



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2024 – 10:42 PM EST

PDB ID : 3JWJ
Title : Crystal structure analysis of the methyltransferase domain of bacterial-AvHen1-CN
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Deposited on : 2009-09-18
Resolution : 2.50 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

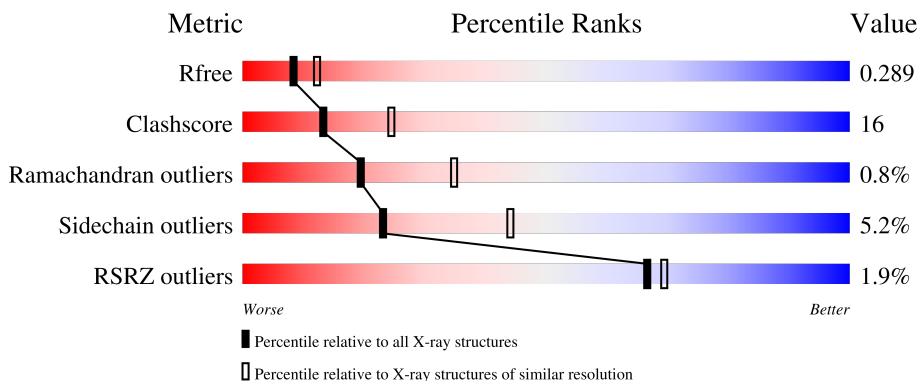
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 $\leq 5\%$. 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	A	202	
1	B	202	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3150 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 Hen1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	188	1542	982	280	277	3	0	0	0
1	B	189	1550	988	281	278	3	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	462	GLY	-	expression tag	UNP Q3MCR9
A	463	HIS	-	expression tag	UNP Q3MCR9
B	462	GLY	-	expression tag	UNP Q3MCR9
B	463	HIS	-	expression tag	UNP Q3MCR9

- Molecule 2 is water.

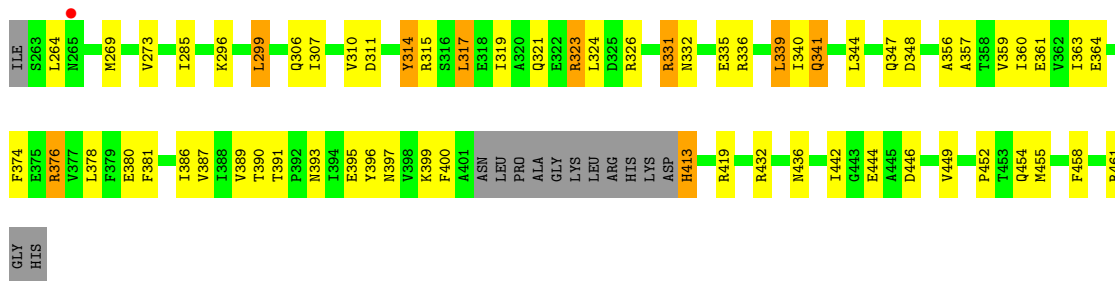
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	29	Total	O	0	0
			29	29		
2	B	29	Total	O	0	0
			29	29		

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.

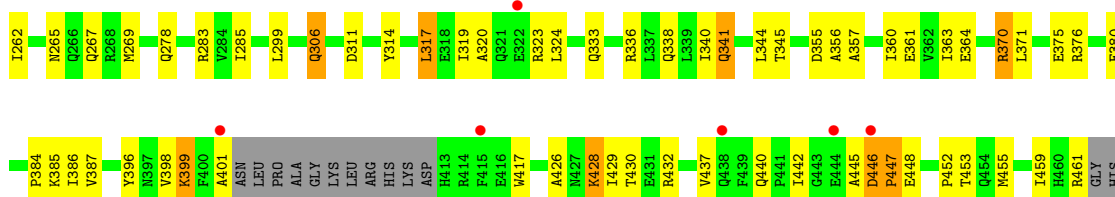
- Molecule 1: Hen1

Chain A: 



