



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 18, 2024 – 03:17 PM EDT

PDB ID : 4DCJ
Title : Crystal structure of caspase 3, L168D mutant
Authors : Chung, S.J.; Kang, H.J.; Kim, S.J.
Deposited on : 2012-01-17
Resolution : 1.70 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
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 1.20.1
EDS : 2.37.1
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) : 2.37.1

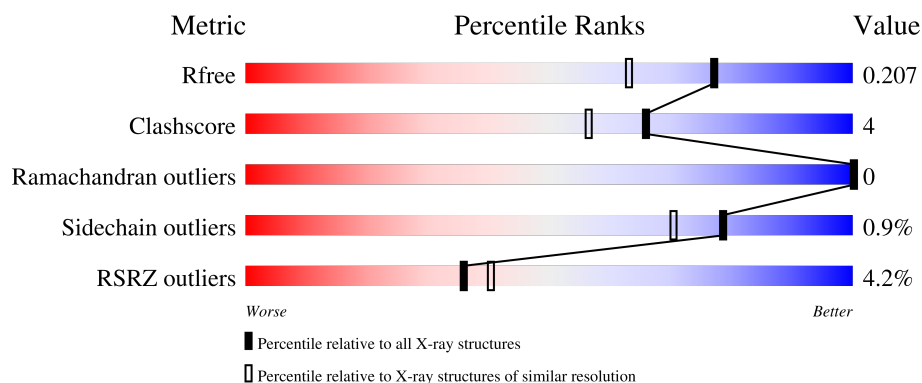
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	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 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	147	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> % </div> <div style="display: flex; justify-content: space-between; align-items: center;"> 84% 12% . </div> </div>
1	D	147	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> 6% </div> <div style="display: flex; justify-content: space-between; align-items: center;"> 84% 12% . </div> </div>
2	B	108	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> 5% </div> <div style="display: flex; justify-content: space-between; align-items: center;"> 81% 6% 12% </div> </div>
2	E	108	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> 4% </div> <div style="display: flex; justify-content: space-between; align-items: center;"> 79% 6% . 14% </div> </div>
3	C	5	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> 80% </div> <div style="display: flex; justify-content: space-between; align-items: center;"> 80% 20% </div> </div>

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Mol	Chain	Length	Quality of chain
3	F	5	<div><div></div><div>80%</div><div>20%</div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4287 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 Caspase-3 subunit p17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	141	Total	C	N	O	S	0	2	0
			1126	692	205	219	10			
1	D	141	Total	C	N	O	S	0	2	0
			1127	694	204	219	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	168	ASP	LEU	ENGINEERED MUTATION	UNP P42574
D	1168	ASP	LEU	ENGINEERED MUTATION	UNP P42574

- Molecule 2 is a protein called Caspase-3 subunit p12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	95	Total	C	N	O	S	0	0	0
			797	523	131	138	5			
2	E	93	Total	C	N	O	S	0	0	0
			777	511	125	136	5			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	278	HIS	-	EXPRESSION TAG	UNP P42574
B	279	HIS	-	EXPRESSION TAG	UNP P42574
B	280	HIS	-	EXPRESSION TAG	UNP P42574
B	281	HIS	-	EXPRESSION TAG	UNP P42574
B	282	HIS	-	EXPRESSION TAG	UNP P42574
B	283	HIS	-	EXPRESSION TAG	UNP P42574
E	1278	HIS	-	EXPRESSION TAG	UNP P42574
E	1279	HIS	-	EXPRESSION TAG	UNP P42574
E	1280	HIS	-	EXPRESSION TAG	UNP P42574
E	1281	HIS	-	EXPRESSION TAG	UNP P42574

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Chain	Residue	Modelled	Actual	Comment	Reference
E	1282	HIS	-	EXPRESSION TAG	UNP P42574
E	1283	HIS	-	EXPRESSION TAG	UNP P42574

- Molecule 3 is a protein called Caspase Inhibitor AC-DEVD-CHO.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	5	Total	C	N	O	0	0	0
			35	20	4	11			
3	F	5	Total	C	N	O	0	0	0
			35	20	4	11			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	127	Total	O	0	0
			127	127		
4	B	79	Total	O	0	0
			79	79		
4	C	8	Total	O	0	0
			8	8		
4	D	98	Total	O	0	0
			98	98		
4	E	69	Total	O	0	0
			69	69		
4	F	9	Total	O	0	0
			9	9		

