



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:42 PM GMT

PDB ID : 4A7X
Title : Crystal structure of uridylylate kinase from Helicobacter pylori
Authors : Chu, C.H.; Liu, M.H.; Chen, P.C.; Sun, Y.J.
Deposited on : 2011-11-15
Resolution : 2.49 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

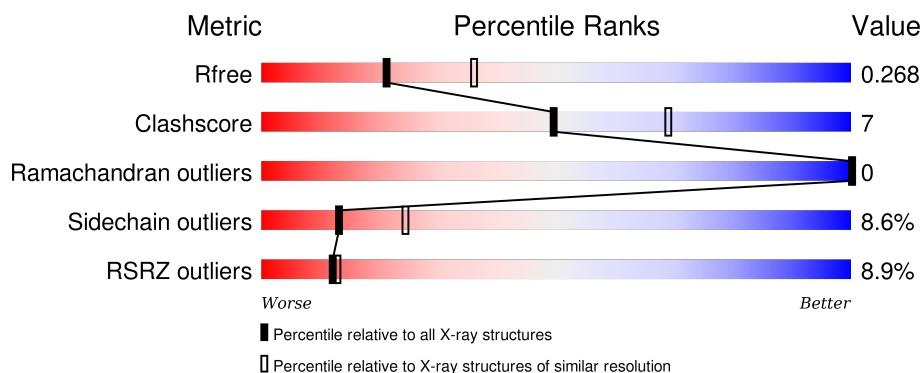
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	240	<div> <div>15%</div> <div>83%</div> <div>13%</div> <div>• •</div> </div>
1	B	240	<div> <div>5%</div> <div>77%</div> <div>16%</div> <div>• 5%</div> </div>
1	C	240	<div> <div>7%</div> <div>78%</div> <div>18%</div> <div>• •</div> </div>
1	D	240	<div> <div>5%</div> <div>77%</div> <div>15%</div> <div>• 5%</div> </div>
1	E	240	<div> <div>8%</div> <div>76%</div> <div>14%</div> <div>• 5%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	240	<div><div></div><div>13%</div><div>81%</div><div>13%</div><div></div><div></div></div>

2 Entry composition [i](#)

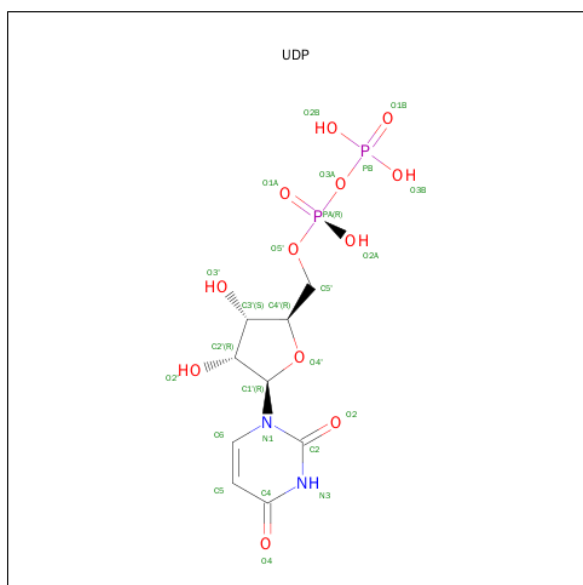
There are 3 unique types of molecules in this entry. The entry contains 10957 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 URIDYLATE KINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	234	Total	C	N	O	S	0	0	0
			1763	1115	305	335	8			
1	B	228	Total	C	N	O	S	0	0	0
			1727	1092	298	329	8			
1	C	234	Total	C	N	O	S	0	0	0
			1783	1129	308	338	8			
1	D	228	Total	C	N	O	S	0	0	0
			1727	1092	298	329	8			
1	E	227	Total	C	N	O	S	0	0	0
			1726	1092	298	328	8			
1	F	230	Total	C	N	O	S	0	0	0
			1738	1100	301	329	8			

- Molecule 2 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: $C_9H_{14}N_2O_{12}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
2	C	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
2	D	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
2	E	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

