



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:32 PM GMT

PDB ID : 4IOL
Title : N10-formyltetrahydrofolate synthetase from Moorella thermoacetica with ADP/ZD9 and XPO
Authors : Stec, B.
Deposited on : 2013-01-08
Resolution : 2.56 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

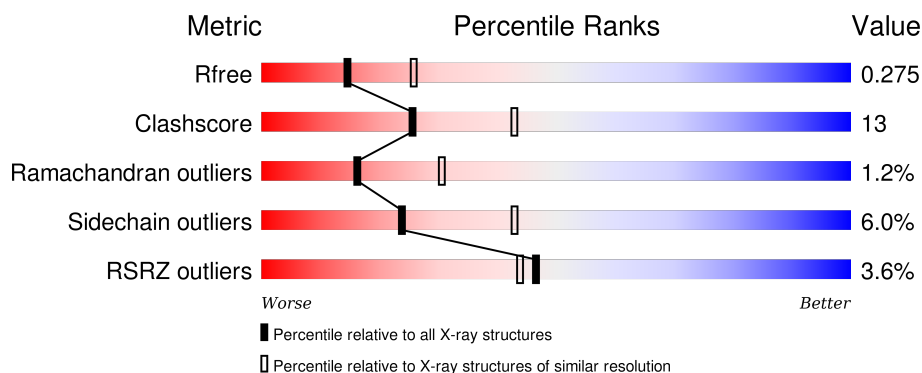
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.56 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	3324 (2.60-2.52)
Clashscore	102246	3729 (2.60-2.52)
Ramachandran outliers	100387	3673 (2.60-2.52)
Sidechain outliers	100360	3673 (2.60-2.52)
RSRZ outliers	91569	3333 (2.60-2.52)

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	559	<div> <div>2%</div> <div>79%</div> <div>18%</div> <div>.</div> </div>
1	B	559	<div> <div>6%</div> <div>63%</div> <div>35%</div> <div>.</div> </div>

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
3	ADP	B	602	-	-	-	X

2 Entry composition [i](#)

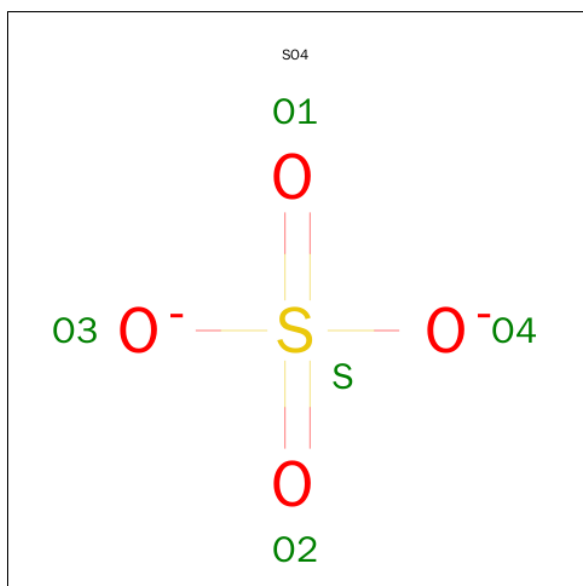
There are 6 unique types of molecules in this entry. The entry contains 8654 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Formate--tetrahydrofolate ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	557	Total	C	N	O	S	0	0	0
			4195	2661	721	792	21			
1	B	557	Total	C	N	O	S	0	0	0
			4195	2661	721	792	21			

- Molecule 2 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).

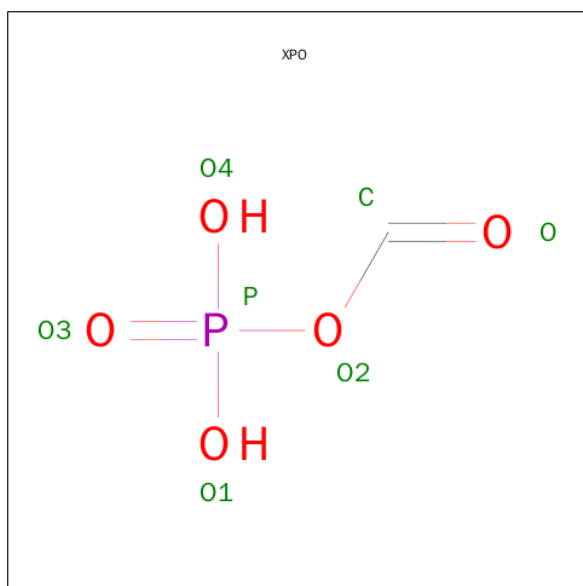


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

Chemical structure of ZD9, a benzimidazole derivative. The structure features a benzimidazole core with a 2-ethynyl-2-methylpropyl group at position 2 and a carbonyl group at position 6. Atoms are labeled with green text: C1-C16 for carbons, N1-N4 for nitrogens, and O1 for oxygen. The ethynyl group consists of C14-C15-C16, and the isopropyl group consists of C12-C13-C14. The benzimidazole ring system includes C1-C9, N1-N4, and O1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	1
			18	14	3	1		

