6-O6
3	С	1	NAG	C8-C7-N2-C2
3	С	1	NAG	O7-C7-N2-C2
3	С	3	NAG	C4-C5-C6-O6
4	D	1	NAG	C3-C2-N2-C7

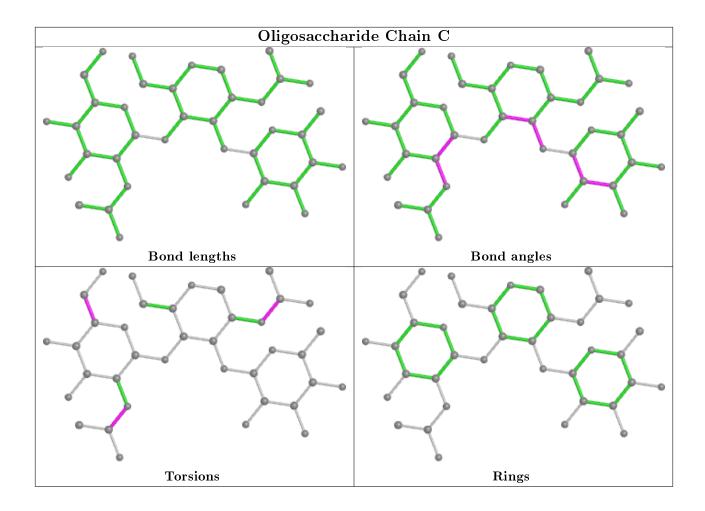
There are no ring outliers.

3 monomers are involved in 2 short contacts:

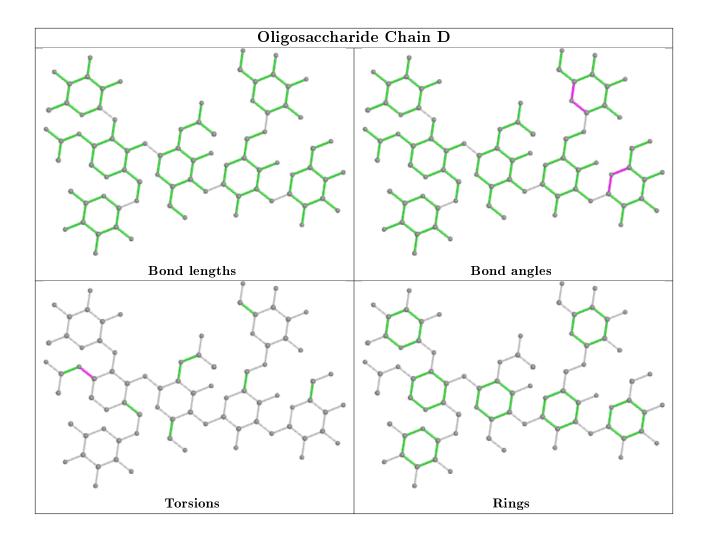
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	С	3	NAG	1	0
3	С	1	NAG	1	0
4	D	2	NAG	1	0

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









5.6 Ligand geometry (i)

Of 16 ligands modelled in this entry, 10 are monoatomic - leaving 6 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	Tuno	Chain	Dog	Res Link Bond lengths			ths	Bond angles		
MIGI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	$\mid \# Z > 2$
7	EDO	В	204	-	3,3,3	0.49	0	2,2,2	0.14	0
7	EDO	A	414	-	3,3,3	0.50	0	2,2,2	0.87	0
7	EDO	A	413	-	3,3,3	0.60	0	2,2,2	0.29	0
9	P50	A	420	-	49,50,50	1.31	4 (8%)	50,56,56	1.10	5 (10%)
6	NAG	A	402	1	14,14,15	0.38	0	17,19,21	0.90	1 (5%)



Mol Type		Chain	Res	Link	Bond lengths			Bond angles		
Wioi I y	туре	Chain	nes	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	CUY	A	401	-	34,34,43	0.68	1 (2%)	34,34,43	0.92	1 (2%)

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
7	EDO	В	204	_	-	1/1/1/1	-
7	EDO	A	414	_	-	1/1/1/1	-
7	EDO	A	413	_	-	1/1/1/1	-
9	P50	A	420	-	-	27/52/55/55	-
6	NAG	A	402	1	-	0/6/23/26	0/1/1/1
5	CUY	A	401	_	_	15/33/33/42	_

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	$\mathbf{Ideal}(\mathbf{\AA})$
9	A	420	P50	P12-O15	4.25	1.66	1.50
9	A	420	P50	O19-C17	3.61	1.43	1.33
9	A	420	P50	O37-C38	3.21	1.43	1.34
5	A	401	CUY	OAQ-CAP	2.67	1.41	1.33
9	A	420	P50	P12-O16	2.19	1.68	1.59

