



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:58 PM UTC

PDB ID : 3Q4J / pdb_00003q4j
Title : Structure of a small peptide ligand bound to E.coli DNA sliding clamp
Authors : Wolff, P.; Olieric, V.; Briand, J.P.; Chaloin, O.; Dejaegere, A.; Dumas, P.;
Ennifar, E.; Guichard, G.; Wagner, J.; Burnouf, D.
Deposited on : 2010-12-23
Resolution : 2.30 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

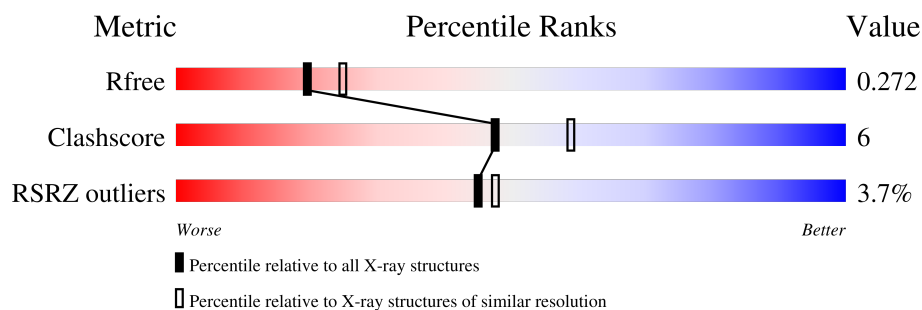
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	366	4% 84% 14% .
1	B	366	2% 87% 12% ..
1	C	366	5% 86% 13% .
1	D	366	4% 81% 18% .
1	E	366	3% 87% 11% ..
1	F	366	2% 84% 15% .
2	H	6	83% 17%

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Mol	Chain	Length	Quality of chain
2	I	6	50%100%
2	J	6	50%83%17%
2	K	6	67%33%
2	L	6	17%67%33%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17322 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 DNA polymerase III subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	366	Total	C	N	O	S	0	1	0
			2774	1747	479	530	18			
1	B	364	Total	C	N	O	S	0	0	0
			2792	1756	485	532	19			
1	C	365	Total	C	N	O	S	0	0	0
			2772	1748	472	533	19			
1	D	365	Total	C	N	O	S	0	0	0
			2779	1748	482	530	19			
1	E	364	Total	C	N	O	S	0	0	0
			2777	1744	482	532	19			
1	F	365	Total	C	N	O	S	0	0	0
			2800	1762	486	533	19			

- Molecule 2 is a protein called peptide ligand.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	H	6	Total	C	N	O	0	0	0
			48	32	6	10			
2	I	6	Total	C	N	O	0	0	0
			48	32	6	10			
2	J	6	Total	C	N	O	0	0	0
			48	32	6	10			
2	K	6	Total	C	N	O	0	0	0
			48	32	6	10			
2	L	6	Total	C	N	O	0	0	0
			48	32	6	10			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	48	Total	O	0	0
			48	48		

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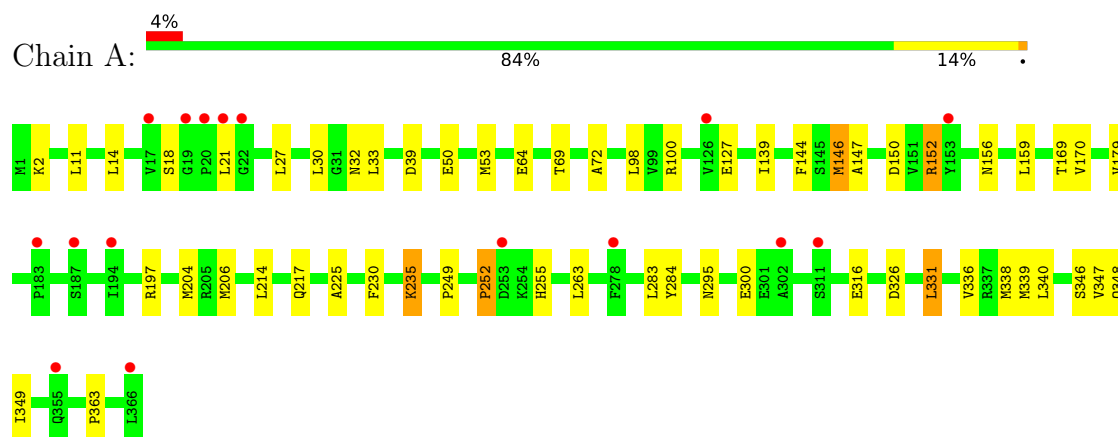
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	98	Total 98	O 98	0	0
3	C	44	Total 44	O 44	0	0
3	D	46	Total 46	O 46	0	0
3	E	75	Total 75	O 75	0	0
3	F	72	Total 72	O 72	0	0
3	H	3	Total 3	O 3	0	0
3	K	2	Total 2	O 2	0	0