- Molecule 5 is FORMYL PHOSPHATE (three-letter code: XPO) (formula: $\text{CH}_3\text{O}_5\text{P}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	O	P	0	1
			7	1	5	1		

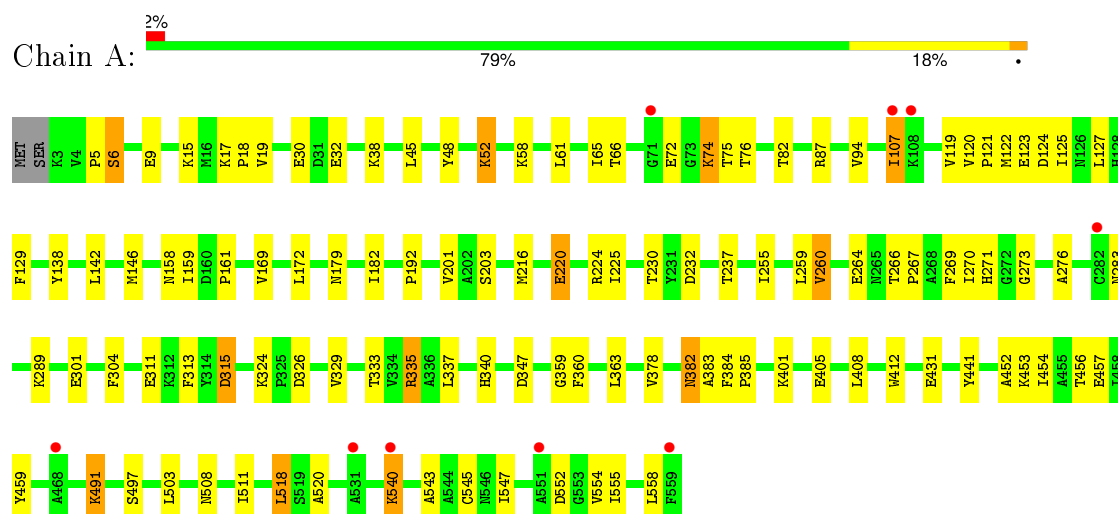
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	118	Total	O	0	0
			118	118		
6	B	47	Total	O	0	0
			47	47		

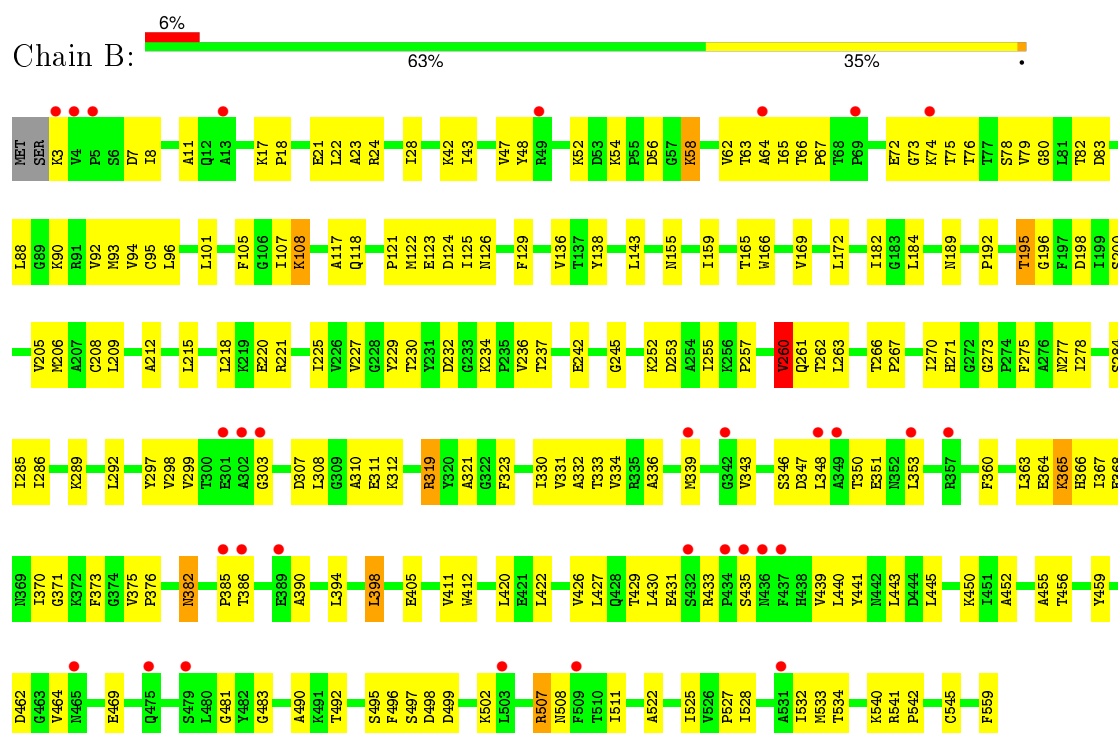
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Formate--tetrahydrofolate ligase



