



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 06:35 pm GMT

PDB ID : 4O9H
Title : Structure of Interleukin-6 in complex with a Camelid Fab fragment
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Deposited on : 2014-01-02
Resolution : 2.42 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

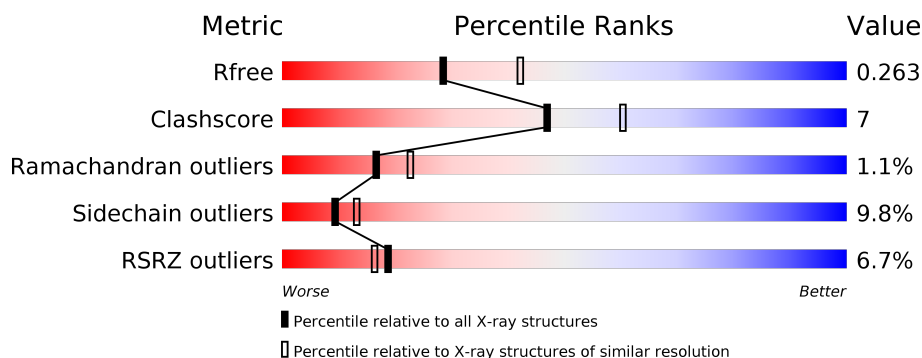
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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}	100719	3709 (2.44-2.40)
Clashscore	112137	4241 (2.44-2.40)
Ramachandran outliers	110173	4178 (2.44-2.40)
Sidechain outliers	110143	4179 (2.44-2.40)
RSRZ outliers	101464	3740 (2.44-2.40)

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. 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	186	
2	H	223	
3	L	215	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4489 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 Interleukin-6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	152	Total	C	N	O	S	0	1	0
			1203	759	209	227	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	EXPRESSION TAG	UNP P05231

- Molecule 2 is a protein called Heavy Chain of the Camelid Fab fragment 61H7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	200	Total	C	N	O	S	0	0	0
			1490	943	253	287	7			

- Molecule 3 is a protein called Light Chain of the Camelid Fab fragment 61H7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	211	Total	C	N	O	S	0	0	0
			1518	949	247	318	4			

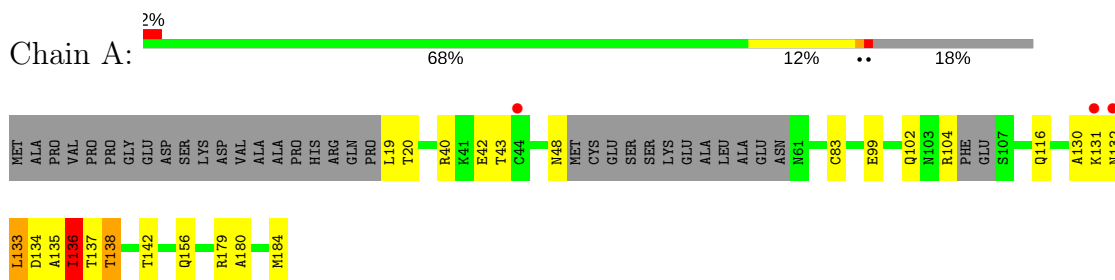
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	127	Total	O	0	0
			127	127		
4	H	76	Total	O	0	0
			76	76		
4	L	75	Total	O	0	0
			75	75		

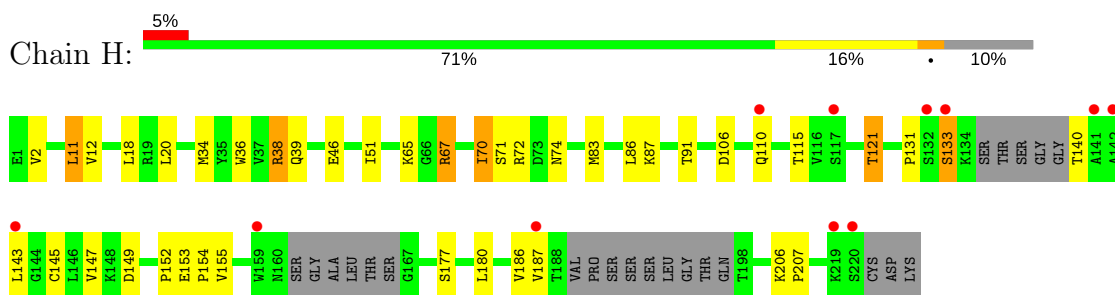
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