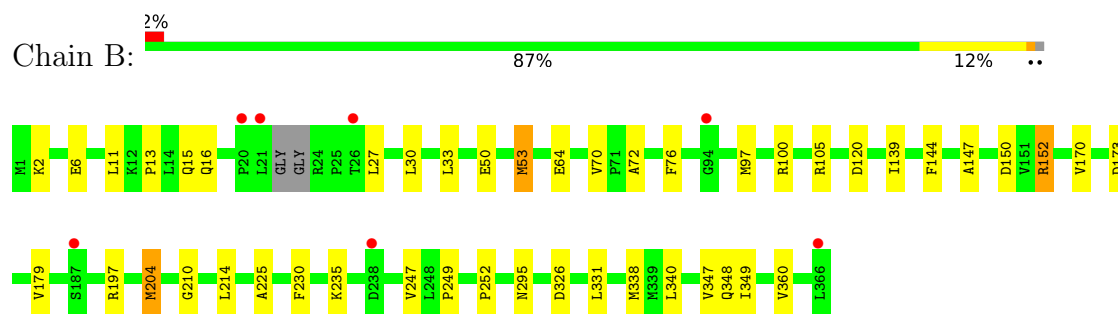
3 Residue-property plots [i](#)

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.

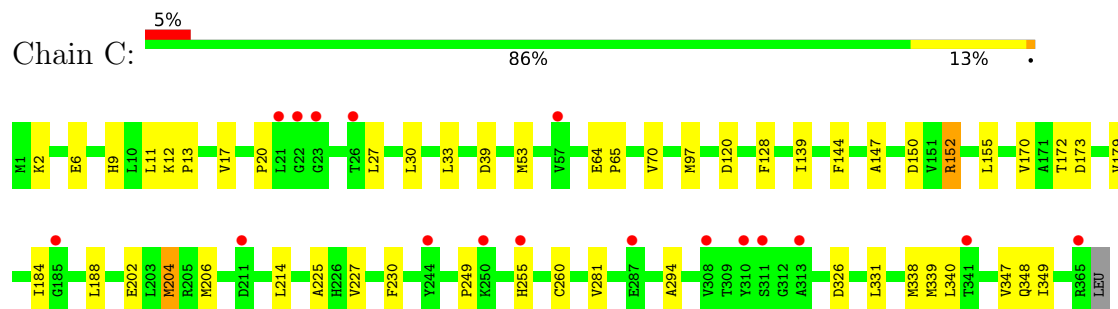
- Molecule 1: DNA polymerase III subunit beta



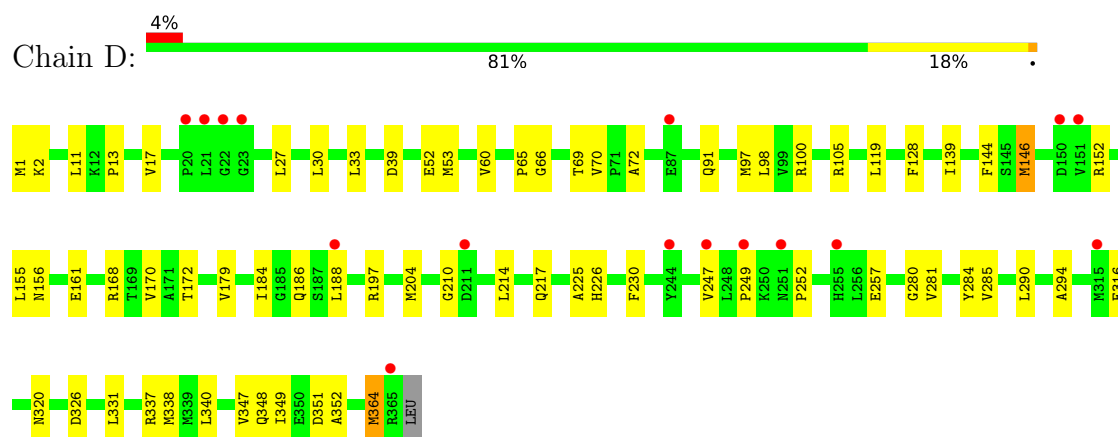
- Molecule 1: DNA polymerase III subunit beta



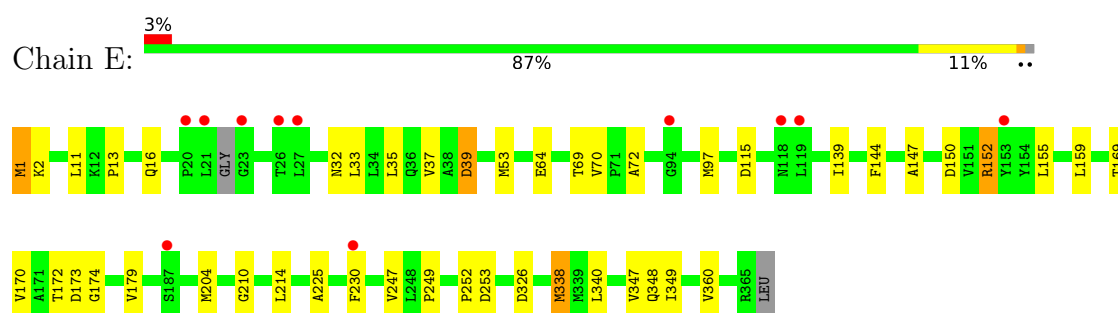
- Molecule 1: DNA polymerase III subunit beta