• Molecule 1: Formate--tetrahydrofolate ligase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	160.37Å 160.37Å 258.07Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	122.30 – 2.56 31.62 – 2.56	Depositor EDS
% Data completeness (in resolution range)	88.2 (122.30-2.56) 88.3 (31.62-2.56)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.29 (at 2.57Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.190 , 0.279 0.190 , 0.275	Depositor DCC
R_{free} test set	1822 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	60.2	Xtriage
Anisotropy	0.133	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 36.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 36317 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8654	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.40% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: XPO, ZD9, SO4, ADP

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	0.62	1/4266 (0.0%)	0.76	1/5778 (0.0%)
1	B	0.54	1/4266 (0.0%)	0.74	1/5778 (0.0%)
All	All	0.58	2/8532 (0.0%)	0.75	2/11556 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	412	TRP	CD2-CE2	5.30	1.47	1.41
1	B	412	TRP	CD2-CE2	5.02	1.47	1.41

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	445	LEU	CA-CB-CG	5.50	127.94	115.30
1	A	315	ASP	CB-CG-OD2	5.01	122.81	118.30

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	4195	0	4278	76	0
1	B	4195	0	4278	144	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	10	0	0	0	0
2	B	10	0	0	1	0
3	A	27	0	12	5	0
3	B	27	0	12	2	0
4	A	18	0	15	3	0
5	A	7	0	3	1	0
6	A	118	0	0	7	0
6	B	47	0	0	6	0
All	All	8654	0	8598	214	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 13.

All (214) 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:B:308:LEU:O	1:B:312:LYS:HG3	1.51	1.10
1:B:376:PRO:HD3	1:B:435:SER:HB2	1.44	0.98
1:A:405:GLU:HG2	6:A:791:HOH:O	1.66	0.96
1:B:422:LEU:O	1:B:426:VAL:HG23	1.73	0.88
1:A:383:ALA:HB2	1:A:408:LEU:HD11	1.57	0.86
1:B:126:ASN:HD22	1:B:559:PHE:HE1	1.22	0.86
1:A:491:LYS:HE3	1:A:497:SER:O	1.77	0.83
1:B:360:PHE:CE1	1:B:364:GLU:HB2	2.15	0.81
1:B:376:PRO:HD3	1:B:435:SER:CB	2.12	0.80
1:B:125:ILE:HD13	1:B:129:PHE:CE1	2.17	0.79
1:B:42:LYS:HD3	1:B:253:ASP:O	1.83	0.79
1:B:8:ILE:HA	1:B:122:MET:HE1	1.64	0.78
1:B:343:VAL:HG13	1:B:347:ASP:HB2	1.69	0.72
1:B:195:THR:OG1	1:B:196:GLY:N	2.22	0.72
1:B:126:ASN:ND2	1:B:559:PHE:HE1	1.87	0.72
1:B:455:ALA:HA	1:B:459:TYR:HD2	1.54	0.71
1:B:353:LEU:HD21	1:B:390:ALA:HB1	1.71	0.71
1:B:308:LEU:O	1:B:312:LYS:CG	2.35	0.70
1:B:339:MET:HG3	1:B:348:LEU:HD21	1.75	0.69
1:B:138:TYR:HB2	6:B:739:HOH:O	1.93	0.68
1:B:78:SER:O	1:B:299:VAL:HG11	1.93	0.67
1:B:93:MET:HG2	1:B:267:PRO:HB2	1.74	0.67
1:B:455:ALA:HA	1:B:459:TYR:CD2	2.29	0.67
1:B:429:THR:HG23	1:B:433:ARG:HD3	1.76	0.66
1:B:129:PHE:HD2	6:B:740:HOH:O	1.79	0.66