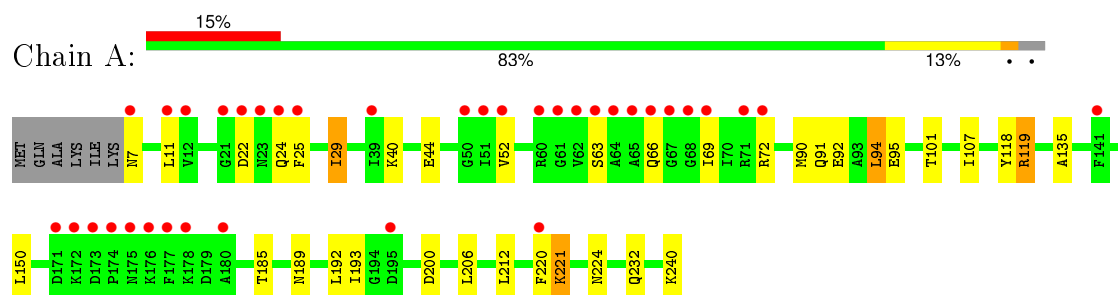
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	59	Total	O	0	0
			59	59		
3	B	74	Total	O	0	0
			74	74		
3	C	89	Total	O	0	0
			89	89		
3	D	76	Total	O	0	0
			76	76		
3	E	34	Total	O	0	0
			34	34		
3	F	61	Total	O	0	0
			61	61		

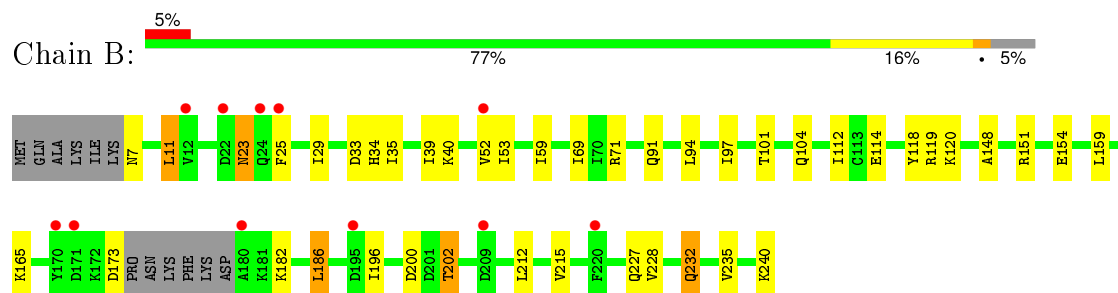
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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.