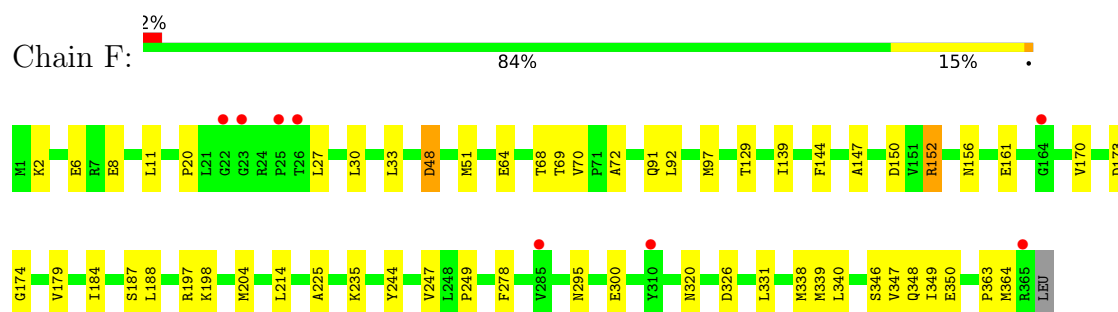
- Molecule 1: DNA polymerase III subunit beta



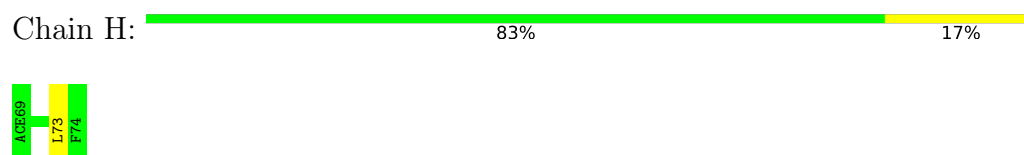
- Molecule 1: DNA polymerase III subunit beta



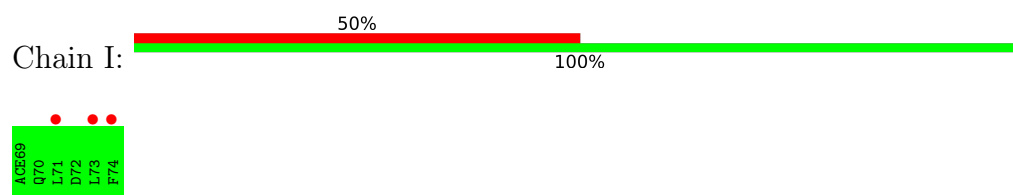
- Molecule 1: DNA polymerase III subunit beta



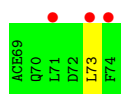
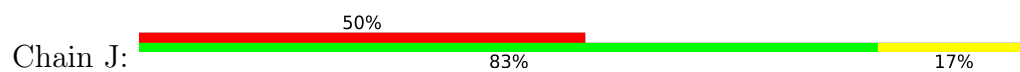
- Molecule 2: peptide ligand



- Molecule 2: peptide ligand



- Molecule 2: peptide ligand



- Molecule 2: peptide ligand



- Molecule 2: peptide ligand



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	35.09Å 132.87Å 137.27Å 62.73° 88.51° 89.77°	Depositor
Resolution (Å)	29.52 – 2.30 29.52 – 2.30	Depositor EDS
% Data completeness (in resolution range)	(Not available) (29.52-2.30) 98.5 (29.52-2.30)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.50 (at 2.31Å)	Xtriage
Refinement program	BUSTER-TNT BUSTER 2.8.0, BUSTER 2.8.0	Depositor
R, R_{free}	0.214 , 0.254 (Not available) , 0.272	Depositor DCC
R_{free} test set	1959 reflections (2.03%)	wwPDB-VP
Wilson B-factor (Å ²)	42.4	Xtriage
Anisotropy	0.487	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 35.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.024 for h,-k,-l 0.000 for -h,k,k-l 0.000 for -h,-k,-k+l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17322	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.06% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.75	1/2827 (0.0%)	1.18	8/3841 (0.2%)
1	B	0.85	1/2840 (0.0%)	1.17	12/3851 (0.3%)
1	C	0.80	0/2821	1.21	9/3829 (0.2%)
1	D	0.79	2/2826 (0.1%)	1.20	11/3837 (0.3%)
1	E	0.83	2/2825 (0.1%)	1.19	8/3831 (0.2%)
1	F	0.85	0/2847	1.19	8/3860 (0.2%)
2	H	0.84	0/46	0.96	0/60
2	I	0.77	0/46	1.02	0/60
2	J	0.69	0/46	1.06	0/60
2	K	0.76	0/46	0.96	0/60
2	L	0.75	0/46	0.97	0/60
All	All	0.81	6/17216 (0.0%)	1.19	56/23349 (0.2%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	146	MET	SD-CE	-8.85	1.57	1.79
1	E	338	MET	SD-CE	-7.50	1.60	1.79
1	A	146	MET	SD-CE	-6.61	1.63	1.79
1	E	1	MET	SD-CE	-6.09	1.64	1.79
1	D	364	MET	SD-CE	-5.48	1.65	1.79
1	B	53	MET	SD-CE	-5.16	1.66	1.79

