



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:44 AM GMT

PDB ID : 3MVK
Title : The Crystal Structure of FucU from Bifidobacterium longum to 1.65Å
Authors : Stein, A.J.; Xu, X.; Cui, H.; Savchenko, A.; Joachimiak, A.; Midwest Center
for Structural Genomics (MCSG)
Deposited on : 2010-05-04
Resolution : 1.65 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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

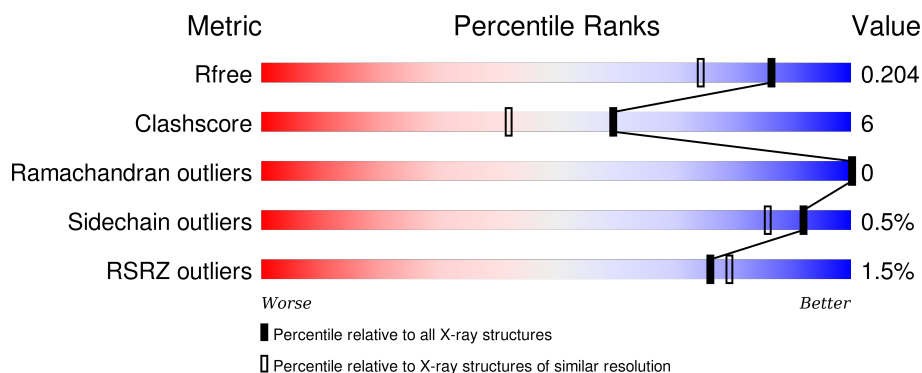
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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}	91344	1226 (1.66-1.66)
Clashscore	102246	1323 (1.66-1.66)
Ramachandran outliers	100387	1295 (1.66-1.66)
Sidechain outliers	100360	1295 (1.66-1.66)
RSRZ outliers	91569	1227 (1.66-1.66)

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	148	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between;"> % 85% 9% • 5% </div> </div>
1	B	148	<div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="display: flex; justify-content: space-between;"> 83% 11% • 5% </div> </div>
1	C	148	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between;"> 4% 79% 15% • 5% </div> </div>
1	D	148	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between;"> % 89% 5% • 5% </div> </div>
1	E	148	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between;"> % 88% 7% • 5% </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	148	 84% 9% • 5%
1	G	148	 88% 8% • •
1	H	148	 89% 5% • 5%
1	I	148	 86% 8% • 5%
1	J	148	 86% 9% • 5%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NA	A	149	-	-	-	X
2	NA	B	149	-	-	-	X
3	PGE	C	150	-	-	-	X
3	PGE	C	151	-	-	X	X
3	PGE	E	151	-	-	-	X
3	PGE	E	152	-	-	-	X
3	PGE	G	150	-	-	-	X
3	PGE	H	151	-	-	-	X
3	PGE	J	150	-	-	X	X
4	GOL	A	152	-	-	-	X
4	GOL	B	152	-	-	X	X
4	GOL	E	153	-	-	X	X
4	GOL	F	152	-	-	-	X
4	GOL	H	152	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12068 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 protein FucU.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	140	Total	C	N	O	S	Se	0	2	0
			1054	679	171	198	3	3			
1	B	140	Total	C	N	O	S	Se	0	1	0
			1053	676	171	200	3	3			
1	C	141	Total	C	N	O	S	Se	0	1	0
			1064	685	172	201	3	3			
1	D	140	Total	C	N	O	S	Se	0	1	0
			1055	681	172	196	3	3			
1	E	141	Total	C	N	O	S	Se	0	1	0
			1061	684	172	199	3	3			
1	F	140	Total	C	N	O	S	Se	0	1	0
			1049	674	171	198	3	3			
1	G	143	Total	C	N	O	S	Se	0	2	0
			1068	687	173	202	3	3			
1	H	141	Total	C	N	O	S	Se	0	2	0
			1063	685	172	200	3	3			
1	I	141	Total	C	N	O	S	Se	0	2	0
			1059	683	171	199	3	3			
1	J	141	Total	C	N	O	S	Se	0	1	0
			1064	686	173	199	3	3			