- Molecule 1: Hen1

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	33.64Å 175.22Å 146.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 37.58 – 2.51	Depositor EDS
% Data completeness (in resolution range)	90.0 (50.00-2.50) 94.9 (37.58-2.51)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.91 (at 2.51Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.215 , 0.280 0.224 , 0.289	Depositor DCC
R_{free} test set	1199 reflections (7.90%)	wwPDB-VP
Wilson B-factor (Å ²)	24.8	Xtrriage
Anisotropy	0.930	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3150	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.47% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.40	0/1573	0.64	0/2125
1	B	0.39	0/1581	0.62	1/2136 (0.0%)
All	All	0.39	0/3154	0.63	1/4261 (0.0%)

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	A	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	399	LYS	N-CA-C	-5.21	96.95	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	314	TYR	Sidechain

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	A	1542	0	1534	47	0
1	B	1550	0	1545	53	0
2	A	29	0	0	0	0
2	B	29	0	0	1	0
All	All	3150	0	3079	100	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 16.

All (100) 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:B:314:TYR:CZ	1:B:341:GLN:HG2	2.08	0.88
1:A:432:ARG:HG3	1:A:432:ARG:HH11	1.51	0.76
1:B:428:LYS:HE3	1:B:428:LYS:HA	1.74	0.70
1:B:442:ILE:HG13	1:B:453:THR:HB	1.75	0.69
1:A:359:VAL:HB	1:A:389:VAL:HG22	1.76	0.68
1:A:331:ARG:HD3	1:A:332:ASN:OD1	1.94	0.66
1:B:283:ARG:HD3	2:B:20:HOH:O	1.96	0.65
1:A:376:ARG:HG2	1:A:376:ARG:HH11	1.60	0.64
1:A:397:ASN:HB3	1:A:413:HIS:CE1	2.37	0.59
1:B:262:ILE:CG2	1:B:267:GLN:HG3	2.32	0.59
1:B:440:GLN:O	1:B:455:MET:HB3	2.03	0.59
1:A:296:LYS:HA	1:A:324:LEU:HD21	1.85	0.58
1:A:299:LEU:O	1:A:336:ARG:NH2	2.37	0.57
1:B:283:ARG:HH11	1:B:306:GLN:HB2	1.70	0.57
1:B:262:ILE:HG22	1:B:267:GLN:HG3	1.86	0.57
1:B:314:TYR:CE1	1:B:341:GLN:HG2	2.40	0.56
1:B:399:LYS:NZ	1:B:448:GLU:OE1	2.40	0.55
1:A:319:ILE:O	1:A:323:ARG:HG2	2.06	0.54
1:B:283:ARG:HG2	1:B:306:GLN:HB2	1.88	0.54
1:B:446:ASP:HB3	1:B:448:GLU:H	1.72	0.54
1:B:446:ASP:H	1:B:447:PRO:HA	1.70	0.54
1:A:363:ILE:HD11	1:A:391:THR:CB	2.37	0.54
1:A:390:THR:HG22	1:A:455:MET:HG3	1.89	0.54
1:A:376:ARG:HG2	1:A:376:ARG:NH1	2.24	0.53
1:B:311:ASP:HB3	1:B:317:LEU:HD13	1.90	0.53
1:B:376:ARG:HD2	1:B:380:GLU:OE1	2.08	0.53
1:B:432:ARG:O	1:B:432:ARG:HD2	2.09	0.53
1:B:265:ASN:O	1:B:269:MET:HG2	2.09	0.53
1:A:332:ASN:O	1:A:335:GLU:HG2	2.09	0.52
1:B:459:ILE:HD12	1:B:459:ILE:N	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:LEU:HG	1:A:374:PHE:CE1	2.44	0.52
1:A:347:GLN:HG3	1:A:381:PHE:CG	2.44	0.52
1:A:347:GLN:HG3	1:A:381:PHE:CD2	2.46	0.51
1:B:356:ALA:HA	1:B:386:ILE:O	2.10	0.51
1:B:363:ILE:HG13	1:B:364:GLU:N	2.25	0.51
1:A:387:VAL:HB	1:A:458:PHE:HB2	1.93	0.51
1:B:338:GLN:NE2	1:B:340:ILE:HD11	2.25	0.51