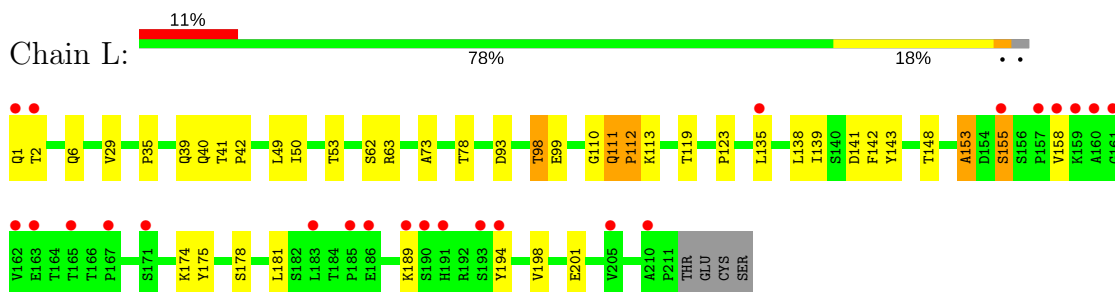
• Molecule 1: Interleukin-6



• Molecule 2: Heavy Chain of the Camelid Fab fragment 61H7



• Molecule 3: Light Chain of the Camelid Fab fragment 61H7



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	108.21Å 47.47Å 148.32Å 90.00° 97.00° 90.00°	Depositor
Resolution (Å)	38.65 – 2.42 38.65 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.7 (38.65-2.42) 99.5 (38.65-2.38)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 2.37Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.216 , 0.244 0.227 , 0.263	Depositor DCC
R_{free} test set	1447 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	43.3	Xtriage
Anisotropy	0.853	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 50.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4489	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.48% 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	0.52	0/1219	0.78	2/1636 (0.1%)
2	H	0.43	0/1526	0.75	0/2073
3	L	0.46	0/1556	0.80	1/2139 (0.0%)
All	All	0.47	0/4301	0.78	3/5848 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	135	ALA	C-N-CA	6.88	138.90	121.70
3	L	110	GLY	C-N-CA	6.16	137.10	121.70
1	A	136	ILE	C-N-CA	5.88	136.40	121.70

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	A	1203	0	1223	14	0
2	H	1490	0	1438	21	0
3	L	1518	0	1422	26	0
4	A	127	0	0	0	0
4	H	76	0	0	0	0
4	L	75	0	0	2	0
All	All	4489	0	4083	59	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 7.