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

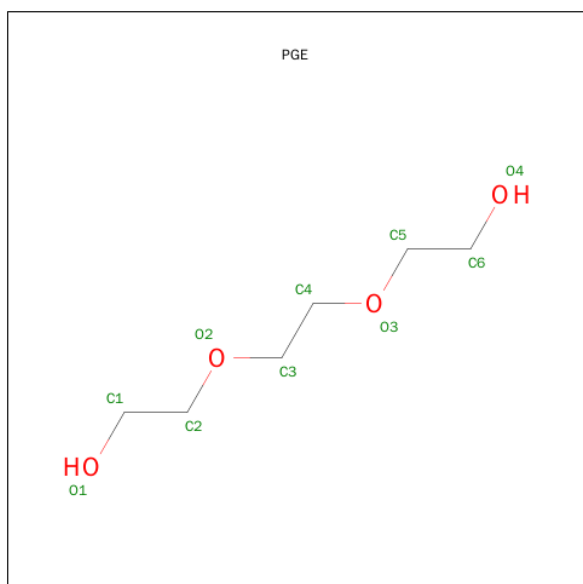
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	1	Total	Na	0	0
			1	1		
2	J	1	Total	Na	0	0
			1	1		
2	D	1	Total	Na	0	0
			1	1		
2	E	2	Total	Na	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	H	2	Total	Na	0	0
			2	2		
2	B	2	Total	Na	0	0
			2	2		
2	I	1	Total	Na	0	0
			1	1		
2	C	1	Total	Na	0	0
			1	1		
2	A	2	Total	Na	0	0
			2	2		
2	F	2	Total	Na	0	0
			2	2		

- Molecule 3 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



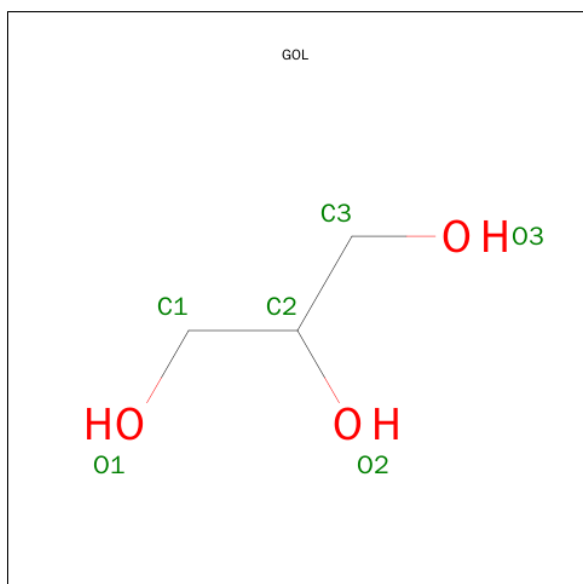
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			10	6	4		
3	B	1	Total	C	O	0	0
			10	6	4		
3	C	1	Total	C	O	0	0
			10	6	4		
3	C	1	Total	C	O	0	0
			10	6	4		
3	E	1	Total	C	O	0	0
			10	6	4		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	E	1	Total	C	O	0	0
			10	6	4		
3	F	1	Total	C	O	0	0
			10	6	4		
3	G	1	Total	C	O	0	0
			10	6	4		
3	H	1	Total	C	O	0	0
			10	6	4		
3	J	1	Total	C	O	0	0
			10	6	4		
3	J	1	Total	C	O	0	0
			10	6	4		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	E	1	Total	C	O	0	0
			6	3	3		
4	F	1	Total	C	O	0	0
			6	3	3		
4	G	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	H	1	Total	C	O	0	0
			6	3	3		

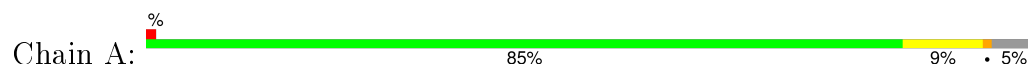
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	137	Total	O	0	0
			137	137		
5	B	139	Total	O	0	0
			139	139		
5	C	116	Total	O	0	0
			116	116		
5	D	121	Total	O	0	0
			121	121		
5	E	136	Total	O	0	0
			136	136		
5	F	120	Total	O	0	0
			120	120		
5	G	156	Total	O	0	0
			156	156		
5	H	138	Total	O	0	0
			138	138		
5	I	121	Total	O	0	0
			121	121		
5	J	133	Total	O	0	0
			133	133		

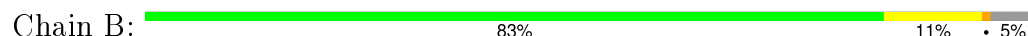
3 Residue-property plots [i](#)

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

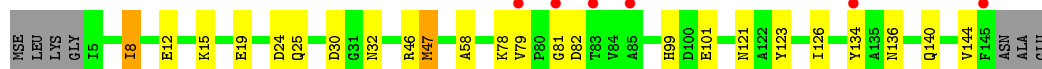
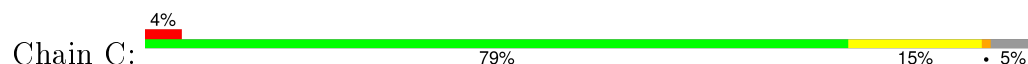
- Molecule 1: protein FucU