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:23:ALA:HA	1:B:28:ILE:HD12	1.78	0.66
1:B:262:THR:HG22	1:B:263:LEU:N	2.11	0.65
1:B:330:ILE:HD11	1:B:370:ILE:HD13	1.78	0.65
1:A:383:ALA:HB2	1:A:408:LEU:CD1	2.28	0.64
1:B:360:PHE:HE1	1:B:364:GLU:HB2	1.62	0.64
1:B:427:LEU:O	1:B:431:GLU:HG2	1.98	0.64
1:B:373:PHE:CE2	1:B:440:LEU:HB2	2.33	0.62
1:B:136:VAL:HG13	1:B:205:VAL:HG12	1.81	0.62
1:B:364:GLU:O	1:B:368:GLU:HG3	2.00	0.61
1:A:337:LEU:O	1:A:359:GLY:HA3	2.00	0.61
1:B:330:ILE:CD1	1:B:370:ILE:HD13	2.31	0.61
1:B:271:HIS:ND1	1:B:284:SER:OG	2.29	0.61
1:A:6:SER:OG	1:A:9:GLU:HG3	2.01	0.60
1:A:543:ALA:O	1:A:547:ILE:HG13	2.01	0.60
1:B:8:ILE:HA	1:B:122:MET:CE	2.31	0.60
1:B:8:ILE:HD12	1:B:122:MET:CE	2.32	0.59
1:A:216:MET:HG2	6:A:806:HOH:O	2.02	0.59
1:B:11:ALA:HB2	1:B:118:GLN:OE1	2.03	0.59
1:B:286:ILE:HD12	6:B:728:HOH:O	2.01	0.59
1:A:179:ASN:ND2	1:B:189:ASN:HD22	2.01	0.59
1:B:182:ILE:O	1:B:192:PRO:HA	2.03	0.59
1:B:67:PRO:HD3	6:B:743:HOH:O	2.04	0.58
1:A:61:LEU:HD13	1:A:313:PHE:CD2	2.39	0.58
1:B:440:LEU:HG	1:B:440:LEU:O	2.03	0.58
1:B:278:ILE:HG23	1:B:532:ILE:HG21	1.86	0.58
1:B:58:LYS:HD3	1:B:430:LEU:HD22	1.84	0.58
1:B:169:VAL:HB	1:B:200:SER:HA	1.85	0.57
1:B:376:PRO:CD	1:B:435:SER:CB	2.81	0.57
1:B:495:SER:OG	1:B:502:LYS:O	2.16	0.57
1:A:120:VAL:HB	1:A:121:PRO:HA	1.87	0.57
1:B:376:PRO:CD	1:B:435:SER:HB2	2.29	0.56
1:B:260:VAL:HG21	1:B:270:ILE:HD12	1.86	0.56
1:B:289:LYS:HB3	1:B:323:PHE:HZ	1.70	0.56
1:B:303:GLY:HA2	2:B:601:SO4:O2	2.06	0.56
1:B:363:LEU:O	1:B:367:ILE:HG13	2.05	0.56
1:A:87:ARG:HH21	1:A:264:GLU:CD	2.09	0.56
1:A:74:LYS:H	3:A:602[A]:ADP:PB	2.29	0.56
1:B:121:PRO:O	1:B:125:ILE:HG12	2.06	0.56
1:B:292:LEU:HD23	1:B:298:VAL:HG21	1.88	0.55
1:B:88:LEU:HD21	1:B:420:LEU:CD2	2.36	0.55
1:A:452:ALA:O	1:A:456:THR:HG23	2.06	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:79:VAL:HG11	1:B:117:ALA:O	2.06	0.55
1:A:19:VAL:HG22	1:A:38:LYS:O	2.07	0.55
1:A:172:LEU:HG	1:B:138:TYR:CZ	2.42	0.55
1:B:278:ILE:CG2	1:B:532:ILE:HG21	2.37	0.55
1:A:220:GLU:OE1	1:A:224:ARG:NH2	2.41	0.54
1:A:453:LYS:HE2	6:A:710:HOH:O	2.07	0.54
1:B:62:VAL:HG21	1:B:78:SER:HA	1.90	0.53
1:A:311:GLU:OE1	1:A:315:ASP:OD1	2.27	0.53
1:B:65:ILE:HA	1:B:366:HIS:CD2	2.44	0.53
1:B:125:ILE:HD12	1:B:270:ILE:HD13	1.91	0.53
1:B:166:TRP:HB2	1:B:227:VAL:HA	1.91	0.53
1:B:394:LEU:O	1:B:398:LEU:HD12	2.09	0.52
1:A:179:ASN:HD22	1:B:189:ASN:HD22	1.58	0.52
1:B:143:LEU:HD23	1:B:166:TRP:CE2	2.44	0.52
1:B:80:GLY:HA3	1:B:411:VAL:HG11	1.92	0.52
1:A:545:CYS:HA	1:B:242:GLU:O	2.09	0.52
1:B:215:LEU:HA	1:B:255:ILE:CD1	2.39	0.52
1:B:95:CYS:O	1:B:96:LEU:HD23	2.10	0.52
1:B:83:ASP:OD1	1:B:262:THR:HG21	2.10	0.52
1:B:277:ASN:OD1	1:B:278:ILE:N	2.39	0.52
1:B:143:LEU:HD23	1:B:166:TRP:NE1	2.26	0.51
1:A:142:LEU:O	1:A:146:MET:HG3	2.10	0.51