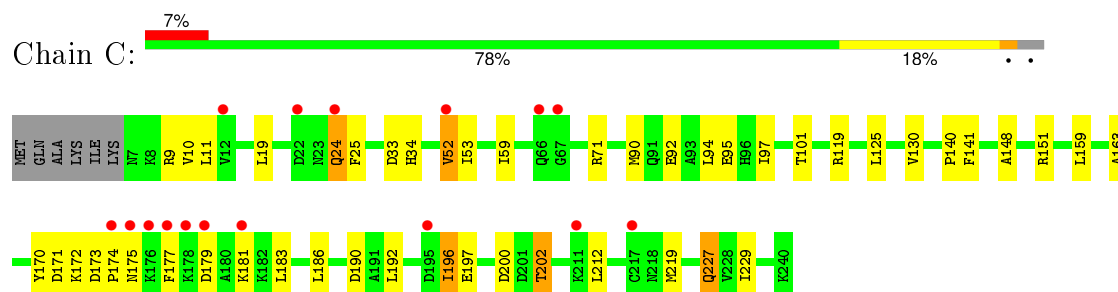
• Molecule 1: URIDYLATE KINASE



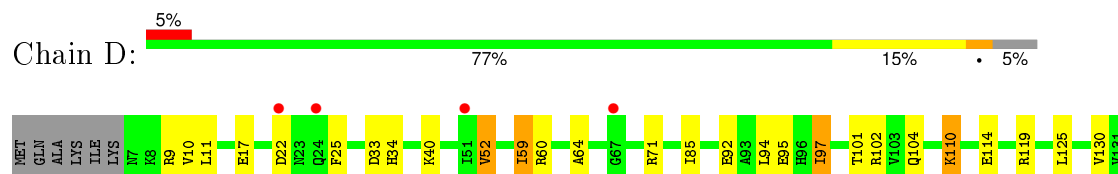
• Molecule 1: URIDYLATE KINASE

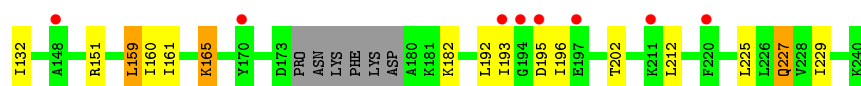


• Molecule 1: URIDYLATE KINASE

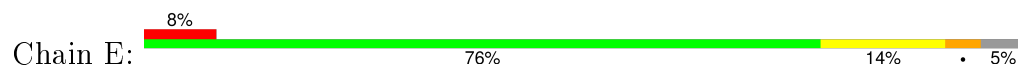


• Molecule 1: URIDYLATE KINASE

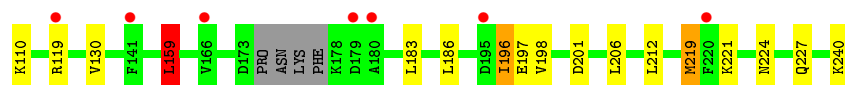
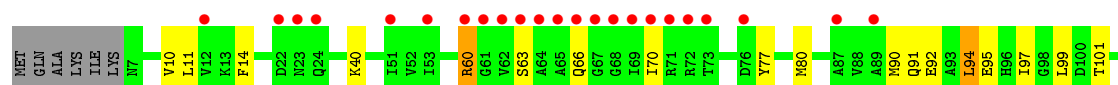
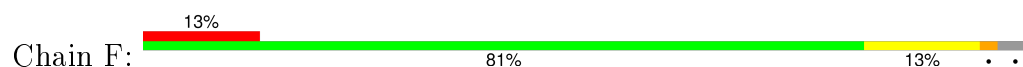




• Molecule 1: URIDYLATE KINASE



• Molecule 1: URIDYLATE KINASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	147.14Å 127.63Å 93.18Å 90.00° 91.46° 90.00°	Depositor
Resolution (Å)	96.40 – 2.49 26.66 – 2.49	Depositor EDS
% Data completeness (in resolution range)	98.9 (96.40-2.49) 99.0 (26.66-2.49)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.93 (at 2.50Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.213 , 0.276 0.209 , 0.268	Depositor DCC
R_{free} test set	2994 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	44.7	Xtriage
Anisotropy	0.211	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 45.7	EDS
Estimated twinning fraction	0.000 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Outliers	1 of 59430 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10957	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.15 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.3776e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.41	0/1782	0.58	0/2402
1	B	0.42	0/1744	0.55	0/2348
1	C	0.55	0/1803	0.62	0/2427
1	D	0.50	0/1744	0.61	0/2348
1	E	0.41	0/1743	0.57	1/2345 (0.0%)
1	F	0.50	0/1755	0.62	1/2362 (0.0%)
All	All	0.47	0/10571	0.59	2/14232 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	159	LEU	CA-CB-CG	5.54	128.05	115.30
1	E	159	LEU	CA-CB-CG	5.32	127.53	115.30