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\mathbf{Ideal}(^{o})$
6	A	402	NAG	C1-O5-C5	3.23	116.57	112.19
9	A	420	P50	O19-C17-C20	2.66	120.24	111.91
9	A	420	P50	C51-C50-C49	-2.61	102.40	113.79
5	A	401	CUY	OAQ-CAP-CAO	2.60	120.08	111.91
9	A	420	P50	O37-C38-C39	2.37	116.61	111.50
9	A	420	P50	P12-OG-CB	-2.02	109.83	121.68
9	A	420	P50	C43-C44-C45	-2.01	105.03	113.79

There are no chirality outliers.

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	A	420	P50	N-CA-CB-OG

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Mol	$\frac{\textit{nued fron}}{\mathbf{Chain}}$	Res	$\overline{ \mathbf{Type} }$	Atoms
9	A	420	P50	C20-C17-O19-C1
5	A	401	CUY	CAO-CAP-OAQ-CAR
5	A	401	CUY	OBO-CAP-OAQ-CAR
9	A	420	P50	O18-C17-O19-C1
9	A	420	P50	C17-C20-C21-C22
7	A	413	EDO	O1-C1-C2-O2
9	A	420	P50	CB-OG-P12-O16
9	A	420	P50	C22-C23-C24-C25
9	A	420	P50	C39-C40-C41-C42
9	A	420	P50	C21-C22-C23-C24
9	A	420	P50	C24-C25-C26-C27
5	A	401	CUY	CAH-CAI-CAJ-CAK
5	A	401	CUY	CAI-CAJ-CAK-CAL
9	A	420	P50	C40-C41-C42-C43
5	A	401	CUY	CAJ-CAK-CAL-CAM
9	A	420	P50	C29-C30-C31-C32
5	A	401	CUY	C1-CBD-CBE-CBF
9	A	420	P50	C51-C52-C53-C54
9	A	420	P50	C52-C53-C54-C55
9	A	420	P50	C50-C51-C52-C53
9	A	420	P50	C42-C43-C44-C45
5	A	401	CUY	C2-C1-CBD-CBE
5	A	401	CUY	CBI-CBJ-CBK-CBL
5	A	401	CUY	CBD-CBE-CBF-CBG
9	A	420	P50	O19-C1-C2-O37
9	A	420	P50	C27-C28-C29-C30
5	A	401	CUY	CAK-CAL-CAM-CAN
9	A	420	P50	C28-C29-C30-C31
9	A	420	P50	O19-C1-C2-C3
9	A	420	P50	C31-C32-C33-C34
5	A	401	CUY	CAU-CAV-CAW-CBM
9	A	420	P50	C25-C26-C27-C28
9	A	420	P50	CB-OG-P12-O13
7	A	414	EDO	O1-C1-C2-O2
9	A	420	P50	C26-C27-C28-C29
5	A	401	CUY	CAT-CAU-CAV-CAW
5	A	401	CUY	CAL-CAM-CAN-CAO
5	A	401	CUY	CAD-CAE-CAF-CAG
9	A	420	P50	C46-C48-C49-C50
9	A	420	P50	C48-C49-C50-C51
7	В	204	EDO	O1-C1-C2-O2
9	A	420	P50	O19-C17-C20-C21

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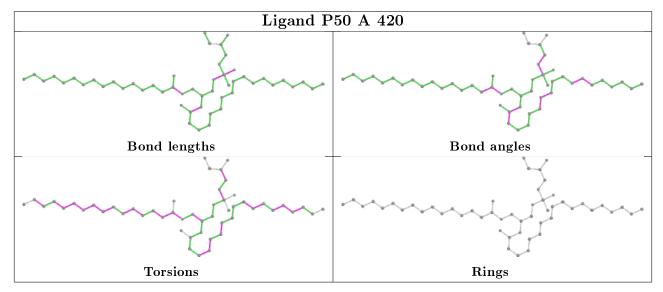
Mol	Chain	Res	Type	Atoms
9	A	420	P50	O18-C17-C20-C21
5	A	401	CUY	CAM-CAN-CAO-CAP

There are no ring outliers.