1:A:385:PRO:HD3	4:A:603[B]:ZD9:H39B	1.92	0.51
1:B:343:VAL:HG13	1:B:347:ASP:CB	2.39	0.51
1:B:225:ILE:O	1:B:237:THR:HA	2.11	0.51
1:B:63:THR:O	1:B:331:VAL:HG23	2.12	0.50
1:A:385:PRO:HD3	4:A:603[B]:ZD9:C39	2.42	0.50
1:A:75:THR:HA	1:A:301:GLU:OE2	2.11	0.50
1:B:229:TYR:N	1:B:229:TYR:CD1	2.78	0.50
1:B:221:ARG:NH1	1:B:522:ALA:O	2.43	0.50
1:B:497:SER:HA	1:B:511:ILE:HD11	1.92	0.50
1:B:441:TYR:CD1	1:B:450:LYS:HD3	2.47	0.49
1:B:88:LEU:HD21	1:B:420:LEU:HD23	1.95	0.49
1:B:307:ASP:OD1	1:B:307:ASP:N	2.45	0.49
1:A:75:THR:N	5:A:604[C]:XPO:O2	2.46	0.49
1:B:94:VAL:HA	1:B:299:VAL:O	2.13	0.48
1:B:92:VAL:HA	1:B:297:TYR:O	2.14	0.48
1:B:285:ILE:HD13	1:B:321:ALA:HB2	1.95	0.48
1:A:335:ARG:HG3	6:A:808:HOH:O	2.12	0.48
1:A:276:ALA:HB3	1:A:304:PHE:CD2	2.49	0.48
1:A:82:THR:HG22	1:A:266:THR:HG21	1.96	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:LEU:HD11	1:B:261:GLN:NE2	2.29	0.48
1:A:125:ILE:HA	1:A:129:PHE:CD1	2.48	0.48
1:B:8:ILE:CD1	1:B:122:MET:CE	2.91	0.48
1:A:453:LYS:O	1:A:457:GLU:HB2	2.14	0.48
1:B:48:TYR:HE1	6:B:729:HOH:O	1.97	0.48
1:B:262:THR:HG22	1:B:263:LEU:H	1.77	0.47
1:A:169:VAL:HG11	1:A:203:SER:HB2	1.95	0.47
1:B:74:LYS:O	1:B:78:SER:HB2	2.14	0.47
1:A:547:ILE:O	1:B:245:GLY:HA3	2.14	0.47
1:A:61:LEU:HD13	1:A:313:PHE:CE2	2.49	0.47
1:A:552:ASP:HB2	1:A:554:VAL:HG23	1.96	0.47
1:B:79:VAL:CG1	1:B:117:ALA:O	2.62	0.47
1:B:101:LEU:CD2	1:B:105:PHE:HE1	2.28	0.47
1:B:11:ALA:HB3	1:B:122:MET:HE2	1.98	0.47
1:B:8:ILE:CD1	1:B:122:MET:HE1	2.45	0.47
1:A:82:THR:OG1	1:A:94:VAL:HB	2.16	0.46
1:A:182:ILE:O	1:A:192:PRO:HA	2.14	0.46
1:B:107:ILE:O	1:B:108:LYS:HB2	2.15	0.46
1:A:382:ASN:ND2	6:A:813:HOH:O	2.48	0.46
1:B:169:VAL:HA	1:B:198:ASP:O	2.16	0.46
1:B:67:PRO:HA	1:B:72:GLU:CD	2.36	0.46
1:A:72:GLU:HB2	3:A:602[A]:ADP:O2B	2.16	0.46
1:B:159:ILE:HA	1:B:230:THR:HA	1.97	0.46
1:B:333:THR:HB	1:B:336:ALA:H	1.80	0.46
1:B:67:PRO:HA	1:B:72:GLU:OE2	2.15	0.46
1:B:75:THR:O	1:B:76:THR:C	2.54	0.46
1:A:441:TYR:CE2	1:A:454:ILE:HD11	2.52	0.45
1:B:198:ASP:OD1	1:B:534:THR:OG1	2.34	0.45
1:B:262:THR:CG2	1:B:263:LEU:N	2.80	0.45
1:A:382:ASN:OD1	3:A:602[A]:ADP:N7	2.50	0.45
1:B:56:ASP:HB3	1:B:292:LEU:HD13	1.97	0.45
1:A:459:TYR:HD2	1:A:511:ILE:HD11	1.82	0.45
1:A:48:TYR:CD1	1:A:289:LYS:HG3	2.51	0.45
1:B:82:THR:HG21	1:B:94:VAL:HB	1.98	0.45
1:A:45:LEU:HD23	6:A:805:HOH:O	2.15	0.45
1:B:17:LYS:O	1:B:18:PRO:C	2.53	0.45
1:B:490:ALA:HB3	1:B:528:ILE:HA	1.99	0.45
1:B:382:ASN:HD21	3:B:602:ADP:HN61	1.65	0.45
1:A:459:TYR:CD2	1:A:511:ILE:HD11	2.52	0.45
1:B:462:ASP:OD2	1:B:508:ASN:HA	2.16	0.45
1:B:376:PRO:CD	1:B:435:SER:HB3	2.46	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:441:TYR:CE1	1:B:450:LYS:HD3	2.52	0.44
1:A:124:ASP:O	1:A:129:PHE:HA	2.16	0.44
1:A:119:VAL:HG11	1:A:270:ILE:HD11	1.99	0.44