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	1763	0	1803	17	0
1	B	1727	0	1774	23	0
1	C	1783	0	1838	38	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1727	0	1774	34	0
1	E	1726	0	1780	29	0
1	F	1738	0	1787	25	0
2	B	25	0	11	0	0
2	C	25	0	11	1	0
2	D	25	0	11	1	0
2	E	25	0	11	4	0
3	A	59	0	0	3	0
3	B	74	0	0	6	0
3	C	89	0	0	6	0
3	D	76	0	0	4	0
3	E	34	0	0	1	0
3	F	61	0	0	2	0
All	All	10957	0	10800	152	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 (152) 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:E:57:ASN:HD21	1:E:86:ASN:HD21	1.07	0.93
1:D:227:GLN:HA	1:D:227:GLN:HE21	1.37	0.89
1:A:101:THR:HB	3:A:2029:HOH:O	1.71	0.88
1:E:57:ASN:HD22	1:E:57:ASN:H	1.25	0.84
1:C:11:LEU:HD21	1:C:52:VAL:HG13	1.61	0.82
1:C:172:LYS:O	1:C:174:PRO:HD3	1.79	0.82
1:E:11:LEU:HD21	1:E:52:VAL:HG13	1.61	0.81
1:D:11:LEU:HD21	1:D:52:VAL:HG13	1.63	0.80
1:C:171:ASP:HA	1:C:196:ILE:HG21	1.64	0.80
1:D:10:VAL:HG11	1:D:229:ILE:HG12	1.64	0.79
1:D:59:ILE:HD11	1:D:64:ALA:HB2	1.68	0.74
1:E:40:LYS:HD2	1:E:97:ILE:HG23	1.73	0.70
1:C:59:ILE:CD1	1:D:85:ILE:HG21	2.24	0.68
1:C:11:LEU:HD21	1:C:52:VAL:CG1	2.24	0.68
1:E:118:TYR:HE1	1:E:154:GLU:O	1.78	0.67
1:D:165:LYS:HD3	1:D:165:LYS:H	1.59	0.67
1:F:10:VAL:HG12	1:F:159:LEU:CD1	2.25	0.67
1:F:63:SER:HA	1:F:66:GLN:HE21	1.60	0.67
1:A:11:LEU:HD21	1:A:52:VAL:HG13	1.77	0.67
2:E:1241:UDP:O2A	2:E:1241:UDP:H3'	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:221:LYS:HB2	1:F:224:ASN:HB2	1.76	0.65
2:D:1241:UDP:H3'	2:D:1241:UDP:PA	2.36	0.64
1:C:95:GLU:OE1	1:E:119:ARG:NH2	2.30	0.64
1:C:186:LEU:HD22	1:C:190:ASP:HB3	1.81	0.63
1:B:151:ARG:NH2	3:B:2052:HOH:O	2.30	0.63
1:E:101:THR:HG22	1:E:130:VAL:HB	1.81	0.63
1:D:40:LYS:HD2	1:D:97:ILE:HG23	1.82	0.62
1:D:119:ARG:HG3	3:D:2047:HOH:O	1.99	0.62
1:A:7:ASN:N	3:A:2001:HOH:O	2.34	0.61
1:C:59:ILE:HD11	1:D:85:ILE:HG21	1.82	0.61
1:B:196:ILE:O	1:B:196:ILE:HG13	2.01	0.60
2:C:1241:UDP:O3B	2:C:1241:UDP:O1A	2.18	0.60
1:F:40:LYS:HG3	1:F:99:LEU:HD21	1.84	0.59
1:F:40:LYS:HD2	1:F:97:ILE:HG23	1.84	0.59
1:C:95:GLU:CD	1:E:119:ARG:HH22	2.07	0.59
1:C:170:TYR:CG	1:C:174:PRO:HG3	2.38	0.58
1:F:10:VAL:HG12	1:F:159:LEU:HD13	1.85	0.58
1:D:104:GLN:HG2	1:D:114:GLU:HG3	1.85	0.57
1:C:19:LEU:HD11	1:C:53:ILE:HD11	1.86	0.57
1:C:227:GLN:HA	1:C:227:GLN:HE21	1.68	0.57
1:E:57:ASN:HD22	1:E:57:ASN:N	1.99	0.57
1:B:186:LEU:HD21	1:B:196:ILE:HD11	1.87	0.57
1:C:140:PRO:O	1:C:141:PHE:HB2	2.04	0.57
1:C:183:LEU:HD13	1:C:186:LEU:HD21	1.85	0.57
1:E:58:ILE:HG22	1:E:59:ILE:HG23	1.87	0.57
1:B:104:GLN:HG2	1:B:114:GLU:HG3	1.86	0.57
1:F:14:PHE:HE1	1:F:219:MET:HE1	1.70	0.56
1:B:23:ASN:HB2	3:B:2009:HOH:O	2.05	0.56
1:F:101:THR:HG22	1:F:130:VAL:HB	1.87	0.56
1:C:151:ARG:NH1	3:C:2067:HOH:O	2.39	0.56
1:B:7:ASN:N	3:B:2001:HOH:O	2.39	0.56
1:C:163:ALA:HB1	1:C:219:MET:HB3	1.88	0.55
1:D:227:GLN:HA	1:D:227:GLN:NE2	2.16	0.55
1:F:196:ILE:HG23	1:F:198:VAL:HG23	1.87	0.55
1:D:17:GLU:OE2	1:D:60:ARG:NH1	2.40	0.54
1:F:91:GLN:O	1:F:95:GLU:HG3	2.07	0.54
1:A:118:TYR:HB3	1:A:119:ARG:NH2	2.22	0.54
1:E:55:GLY:H	2:E:1241:UDP:PA	2.30	0.54
1:B:215:VAL:HG11	1:B:228:VAL:HG11	1.90	0.53
1:C:177:PHE:C	1:C:179:ASP:H	2.12	0.53
1:C:10:VAL:HG11	1:C:229:ILE:HG12	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:2056:HOH:O	1:E:151:ARG:NH2	2.41	0.53
1:C:52:VAL:HG11	1:C:148:ALA:HA	1.90	0.52
1:A:63:SER:O	1:A:66:GLN:HG2	2.10	0.52
1:B:40:LYS:HD2	1:B:97:ILE:HD12	1.90	0.52
1:D:101:THR:HG22	1:D:130:VAL:HB	1.92	0.52
1:A:94:LEU:HB3	1:A:101:THR:HG21	1.91	0.52
1:E:240:LYS:HE2	3:E:2034:HOH:O	2.10	0.51
1:D:193:ILE:O	1:D:193:ILE:HG22	2.10	0.51
