

Full wwPDB X-ray Structure Validation Report (i)

Aug 27, 2023 – 11:35 AM EDT

PDB ID	:	3HNV
Title	:	CS-35 Fab Complex with Oligoarabinofuranosyl Tetrasaccharide (branch part
		of Hexasaccharide)
Authors	:	Murase, T.; Zheng, R.B.; Joe, M.; Bai, Y.; Marcus, S.L.; Lowary, T.L.; Ng,
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Deposited on	:	2009-06-01
Resolution	:	2.00 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 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:

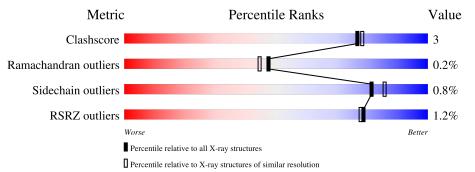
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 2.00 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	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	Н	220	^{2%} 97%	•
2	L	214	88%	12%
3	А	4	100%	

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
3	BXX	А	4	-	-	-	Х



3HNV

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3827 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 CS-35 Fab Heavy Chain.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	Н	220	Total 1675	C 1069	N 277	0 321	S 8	0	0	0

• Molecule 2 is a protein called CS-35 Fab Light Chain.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
2	L	214	Total 1665	C 1032	N 285	0 341	${ m S} 7$	0	0	0

• Molecule 3 is an oligosaccharide called beta-D-arabinofuranose-(1-2)-alpha-D-arabinofurano se-(1-3)-alpha-D-arabinofuranose-(1-5)-methyl alpha-D-arabinofuranoside.

$$\bigstar_{\beta 2} \bigstar_{\alpha 3} \bigstar_{\alpha 5} \bigstar_{\alpha}$$

1Me

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace	
3	А	4	Total 38	C 21	O 17	0	0	0

• Molecule 4 is water.

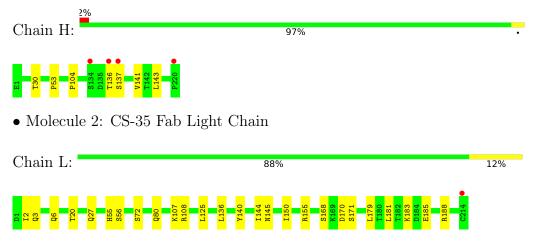
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	Н	242	Total O 242 242	0	0
4	L	207	Total O 207 207	0	0



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: CS-35 Fab Heavy Chain



 \bullet Molecule 3: beta-D-arabinofuranose-(1-2)-alpha-D-arabinofuranose-(1-3)-alpha-D-arabinofurano ose-(1-5)-methyl alpha-D-arabinofuranoside

Chain A:

100%

AXR1 BXY2 BXY3 BXY3 BXX4



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	145.78Å 95.11Å 35.67Å	Depositor
a, b, c, α , β , γ	90.00° 102.51° 90.00°	Depositor
Resolution (Å)	39.53 - 2.00	Depositor
Resolution (A)	39.54 $ 2.00$	EDS
% Data completeness	100.0 (39.53-2.00)	Depositor
(in resolution range)	96.6(39.54-2.00)	EDS
R _{merge}	0.10	Depositor
R _{sym}	0.10	Depositor
$< I/\sigma(I) > 1$	$3.12 (at 2.00 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
D D.	0.174 , 0.230	Depositor
R, R_{free}	0.178 , (Not available)	DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor $(Å^2)$	23.6	Xtriage
Anisotropy	0.484	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.37, 42.0	EDS
L-test for twinning ²	$< L > = 0.49, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.029 for -h-2*l,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3827	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 6.14% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



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: AXR, BXY, BXX

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		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	Н	0.38	0/1721	0.54	0/2352	
2	L	0.38	0/1702	0.56	0/2309	
All	All	0.38	0/3423	0.55	0/4661	