1:A:271:HIS:O	1:A:283:ASN:HB2	2.17	0.44
1:B:21:GLU:O	1:B:24:ARG:HB2	2.18	0.44
1:A:201:VAL:HB	1:A:273:GLY:O	2.17	0.44
1:B:339:MET:CG	1:B:348:LEU:HD21	2.47	0.44
1:A:138:TYR:CZ	1:B:172:LEU:HG	2.52	0.44
1:B:507:ARG:O	1:B:508:ASN:HB2	2.17	0.44
1:B:499:ASP:HB3	1:B:502:LYS:HG3	2.00	0.43
1:A:124:ASP:HB3	6:A:758:HOH:O	2.17	0.43
1:B:292:LEU:HD23	1:B:298:VAL:CG2	2.48	0.43
1:B:107:ILE:HG13	1:B:108:LYS:H	1.84	0.43
1:A:518:LEU:HD21	1:A:520:ALA:HB2	1.99	0.43
1:A:87:ARG:NH2	1:A:264:GLU:OE2	2.49	0.43
1:A:138:TYR:CZ	1:B:101:LEU:HD13	2.53	0.43
1:B:125:ILE:HD13	1:B:129:PHE:CD1	2.54	0.43
1:A:333:THR:HG22	1:A:382:ASN:HB3	2.01	0.43
1:B:375:VAL:HA	1:B:376:PRO:HD3	1.90	0.43
1:B:8:ILE:CA	1:B:122:MET:HE1	2.44	0.43
3:A:602[A]:ADP:H8	3:A:602[A]:ADP:O1A	2.01	0.43
1:A:555:ILE:HG21	1:A:558:LEU:HD23	2.01	0.43
1:A:540:LYS:HG3	1:A:540:LYS:H	1.56	0.43
1:A:329:VAL:HA	1:A:378:VAL:O	2.19	0.43
1:A:382:ASN:HA	1:A:382:ASN:HD22	1.68	0.42
1:B:73:GLY:HA2	3:B:602:ADP:O5'	2.18	0.42
1:A:17:LYS:O	1:A:18:PRO:C	2.55	0.42
1:B:79:VAL:O	1:B:82:THR:HB	2.18	0.42
1:B:90:LYS:HB3	1:B:297:TYR:CD2	2.53	0.42
1:A:360:PHE:O	1:A:363:LEU:N	2.53	0.42
1:A:123:GLU:O	1:A:127:LEU:HG	2.19	0.42
1:A:30:GLU:C	1:A:32:GLU:H	2.23	0.42
1:B:365:LYS:HG3	1:B:496:PHE:CE1	2.55	0.42
1:B:260:VAL:HB	1:B:261:GLN:H	1.68	0.42
1:A:74:LYS:HB3	3:A:602[A]:ADP:O1B	2.19	0.42
1:A:255:ILE:HG13	1:A:255:ILE:O	2.20	0.42
1:B:464:VAL:HA	1:B:511:ILE:O	2.20	0.42
1:B:208:CYS:O	1:B:209:LEU:C	2.59	0.41
1:B:370:ILE:O	1:B:371:GLY:C	2.58	0.41
1:B:319:ARG:HE	1:B:443:LEU:HD21	1.85	0.41
1:B:257:PRO:HA	1:B:271:HIS:CD2	2.56	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:236:VAL:HG12	1:B:237:THR:N	2.34	0.41
1:A:266:THR:CG2	1:A:267:PRO:HD2	2.50	0.41
1:B:155:ASN:N	6:B:706:HOH:O	2.52	0.41
1:B:125:ILE:CD1	1:B:129:PHE:CE1	2.97	0.41
1:B:8:ILE:HD12	1:B:122:MET:HE1	2.02	0.41
1:B:64:ALA:HA	1:B:331:VAL:H	1.86	0.41
1:A:52:LYS:HE2	1:A:52:LYS:HB3	1.73	0.41
1:B:481:GLY:C	1:B:483:GLY:H	2.23	0.41
1:B:525:ILE:HG22	1:B:527:PRO:HD3	2.02	0.41
1:B:212:ALA:HB2	1:B:218:LEU:HD13	2.03	0.41
1:A:124:ASP:O	1:A:129:PHE:CA	2.69	0.41
1:A:259:LEU:O	1:A:260:VAL:HG13	2.20	0.41
1:B:184:LEU:CD2	1:B:192:PRO:HB3	2.51	0.41
1:B:452:ALA:O	1:B:456:THR:HG22	2.21	0.41
1:B:331:VAL:HG12	1:B:332:ALA:H	1.86	0.40
1:B:43:ILE:CG2	1:B:47:VAL:HG21	2.50	0.40
1:A:66:THR:HB	1:A:340:HIS:CE1	2.56	0.40
1:B:310:ALA:O	1:B:311:GLU:C	2.60	0.40
1:A:159:ILE:O	1:A:161:PRO:HD3	2.21	0.40
1:A:158:ASN:O	1:A:230:THR:HA	2.21	0.40
1:A:225:ILE:O	1:A:237:THR:HA	2.21	0.40
1:A:58:LYS:HA	1:A:326:ASP:OD2	2.21	0.40
1:B:542:PRO:HD2	1:B:545:CYS:SG	2.61	0.40
1:A:611:LEU:HD13	1:A:313:PHE:CG	2.57	0.40
1:A:385:PRO:HG3	4:A:603[B]:ZD9:H39	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	555/559 (99%)	503 (91%)	46 (8%)	6 (1%)	17 33