1:C:173:ASP:OD1	1:C:175:ASN:N	2.29	0.50
1:E:57:ASN:ND2	1:E:86:ASN:HD21	1.91	0.50
1:E:71:ARG:HD3	1:F:92:GLU:OE1	2.11	0.50
1:C:171:ASP:HA	1:C:196:ILE:CG2	2.39	0.50
1:E:57:ASN:ND2	1:E:57:ASN:H	2.02	0.49
1:A:95:GLU:OE1	1:C:119:ARG:NH2	2.46	0.49
1:E:187:SER:HA	1:E:240:LYS:HD2	1.94	0.49
1:F:14:PHE:HE1	1:F:219:MET:CE	2.24	0.49
1:C:170:TYR:CD2	1:C:174:PRO:HG3	2.47	0.49
1:C:71:ARG:HD3	1:D:92:GLU:OE1	2.12	0.49
1:D:193:ILE:O	1:D:193:ILE:CG2	2.58	0.48
1:B:119:ARG:NH2	1:D:95:GLU:OE1	2.46	0.48
1:A:189:ASN:O	1:A:193:ILE:HG12	2.13	0.48
1:C:179:ASP:HB2	3:C:2075:HOH:O	2.13	0.48
1:A:92:GLU:OE1	1:B:71:ARG:HD3	2.13	0.48
1:F:60:ARG:HB2	1:F:63:SER:HB2	1.96	0.48
1:A:40:LYS:O	1:A:44:GLU:HG3	2.14	0.48
1:A:221:LYS:HB3	1:A:224:ASN:HB2	1.96	0.48
1:E:54:GLY:HA2	2:E:1241:UDP:O1A	2.14	0.48
1:B:91:GLN:HG3	1:B:101:THR:OG1	2.14	0.47
1:F:183:LEU:HD13	1:F:186:LEU:HD21	1.95	0.47
1:B:119:ARG:HH22	1:D:95:GLU:CD	2.17	0.47
1:B:227:GLN:NE2	3:B:2070:HOH:O	2.39	0.47
1:A:29:ILE:HG23	1:B:69:ILE:HG13	1.96	0.47
1:F:227:GLN:NE2	3:F:2060:HOH:O	2.42	0.47
1:D:165:LYS:CD	1:D:165:LYS:H	2.25	0.47
1:E:92:GLU:HG3	1:F:70:ILE:HG23	1.96	0.47
1:C:25:PHE:HB3	1:D:25:PHE:CE2	2.50	0.47
1:A:91:GLN:O	1:A:95:GLU:HG3	2.15	0.47
1:C:34:HIS:HB3	3:C:2008:HOH:O	2.13	0.47
1:E:17:GLU:CD	1:E:17:GLU:H	2.18	0.47
1:D:159:LEU:CD2	1:D:161:ILE:HG13	2.44	0.47
1:B:34:HIS:HB2	3:B:2011:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:2038:HOH:O	1:C:119:ARG:HG3	2.15	0.46
1:A:69:ILE:HD13	1:B:33:ASP:OD1	2.15	0.46
1:A:25:PHE:CE1	1:B:25:PHE:HB2	2.51	0.46
1:B:118:TYR:HE1	1:B:154:GLU:O	1.97	0.46
1:B:11:LEU:HD21	1:B:52:VAL:HG23	1.97	0.46
1:F:119:ARG:HG2	3:F:2035:HOH:O	2.14	0.46
1:D:34:HIS:HB3	3:D:2021:HOH:O	2.15	0.46
3:C:2045:HOH:O	1:D:110:LYS:HG3	2.15	0.46
1:F:60:ARG:CB	1:F:63:SER:HB2	2.46	0.45
1:F:94:LEU:HB3	1:F:101:THR:HG21	1.99	0.45
1:C:92:GLU:OE1	1:D:71:ARG:HD3	2.16	0.45
1:C:177:PHE:HB3	3:C:2075:HOH:O	2.16	0.45
1:F:14:PHE:CE1	1:F:219:MET:HE1	2.52	0.45
1:F:91:GLN:HG3	1:F:101:THR:OG1	2.16	0.45
1:D:227:GLN:CA	1:D:227:GLN:HE21	2.13	0.45
1:C:101:THR:HG22	1:C:130:VAL:HB	1.99	0.45
1:B:200:ASP:OD2	1:B:202:THR:HG23	2.17	0.45
1:E:172:LYS:O	1:E:173:ASP:C	2.55	0.44
1:C:59:ILE:CD1	1:D:85:ILE:CG2	2.94	0.44
1:E:10:VAL:HG11	1:E:229:ILE:HG12	1.98	0.44
1:B:52:VAL:HG21	1:B:148:ALA:HA	1.99	0.44
1:D:9:ARG:HD3	1:D:125:LEU:HD13	1.99	0.44
1:F:196:ILE:O	1:F:196:ILE:CG2	2.65	0.43
1:C:200:ASP:OD2	1:C:202:THR:HG23	2.18	0.43
1:C:24:GLN:N	3:C:2010:HOH:O	2.46	0.43
1:B:232:GLN:HB2	1:B:232:GLN:HE21	1.64	0.43
1:E:53:ILE:HD13	1:E:54:GLY:N	2.33	0.43
1:D:94:LEU:HD22	1:D:132:ILE:HD11	2.00	0.43
1:E:219:MET:HE2	1:E:220:PHE:CZ	2.54	0.43
1:B:35:ILE:O	1:B:39:ILE:HG12	2.19	0.43
1:F:90:MET:HG3	1:F:94:LEU:HD22	2.01	0.43
1:A:107:ILE:HD12	1:A:135:ALA:HB1	2.00	0.43
1:E:57:ASN:N	1:E:57:ASN:ND2	2.66	0.42
1:C:196:ILE:O	1:C:196:ILE:HG12	2.19	0.42
1:F:219:MET:HB3	1:F:219:MET:HE3	1.89	0.42
1:D:165:LYS:N	1:D:165:LYS:HD3	2.30	0.42
1:E:144:THR:HG1	2:E:1241:UDP:PA	2.42	0.42
1:D:151:ARG:NH2	3:D:2058:HOH:O	2.51	0.42
1:D:97:ILE:O	1:D:97:ILE:HG23	2.20	0.42
1:C:140:PRO:O	1:C:141:PHE:CB	2.67	0.42
1:E:164:THR:HB	1:E:165:LYS:H	1.74	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:MET:HG3	1:A:94:LEU:HD22	2.02	0.42
1:D:160:ILE:HG22	1:D:212:LEU:CD1	2.50	0.42
1:F:77:TYR:HA	1:F:80:MET:HE3	2.01	0.41
1:D:10:VAL:HG12	1:D:159:LEU:HD13	2.03	0.41
1:C:9:ARG:HD3	1:C:125:LEU:HD13	2.02	0.41
1:E:183:LEU:HD23	1:E:186:LEU:HD11	2.03	0.41
3:B:2046:HOH:O	1:D:102:ARG:NH2	2.54	0.41
1:C:227:GLN:HA	1:C:227:GLN:NE2	2.35	0.41
1:E:10:VAL:HG12	1:E:159:LEU:HD22	2.03	0.41