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)	H(added)	Clashes	Symm-Clashes
1	Н	1675	0	1651	3	0
2	L	1665	0	1587	17	0
3	А	38	0	31	0	0
4	Н	242	0	0	0	0
4	L	207	0	0	2	0
All	All	3827	0	3269	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	Interatomic distance (Å)	Clash overlap (Å)
2:L:185:GLU:HG3	2:L:188:ARG:HH21	1.60	0.65
2:L:155:ARG:HD3	2:L:179:LEU:HD11	1.79	0.63
2:L:150:ILE:HD12	2:L:155:ARG:HD2	1.87	0.56
2:L:55:HIS:CG	2:L:56:SER:H	2.24	0.56
2:L:108:ARG:HG3	2:L:171:SER:HB2	1.88	0.55
1:H:136:THR:HG23	1:H:137:SER:HB2	1.89	0.54
2:L:80:GLN:HG3	2:L:168:SER:O	2.09	0.52
2:L:20:THR:CG2	2:L:72:SER:HB2	2.42	0.50
2:L:3:GLN:HB2	4:L:226:HOH:O	2.11	0.49
2:L:125:LEU:O	2:L:183:LYS:HD3	2.12	0.49
1:H:141:VAL:HG13	1:H:143:LEU:HD21	1.95	0.48
2:L:2:ILE:HG12	2:L:27:GLN:HG2	1.96	0.46
2:L:145:ASN:HB3	4:L:404:HOH:O	2.15	0.46
2:L:107:LYS:HA	2:L:140:TYR:OH	2.17	0.45
2:L:185:GLU:HG3	2:L:188:ARG:NH2	2.29	0.45
2:L:136:LEU:HD23	2:L:144:ILE:CD1	2.47	0.44
2:L:136:LEU:HD23	2:L:144:ILE:HD13	2.00	0.44
2:L:150:ILE:HD11	2:L:179:LEU:HD21	2.00	0.44
1:H:30:THR:HA	1:H:53:PRO:HB2	2.01	0.41
2:L:55:HIS:CG	2:L:56:SER:N	2.90	0.40

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	Percentiles
1	Η	218/220~(99%)	214 (98%)	3~(1%)	1 (0%)	29 23
2	L	212/214~(99%)	208 (98%)	4 (2%)	0	100 100
All	All	430/434~(99%)	422 (98%)	7~(2%)	1 (0%)	47 44

All (1) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	Н	104	PRO

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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles		
1	Н	188/188~(100%)	188 (100%)	0	100 100		
2	L	191/191 (100%)	188~(98%)	3~(2%)	62 67		
All	All	379/379~(100%)	376~(99%)	3~(1%)	81 86		

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

Mol	Chain	Res	Type
2	L	6	GLN
2	L	170	ASP
2	L	181	LEU

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)

4 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 length (or angles).

Mal	Mol Type Chain Re	Dec	Link	Bond lengths			Bond angles			
	туре	Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2
3	AXR	А	1	3	11,11,11	0.76	0	$15,\!15,\!15$	2.06	6 (40%)
3	BXY	А	2	3	9,9,10	1.61	1 (11%)	10,12,14	2.53	4 (40%)
3	BXY	А	3	3	9,9,10	1.63	1 (11%)	10,12,14	2.53	5 (50%)
3	BXX	А	4	3	9,9,10	0.57	0	10,12,14	2.18	2 (20%)

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	AXR	А	1	3	-	0/4/20/20	0/1/1/1
3	BXY	А	2	3	-	0/2/15/18	0/1/1/1
3	BXY	А	3	3	-	0/2/15/18	0/1/1/1
3	BXX	А	4	3	-	2/2/15/18	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
3	А	2	BXY	O5-C5	-4.73	1.22	1.42
3	А	3	BXY	O5-C5	-4.68	1.22	1.42