All (59) 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)
3:L:111:GLN:HB3	3:L:112:PRO:HD2	1.33	1.07
3:L:35:PRO:HD2	3:L:53:THR:HG22	1.41	0.98
1:A:136:ILE:HD13	1:A:137:THR:HA	1.59	0.85
3:L:35:PRO:HD2	3:L:53:THR:CG2	2.13	0.78
3:L:111:GLN:HE21	3:L:112:PRO:HD2	1.51	0.75
3:L:153:ALA:HB2	3:L:194:TYR:CD2	2.25	0.72
2:H:36:TRP:HD1	2:H:70:ILE:HD11	1.54	0.72
3:L:35:PRO:CD	3:L:53:THR:HG22	2.21	0.70
3:L:111:GLN:HB3	3:L:112:PRO:CD	2.18	0.70
1:A:104:ARG:HH22	1:A:156:GLN:NE2	1.91	0.69
3:L:111:GLN:CB	3:L:112:PRO:HD2	2.20	0.66
3:L:113:LYS:HD2	3:L:201:GLU:HG3	1.78	0.65
1:A:180:ALA:HB1	1:A:184:MET:CE	2.26	0.65
2:H:51:ILE:HB	2:H:70:ILE:CG2	2.27	0.64
2:H:51:ILE:HD13	2:H:72:ARG:HB2	1.81	0.62
2:H:51:ILE:HB	2:H:70:ILE:HG21	1.81	0.62
2:H:39:GLN:HE22	3:L:40:GLN:HE22	1.48	0.61
3:L:158:VAL:HG11	3:L:181:LEU:HD11	1.81	0.61
1:A:136:ILE:HD12	1:A:136:ILE:H	1.67	0.59
3:L:39:GLN:HB2	3:L:49:LEU:HD11	1.82	0.59
2:H:36:TRP:CD1	2:H:70:ILE:HD11	2.38	0.59
1:A:136:ILE:CD1	1:A:137:THR:HA	2.31	0.58
1:A:133:LEU:HD23	1:A:134:ASP:H	1.70	0.57
3:L:153:ALA:HB2	3:L:194:TYR:CG	2.41	0.55
3:L:123:PRO:HD3	3:L:135:LEU:HD12	1.89	0.55
2:H:67:ARG:HG2	2:H:67:ARG:HH11	1.71	0.55
1:A:136:ILE:H	1:A:136:ILE:CD1	2.20	0.54
1:A:83:CYS:SG	1:A:184:MET:HE2	2.47	0.54
2:H:20:LEU:HD13	2:H:83:MET:SD	2.47	0.54
1:A:180:ALA:HB1	1:A:184:MET:HE1	1.89	0.53
3:L:53:THR:HG23	4:L:301:HOH:O	2.11	0.51
2:H:38:ARG:HG3	2:H:46:GLU:HB3	1.93	0.50
1:A:83:CYS:SG	1:A:184:MET:CE	3.00	0.49
3:L:141:ASP:HA	3:L:174:LYS:HB3	1.94	0.49
1:A:104:ARG:HH12	1:A:156:GLN:HE22	1.60	0.48
3:L:119:THR:HG23	3:L:138:LEU:HB2	1.95	0.48
2:H:121:THR:HG21	2:H:207:PRO:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:ILE:HG12	1:A:138:THR:HG22	1.95	0.47
3:L:139:ILE:HG12	3:L:198:VAL:HG11	1.97	0.47
2:H:186:VAL:HG21	3:L:138:LEU:HD13	1.97	0.46
2:H:11:LEU:HG	2:H:152:PRO:HG3	1.98	0.46
2:H:149:ASP:HB3	2:H:180:LEU:HD13	1.98	0.46
2:H:153:GLU:HG2	2:H:154:PRO:HA	1.98	0.46
3:L:63:ARG:HB2	3:L:78:THR:O	2.17	0.45
2:H:147:VAL:HG11	2:H:155:VAL:HG11	1.98	0.44
3:L:142:PHE:O	3:L:175:TYR:HB2	2.17	0.44
2:H:131:PRO:HG3	2:H:143:LEU:HB3	2.00	0.44
1:A:130:ALA:O	1:A:132:ASN:N	2.51	0.43
2:H:34:MET:HG2	2:H:72:ARG:HH21	1.84	0.43
1:A:102:GLN:NE2	1:A:116:GLN:NE2	2.67	0.43
3:L:174:LYS:HA	4:L:369:HOH:O	2.19	0.42
3:L:135:LEU:HD21	3:L:194:TYR:HB3	2.01	0.42
3:L:29:VAL:HG11	3:L:73:ALA:HB2	2.01	0.42
3:L:41:THR:HG23	3:L:42:PRO:HD2	2.00	0.42
3:L:93:ASP:OD1	3:L:98:THR:HB	2.21	0.41
2:H:20:LEU:CD1	2:H:83:MET:SD	3.09	0.41
2:H:91:THR:HG23	2:H:115:THR:HA	2.02	0.40
2:H:152:PRO:HD2	2:H:207:PRO:HB2	2.04	0.40
2:H:83:MET:HB3	2:H:86:LEU:HD21	2.03	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	A	147/186 (79%)	138 (94%)	7 (5%)	2 (1%)	13	17
2	H	192/223 (86%)	182 (95%)	9 (5%)	1 (0%)	32	45
3	L	209/215 (97%)	196 (94%)	10 (5%)	3 (1%)	13	17