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	184	ILE	CB-CA-C	8.45	120.22	110.93
1	E	39	ASP	N-CA-C	7.38	126.51	110.80
1	B	331	LEU	N-CA-C	7.12	118.82	111.14
1	D	331	LEU	N-CA-C	6.76	118.30	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	331	LEU	N-CA-C	6.75	118.30	111.07
1	A	331	LEU	N-CA-C	6.72	118.26	111.07
1	D	152	ARG	CA-C-N	6.67	129.51	120.38
1	D	152	ARG	C-N-CA	6.67	129.51	120.38
1	C	331	LEU	N-CA-C	6.42	118.07	111.14
1	C	120	ASP	CA-CB-CG	6.42	119.02	112.60
1	D	11	LEU	N-CA-C	6.39	117.93	110.97
1	A	11	LEU	N-CA-C	6.31	117.85	110.97
1	D	351	ASP	N-CA-C	-6.21	100.39	110.20
1	E	11	LEU	N-CA-C	6.17	117.69	110.97
1	B	11	LEU	N-CA-C	6.13	117.65	110.97
1	A	39	ASP	CA-CB-CG	6.02	118.62	112.60
1	B	210	GLY	CA-C-N	5.83	128.67	120.28
1	B	210	GLY	C-N-CA	5.83	128.67	120.28
1	F	11	LEU	N-CA-C	5.81	117.31	110.97
1	C	11	LEU	N-CA-C	5.77	117.26	110.97
1	E	115	ASP	CA-CB-CG	5.74	118.33	112.60
1	C	204	MET	CA-C-N	5.68	127.89	120.28
1	C	204	MET	C-N-CA	5.68	127.89	120.28
1	E	210	GLY	CA-C-N	5.60	128.34	120.28
1	E	210	GLY	C-N-CA	5.60	128.34	120.28
1	D	210	GLY	CA-C-N	5.57	128.30	120.28
1	D	210	GLY	C-N-CA	5.57	128.30	120.28
1	F	152	ARG	CA-C-N	5.42	127.81	120.38
1	F	152	ARG	C-N-CA	5.42	127.81	120.38
1	B	152	ARG	CA-C-N	5.42	127.81	120.38
1	B	152	ARG	C-N-CA	5.42	127.81	120.38
1	F	129	THR	CB-CA-C	5.37	119.90	109.33
1	E	253	ASP	CA-CB-CG	5.33	117.93	112.60
1	B	204	MET	CA-C-N	5.30	127.39	120.28
1	B	204	MET	C-N-CA	5.30	127.39	120.28
1	E	152	ARG	CA-C-N	5.28	127.62	120.38
1	E	152	ARG	C-N-CA	5.28	127.62	120.38
1	A	252	PRO	CA-C-N	5.22	129.50	121.19
1	A	252	PRO	C-N-CA	5.22	129.50	121.19
1	D	60	VAL	CA-C-N	5.21	128.61	120.68
1	D	60	VAL	C-N-CA	5.21	128.61	120.68
1	F	48	ASP	CA-CB-CG	5.19	117.79	112.60
1	C	152	ARG	CA-C-N	5.17	127.46	120.38
1	C	152	ARG	C-N-CA	5.17	127.46	120.38
1	B	120	ASP	CA-CB-CG	5.15	117.75	112.60
1	B	295	ASN	CA-CB-CG	5.14	117.75	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	320	ASN	N-CA-C	-5.11	101.79	109.15
1	A	152	ARG	CA-C-N	5.10	131.27	121.54
1	A	152	ARG	C-N-CA	5.10	131.27	121.54
1	A	235	LYS	N-CA-C	-5.09	103.32	110.50
1	F	235	LYS	N-CA-C	-5.07	103.36	110.50
1	C	184	ILE	CA-CB-CG2	5.06	119.10	110.50
1	D	320	ASN	N-CA-C	-5.03	101.91	109.15
1	D	184	ILE	CB-CA-C	5.02	115.47	111.06
1	B	6	GLU	CA-C-N	5.01	127.24	120.38
1	B	6	GLU	C-N-CA	5.01	127.24	120.38

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	2774	0	2722	37	0
1	B	2792	0	2771	27	0
1	C	2772	0	2731	27	0
1	D	2779	0	2738	41	0
1	E	2777	0	2726	29	0
1	F	2800	0	2789	30	0
2	H	48	0	46	2	0
2	I	48	0	46	0	0
2	J	48	0	46	1	0
2	K	48	0	46	3	0
2	L	48	0	46	2	0
3	A	48	0	0	0	0
3	B	98	0	0	0	0
3	C	44	0	0	1	0
3	D	46	0	0	0	0
3	E	75	0	0	0	0
3	F	72	0	0	0	0
3	H	3	0	0	0	0
3	K	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	17322	0	16707	188	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 6.