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	555/559 (99%)	475 (86%)	73 (13%)	7 (1%)	15	28
All	All	1110/1118 (99%)	978 (88%)	119 (11%)	13 (1%)	16	30

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	260	VAL
1	B	108	LYS
1	B	252	LYS
1	A	107	ILE
1	A	335	ARG
1	A	122	MET
1	B	7	ASP
1	A	5	PRO
1	B	492	THR
1	B	273	GLY
1	B	385	PRO
1	B	260	VAL
1	A	65	ILE

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	439/441 (100%)	419 (95%)	20 (5%)	33	57
1	B	439/441 (100%)	406 (92%)	33 (8%)	17	30
All	All	878/882 (100%)	825 (94%)	53 (6%)	24	43

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	SER
1	A	15	LYS
1	A	52	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	74	LYS
1	A	76	THR
1	A	107	ILE
1	A	220	GLU
1	A	232	ASP
1	A	269	PHE
1	A	324	LYS
1	A	347	ASP
1	A	382	ASN
1	A	384	PHE
1	A	401	LYS
1	A	431	GLU
1	A	491	LYS
1	A	503	LEU
1	A	508	ASN
1	A	518	LEU
1	A	540	LYS
1	B	3	LYS
1	B	52	LYS
1	B	54	LYS
1	B	58	LYS
1	B	66	THR
1	B	123	GLU
1	B	124	ASP
1	B	165	THR
1	B	195	THR
1	B	206	MET
1	B	220	GLU
1	B	232	ASP
1	B	234	LYS
1	B	260	VAL
1	B	266	THR
1	B	275	PHE
1	B	319	ARG
1	B	334	VAL
1	B	346	SER
1	B	350	THR
1	B	351	GLU
1	B	365	LYS
1	B	382	ASN
1	B	386	THR
1	B	398	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	405	GLU
1	B	439	VAL
1	B	469	GLU
1	B	498	ASP
1	B	507	ARG
1	B	533	MET
1	B	540	LYS
1	B	541	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	179	ASN
1	A	283	ASN
1	A	382	ASN
1	B	29	GLN
1	B	158	ASN
1	B	179	ASN
1	B	265	ASN
1	B	382	ASN
1	B	438	HIS
1	B	465	ASN

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](#)

8 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	SO4	A	601	-	4,4,4	0.64	0	6,6,6	0.64	0
3	ADP	A	602[A]	-	22,29,29	0.93	1 (4%)	27,45,45	1.91	2 (7%)
4	ZD9	A	603[B]	-	18,19,19	1.85	2 (11%)	16,26,26	2.04	6 (37%)
5	XPO	A	604[C]	-	4,6,6	2.27	2 (50%)	5,8,8	0.77	0
2	SO4	A	605	-	4,4,4	0.56	0	6,6,6	0.28	0
2	SO4	B	601	-	4,4,4	0.40	0	6,6,6	0.31	0
3	ADP	B	602	-	22,29,29	1.09	2 (9%)	27,45,45	1.96	5 (18%)
2	SO4	B	603	-	4,4,4	0.43	0	6,6,6	0.16	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
2	SO4	A	601	-	-	0/0/0/0	0/0/0/0
3	ADP	A	602[A]	-	-	0/12/32/32	0/3/3/3
4	ZD9	A	603[B]	-	-	0/3/5/5	0/2/2/2
5	XPO	A	604[C]	-	-	0/0/4/4	0/0/0/0
2	SO4	A	605	-	-	0/0/0/0	0/0/0/0
2	SO4	B	601	-	-	0/0/0/0	0/0/0/0
3	ADP	B	602	-	-	0/12/32/32	0/3/3/3
2	SO4	B	603	-	-	0/0/0/0	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	603[B]	ZD9	C12-N13	-5.28	1.37	1.45
5	A	604[C]	XPO	P-O4	2.19	1.62	1.54
3	B	602	ADP	O4'-C1'	2.31	1.44	1.41
3	B	602	ADP	C5-C4	2.96	1.47	1.40
3	A	602[A]	ADP	C5-C4	3.03	1.47	1.40
5	A	604[C]	XPO	P-O3	3.45	1.62	1.51
4	A	603[B]	ZD9	C1-C6	4.74	1.49	1.41