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	232/240 (97%)	227 (98%)	5 (2%)	0	100	100
1	B	224/240 (93%)	216 (96%)	8 (4%)	0	100	100
1	C	232/240 (97%)	225 (97%)	7 (3%)	0	100	100
1	D	224/240 (93%)	219 (98%)	5 (2%)	0	100	100
1	E	223/240 (93%)	218 (98%)	5 (2%)	0	100	100
1	F	226/240 (94%)	215 (95%)	11 (5%)	0	100	100
All	All	1361/1440 (94%)	1320 (97%)	41 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	186/197 (94%)	170 (91%)	16 (9%)	13	24
1	B	184/197 (93%)	166 (90%)	18 (10%)	10	19
1	C	191/197 (97%)	177 (93%)	14 (7%)	17	32
1	D	184/197 (93%)	169 (92%)	15 (8%)	14	27
1	E	185/197 (94%)	164 (89%)	21 (11%)	7	13
1	F	184/197 (93%)	172 (94%)	12 (6%)	21	39
All	All	1114/1182 (94%)	1018 (91%)	96 (9%)	13	24

All (96) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ASP
1	A	24	GLN
1	A	29	ILE
1	A	72	ARG
1	A	94	LEU
1	A	119	ARG
1	A	150	LEU
1	A	185	THR
1	A	192	LEU
1	A	200	ASP
1	A	206	LEU
1	A	212	LEU
1	A	220	PHE
1	A	221	LYS
1	A	232	GLN
1	A	240	LYS
1	B	11	LEU
1	B	23	ASN
1	B	29	ILE
1	B	53	ILE
1	B	59	ILE
1	B	94	LEU
1	B	112	ILE
1	B	120	LYS
1	B	159	LEU
1	B	165	LYS
1	B	173	ASP

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Mol	Chain	Res	Type
1	B	182	LYS
1	B	186	LEU
1	B	202	THR
1	B	212	LEU
1	B	232	GLN
1	B	235	VAL
1	B	240	LYS
1	C	24	GLN
1	C	33	ASP
1	C	52	VAL
1	C	90	MET
1	C	94	LEU
1	C	97	ILE
1	C	159	LEU
1	C	181	LYS
1	C	192	LEU
1	C	196	ILE
1	C	197	GLU
1	C	202	THR
1	C	212	LEU
1	C	227	GLN
1	D	22	ASP
1	D	33	ASP
1	D	52	VAL
1	D	59	ILE
1	D	97	ILE
1	D	110	LYS
1	D	159	LEU
1	D	165	LYS
1	D	182	LYS
1	D	192	LEU
1	D	195	ASP
1	D	196	ILE
1	D	202	THR
1	D	225	LEU
1	D	227	GLN
1	E	7	ASN
1	E	8	LYS
1	E	11	LEU
1	E	23	ASN
1	E	52	VAL
1	E	53	ILE