All (188) 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:32:ASN:HB3	1:A:69:THR:HG22	1.28	1.13
1:D:280:GLY:HA2	1:D:364:MET:HE1	1.31	1.11
1:E:32:ASN:HB3	1:E:69:THR:HG22	1.33	1.05
1:D:280:GLY:CA	1:D:364:MET:HE1	2.02	0.89
1:D:1:MET:HB3	1:D:66:GLY:HA3	1.57	0.86
1:E:1:MET:HE1	1:E:35:LEU:HB3	1.66	0.78
1:C:260:CYS:SG	3:C:415:HOH:O	2.42	0.77
1:B:16:GLN:HE21	1:B:53:MET:HE2	1.49	0.77
1:A:127:GLU:HB2	1:C:65:PRO:HG3	1.70	0.73
1:A:32:ASN:CB	1:A:69:THR:HG22	2.14	0.73
1:D:280:GLY:HA2	1:D:364:MET:CE	2.15	0.73
1:A:98:LEU:HD23	1:A:100:ARG:HH12	1.50	0.72
1:E:249:PRO:HD2	1:E:348:GLN:HE21	1.54	0.71
1:F:2:LYS:HB2	1:F:64:GLU:HB2	1.73	0.70
1:A:249:PRO:HD2	1:A:348:GLN:HE21	1.56	0.70
1:A:206:MET:HE3	1:A:230:PHE:HD2	1.57	0.69
1:B:249:PRO:HD2	1:B:348:GLN:HE21	1.57	0.69
1:C:214:LEU:HD11	1:C:225:ALA:HB1	1.75	0.68
1:A:206:MET:HE3	1:A:230:PHE:CD2	2.28	0.68
1:C:249:PRO:HD2	1:C:348:GLN:HE21	1.58	0.67
1:E:214:LEU:HD11	1:E:225:ALA:HB1	1.76	0.67
1:E:1:MET:HE3	1:E:37:VAL:HG23	1.76	0.66
1:E:32:ASN:CB	1:E:69:THR:HG22	2.19	0.66
1:D:214:LEU:HD11	1:D:225:ALA:HB1	1.77	0.66
1:D:98:LEU:HD23	1:D:100:ARG:HH12	1.59	0.66
1:A:249:PRO:HB2	1:A:252:PRO:HG3	1.78	0.66
1:F:139:ILE:HG21	1:F:204:MET:HG2	1.76	0.66
1:A:214:LEU:HD11	1:A:225:ALA:HB1	1.78	0.65
1:B:100:ARG:HG2	1:B:105:ARG:HG2	1.78	0.65
1:D:1:MET:HB3	1:D:66:GLY:CA	2.26	0.65
1:C:338:MET:HG2	1:C:349:ILE:HG12	1.79	0.65
1:D:249:PRO:HD2	1:D:348:GLN:HE21	1.62	0.64
1:F:214:LEU:HD11	1:F:225:ALA:HB1	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:249:PRO:HD2	1:F:348:GLN:HE21	1.64	0.62
1:B:53:MET:CE	1:B:230:PHE:HB3	2.30	0.62
1:C:206:MET:HE2	1:C:227:VAL:HG12	1.83	0.61
1:D:139:ILE:HG21	1:D:204:MET:HG2	1.82	0.61
1:E:139:ILE:HG21	1:E:204:MET:HG2	1.83	0.60
1:B:139:ILE:HG21	1:B:204:MET:HG2	1.84	0.60
1:C:281:VAL:HG12	1:C:294:ALA:HB2	1.84	0.60
1:B:214:LEU:HD11	1:B:225:ALA:HB1	1.82	0.59
1:C:139:ILE:HG21	1:C:204:MET:HG2	1.84	0.59
1:F:20:PRO:O	1:F:48:ASP:HB3	2.03	0.58
1:F:70:VAL:HG11	1:F:97:MET:SD	2.44	0.58
1:E:39:ASP:CG	1:E:39:ASP:O	2.47	0.58
1:E:247:VAL:HG13	2:K:73:LEU:HD22	1.86	0.58
1:C:206:MET:HE3	1:C:230:PHE:CD2	2.39	0.57
1:D:128:PHE:HA	1:D:186:GLN:HE22	1.71	0.56
1:F:68:THR:HG21	1:F:92:LEU:HD21	1.87	0.56
1:A:139:ILE:HG21	1:A:204:MET:HG2	1.87	0.56
1:F:6:GLU:OE2	1:F:8:GLU:HG2	2.06	0.55
1:D:338:MET:HE3	1:D:347:VAL:HG21	1.89	0.55
1:D:65:PRO:HD2	1:F:187:SER:O	2.07	0.55
1:E:340:LEU:HG	1:E:347:VAL:HG23	1.89	0.55
1:C:20:PRO:HG2	1:C:53:MET:HG2	1.90	0.54
1:E:1:MET:CE	1:E:37:VAL:HG23	2.38	0.54
1:A:284:TYR:CE1	1:A:316:GLU:HG3	2.43	0.54
1:B:170:VAL:HG22	1:B:179:VAL:HG23	1.89	0.54
1:F:247:VAL:HG13	2:L:73:LEU:HD22	1.90	0.54
1:B:338:MET:HE3	1:B:347:VAL:HG21	1.90	0.53
1:A:32:ASN:HB3	1:A:69:THR:CG2	2.20	0.53
1:C:128:PHE:HB3	1:C:188:LEU:HD21	1.91	0.53
1:E:338:MET:HG2	1:E:349:ILE:HG12	1.91	0.53
1:D:146:MET:HE1	1:D:197:ARG:HA	1.90	0.52
1:B:53:MET:HE1	1:B:230:PHE:HB3	1.89	0.52
1:D:249:PRO:HB2	1:D:252:PRO:HG3	1.91	0.52
1:D:284:TYR:CE1	1:D:316:GLU:HG3	2.44	0.52
1:A:127:GLU:CD	1:C:39:ASP:H	2.17	0.52
