



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 06:38 AM GMT

PDB ID : 2D1L
Title : Structure of F-actin binding domain IMD of MIM (Missing In Metastasis)
Authors : Lee, S.H.; Kerff, F.; Chereau, D.; Ferron, F.; Dominguez, R.
Deposited on : 2005-08-27
Resolution : 1.85 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

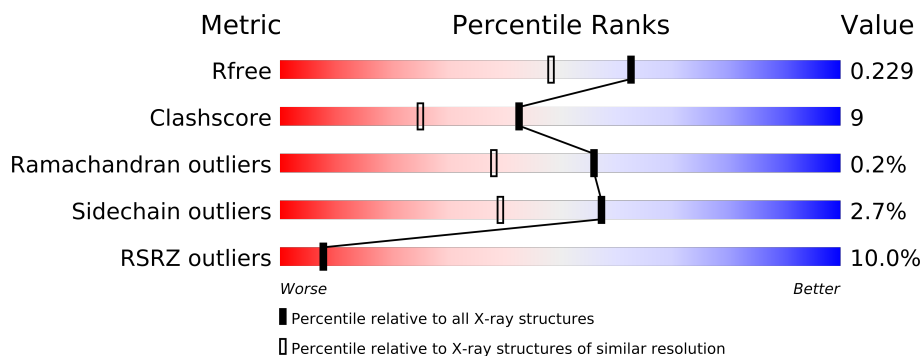
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.85 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}	66092	1269 (1.86-1.86)
Clashscore	79885	1470 (1.86-1.86)
Ramachandran outliers	78287	1451 (1.86-1.86)
Sidechain outliers	78261	1451 (1.86-1.86)
RSRZ outliers	66119	1269 (1.86-1.86)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	253	
1	B	253	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4390 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Metastasis suppressor protein 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	Se	0	19	0
			2058	1290	358	394	4	12			
1	B	231	Total	C	N	O	S	Se	0	20	0
			1940	1224	339	362	4	11			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	ALA	-	CLONING ARTIFACT	UNP Q8R1S4
A	-1	GLY	-	CLONING ARTIFACT	UNP Q8R1S4
A	0	HIS	-	CLONING ARTIFACT	UNP Q8R1S4
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
A	23	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
A	64	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
A	82	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
A	84	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
A	112	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
A	139	LYS	ASN	MODIFIED RESIDUE	UNP Q8R1S4
A	205	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
A	216	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
A	235	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
B	-2	ALA	-	CLONING ARTIFACT	UNP Q8R1S4
B	-1	GLY	-	CLONING ARTIFACT	UNP Q8R1S4
B	0	HIS	-	CLONING ARTIFACT	UNP Q8R1S4
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
B	23	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
B	64	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
B	82	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
B	84	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
B	112	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
B	139	LYS	ASN	MODIFIED RESIDUE	UNP Q8R1S4
B	205	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4
B	216	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	235	MSE	MET	MODIFIED RESIDUE	UNP Q8R1S4

- Molecule 2 is water.

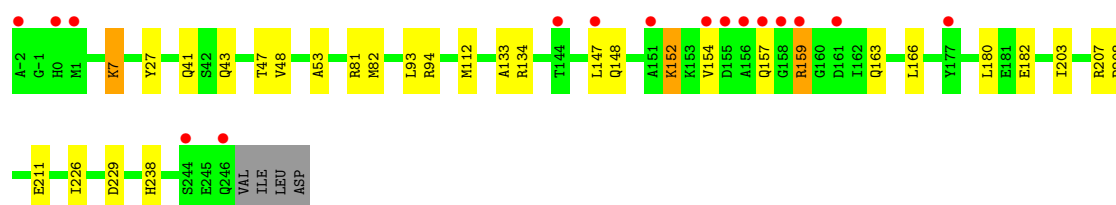
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	208	Total 208	O 208	0	0
2	B	184	Total 184	O 184	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