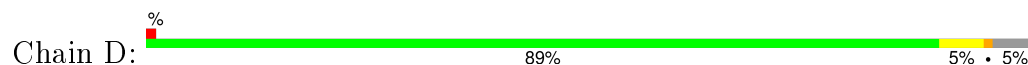
- Molecule 1: protein FucU



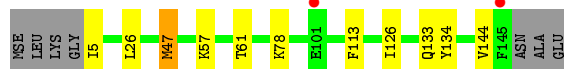
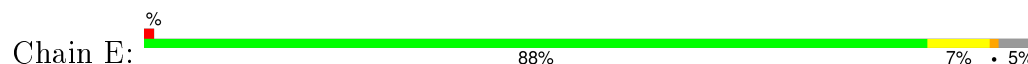
- Molecule 1: protein FucU



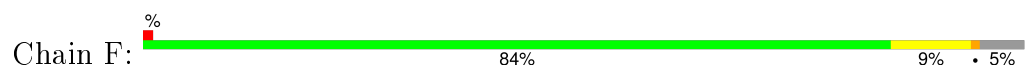
- Molecule 1: protein FucU



- Molecule 1: protein FucU

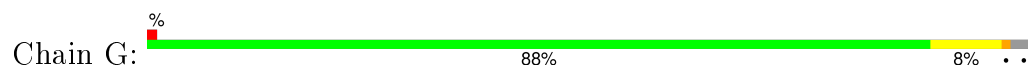


- Molecule 1: protein FucU

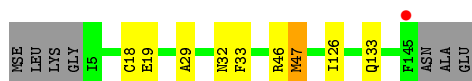
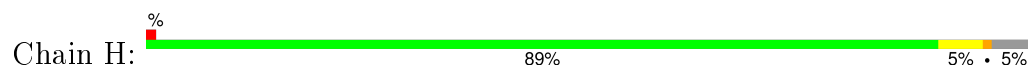




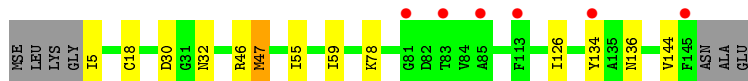
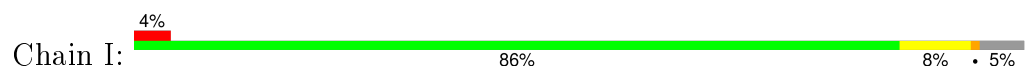
- Molecule 1: protein FucU



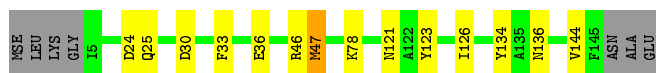
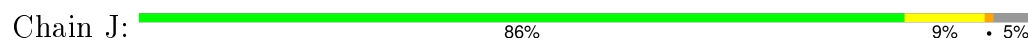
- Molecule 1: protein FucU



- Molecule 1: protein FucU



- Molecule 1: protein FucU



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.05Å 83.86Å 144.81Å 90.00° 95.11° 90.00°	Depositor
Resolution (Å)	40.72 – 1.65 40.72 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.8 (40.72-1.65) 99.8 (40.72-1.65)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.182 , 0.205 0.181 , 0.204	Depositor DCC
R_{free} test set	9474 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	20.2	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 189234 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12068	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: GOL, PGE, NA

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	0.54	0/1078	0.68	1/1462 (0.1%)
1	B	0.56	0/1074	0.67	1/1456 (0.1%)
1	C	0.55	0/1086	0.66	1/1472 (0.1%)
1	D	0.52	0/1076	0.65	1/1454 (0.1%)
1	E	0.57	0/1083	0.63	1/1468 (0.1%)
1	F	0.53	0/1070	0.64	1/1451 (0.1%)
1	G	0.55	0/1093	0.65	1/1483 (0.1%)
1	H	0.57	0/1088	0.63	1/1475 (0.1%)
1	I	0.55	0/1084	0.65	1/1471 (0.1%)
1	J	0.58	0/1086	0.66	1/1471 (0.1%)
All	All	0.55	0/10818	0.65	10/14663 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	47	MSE	CG-SE-CE	-9.48	78.03	98.90
1	D	47	MSE	CG-SE-CE	-9.34	78.35	98.90
1	A	47	MSE	CG-SE-CE	-9.16	78.75	98.90
1	J	47	MSE	CG-SE-CE	-8.97	79.16	98.90
1	G	47	MSE	CG-SE-CE	-8.69	79.79	98.90
1	C	47	MSE	CG-SE-CE	-8.63	79.90	98.90
1	F	47	MSE	CG-SE-CE	-8.26	80.72	98.90
1	I	47	MSE	CG-SE-CE	-8.20	80.86	98.90
1	H	47	MSE	CG-SE-CE	-7.61	82.16	98.90
1	E	47	MSE	CG-SE-CE	-5.51	86.77	98.90

There are no chirality outliers.