1:D:340:LEU:HG	1:D:347:VAL:HG23	1.89	0.52
1:F:340:LEU:HG	1:F:347:VAL:HG23	1.91	0.52
1:A:146:MET:HE1	1:A:197:ARG:HA	1.91	0.52
1:A:340:LEU:HG	1:A:347:VAL:HG23	1.91	0.51
1:B:150:ASP:OD2	1:B:152:ARG:HD3	2.11	0.51
1:A:2:LYS:HB3	1:A:64:GLU:HB3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:39:ASP:OD2	1:D:39:ASP:CB	2.58	0.51
1:E:249:PRO:HB2	1:E:252:PRO:HG3	1.93	0.51
1:B:249:PRO:HB2	1:B:252:PRO:HG3	1.92	0.51
1:C:340:LEU:HG	1:C:347:VAL:HG23	1.93	0.51
1:E:170:VAL:HG22	1:E:179:VAL:HG23	1.92	0.51
1:F:278:PHE:O	1:F:364:MET:HE1	2.11	0.51
1:B:340:LEU:HG	1:B:347:VAL:HG23	1.92	0.51
1:A:170:VAL:HG22	1:A:179:VAL:HG23	1.92	0.50
1:C:2:LYS:HB3	1:C:64:GLU:HB3	1.92	0.50
1:C:6:GLU:HB3	1:C:9:HIS:HD2	1.76	0.50
1:D:52:GLU:HG2	1:D:119:LEU:HD22	1.93	0.50
1:B:144:PHE:CD2	1:B:326:ASP:HB3	2.47	0.50
1:F:338:MET:HE3	1:F:347:VAL:HG21	1.93	0.50
1:B:13:PRO:HA	1:B:230:PHE:HE1	1.76	0.50
1:D:13:PRO:HA	1:D:230:PHE:HE1	1.77	0.50
1:D:280:GLY:CA	1:D:364:MET:CE	2.82	0.50
1:D:33:LEU:HG	1:D:72:ALA:HB2	1.94	0.50
1:D:338:MET:HG2	1:D:349:ILE:HG12	1.93	0.50
1:E:360:VAL:HG12	2:K:73:LEU:HD21	1.94	0.50
1:F:338:MET:HG2	1:F:349:ILE:HG12	1.94	0.50
1:B:2:LYS:HB3	1:B:64:GLU:HB3	1.94	0.49
1:F:184:ILE:HD11	1:F:188:LEU:HD11	1.95	0.49
1:E:13:PRO:HA	1:E:230:PHE:HE1	1.77	0.49
1:D:100:ARG:HG2	1:D:105:ARG:HG3	1.94	0.49
1:B:53:MET:HE3	1:B:230:PHE:HB3	1.94	0.49
1:E:2:LYS:HB3	1:E:64:GLU:HB2	1.94	0.49
1:E:150:ASP:OD2	1:E:152:ARG:HD3	2.12	0.49
1:A:150:ASP:OD1	1:A:152:ARG:HD3	2.12	0.49
1:A:144:PHE:CD2	1:A:326:ASP:HB3	2.48	0.48
1:B:247:VAL:HG13	2:H:73:LEU:HD22	1.95	0.48
1:C:255:HIS:HD2	1:C:339:MET:HG2	1.78	0.48
1:D:70:VAL:HG11	1:D:97:MET:SD	2.53	0.48
1:C:170:VAL:HG22	1:C:179:VAL:HG23	1.96	0.48
1:A:53:MET:HB3	1:A:230:PHE:CZ	2.49	0.48
1:F:156:ASN:O	1:F:197:ARG:HB2	2.14	0.48
1:D:257:GLU:HG2	1:D:337:ARG:HG2	1.96	0.48
1:E:16:GLN:HB3	1:E:53:MET:HE2	1.95	0.47
1:D:170:VAL:HG22	1:D:179:VAL:HG23	1.95	0.47
1:C:338:MET:HE3	1:C:347:VAL:HG21	1.96	0.47
1:C:144:PHE:CD2	1:C:326:ASP:HB3	2.49	0.47
1:F:150:ASP:OD2	1:F:152:ARG:HD3	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:LEU:H	1:A:21:LEU:HD12	1.80	0.47
1:F:33:LEU:HG	1:F:72:ALA:HB2	1.96	0.47
1:F:147:ALA:HB2	1:F:173:ASP:HA	1.97	0.47
1:C:70:VAL:HG11	1:C:97:MET:SD	2.55	0.47
1:F:161:GLU:OE1	1:F:244:TYR:OH	2.25	0.46
1:F:170:VAL:HG22	1:F:179:VAL:HG23	1.98	0.46
1:B:360:VAL:HG12	2:H:73:LEU:HD21	1.96	0.46
1:C:17:VAL:HG23	1:C:33:LEU:HD11	1.97	0.46
1:A:127:GLU:HG3	1:A:217:GLN:HG2	1.98	0.46
1:D:247:VAL:HG13	2:J:73:LEU:HD22	1.98	0.46
1:F:2:LYS:HD2	1:F:91:GLN:HB2	1.97	0.46
1:A:338:MET:HG2	1:A:349:ILE:HG12	1.97	0.46
1:B:33:LEU:HG	1:B:72:ALA:HB2	1.98	0.46
1:D:144:PHE:CD2	1:D:326:ASP:HB3	2.50	0.46
1:A:147:ALA:HB1	1:A:150:ASP:HB2	1.98	0.46
1:E:33:LEU:HG	1:E:72:ALA:HB2	1.98	0.45
1:E:155:LEU:HD22	1:E:172:THR:HG23	1.97	0.45
1:C:20:PRO:HG3	1:C:202:GLU:OE1	2.16	0.45
1:D:281:VAL:HG12	1:D:294:ALA:HB2	1.98	0.45
1:D:17:VAL:HG12	1:D:53:MET:HG3	1.97	0.45
1:F:33:LEU:O	1:F:69:THR:HA	2.16	0.45
1:E:144:PHE:CD2	1:E:326:ASP:HB3	2.51	0.45