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Mol	Chain	Res	Type
1	E	57	ASN
1	E	59	ILE
1	E	63	SER
1	E	66	GLN
1	E	94	LEU
1	E	97	ILE
1	E	119	ARG
1	E	159	LEU
1	E	164	THR
1	E	166	VAL
1	E	183	LEU
1	E	185	THR
1	E	202	THR
1	E	235	VAL
1	E	240	LYS
1	F	11	LEU
1	F	60	ARG
1	F	94	LEU
1	F	110	LYS
1	F	159	LEU
1	F	196	ILE
1	F	197	GLU
1	F	201	ASP
1	F	206	LEU
1	F	212	LEU
1	F	219	MET
1	F	240	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	B	23	ASN
1	B	139	ASN
1	B	232	GLN
1	C	23	ASN
1	C	227	GLN
1	C	233	GLN
1	D	227	GLN
1	E	30	HIS
1	E	57	ASN
1	E	233	GLN

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Mol	Chain	Res	Type
1	F	66	GLN
1	F	233	GLN

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 ⓘ

4 ligands 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	$\# Z > 2$
2	UDP	B	1241	-	18,26,26	1.25	2 (11%)	26,40,40	2.46	8 (30%)
2	UDP	C	1241	-	18,26,26	1.11	1 (5%)	26,40,40	2.08	3 (11%)
2	UDP	D	1241	-	18,26,26	1.20	1 (5%)	26,40,40	1.86	4 (15%)
2	UDP	E	1241	-	18,26,26	1.26	1 (5%)	26,40,40	1.50	2 (7%)

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
2	UDP	B	1241	-	-	0/12/32/32	0/2/2/2
2	UDP	C	1241	-	-	0/12/32/32	0/2/2/2
2	UDP	D	1241	-	-	0/12/32/32	0/2/2/2
2	UDP	E	1241	-	-	0/12/32/32	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1241	UDP	C6-N1	2.16	1.38	1.35
2	C	1241	UDP	C4-N3	2.64	1.38	1.33
2	D	1241	UDP	C4-N3	3.00	1.38	1.33
2	B	1241	UDP	C4-N3	3.14	1.38	1.33
2	E	1241	UDP	C4-N3	3.16	1.39	1.33

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1241	UDP	PA-O3A-PB	-7.13	108.75	132.67
2	B	1241	UDP	PA-O3A-PB	-6.83	109.75	132.67
2	D	1241	UDP	PA-O3A-PB	-5.36	114.70	132.67
2	B	1241	UDP	C6-N1-C2	-2.94	116.52	121.28
2	D	1241	UDP	O3A-PA-O5'	-2.04	97.52	102.94
2	D	1241	UDP	O3B-PB-O2B	2.06	115.23	107.38
2	E	1241	UDP	O4'-C1'-N1	2.16	112.64	108.08
2	B	1241	UDP	C2'-C3'-C4'	2.26	107.27	102.61
2	B	1241	UDP	O5'-C5'-C4'	2.35	117.80	109.12
2	C	1241	UDP	O3A-PA-O5'	2.63	109.92	102.94
2	B	1241	UDP	O4'-C1'-N1	2.83	114.06	108.08
2	B	1241	UDP	O3A-PA-O5'	3.72	112.81	102.94
2	B	1241	UDP	C4'-O4'-C1'	3.97	114.08	109.72
2	E	1241	UDP	C4-N3-C2	5.84	119.93	114.14
2	D	1241	UDP	C4-N3-C2	6.07	120.15	114.14
2	C	1241	UDP	C4-N3-C2	6.31	120.39	114.14
2	B	1241	UDP	C4-N3-C2	6.36	120.44	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1241	UDP	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1241	UDP	1	0
2	E	1241	UDP	4	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	234/240 (97%)	0.76	36 (15%) 3 3	32, 54, 86, 98	0
1	B	228/240 (95%)	0.25	11 (4%) 34 39	26, 51, 77, 94	0
1	C	234/240 (97%)	0.34	16 (6%) 20 23	23, 47, 71, 84	0
1	D	228/240 (95%)	0.15	12 (5%) 30 34	24, 45, 66, 73	0
1	E	227/240 (94%)	0.42	18 (7%) 15 17	25, 51, 74, 83	0
1	F	230/240 (95%)	0.60	30 (13%) 5 4	32, 53, 80, 97	0
All	All	1381/1440 (95%)	0.42	123 (8%) 12 13	23, 50, 77, 98	0