1:A:400:PHE:HZ	1:A:446:ASP:OD2	1.93	0.50
1:B:283:ARG:NH1	1:B:306:GLN:CG	2.75	0.50
1:A:323:ARG:HB3	1:A:326:ARG:NH2	2.26	0.50
1:B:430:THR:O	1:B:461:ARG:NH2	2.44	0.49
1:A:269:MET:O	1:A:273:VAL:HG23	2.12	0.49
1:A:321:GLN:HG2	1:A:339:LEU:HD23	1.93	0.49
1:B:283:ARG:NH1	1:B:306:GLN:HB2	2.27	0.49
1:A:432:ARG:HG3	1:A:432:ARG:NH1	2.24	0.49
1:B:426:ALA:O	1:B:429:ILE:HB	2.13	0.49
1:B:375:GLU:OE2	1:B:429:ILE:HD11	2.13	0.48
1:A:363:ILE:HD12	1:A:364:GLU:N	2.29	0.48
1:A:436:ASN:ND2	1:A:461:ARG:HE	2.12	0.48
1:B:299:LEU:O	1:B:336:ARG:NH2	2.44	0.48
1:A:307:ILE:N	1:A:307:ILE:HD12	2.30	0.47
1:B:333:GLN:HB3	1:B:336:ARG:HH21	1.80	0.47
1:B:319:ILE:O	1:B:323:ARG:HG2	2.14	0.47
1:B:396:TYR:CE2	1:B:452:PRO:HB3	2.50	0.47
1:A:432:ARG:HH11	1:A:432:ARG:CG	2.25	0.46
1:B:333:GLN:OE1	1:B:336:ARG:CZ	2.63	0.46
1:B:345:THR:HG21	1:B:370:ARG:CG	2.46	0.46
1:A:378:LEU:HD21	1:A:387:VAL:HG11	1.98	0.46
1:A:344:LEU:HB3	1:A:374:PHE:CD1	2.51	0.45
1:B:446:ASP:HB3	1:B:448:GLU:N	2.30	0.45
1:B:320:ALA:O	1:B:324:LEU:HG	2.17	0.45
1:B:345:THR:HG21	1:B:370:ARG:HG3	1.98	0.45
1:B:398:VAL:O	1:B:401:ALA:HB2	2.16	0.45
1:A:363:ILE:HD11	1:A:391:THR:OG1	2.17	0.45
1:A:348:ASP:O	1:A:381:PHE:HE2	2.00	0.45
1:B:442:ILE:CG1	1:B:453:THR:HB	2.47	0.45
1:A:311:ASP:O	1:A:341:GLN:HA	2.17	0.44
1:A:285:ILE:O	1:A:285:ILE:HG23	2.18	0.44
1:A:442:ILE:HD11	1:A:455:MET:SD	2.57	0.44
1:B:363:ILE:HD12	1:B:417:TRP:CZ3	2.53	0.44
1:A:356:ALA:HA	1:A:386:ILE:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:419:ARG:HA	1:A:454:GLN:OE1	2.18	0.44
1:A:376:ARG:HG3	1:A:380:GLU:OE1	2.18	0.43
1:B:344:LEU:N	1:B:344:LEU:HD12	2.33	0.43
1:B:446:ASP:CB	1:B:448:GLU:H	2.31	0.43
1:A:360:ILE:HG23	1:A:390:THR:OG1	2.17	0.43
1:A:396:TYR:CE2	1:A:452:PRO:HB3	2.54	0.43
1:A:296:LYS:HA	1:A:324:LEU:CD2	2.46	0.43
1:A:399:LYS:HE3	1:A:449:VAL:CG1	2.49	0.43
1:B:384:PRO:HG2	1:B:387:VAL:HG22	2.00	0.43
1:A:341:GLN:CD	1:A:341:GLN:C	2.77	0.43
1:A:310:VAL:HA	1:A:340:ILE:O	2.19	0.43
1:B:262:ILE:HG21	1:B:267:GLN:HG3	2.01	0.43
1:B:299:LEU:HD23	1:B:299:LEU:HA	1.81	0.43
1:B:355:ASP:O	1:B:385:LYS:HB3	2.18	0.43
1:A:314:TYR:CD1	1:A:315:ARG:N	2.86	0.42
1:B:283:ARG:HH11	1:B:306:GLN:CG	2.32	0.42
1:B:285:ILE:O	1:B:357:ALA:HA	2.18	0.42
1:B:283:ARG:HH11	1:B:306:GLN:CB	2.31	0.42
1:B:384:PRO:HG2	1:B:387:VAL:CG2	2.50	0.42
1:A:285:ILE:O	1:A:357:ALA:HA	2.20	0.41
1:A:393:ASN:OD1	1:A:395:GLU:HB2	2.19	0.41
1:A:360:ILE:HG22	1:A:361:GLU:HG2	2.02	0.41
1:B:314:TYR:OH	1:B:341:GLN:HG2	2.19	0.41
1:B:317:LEU:HA	1:B:317:LEU:HD12	1.80	0.41
1:B:360:ILE:HG22	1:B:361:GLU:HG2	2.03	0.41
1:B:371:LEU:O	1:B:375:GLU:HG3	2.21	0.41
1:B:426:ALA:HB1	1:B:437:VAL:HG21	2.03	0.41
1:A:399:LYS:HD2	1:A:400:PHE:CZ	2.56	0.40
1:A:311:ASP:HB3	1:A:317:LEU:HD13	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	A	184/202 (91%)	175 (95%)	8 (4%)	1 (0%)	29	48
1	B	185/202 (92%)	176 (95%)	7 (4%)	2 (1%)	14	26
All	All	369/404 (91%)	351 (95%)	15 (4%)	3 (1%)	19	35