- Molecule 1: Caspase-3 subunit p17



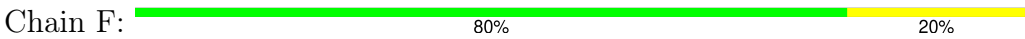
- | SER | GLY | ILE | SER | SER | LEU |
|-------|-----|-----|-----|-----|-----|
| D1034 | | | | | |
| K1038 | | | | | |
| M1039 | | | | | |
| D1040 | | | | | |
| Y1041 | | | | | |
| M1051 | | | | | |
| H1056 | | | | | |
| K1057 | | | | | |
| S1058 | | | | | |
| T1059 | | | | | |
| G1060 | | | | | |
| M1061 | | | | | |
| L1074 | | | | | |
| M1089 | | | | | |
| R1093 | | | | | |
| L1099 | | | | | |
| M1100 | | | | | |
| V1103 | | | | | |
| S1112 | | | | | |
| V1117 | | | | | |
| G1129 | | | | | |
| V1134 | | | | | |
| L1139 | | | | | |
| I1159 | | | | | |
| E1173 | | | | | |
| T1174 | | | | | |
| ASP | | | | | |

- SER
GLY
VAL
ASP
ASP
ASP
MET
ALA
CYS
H185
L194
R207
S213
M268
Y276
H277
H278
H279
HIS
HIS
HIS
HIS

- [illegible]

- ACE501
V504
ASJ505

- 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.83Å 68.55Å 93.81Å 90.00° 101.88° 90.00°	Depositor
Resolution (Å)	32.11 – 1.70 32.11 – 1.70	Depositor EDS
% Data completeness (in resolution range)	95.3 (32.11-1.70) 95.4 (32.11-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 1.70Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.190 , 0.213 0.186 , 0.207	Depositor DCC
R_{free} test set	3457 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	14.5	Xtriage
Anisotropy	0.486	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4287	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.58% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ASJ, ACE

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	0.30	0/1152	0.60	1/1542 (0.1%)
1	D	0.29	0/1152	0.58	1/1543 (0.1%)
2	B	0.34	0/824	0.62	0/1113
2	E	0.33	0/802	0.63	0/1083
3	C	0.50	0/24	0.59	0/32
3	F	0.48	0/24	0.58	0/32
All	All	0.32	0/3978	0.60	2/5345 (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
3	C	0	1
3	F	0	1
All	All	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	129	GLY	N-CA-C	-5.93	98.27	113.10
1	D	1129	GLY	N-CA-C	-5.10	100.34	113.10

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	504	VAL	Mainchain
3	F	1504	VAL	Mainchain

5.2 Too-close contacts

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	1126	0	1112	14	0
1	D	1127	0	1110	9	0
2	B	797	0	768	6	0
2	E	777	0	754	5	0
3	C	35	0	27	0	0
3	F	35	0	27	0	0
4	A	127	0	0	1	0
4	B	79	0	0	0	0
4	C	8	0	0	0	0
4	D	98	0	0	0	0
4	E	69	0	0	0	0
4	F	9	0	0	0	0
All	All	4287	0	3798	31	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 4.