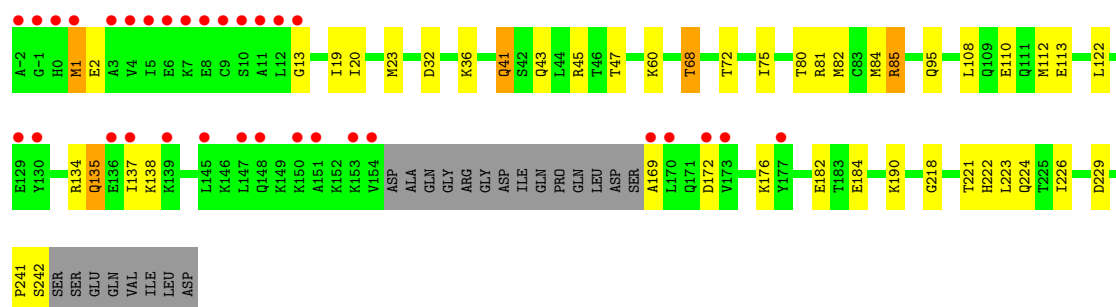
- Molecule 1: Metastasis suppressor protein 1

Chain A: 



- Molecule 1: Metastasis suppressor protein 1

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	53.50Å 37.32Å 129.02Å 90.00° 94.07° 90.00°	Depositor
Resolution (Å)	48.10 – 1.85 48.10 – 1.80	Depositor EDS
% Data completeness (in resolution range)	98.7 (48.10-1.85) 96.2 (48.10-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.183 , 0.228 0.184 , 0.229	Depositor DCC
R_{free} test set	2197 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.295	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 70.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 46045 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4390	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.41% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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	0.52	0/2127	0.55	0/2825
1	B	0.56	1/2012 (0.0%)	0.57	0/2670
All	All	0.54	1/4139 (0.0%)	0.56	0/5495