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	331	ARG
1	B	445	ALA
1	B	446	ASP

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	164/175 (94%)	154 (94%)	10 (6%)	18	36
1	B	165/175 (94%)	158 (96%)	7 (4%)	30	54
All	All	329/350 (94%)	312 (95%)	17 (5%)	23	44

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	264	LEU
1	A	299	LEU
1	A	306	GLN
1	A	317	LEU
1	A	323	ARG
1	A	339	LEU
1	A	341	GLN
1	A	376	ARG
1	A	413	HIS
1	A	444	GLU
1	B	278	GLN
1	B	306	GLN
1	B	317	LEU

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Mol	Chain	Res	Type
1	B	341	GLN
1	B	370	ARG
1	B	428	LYS
1	B	447	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	347	GLN
1	A	352	HIS
1	A	421	GLN
1	A	427	ASN
1	A	436	ASN
1	B	338	GLN
1	B	383	GLN
1	B	421	GLN
1	B	436	ASN
1	B	440	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](#)

There are no monosaccharides in this entry.

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

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, 95th 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>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	188/202 (93%)	-0.06	1 (0%) 91 91	11, 22, 40, 48	0
1	B	189/202 (93%)	0.08	6 (3%) 47 51	13, 27, 42, 54	0
All	All	377/404 (93%)	0.01	7 (1%) 66 69	11, 25, 41, 54	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	444	GLU	3.0
1	B	401	ALA	2.5
1	B	322	GLU	2.3
1	A	265	ASN	2.2
1	B	415	PHE	2.1
1	B	446	ASP	2.1
1	B	438	GLN	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](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.