There are no planarity outliers.

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	1054	0	1057	12	0
1	B	1053	0	1050	20	0
1	C	1064	0	1059	26	0
1	D	1055	0	1056	7	0
1	E	1061	0	1057	14	0
1	F	1049	0	1046	14	0
1	G	1068	0	1054	13	0
1	H	1063	0	1060	8	0
1	I	1059	0	1048	12	0
1	J	1064	0	1066	18	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
2	G	1	0	0	0	0
2	H	2	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
3	A	10	0	14	0	0
3	B	10	0	14	0	0
3	C	20	0	28	7	0
3	E	20	0	28	2	0
3	F	10	0	14	0	0
3	G	10	0	14	0	0
3	H	10	0	14	0	0
3	J	20	0	28	9	0
4	A	6	0	8	0	0
4	B	6	0	7	5	0
4	E	6	0	7	4	0
4	F	6	0	8	1	0
4	G	6	0	8	0	0
4	H	6	0	8	1	0
5	A	137	0	0	1	0
5	B	139	0	0	3	0
5	C	116	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	121	0	0	2	0
5	E	136	0	0	0	0
5	F	120	0	0	2	0
5	G	156	0	0	1	0
5	H	138	0	0	0	0
5	I	121	0	0	0	0
5	J	133	0	0	0	0
All	All	12068	0	10753	122	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 (122) 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:C:25:GLN:H	3:C:151:PGE:H62	1.05	1.12
1:B:133:GLN:HE21	4:B:152:GOL:C3	1.80	0.95
1:J:123:TYR:HB2	3:J:150:PGE:H3	1.47	0.94
1:H:32:ASN:HD21	1:I:144:VAL:H	1.13	0.91
1:B:32:ASN:HD21	1:E:144:VAL:H	1.18	0.91
1:B:133:GLN:HE21	4:B:152:GOL:H32	1.35	0.89
1:C:144:VAL:H	1:F:32:ASN:HD21	1.15	0.88
1:G:32:ASN:HD21	1:J:144:VAL:H	1.22	0.88
1:F:99:HIS:HE1	5:F:224:HOH:O	1.57	0.87
1:C:25:GLN:HE21	1:C:140:GLN:HE21	1.22	0.85
1:A:144:VAL:H	1:I:32:ASN:HD21	1.25	0.84
1:B:144:VAL:H	1:C:32:ASN:HD21	1.24	0.82
1:C:25:GLN:N	3:C:151:PGE:H62	1.90	0.81
1:D:32:ASN:HD21	1:F:144:VAL:H	1.31	0.77
1:E:133:GLN:HE21	4:E:153:GOL:C3	2.00	0.75
1:C:25:GLN:H	3:C:151:PGE:C6	1.94	0.74
1:A:144:VAL:H	1:I:32:ASN:ND2	1.85	0.74
1:C:8:ILE:HD11	1:C:58:ALA:CB	2.18	0.73
1:A:5:ILE:HD11	1:E:5:ILE:HD11	1.71	0.70
1:C:12:GLU:HG3	5:C:405:HOH:O	1.90	0.70
1:I:30:ASP:H	1:I:136:ASN:ND2	1.90	0.70
1:J:121:ASN:O	3:J:150:PGE:H2	1.92	0.70
1:B:144:VAL:H	1:C:32:ASN:ND2	1.90	0.69
1:F:12:GLU:HG3	5:F:1307:HOH:O	1.92	0.69
1:C:30:ASP:H	1:C:136:ASN:ND2	1.90	0.69
1:A:30:ASP:H	1:A:136:ASN:ND2	1.90	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:144:VAL:H	1:F:32:ASN:ND2	1.90	0.68
1:J:121:ASN:O	3:J:150:PGE:H42	1.93	0.68
1:C:24:ASP:OD1	3:C:151:PGE:H5	1.94	0.67
1:B:32:ASN:ND2	1:E:144:VAL:H	1.90	0.67
1:C:81:GLY:HA3	5:C:1305:HOH:O	1.96	0.65
1:D:32:ASN:ND2	1:F:144:VAL:H	1.95	0.65
1:C:99:HIS:HE1	5:C:162:HOH:O	1.80	0.65
1:G:99:HIS:HE1	5:G:291:HOH:O	1.78	0.64