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	А	2	BXY	C5-C4-C3	-5.65	101.48	115.09
3	А	3	BXY	C5-C4-C3	-5.19	102.57	115.09
3	А	4	BXX	C5-C4-C3	-5.11	102.76	115.09
3	А	1	AXR	C5-C4-C3	-3.78	105.97	115.09
3	А	3	BXY	O2-C2-C3	-3.52	104.60	111.27
3	А	2	BXY	O2-C2-C3	-3.39	104.84	111.27
3	А	4	BXX	O2-C2-C3	-3.26	105.09	111.27
3	А	1	AXR	C2-C3-C4	-2.90	97.00	102.64
3	А	3	BXY	O3-C3-C2	-2.80	105.30	112.04
3	А	3	BXY	O4-C4-C3	2.61	107.01	104.70
3	А	1	AXR	O1-C1-C2	-2.47	105.06	108.00

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Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	А	1	AXR	O4-C4-C3	-2.45	100.27	105.11
3	А	1	AXR	O5-C5-C4	-2.44	102.92	111.29
3	А	2	BXY	O3-C3-C2	-2.28	106.54	112.04
3	А	2	BXY	O4-C1-C2	-2.10	101.93	105.99
3	А	1	AXR	O2-C2-C1	-2.06	106.07	111.83
3	А	3	BXY	O3-C3-C4	-2.01	105.24	111.05

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There are no chirality outliers.

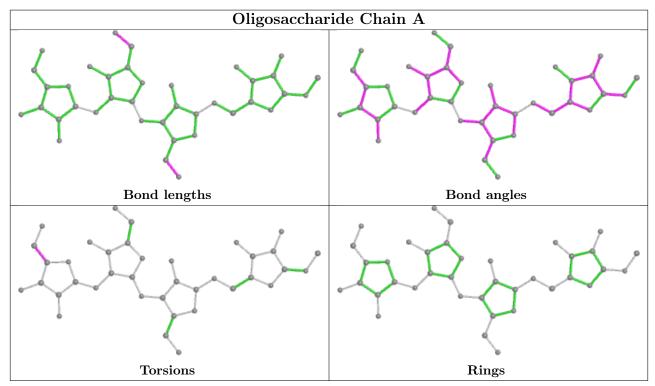
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	А	4	BXX	O4-C4-C5-O5
3	А	4	BXX	C3-C4-C5-O5

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)

There are no ligands in this entry.

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	$\langle RSRZ \rangle$	#RSRZ>2		$\mathbf{OWAB}(\mathbf{A}^2)$	Q < 0.9
1	Н	220/220~(100%)	-0.16	4 (1%) 68 6	66	15, 23, 44, 68	0
2	L	214/214~(100%)	-0.23	1 (0%) 91 9	90	16, 25, 37, 62	0
All	All	434/434~(100%)	-0.19	5 (1%) 79 7	78	15, 23, 37, 68	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	Н	136	THR	7.3
1	Н	137	SER	6.0
2	L	214	CYS	6.0
1	Н	220	PRO	4.0
1	Н	134	SER	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.

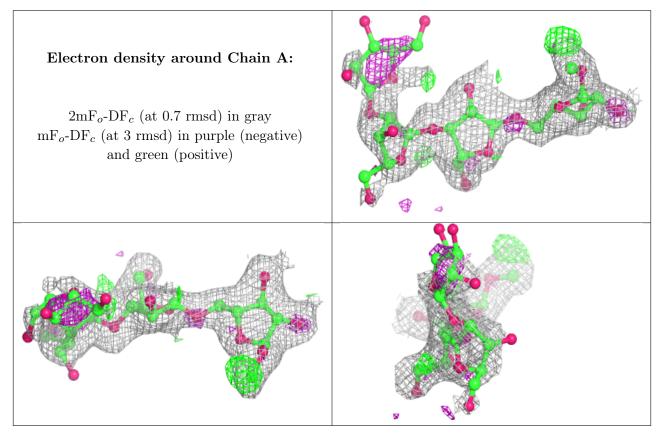
6.3 Carbohydrates (i)

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	B-factors(Å ²)	Q<0.9
3	BXX	А	4	9/10	0.33	0.55	83,84,86,86	0
3	BXY	А	3	9/10	0.78	0.33	69,74,75,79	0
3	BXY	А	2	9/10	0.78	0.14	46,52,55,58	0
3	AXR	А	1	11/11	0.93	0.09	23,27,36,41	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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.