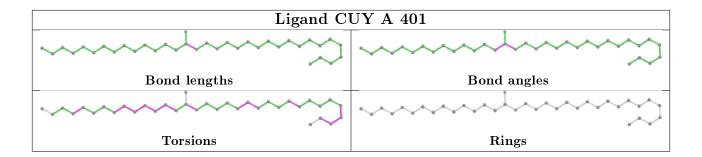
3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	414	EDO	3	0
7	A	413	EDO	1	0
6	A	402	NAG	1	0

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

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	<RSRZ $>$	$\#\mathrm{RSRZ}{>}2$	$OWAB(Å^2)$	Q < 0.9
1	A	279/300 (93%)	0.10	11 (3%) 39 44	11, 22, 40, 63	1 (0%)
2	В	99/99 (100%)	0.08	4 (4%) 38 42	11, 23, 46, 63	0
All	All	378/399 (94%)	0.09	15 (3%) 38 42	11, 22, 42, 63	1 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	152	GLN	4.0
1	A	279	GLY	4.0
1	A	282	LEU	3.9
2	В	101	MET	3.7
1	A	280	SER	2.8
1	A	128	ASN	2.8
1	A	270	GLY	2.7
2	В	3	ILE	2.6
1	A	151	TYR	2.6
2	В	21	LYS	2.5
1	A	253[A]	ASP	2.5
2	В	100	ASP	2.4
1	A	4	PHE	2.3
1	A	75	PHE	2.1
1	A	200	GLY	2.0

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

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



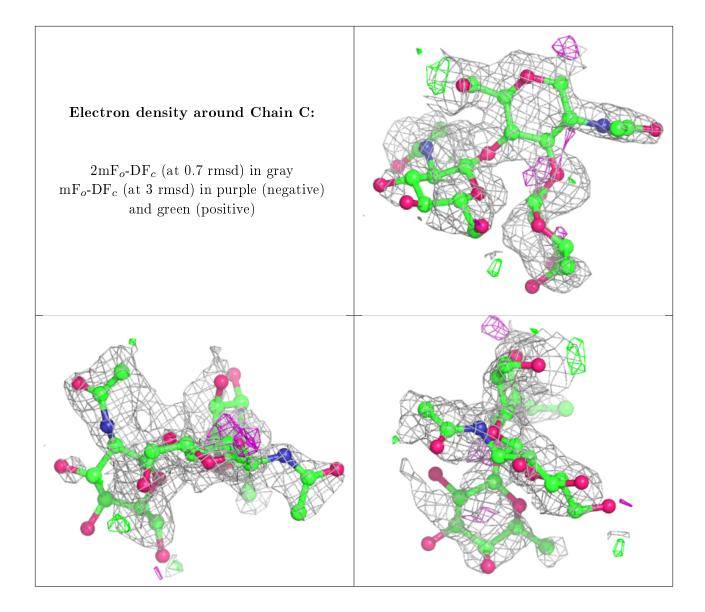
6.3 Carbohydrates (i)

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.

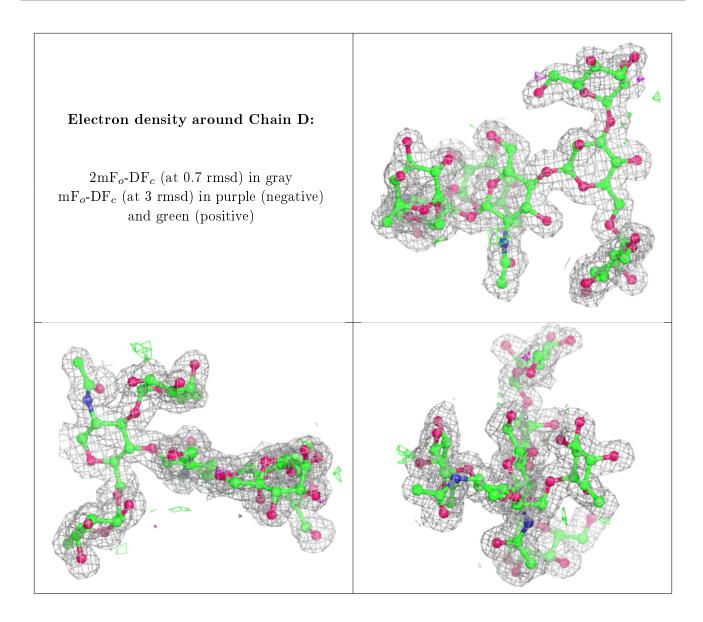
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
3	FUC	С	2	10/11	0.21	0.40	57,58,61,63	3
3	NAG	С	1	14/15	0.55	0.33	50,53,56,59	0
3	NAG	С	3	14/15	0.76	0.29	60,62,65,65	2
4	MAN	D	4	11/12	0.90	0.11	24,28,32,32	0
4	FUC	D	7	10/11	0.90	0.13	21,25,30,30	0
4	FUC	D	6	10/11	0.94	0.08	19,23,27,27	0
4	MAN	D	5	11/12	0.95	0.07	17,19,23,25	0
4	NAG	D	2	14/15	0.95	0.09	13,17,19,21	0
4	BMA	D	3	11/12	0.96	0.07	18,19,22,24	0
4	NAG	D	1	14/15	0.96	0.07	12,15,17,18	0