1:C:25:GLN:HE21	1:C:140:GLN:NE2	1.93	0.64
5:D:645:HOH:O	1:I:18[A]:CYS:SG	2.40	0.64
1:C:47:MSE:HG3	1:C:126:ILE:HG23	1.81	0.63
1:D:30:ASP:H	1:D:136:ASN:ND2	1.95	0.63
1:G:30:ASP:H	1:G:136:ASN:ND2	1.96	0.63
1:B:32:ASN:HD21	1:E:144:VAL:N	1.94	0.62
1:G:32:ASN:ND2	1:J:144:VAL:H	1.97	0.62
1:H:32:ASN:ND2	1:I:144:VAL:H	1.93	0.61
1:J:30:ASP:H	1:J:136:ASN:ND2	1.98	0.61
1:C:8:ILE:HD11	1:C:58:ALA:HB2	1.82	0.61
1:B:5:ILE:HD11	1:G:5:ILE:HD11	1.82	0.60
1:I:47:MSE:HG3	1:I:126:ILE:HG23	1.84	0.60
1:B:12:GLU:HG3	5:B:620:HOH:O	2.00	0.60
1:A:57:LYS:O	1:A:61:THR:HG23	2.02	0.59
1:B:133:GLN:NE2	4:B:152:GOL:H32	2.13	0.58
1:A:46:ARG:C	1:A:47:MSE:HG2	2.24	0.58
1:C:144:VAL:N	1:F:32:ASN:HD21	1.95	0.58
1:G:37:SER:HB2	3:J:150:PGE:H52	1.86	0.57
1:B:47:MSE:HG3	1:B:126:ILE:HG23	1.86	0.57
1:E:133:GLN:HE21	4:E:153:GOL:H32	1.69	0.57
1:J:47:MSE:HG3	1:J:126:ILE:HG23	1.86	0.56
1:F:22:HIS:HE1	1:F:66:ASP:OD1	1.86	0.56
1:B:22:HIS:HE1	1:B:66:ASP:OD1	1.88	0.56
1:C:121:ASN:O	3:C:151:PGE:H42	2.05	0.56
1:I:46:ARG:C	1:I:47:MSE:HG2	2.25	0.56
1:H:133:GLN:HE21	4:H:152:GOL:C3	2.18	0.56
1:B:46:ARG:C	1:B:47:MSE:HG2	2.27	0.55
1:A:22:HIS:HE1	1:A:66:ASP:OD1	1.90	0.55
1:B:57:LYS:O	1:B:61:THR:HG23	2.08	0.54
1:D:46:ARG:C	1:D:47:MSE:HG2	2.26	0.54
1:A:30:ASP:H	1:A:136:ASN:HD22	1.52	0.54
1:C:24:ASP:HA	3:C:151:PGE:H5	1.90	0.53
1:C:46:ARG:C	1:C:47:MSE:HG2	2.29	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:25:GLN:HB2	3:J:150:PGE:H22	1.91	0.52
1:J:46:ARG:C	1:J:47:MSE:HG2	2.28	0.52
1:C:123:TYR:N	3:C:151:PGE:H6	2.24	0.52
1:E:26:LEU:C	1:E:26:LEU:HD12	2.30	0.52
1:A:47:MSE:HG3	1:A:126:ILE:HG23	1.93	0.51
1:F:5:ILE:HG13	1:F:5:ILE:O	2.10	0.51
1:E:57:LYS:O	1:E:61:THR:HG23	2.10	0.51
1:D:57:LYS:O	1:D:61:THR:HG23	2.12	0.50
1:H:47:MSE:HG3	1:H:126:ILE:HG23	1.93	0.49
1:F:133:GLN:HE21	4:F:152:GOL:H31	1.76	0.49
1:J:123:TYR:CB	3:J:150:PGE:H3	2.31	0.49
1:G:32:ASN:HD21	1:J:144:VAL:N	2.02	0.49
1:G:47:MSE:HG3	1:G:126:ILE:HG23	1.94	0.49
1:J:24:ASP:OD1	3:J:150:PGE:H32	2.12	0.49
1:F:11:PRO:HG3	1:H:18[B]:CYS:SG	2.53	0.49
1:A:54:GLU:HG3	5:A:156:HOH:O	2.12	0.49
1:C:15:LYS:O	1:C:19:GLU:HG3	2.13	0.48
1:G:46:ARG:C	1:G:47:MSE:HG2	2.34	0.48
1:J:30:ASP:H	1:J:136:ASN:HD22	1.60	0.48
1:B:133:GLN:HE21	4:B:152:GOL:H31	1.74	0.48
1:B:79:VAL:CG2	1:B:82:ASP:HB2	2.44	0.47
1:G:30:ASP:H	1:G:136:ASN:HD22	1.61	0.47
1:D:47:MSE:HG3	1:D:126:ILE:HG23	1.96	0.47
1:E:133:GLN:NE2	4:E:153:GOL:C3	2.75	0.47