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	548/624 (88%)	516 (94%)	26 (5%)	6 (1%)	17	23

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	131	LYS
1	A	136	ILE
2	H	133	SER
3	L	153	ALA
3	L	112	PRO
3	L	155	SER

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	133/167 (80%)	121 (91%)	12 (9%)	11	16
2	H	161/183 (88%)	141 (88%)	20 (12%)	5	6
3	L	165/178 (93%)	152 (92%)	13 (8%)	14	21
All	All	459/528 (87%)	414 (90%)	45 (10%)	9	13

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	20	THR
1	A	40	ARG
1	A	42	GLU
1	A	43	THR
1	A	48	ASN
1	A	99	GLU
1	A	133	LEU
1	A	136	ILE
1	A	138	THR

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Mol	Chain	Res	Type
1	A	142	THR
1	A	179	ARG
2	H	2	VAL
2	H	11	LEU
2	H	12	VAL
2	H	18	LEU
2	H	38	ARG
2	H	65	LYS
2	H	67	ARG
2	H	70	ILE
2	H	71	SER
2	H	74	ASN
2	H	87	LYS
2	H	106	ASP
2	H	110	GLN
2	H	121	THR
2	H	133	SER
2	H	140	THR
2	H	145	CYS
2	H	177	SER
2	H	187	VAL
2	H	206	LYS
3	L	1	GLN
3	L	2	THR
3	L	6	GLN
3	L	50	ILE
3	L	62	SER
3	L	98	THR
3	L	99	GLU
3	L	111	GLN
3	L	143	TYR
3	L	148	THR
3	L	155	SER
3	L	178	SER
3	L	189	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	102	GLN
1	A	116	GLN

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Mol	Chain	Res	Type
1	A	127	GLN
1	A	156	GLN
1	A	164	HIS
3	L	40	GLN
3	L	111	GLN
3	L	191	HIS

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 carbohydrates 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	152/186 (81%)	-0.11	3 (1%) 65 62	32, 48, 99, 128	0
2	H	200/223 (89%)	0.26	11 (5%) 26 24	36, 66, 104, 161	0
3	L	211/215 (98%)	0.48	24 (11%) 6 5	33, 70, 122, 134	0
All	All	563/624 (90%)	0.24	38 (6%) 19 16	32, 60, 111, 161	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	L	158	VAL	7.5
3	L	1	GLN	6.2
3	L	186	GLU	5.2
1	A	132	ASN	5.1
3	L	160	ALA	4.7
3	L	183	LEU	4.7
3	L	157	PRO	4.5
3	L	193	SER	4.5
2	H	219	LYS	4.4
3	L	205	VAL	3.9
2	H	187	VAL	3.8
2	H	141	ALA	3.7
2	H	142	ALA	3.6
1	A	44	CYS	3.5
3	L	2	THR	3.5
3	L	155	SER	3.5
3	L	190	SER	3.3
3	L	210	ALA	3.3
2	H	220	SER	3.1
3	L	162	VAL	2.9
3	L	161	GLY	2.9
2	H	133	SER	2.8
3	L	194	TYR	2.8

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Mol	Chain	Res	Type	RSRZ
2	H	159	TRP	2.7
3	L	189	LYS	2.7
3	L	163	GLU	2.7
1	A	131	LYS	2.7
2	H	110	GLN	2.6
2	H	132	SER	2.6
3	L	167	PRO	2.5
2	H	143	LEU	2.5
3	L	185	PRO	2.4
3	L	135	LEU	2.3
3	L	159	LYS	2.3
3	L	171	SER	2.3
3	L	165	THR	2.3
3	L	191	HIS	2.2
2	H	117	SER	2.2

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 carbohydrates 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.