All (123) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	66	GLN	8.4
1	B	180	ALA	7.3
1	A	68	GLY	7.1
1	F	141	PHE	6.9
1	F	63	SER	6.9
1	A	67	GLY	6.7
1	C	177	PHE	6.6
1	F	65	ALA	6.4
1	F	64	ALA	6.2
1	F	62	VAL	6.1
1	A	63	SER	6.0
1	A	65	ALA	6.0
1	A	64	ALA	5.9
1	E	193	ILE	5.7
1	A	62	VAL	5.4
1	A	220	PHE	5.3
1	A	23	ASN	5.2
1	A	141	PHE	5.2
1	B	22	ASP	5.1

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Mol	Chain	Res	Type	RSRZ
1	A	61	GLY	5.1
1	E	22	ASP	5.1
1	F	72	ARG	5.1
1	F	67	GLY	5.0
1	A	22	ASP	4.9
1	F	68	GLY	4.9
1	E	171	ASP	4.8
1	E	195	ASP	4.7
1	A	72	ARG	4.7
1	C	175	ASN	4.7
1	C	22	ASP	4.6
1	F	66	GLN	4.6
1	E	194	GLY	4.3
1	A	60	ARG	4.3
1	B	195	ASP	4.1
1	E	170	TYR	4.0
1	A	173	ASP	3.9
1	A	174	PRO	3.9
1	E	12	VAL	3.8
1	E	23	ASN	3.8
1	C	176	LYS	3.8
1	F	61	GLY	3.7
1	F	22	ASP	3.7
1	B	25	PHE	3.6
1	E	196	ILE	3.6
1	A	175	ASN	3.5
1	C	178	LYS	3.4
1	A	195	ASP	3.3
1	D	22	ASP	3.2
1	A	177	PHE	3.2
1	E	173	ASP	3.1
1	E	25	PHE	3.1
1	D	24	GLN	3.1
1	B	220	PHE	3.0
1	C	174	PRO	3.0
1	F	179	ASP	3.0
1	A	180	ALA	3.0
1	F	24	GLN	3.0
1	A	11	LEU	3.0
1	C	181	LYS	3.0
1	A	24	GLN	2.9
1	C	195	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	220	PHE	2.9
1	B	171	ASP	2.9
1	C	211	LYS	2.9
1	D	194	GLY	2.9
1	F	220	PHE	2.9
1	E	189	ASN	2.8
1	F	23	ASN	2.8
1	B	170	TYR	2.8
1	F	76	ASP	2.8
1	F	71	ARG	2.8
1	E	172	LYS	2.8
1	F	73	THR	2.7
1	A	176	LYS	2.7
1	F	51	ILE	2.7
1	A	12	VAL	2.7
1	A	25	PHE	2.7
1	C	24	GLN	2.7
1	E	14	PHE	2.6
1	E	220	PHE	2.6
1	F	195	ASP	2.6
1	A	172	LYS	2.6
1	D	170	TYR	2.6
1	F	180	ALA	2.6
1	F	60	ARG	2.5
1	F	87	ALA	2.5
1	A	51	ILE	2.5
1	F	53	ILE	2.5
1	A	69	ILE	2.5
1	A	21	GLY	2.5
1	D	195	ASP	2.4
1	E	52	VAL	2.4
1	A	7	ASN	2.4
1	B	12	VAL	2.3
1	F	166	VAL	2.3
1	A	178	LYS	2.3
1	A	171	ASP	2.3
1	C	179	ASP	2.3
1	A	39	ILE	2.3
1	F	70	ILE	2.3
1	A	50	GLY	2.3
1	D	51	ILE	2.3
1	C	52	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	217	CYS	2.2
1	D	197	GLU	2.2
1	C	66	GLN	2.2
1	E	24	GLN	2.2
1	B	209	ASP	2.2
1	D	67	GLY	2.2
1	F	12	VAL	2.2
1	D	148	ALA	2.2
1	A	71	ARG	2.2
1	C	67	GLY	2.2
1	D	193	ILE	2.2
1	B	24	GLN	2.2
1	D	211	LYS	2.2
1	F	89	ALA	2.1
1	F	119	ARG	2.1
1	B	52	VAL	2.1
1	C	12	VAL	2.1
1	F	69	ILE	2.1
1	E	163	ALA	2.0
1	A	52	VAL	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
2	UDP	D	1241	25/25	0.81	0.20	1.24	41,47,55,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	UDP	B	1241	25/25	0.83	0.20	1.18	44,47,56,56	0
2	UDP	E	1241	25/25	0.81	0.21	1.05	44,47,53,55	0
2	UDP	C	1241	25/25	0.86	0.18	0.36	42,46,54,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.