1:A:50:GLU:HA	1:A:235:LYS:HD2	1.99	0.44
1:B:147:ALA:O	1:B:197:ARG:NH1	2.46	0.44
1:C:150:ASP:OD2	1:C:152:ARG:HD3	2.18	0.44
1:C:155:LEU:HD22	1:C:172:THR:HG23	1.99	0.44
1:D:249:PRO:HD2	1:D:348:GLN:NE2	2.32	0.44
1:E:147:ALA:HB2	1:E:173:ASP:HA	1.98	0.44
1:E:174:GLY:HA2	2:K:74:PHE:CZ	2.52	0.44
1:B:147:ALA:HB2	1:B:173:ASP:HA	1.98	0.44
1:A:33:LEU:HG	1:A:72:ALA:HB2	2.00	0.44
1:F:174:GLY:HA2	2:L:74:PHE:CZ	2.53	0.43
1:A:255[A]:HIS:NE2	1:A:339:MET:HG2	2.33	0.43
1:D:2:LYS:HG2	1:D:91:GLN:CB	2.48	0.43
1:C:147:ALA:HB2	1:C:173:ASP:HA	2.00	0.43
1:D:155:LEU:HD22	1:D:172:THR:HG23	2.01	0.43
1:E:70:VAL:HG11	1:E:97:MET:SD	2.58	0.43
1:E:147:ALA:HB1	1:E:150:ASP:HB2	2.01	0.43
1:A:263:LEU:HD23	1:A:336:VAL:HG21	1.99	0.43
1:F:346:SER:HA	1:F:363:PRO:HD3	2.01	0.43
1:B:70:VAL:HG11	1:B:97:MET:SD	2.58	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:147:ALA:HB1	1:B:150:ASP:HB2	2.00	0.43
1:D:27:LEU:HB2	1:D:30:LEU:HG	2.01	0.43
1:F:144:PHE:CD2	1:F:326:ASP:HB3	2.54	0.43
1:A:331:LEU:HD13	1:A:336:VAL:HG12	2.00	0.42
1:B:27:LEU:HB2	1:B:30:LEU:HG	2.01	0.42
1:A:338:MET:HE3	1:A:347:VAL:HG21	2.02	0.42
1:B:338:MET:HG2	1:B:349:ILE:HG12	2.01	0.42
1:D:161:GLU:OE1	1:D:168:ARG:NH2	2.53	0.42
1:D:33:LEU:O	1:D:69:THR:HA	2.19	0.42
1:D:337:ARG:HG3	1:D:352:ALA:HA	2.02	0.42
1:A:14:LEU:O	1:A:18:SER:HB3	2.20	0.42
1:A:346:SER:HA	1:A:363:PRO:HD3	2.02	0.42
1:F:51:MET:HE1	1:F:198:LYS:HB3	2.02	0.42
1:B:15:GLN:HG3	1:B:76:PHE:CE1	2.55	0.41
1:C:27:LEU:HB2	1:C:30:LEU:HG	2.01	0.41
1:C:12:LYS:HB3	1:C:13:PRO:HD3	2.01	0.41
1:A:27:LEU:HB2	1:A:30:LEU:HG	2.01	0.41
1:A:159:LEU:O	1:A:169:THR:HA	2.21	0.41
1:D:217:GLN:NE2	1:D:226:HIS:NE2	2.68	0.41
1:A:295:ASN:HA	1:A:300:GLU:O	2.21	0.41
1:D:156:ASN:O	1:D:197:ARG:HB2	2.21	0.41
1:D:285:VAL:HG12	1:D:290:LEU:HD13	2.03	0.41
1:E:53:MET:HE1	1:E:230:PHE:HB3	2.03	0.41
1:A:283:LEU:O	1:A:316:GLU:HA	2.21	0.41
1:F:27:LEU:HB2	1:F:30:LEU:HG	2.03	0.41
1:E:159:LEU:O	1:E:169:THR:HA	2.22	0.40
1:F:339:MET:CE	1:F:350:GLU:HG2	2.51	0.40
1:E:249:PRO:HD2	1:E:348:GLN:NE2	2.30	0.40
1:A:156:ASN:O	1:A:197:ARG:HB2	2.21	0.40
1:B:50:GLU:HA	1:B:235:LYS:HD2	2.03	0.40
1:D:128:PHE:HB3	1:D:188:LEU:HD21	2.04	0.40
1:F:295:ASN:HA	1:F:300:GLU:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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 oligosaccharides 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	366/366 (100%)	0.64	16 (4%) 39 41	27, 56, 85, 104	2 (0%)
1	B	364/366 (99%)	0.21	7 (1%) 66 68	23, 45, 72, 97	2 (0%)
1	C	365/366 (99%)	0.60	17 (4%) 36 38	31, 55, 80, 101	1 (0%)
1	D	365/366 (99%)	0.61	16 (4%) 39 41	32, 54, 83, 95	1 (0%)
1	E	364/366 (99%)	0.49	11 (3%) 52 54	26, 51, 76, 107	0
1	F	365/366 (99%)	0.25	8 (2%) 62 64	29, 47, 75, 99	1 (0%)
2	H	5/6 (83%)	0.48	0 100 100	44, 45, 58, 59	0
2	I	5/6 (83%)	1.80	3 (60%) 0 0	62, 67, 74, 78	0
2	J	5/6 (83%)	2.11	3 (60%) 0 0	78, 90, 100, 108	0
2	K	5/6 (83%)	0.64	0 100 100	54, 54, 65, 66	0
2	L	5/6 (83%)	1.52	1 (20%) 3 3	66, 72, 81, 86	0
All	All	2214/2226 (99%)	0.48	82 (3%) 45 48	23, 52, 81, 108	7 (0%)