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	Type	Chain	Res	Atoms	RSCC	RSR	${f B\text{-factors}}({f \AA}^2)$	Q < 0.9
7	EDO	В	204	4/4	0.55	0.30	29,30,34,36	2
6	NAG	A	402	14/15	0.59	0.24	44,48,54,55	0
9	P50	A	420	51/51	0.63	0.25	31,42,53,54	2
5	CUY	A	401	35/44	0.73	0.23	36,44,50,55	0
8	CL	В	205	1/1	0.82	0.14	66,66,66,66	0
8	CL	A	417	1/1	0.89	0.10	42,42,42,42	0
7	EDO	A	414	4/4	0.91	0.18	11,25,25,27	0

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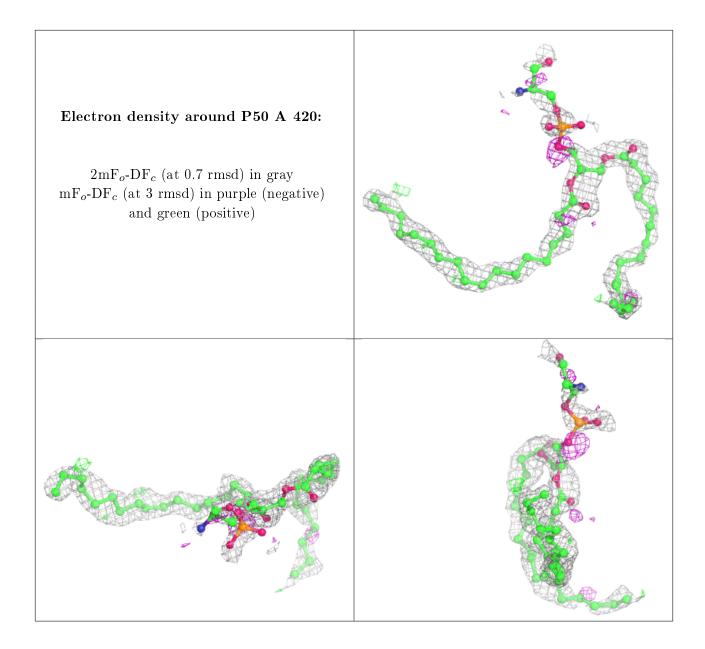


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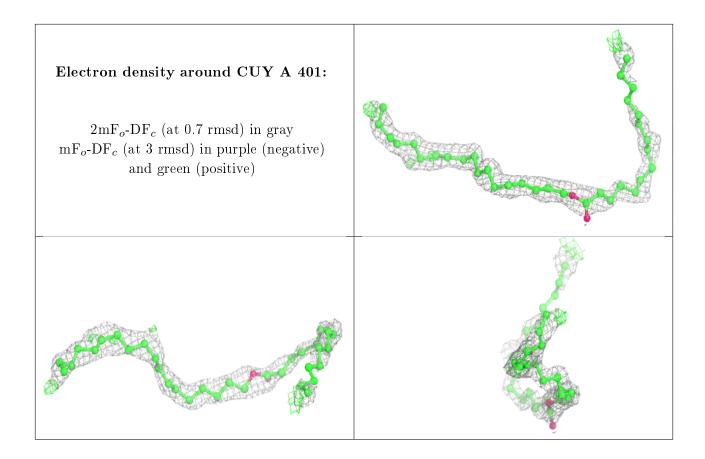
Mol	Type	Chain	Res	Atoms	RSCC	RSR	${f B-factors}({f \AA}^2)$	Q<0.9
8	CL	A	418	1/1	0.91	0.07	48,48,48,48	0
8	CL	A	419	1/1	0.91	0.08	53,53,53,53	0
10	NA	В	201	1/1	0.93	0.24	48,48,48,48	0
7	EDO	A	413	4/4	0.94	0.13	23,29,32,33	0
10	NA	В	202	1/1	0.94	0.09	45,45,45,45	0
8	CL	В	206	1/1	0.94	0.18	76,76,76,76	0
10	NA	В	203	1/1	0.96	0.23	38,38,38,38	0
8	CL	A	415	1/1	0.97	0.10	30,30,30,30	0
8	CL	A	416	1/1	0.99	0.05	26,26,26,26	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.