All (31) 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:A:93:ARG:HB2	1:A:134:VAL:CG1	2.06	0.86
1:A:93:ARG:HB2	1:A:134:VAL:HG13	1.71	0.73
1:A:40:ASP:HB3	2:B:277:HIS:NE2	2.07	0.69
1:D:1093:ARG:HB2	1:D:1134:VAL:HG22	1.78	0.65
1:D:1100:MET:HG3	1:D:1139:ILE:HG23	1.81	0.63
2:B:207:ARG:HA	2:B:213:SER:HA	1.81	0.62
2:E:1207:ARG:HA	2:E:1213:SER:HA	1.84	0.59
1:A:40:ASP:HB3	2:B:277:HIS:CD2	2.36	0.59
1:A:93:ARG:HB2	1:A:134:VAL:HG11	1.80	0.59
1:D:1173:GLU:H	1:D:1173:GLU:CD	2.06	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1093:ARG:HB2	1:D:1134:VAL:CG2	2.35	0.56
1:A:74:LEU:HD13	1:A:117:VAL:HG11	1.89	0.55
1:A:149:ARG:NH1	1:A:152:THR:OG1	2.40	0.54
1:A:98:GLU:OE1	1:A:101[A]:ARG:NH1	2.42	0.52
1:A:40:ASP:HB3	2:B:277:HIS:CE1	2.45	0.50
1:A:100:MET:HG3	1:A:139:ILE:HG23	1.92	0.50
1:D:1099:LEU:O	1:D:1103:VAL:HG23	2.13	0.48
2:E:1194:LEU:HD13	2:E:1233:MET:CE	2.46	0.46
2:E:1194:LEU:HD13	2:E:1233:MET:HE3	1.98	0.46
1:D:1074:LEU:HD13	1:D:1117:VAL:HG11	1.97	0.46
2:B:194:LEU:C	2:B:194:LEU:HD13	2.38	0.45
1:A:94:GLU:O	1:A:98:GLU:HG2	2.18	0.44
1:A:144:ARG:HD2	4:A:2045:HOH:O	2.17	0.43
1:A:101[B]:ARG:NH1	1:A:102:ASP:OD1	2.51	0.43
1:D:1041:TYR:HB2	1:D:1112:SER:OG	2.17	0.43
1:D:1051:ASN:HD22	1:D:1089:ASN:HD22	1.65	0.43
1:D:1038:LYS:HG3	1:D:1040:ASP:OD1	2.19	0.42
2:E:1194:LEU:HD22	2:E:1232:PHE:HE2	1.84	0.42
1:A:51:ASN:HB3	1:A:89:ASN:HD22	1.85	0.41
2:B:276:TYR:HE1	2:B:278:HIS:CE1	2.37	0.41
2:E:1240:ASN:OD1	2:E:1263:PRO:HB2	2.21	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	141/147 (96%)	138 (98%)	3 (2%)	0	100	100
1	D	141/147 (96%)	137 (97%)	4 (3%)	0	100	100
2	B	93/108 (86%)	92 (99%)	1 (1%)	0	100	100
2	E	91/108 (84%)	89 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	3/5 (60%)	3 (100%)	0	0	100	100
3	F	3/5 (60%)	3 (100%)	0	0	100	100
All	All	472/520 (91%)	462 (98%)	10 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	129/132 (98%)	129 (100%)	0	100	100
1	D	129/132 (98%)	128 (99%)	1 (1%)	81	74
2	B	85/96 (88%)	84 (99%)	1 (1%)	71	59
2	E	83/96 (86%)	81 (98%)	2 (2%)	49	31
3	C	3/3 (100%)	3 (100%)	0	100	100
3	F	3/3 (100%)	3 (100%)	0	100	100
All	All	432/462 (94%)	428 (99%)	4 (1%)	78	70

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	268	MET
1	D	1034	ASP
2	E	1194	LEU
2	E	1268	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	89	ASN
2	B	277	HIS
1	D	1089	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	ASJ	F	1505	1,3	7,7,7	0.81	0	5,8,8	1.06	0
3	ASJ	C	505	1,3	7,7,7	0.80	0	5,8,8	1.03	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	ASJ	F	1505	1,3	-	3/6/6/6	-
3	ASJ	C	505	1,3	-	2/6/6/6	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	F	1505	ASJ	N-CA-CB-CG
3	C	505	ASJ	N-CA-CB-CG
3	C	505	ASJ	O-C-CA-N
3	F	1505	ASJ	O-C-CA-N
3	F	1505	ASJ	C-CA-CB-CG

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	141/147 (95%)	-0.10	2 (1%) 75 79	10, 17, 28, 44	0
1	D	141/147 (95%)	0.31	9 (6%) 19 21	13, 21, 43, 70	0
2	B	95/108 (87%)	-0.02	5 (5%) 26 29	9, 13, 30, 54	0
2	E	93/108 (86%)	0.08	4 (4%) 35 39	9, 16, 30, 57	0
3	C	3/5 (60%)	-0.35	0 100 100	17, 17, 17, 18	0
3	F	3/5 (60%)	0.16	0 100 100	20, 20, 21, 23	0
All	All	476/520 (91%)	0.07	20 (4%) 36 40	9, 18, 35, 70	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	34	ASP	5.9
2	B	279	HIS	5.7
2	E	1277	HIS	4.7
1	D	1034	ASP	4.4
1	D	1057	LYS	4.3
1	D	1174	THR	4.3
2	E	1185	HIS	4.2
2	B	185	HIS	4.1
1	D	1056	HIS	3.9
2	B	277	HIS	3.9
1	D	1061	MET	3.8
2	B	278	HIS	3.6
1	D	1058	SER	3.6
2	E	1229	LYS	2.7
2	B	276	TYR	2.4
1	D	1159	ILE	2.4
2	E	1194	LEU	2.3
1	D	1059	THR	2.3
1	A	174	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	1060	GLY	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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, 95th 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(\AA^2)	Q<0.9
3	ASJ	C	505	8/8	0.93	0.10	13,14,16,19	0
3	ASJ	F	1505	8/8	0.93	0.09	17,18,19,21	0

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.