1:I:55:ILE:O	1:I:59[A]:ILE:HG12	2.15	0.47
1:E:78:LYS:HA	1:E:134:TYR:CE1	2.50	0.47
1:B:133:GLN:NE2	4:B:152:GOL:C3	2.64	0.45
1:F:47:MSE:HG3	1:F:126:ILE:HG23	1.97	0.45
1:C:78:LYS:HA	1:C:134:TYR:CE1	2.52	0.45
1:G:78:LYS:HA	1:G:134:TYR:CE1	2.52	0.45
1:H:46:ARG:C	1:H:47:MSE:HG2	2.37	0.45
1:J:25:GLN:H	3:J:150:PGE:H22	1.82	0.44
1:B:144:VAL:HG22	1:C:32:ASN:ND2	2.32	0.44
1:D:57:LYS:HD2	5:D:959:HOH:O	2.17	0.44
1:E:47:MSE:HG3	1:E:126:ILE:HG23	1.99	0.44
1:G:26:LEU:C	1:G:26:LEU:HD12	2.37	0.44
1:E:113:PHE:CE2	3:E:151:PGE:H52	2.53	0.44
1:E:133:GLN:NE2	4:E:153:GOL:H32	2.32	0.43
1:F:46:ARG:C	1:F:47:MSE:HG2	2.39	0.43
1:A:26:LEU:C	1:A:26:LEU:HD12	2.39	0.43
1:B:54:GLU:OE1	5:B:1303:HOH:O	2.21	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59[A]:ILE:HD11	1:A:126:ILE:CD1	2.49	0.43
1:J:25:GLN:H	3:J:150:PGE:C2	2.32	0.42
5:B:1087:HOH:O	1:G:18[B]:CYS:SG	2.60	0.42
1:C:79:VAL:HG23	1:C:82:ASP:HB2	2.01	0.42
1:F:78:LYS:HA	1:F:134:TYR:CE1	2.53	0.42
1:E:113:PHE:HE2	3:E:151:PGE:H52	1.84	0.42
1:B:77:GLU:HG3	1:B:109:SER:HB2	2.01	0.42
1:J:78:LYS:HA	1:J:134:TYR:CE1	2.55	0.42
1:I:30:ASP:H	1:I:136:ASN:HD22	1.67	0.42
1:H:29:ALA:HB1	1:H:33:PHE:HB3	2.02	0.41
1:J:33:PHE:HB2	1:J:136:ASN:HD21	1.85	0.41
1:I:78:LYS:HA	1:I:134:TYR:CE1	2.54	0.41
1:I:47:MSE:HG3	1:I:126:ILE:CG2	2.51	0.40
1:H:19:GLU:HB2	1:J:36:GLU:HG3	2.04	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	140/148 (95%)	140 (100%)	0	0	100	100
1	B	139/148 (94%)	138 (99%)	1 (1%)	0	100	100
1	C	140/148 (95%)	137 (98%)	3 (2%)	0	100	100
1	D	137/148 (93%)	136 (99%)	1 (1%)	0	100	100
1	E	140/148 (95%)	139 (99%)	1 (1%)	0	100	100
1	F	139/148 (94%)	138 (99%)	1 (1%)	0	100	100
1	G	143/148 (97%)	142 (99%)	1 (1%)	0	100	100
1	H	141/148 (95%)	140 (99%)	1 (1%)	0	100	100
1	I	141/148 (95%)	141 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	J	140/148 (95%)	140 (100%)	0	0	100	100
All	All	1400/1480 (95%)	1391 (99%)	9 (1%)	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	110/113 (97%)	110 (100%)	0	100	100
1	B	110/113 (97%)	110 (100%)	0	100	100
1	C	111/113 (98%)	109 (98%)	2 (2%)	66	43
1	D	109/113 (96%)	109 (100%)	0	100	100
1	E	110/113 (97%)	110 (100%)	0	100	100
1	F	109/113 (96%)	107 (98%)	2 (2%)	66	43
1	G	110/113 (97%)	110 (100%)	0	100	100
1	H	111/113 (98%)	111 (100%)	0	100	100
1	I	109/113 (96%)	108 (99%)	1 (1%)	84	72
1	J	111/113 (98%)	111 (100%)	0	100	100
All	All	1100/1130 (97%)	1095 (100%)	5 (0%)	92	86