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	602	ADP	N3-C2-N1	-7.96	122.80	128.89
3	A	602[A]	ADP	N3-C2-N1	-7.41	123.22	128.89
3	A	602[A]	ADP	C4-C5-N7	-3.74	106.04	109.48
4	A	603[B]	ZD9	C9-C8-C6	-2.67	119.86	122.89
3	B	602	ADP	C4-C5-N7	-2.13	107.52	109.48
3	B	602	ADP	C4'-O4'-C1'	2.02	111.94	109.72
4	A	603[B]	ZD9	C8-C6-C5	2.11	120.66	118.14
3	B	602	ADP	C2'-C3'-C4'	2.23	107.19	102.61
4	A	603[B]	ZD9	C39-C3-N4	2.35	121.20	117.21
3	B	602	ADP	O2B-PB-O1B	2.44	118.42	110.58
4	A	603[B]	ZD9	C9-C12-N13	2.75	118.65	113.08
4	A	603[B]	ZD9	C11-C10-C9	3.07	120.27	117.92
4	A	603[B]	ZD9	C3-N4-C5	4.82	121.30	115.86

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602[A]	ADP	5	0
4	A	603[B]	ZD9	3	0
5	A	604[C]	XPO	1	0
2	B	601	SO4	1	0
3	B	602	ADP	2	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	557/559 (99%)	-0.18	9 (1%) 74 73	33, 49, 79, 129	0
1	B	557/559 (99%)	0.19	31 (5%) 28 26	31, 72, 107, 149	0
All	All	1114/1118 (99%)	0.00	40 (3%) 46 44	31, 60, 102, 149	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	349	ALA	5.4
1	B	3	LYS	5.2
1	B	503	LEU	3.9
1	A	559	PHE	3.9
1	A	108	LYS	3.4
1	A	540	LYS	3.4
1	B	437	PHE	3.4
1	A	107	ILE	3.3
1	B	4	VAL	3.2
1	B	357	ARG	3.2
1	B	342	GLY	3.1
1	B	389	GLU	3.1
1	B	475	GLN	3.1
1	B	479	SER	3.0
1	B	434	PRO	2.9
1	B	509	PHE	2.9
1	B	436	ASN	2.8
1	B	386	THR	2.8
1	A	551	ALA	2.8
1	B	302	ALA	2.8
1	B	435	SER	2.7
1	B	5	PRO	2.7
1	B	385	PRO	2.7
1	B	432	SER	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	13	ALA	2.6
1	B	64	ALA	2.5
1	B	69	PRO	2.5
1	B	303	GLY	2.4
1	B	339	MET	2.4
1	B	348	LEU	2.3
1	A	71	GLY	2.2
1	B	74	LYS	2.2
1	B	531	ALA	2.2
1	B	353	LEU	2.2
1	A	531	ALA	2.2
1	A	282	CYS	2.1
1	B	301	GLU	2.1
1	A	468	ALA	2.0
1	B	49	ARG	2.0
1	B	465	ASN	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	ADP	B	602	27/27	0.77	0.35	2.22	59,71,80,81	27
4	ZD9	A	603[B]	18/18	0.95	0.16	0.88	53,56,67,69	18
2	SO4	B	603	5/5	0.95	0.14	-0.01	92,98,103,104	0
2	SO4	A	605	5/5	0.93	0.12	-0.50	88,94,109,113	0
3	ADP	A	602[A]	27/27	0.93	0.14	-0.54	48,64,76,81	27

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	XPO	A	604[C]	7/7	0.97	0.14	-0.79	57,60,66,68	7
2	SO4	A	601	5/5	0.97	0.12	-1.47	50,64,70,70	0
2	SO4	B	601	5/5	0.95	0.22	-	76,87,93,95	0

6.5 Other polymers [i](#)

There are no such residues in this entry.