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1	MSE	SE-CE	8.71	2.46	1.95

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2058	0	2157	31	0
1	B	1940	0	2058	47	0
2	A	208	0	0	1	0
2	B	184	0	0	9	0
All	All	4390	0	4215	73	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (73) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:224[B]:GLN:HG3	2:B:340:HOH:O	1.47	1.14
1:B:1:MSE:SE	1:B:1:MSE:CE	2.46	1.13
1:A:48:VAL:HG11	1:A:94:ARG:HG2	1.46	0.98
1:B:85[A]:ARG:HD3	1:B:222:HIS:HB3	1.45	0.95
1:B:85[A]:ARG:HD3	1:B:222:HIS:CB	1.98	0.93
1:B:84[A]:MSE:SE	2:B:420:HOH:O	2.40	0.89
1:A:152:LYS:HE3	1:A:152:LYS:HA	1.57	0.84
1:A:148:GLN:HG3	1:A:166:LEU:HD11	1.62	0.82
1:A:48:VAL:HG22	1:A:93:LEU:HD23	1.62	0.81
1:B:134:ARG:HD2	1:B:138:LYS:HE2	1.62	0.80
1:A:207[B]:ARG:HG3	1:A:208:PRO:HD3	1.65	0.77
1:B:81[A]:ARG:HA	1:B:84[A]:MSE:HE3	1.67	0.77
1:B:113[B]:GLU:HB2	2:B:316:HOH:O	1.86	0.75
1:A:154:VAL:HG13	1:A:159:ARG:HG3	1.72	0.72
1:B:95:GLN:NE2	2:B:432:HOH:O	2.23	0.72
1:A:48:VAL:HG11	1:A:94:ARG:CG	2.20	0.72
1:A:27:TYR:HD1	1:A:112[A]:MSE:HE1	1.57	0.70
1:B:80:THR:HG22	1:B:84[A]:MSE:HE2	1.74	0.69
1:B:81[A]:ARG:NH2	1:B:229[A]:ASP:OD2	2.20	0.68
1:A:27:TYR:HD1	1:A:112[A]:MSE:CE	2.08	0.66
1:B:122:LEU:HD13	1:B:190:LYS:HE2	1.76	0.66
1:A:112[B]:MSE:HE2	1:A:112[B]:MSE:HA	1.77	0.66
1:B:41:GLN:HE22	1:B:45[A]:ARG:HH11	1.45	0.63
1:B:68:THR:OG1	1:B:72:THR:HB	2.01	0.61
1:A:112[B]:MSE:CE	1:A:112[B]:MSE:HA	2.31	0.61
1:A:7:LYS:N	1:A:7:LYS:HD3	2.17	0.60
1:A:207[B]:ARG:HG3	1:A:208:PRO:CD	2.33	0.59
1:B:110[B]:GLU:OE2	2:B:292:HOH:O	2.17	0.59
1:A:133:ALA:HB3	1:A:180:LEU:HD13	1.83	0.59
1:B:80:THR:O	1:B:84[A]:MSE:HG3	2.04	0.58
1:B:19:ILE:HG22	1:B:23:MSE:HE2	1.85	0.58
1:A:43[B]:GLN:HG3	1:B:47:THR:OG1	2.03	0.58
1:B:85[A]:ARG:HD3	1:B:222:HIS:HB2	1.81	0.58
1:B:113[A]:GLU:HB2	2:B:316:HOH:O	2.07	0.55
1:A:238:HIS:H	1:A:238:HIS:CD2	2.24	0.55
1:B:134:ARG:O	1:B:137:ILE:HG12	2.07	0.54
1:B:218:GLY:O	1:B:221:THR:HG22	2.08	0.54
1:A:203:ILE:HD11	1:B:75:ILE:HG23	1.89	0.53
1:B:32:ASP:OD2	2:B:408:HOH:O	2.18	0.53
1:A:152:LYS:CE	1:A:152:LYS:HA	2.34	0.53
1:B:108:LEU:O	1:B:112:MSE:HG2	2.09	0.52
1:B:241:PRO:O	1:B:242:SER:C	2.47	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:43[B]:GLN:OE1	1:B:43[B]:GLN:NE2	2.40	0.52
1:B:13:GLY:HA3	1:B:184:GLU:OE2	2.10	0.50
1:B:85[A]:ARG:HG3	1:B:226:ILE:CD1	2.41	0.50
1:A:207[B]:ARG:NH1	1:A:211:GLU:HB2	2.27	0.50
1:B:113[A]:GLU:HG2	2:B:369:HOH:O	2.11	0.50
1:A:47:THR:HA	1:B:43[A]:GLN:HG2	1.94	0.50
1:A:133:ALA:CB	1:A:180:LEU:HD13	2.41	0.50
1:B:85[A]:ARG:HG3	1:B:226:ILE:HD12	1.95	0.48
1:B:137:ILE:HG22	1:B:176:LYS:CB	2.43	0.48
1:B:82[A]:MSE:SE	1:B:223:LEU:HD22	2.64	0.47
1:A:147:LEU:HB3	1:A:166:LEU:HD13	1.96	0.47
1:A:27:TYR:CD1	1:A:112[A]:MSE:CE	2.94	0.46
1:B:122:LEU:CD1	1:B:190:LYS:HE2	2.44	0.46
1:B:137:ILE:CG2	1:B:176:LYS:HB3	2.46	0.46
1:B:137:ILE:HG22	1:B:176:LYS:HB2	1.99	0.45
1:B:224[B]:GLN:CG	2:B:340:HOH:O	2.30	0.45
1:A:81[B]:ARG:NH1	1:A:229[B]:ASP:OD2	2.49	0.45
1:B:32:ASP:OD1	1:B:36[A]:LYS:NZ	2.47	0.44
1:B:169:ALA:HA	1:B:172:ASP:OD2	2.17	0.44
1:B:20:ILE:HD13	1:B:23:MSE:HE3	1.99	0.44
1:B:135:GLN:O	1:B:135:GLN:HG3	2.19	0.43
1:A:207[B]:ARG:NE	2:A:368:HOH:O	2.52	0.43
1:A:82[B]:MSE:HE2	1:A:226:ILE:HG22	2.00	0.43
1:B:134:ARG:HD2	1:B:138:LYS:CE	2.38	0.42
1:A:159:ARG:HH11	1:A:163:GLN:HG2	1.83	0.42
1:A:154:VAL:HG12	1:A:154:VAL:O	2.20	0.42
1:B:2:GLU:HG3	1:B:138:LYS:HD3	2.01	0.42
1:B:60:LYS:HE2	1:B:60:LYS:HB3	1.76	0.42
1:A:27:TYR:CD1	1:A:112[A]:MSE:HE3	2.55	0.41
1:A:53:ALA:HB1	1:B:36[B]:LYS:HD2	2.03	0.41
1:B:41:GLN:HE22	1:B:45[A]:ARG:NH1	2.14	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	266/253 (105%)	258 (97%)	8 (3%)	0	100	100
1	B	247/253 (98%)	242 (98%)	4 (2%)	1 (0%)	43	24
All	All	513/506 (101%)	500 (98%)	12 (2%)	1 (0%)	56	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	68	THR