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	8	ILE
1	C	101	GLU
1	F	5	ILE
1	F	14	LEU
1	I	5	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	HIS
1	A	136	ASN
1	B	22	HIS
1	B	32	ASN
1	B	133	GLN
1	C	32	ASN
1	C	99	HIS
1	C	136	ASN
1	C	140	GLN
1	D	32	ASN
1	D	99	HIS
1	D	118	GLN
1	D	136	ASN
1	E	133	GLN
1	F	22	HIS
1	F	32	ASN
1	F	99	HIS
1	F	133	GLN
1	G	32	ASN
1	G	99	HIS
1	G	136	ASN
1	H	32	ASN
1	H	118	GLN
1	I	32	ASN
1	I	118	GLN
1	I	136	ASN
1	J	136	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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

Of 32 ligands modelled in this entry, 15 are monoatomic - leaving 17 for Mogul analysis.

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	PGE	A	151	-	9,9,9	0.50	0	8,8,8	0.16	0
4	GOL	A	152	2	5,5,5	0.38	0	5,5,5	0.73	0
3	PGE	B	151	-	9,9,9	0.44	0	8,8,8	0.41	0
4	GOL	B	152	2	5,5,5	0.33	0	5,5,5	0.62	0
3	PGE	C	150	-	9,9,9	0.47	0	8,8,8	0.31	0
3	PGE	C	151	-	9,9,9	0.40	0	8,8,8	0.49	0
3	PGE	E	151	-	9,9,9	0.48	0	8,8,8	0.22	0
3	PGE	E	152	-	9,9,9	0.46	0	8,8,8	0.30	0
4	GOL	E	153	2	5,5,5	0.39	0	5,5,5	0.86	0
3	PGE	F	151	-	9,9,9	0.47	0	8,8,8	0.29	0
4	GOL	F	152	2	5,5,5	0.42	0	5,5,5	0.51	0
3	PGE	G	150	-	9,9,9	0.43	0	8,8,8	0.36	0
4	GOL	G	151	-	5,5,5	0.33	0	5,5,5	0.43	0
3	PGE	H	151	-	9,9,9	0.55	0	8,8,8	0.34	0
4	GOL	H	152	2	5,5,5	0.52	0	5,5,5	1.32	1 (20%)
3	PGE	J	150	-	9,9,9	0.62	0	8,8,8	0.89	0
3	PGE	J	151	-	9,9,9	0.43	0	8,8,8	0.40	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	PGE	A	151	-	-	0/7/7/7	0/0/0/0
4	GOL	A	152	2	-	0/4/4/4	0/0/0/0
3	PGE	B	151	-	-	0/7/7/7	0/0/0/0
4	GOL	B	152	2	-	0/4/4/4	0/0/0/0
3	PGE	C	150	-	-	0/7/7/7	0/0/0/0
3	PGE	C	151	-	-	0/7/7/7	0/0/0/0
3	PGE	E	151	-	-	0/7/7/7	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PGE	E	152	-	-	0/7/7/7	0/0/0/0
4	GOL	E	153	2	-	0/4/4/4	0/0/0/0
3	PGE	F	151	-	-	0/7/7/7	0/0/0/0
4	GOL	F	152	2	-	0/4/4/4	0/0/0/0
3	PGE	G	150	-	-	0/7/7/7	0/0/0/0
4	GOL	G	151	-	-	0/4/4/4	0/0/0/0
3	PGE	H	151	-	-	0/7/7/7	0/0/0/0
4	GOL	H	152	2	-	0/4/4/4	0/0/0/0
3	PGE	J	150	-	-	0/7/7/7	0/0/0/0
3	PGE	J	151	-	-	0/7/7/7	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	152	GOL	O3-C3-C2	-2.11	99.97	110.18

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 29 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	152	GOL	5	0
3	C	151	PGE	7	0
3	E	151	PGE	2	0
4	E	153	GOL	4	0
4	F	152	GOL	1	0
4	H	152	GOL	1	0
3	J	150	PGE	9	0