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	20	PRO	4.7
1	A	21	LEU	4.5
1	A	366	LEU	4.0
1	E	21	LEU	3.8
1	B	20	PRO	3.7
1	D	315	MET	3.7
1	A	20	PRO	3.7
1	D	244	TYR	3.6
1	C	57	VAL	3.6
1	E	23	GLY	3.5
1	B	21	LEU	3.4
2	L	74	PHE	3.3

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Mol	Chain	Res	Type	RSRZ
1	D	247	VAL	3.3
1	B	94	GLY	3.2
1	D	365	ARG	3.1
2	J	74	PHE	3.0
2	J	73	LEU	3.0
1	C	310	TYR	3.0
1	C	23	GLY	3.0
1	C	244	TYR	3.0
1	F	26	THR	2.9
1	A	153	TYR	2.9
1	F	365	ARG	2.9
1	C	211	ASP	2.9
1	A	187	SER	2.8
1	E	187	SER	2.8
1	E	118	ASN	2.7
1	D	23	GLY	2.7
1	E	119	LEU	2.7
1	C	26	THR	2.7
1	D	211	ASP	2.7
1	C	22	GLY	2.7
1	E	94	GLY	2.6
1	B	366	LEU	2.6
1	D	251	ASN	2.6
1	C	255	HIS	2.6
1	B	187	SER	2.5
1	D	20	PRO	2.5
1	D	249	PRO	2.5
1	F	285	VAL	2.5
2	J	71	LEU	2.4
1	E	230	PHE	2.4
1	A	355	GLN	2.4
1	B	26	THR	2.4
1	D	21	LEU	2.4
1	C	365	ARG	2.4
2	I	74	PHE	2.3
1	F	310	TYR	2.3
1	D	22	GLY	2.3
1	D	151	VAL	2.3
1	D	255	HIS	2.3
1	C	21	LEU	2.3
1	B	238	ASP	2.3
1	A	19	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	153	TYR	2.3
1	C	308	VAL	2.3
1	E	27	LEU	2.3
1	A	22	GLY	2.3
1	F	22	GLY	2.2
1	A	183	PRO	2.2
1	A	311	SER	2.2
1	F	164	GLY	2.2
1	A	302	ALA	2.2
1	A	126	VAL	2.1
1	F	23	GLY	2.1
1	C	250	LYS	2.1
2	I	73	LEU	2.1
1	D	188	LEU	2.1
1	A	17	VAL	2.1
1	F	25	PRO	2.1
1	A	278	PHE	2.1
1	E	26	THR	2.1
1	C	185	GLY	2.1
1	C	311	SER	2.0
1	C	287	GLU	2.0
1	A	194	ILE	2.0
1	C	341	THR	2.0
1	D	87	GLU	2.0
1	C	313	ALA	2.0
1	A	253	ASP	2.0
1	D	150	ASP	2.0
2	I	71	LEU	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 oligosaccharides 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.