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	234/210 (111%)	227 (97%)	7 (3%)	53	34
1	B	220/210 (105%)	215 (98%)	5 (2%)	63	46
All	All	454/420 (108%)	442 (97%)	12 (3%)	57	41

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LYS
1	A	41	GLN
1	A	134	ARG
1	A	152	LYS
1	A	157	GLN
1	A	159	ARG
1	A	182	GLU
1	B	41	GLN
1	B	85[A]	ARG
1	B	85[B]	ARG
1	B	135	GLN
1	B	182	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	GLN
1	A	238	HIS
1	B	41	GLN
1	B	95	GLN
1	B	106	ASN
1	B	121	GLN
1	B	135	GLN
1	B	186	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

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 ⓘ

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	249/253 (98%)	0.20	16 (6%) 19 18	16, 26, 69, 84	0
1	B	231/253 (91%)	0.52	32 (13%) 4 3	14, 26, 64, 83	0
All	All	480/506 (94%)	0.36	48 (10%) 8 8	14, 26, 69, 84	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	158	GLY	11.2
1	B	-2	ALA	6.0
1	B	151	ALA	5.7
1	B	147	LEU	5.6
1	B	170	LEU	5.4
1	A	156	ALA	5.3
1	B	8	GLU	5.2
1	A	155	ASP	4.8
1	B	145	LEU	4.7
1	A	151	ALA	4.4
1	B	169	ALA	4.4
1	B	137	ILE	4.1
1	B	177	TYR	4.0
1	B	173	VAL	3.9
1	B	12	LEU	3.9
1	B	9	CYS	3.8
1	B	150	LYS	3.7
1	B	5	ILE	3.5
1	A	157	GLN	3.5
1	B	148	GLN	3.2
1	B	-1	GLY	3.1
1	A	-2	ALA	3.1
1	A	159	ARG	3.1
1	B	6	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	0	HIS	2.9
1	B	11	ALA	2.9
1	A	154	VAL	2.8
1	A	1	MSE	2.8
1	B	172	ASP	2.7
1	B	13	GLY	2.7
1	A	246	GLN	2.7
1	B	10	SER	2.6
1	B	154	VAL	2.6
1	A	161	ASP	2.6
1	A	177	TYR	2.5
1	A	144	THR	2.5
1	B	136	GLU	2.5
1	B	0	HIS	2.4
1	B	129	GLU	2.4
1	B	4	VAL	2.4
1	B	153	LYS	2.3
1	A	147	LEU	2.2
1	A	244	SER	2.2
1	B	139	LYS	2.2
1	B	130	TYR	2.1
1	B	3	ALA	2.0
1	B	1	MSE	2.0
1	B	7	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.