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	137/148 (92%)	-0.24	1 (0%) 89 90	13, 18, 29, 32	0
1	B	137/148 (92%)	-0.33	0 100 100	13, 18, 30, 36	0
1	C	138/148 (93%)	-0.15	6 (4%) 39 38	14, 21, 35, 40	0
1	D	137/148 (92%)	-0.15	2 (1%) 76 79	12, 19, 34, 43	0
1	E	138/148 (93%)	-0.11	2 (1%) 78 80	13, 19, 33, 42	0
1	F	137/148 (92%)	-0.22	1 (0%) 89 90	14, 20, 32, 39	0
1	G	140/148 (94%)	-0.40	1 (0%) 89 90	13, 18, 29, 43	0
1	H	138/148 (93%)	-0.22	1 (0%) 89 90	13, 18, 29, 38	0
1	I	138/148 (93%)	-0.13	6 (4%) 39 38	14, 21, 35, 38	0
1	J	138/148 (93%)	-0.22	0 100 100	13, 19, 29, 34	0
All	All	1378/1480 (93%)	-0.22	20 (1%) 76 79	12, 19, 32, 43	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	145	PHE	5.6
1	C	81	GLY	4.7
1	I	134	TYR	4.3
1	C	83	THR	3.7
1	H	145	PHE	3.6
1	I	85	ALA	3.5
1	I	83	THR	3.5
1	C	85	ALA	3.2
1	D	80	PRO	3.2
1	I	81	GLY	3.0
1	C	145	PHE	3.0
1	E	101	GLU	2.9
1	C	79	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	134	TYR	2.8
1	D	81	GLY	2.6
1	A	85	ALA	2.5
1	I	113	PHE	2.3
1	G	4	GLY	2.1
1	F	105	ASP	2.1
1	I	145	PHE	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 carbohydrates in this entry.

6.4 Ligands [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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	PGE	C	151	10/10	0.78	0.38	15.11	38,41,47,48	0
4	GOL	H	152	6/6	0.79	0.25	13.53	31,39,39,40	0
3	PGE	J	150	10/10	0.79	0.29	11.22	33,35,38,40	0
2	NA	B	149	1/1	0.98	0.15	10.84	22,22,22,22	0
3	PGE	E	152	10/10	0.84	0.16	7.83	44,46,48,49	0
4	GOL	E	153	6/6	0.75	0.20	7.06	34,35,37,39	0
4	GOL	A	152	6/6	0.86	0.16	6.32	27,33,35,39	0
4	GOL	F	152	6/6	0.79	0.19	4.37	29,36,39,42	0
4	GOL	B	152	6/6	0.87	0.18	4.13	34,39,40,41	0
3	PGE	H	151	10/10	0.78	0.12	3.96	43,47,49,49	0
3	PGE	C	150	10/10	0.83	0.11	3.58	40,41,42,43	0
2	NA	A	149	1/1	0.97	0.13	3.27	23,23,23,23	0
3	PGE	G	150	10/10	0.89	0.10	2.65	43,44,44,45	0
3	PGE	E	151	10/10	0.87	0.11	2.55	44,45,45,45	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	PGE	B	151	10/10	0.85	0.12	1.95	47,47,49,49	0
3	PGE	A	151	10/10	0.84	0.13	1.45	48,49,50,51	0
2	NA	F	149	1/1	0.97	0.10	1.29	27,27,27,27	0
2	NA	F	150	1/1	0.96	0.08	0.68	28,28,28,28	0
3	PGE	J	151	10/10	0.92	0.08	0.54	48,50,52,53	0
4	GOL	G	151	6/6	0.95	0.08	0.45	20,25,25,27	0
2	NA	E	149	1/1	0.97	0.11	0.17	24,24,24,24	0
2	NA	G	149	1/1	0.97	0.07	0.01	22,22,22,22	0
2	NA	D	149	1/1	0.97	0.07	-0.16	20,20,20,20	0
3	PGE	F	151	10/10	0.89	0.10	-0.27	51,51,53,54	0
2	NA	H	149	1/1	0.98	0.08	-0.50	23,23,23,23	0
2	NA	C	149	1/1	0.99	0.05	-0.76	24,24,24,24	0
2	NA	I	149	1/1	0.99	0.05	-0.98	23,23,23,23	0
2	NA	J	149	1/1	0.99	0.06	-1.16	22,22,22,22	0
2	NA	H	150	1/1	0.94	0.06	-1.62	25,25,25,25	0
2	NA	A	150	1/1	0.99	0.05	-4.58	27,27,27,27	0
2	NA	E	150	1/1	0.97	0.09	-	27,27,27,27	0
2	NA	B	150	1/1	0.99	0.08	-	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.