



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:01 AM UTC

PDB ID : 2FTY / pdb_00002fty
Title : Crystal structure of dihydropyrimidinase from *Saccharomyces kluyveri*
Authors : Dobritsch, D.; Lohkamp, B.
Deposited on : 2006-01-25
Resolution : 2.40 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

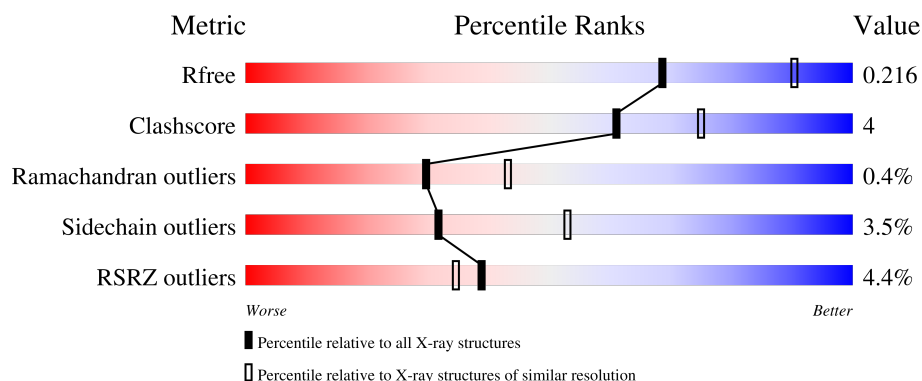
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	559	<div> <div>3%</div> <div>83% 12% 5%</div> </div>
1	B	559	<div> <div>4%</div> <div>82% 13% 5%</div> </div>
1	C	559	<div> <div>6%</div> <div>84% 11% 5%</div> </div>
1	D	559	<div> <div>4%</div> <div>82% 12% 5%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17311 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 dihydropyrimidinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	532	Total	C	N	O	S	0	2	0
			4164	2646	680	811	27			
1	B	532	Total	C	N	O	S	0	3	0
			4169	2651	680	811	27			
1	C	531	Total	C	N	O	S	0	0	0
			4142	2630	677	809	26			
1	D	530	Total	C	N	O	S	0	2	0
			4147	2633	678	809	27			

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	167	KCX	LYS	modified residue	UNP Q9P903
A	543	PRO	-	expression tag	UNP Q9P903
A	544	GLY	-	expression tag	UNP Q9P903
A	545	ASP	-	expression tag	UNP Q9P903
A	546	ASP	-	expression tag	UNP Q9P903
A	547	ASP	-	expression tag	UNP Q9P903
A	548	ASP	-	expression tag	UNP Q9P903
A	549	LYS	-	expression tag	UNP Q9P903
A	550	HIS	-	expression tag	UNP Q9P903
A	551	HIS	-	expression tag	UNP Q9P903
A	552	HIS	-	expression tag	UNP Q9P903
A	553	HIS	-	expression tag	UNP Q9P903
A	554	HIS	-	expression tag	UNP Q9P903
A	555	HIS	-	expression tag	UNP Q9P903
A	556	HIS	-	expression tag	UNP Q9P903
A	557	HIS	-	expression tag	UNP Q9P903
A	558	SER	-	expression tag	UNP Q9P903
A	559	GLY	-	expression tag	UNP Q9P903
A	560	ASP	-	expression tag	UNP Q9P903
B	167	KCX	LYS	modified residue	UNP Q9P903
B	543	PRO	-	expression tag	UNP Q9P903

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Chain	Residue	Modelled	Actual	Comment	Reference
B	544	GLY	-	expression tag	UNP Q9P903
B	545	ASP	-	expression tag	UNP Q9P903
B	546	ASP	-	expression tag	UNP Q9P903
B	547	ASP	-	expression tag	UNP Q9P903
B	548	ASP	-	expression tag	UNP Q9P903
B	549	LYS	-	expression tag	UNP Q9P903
B	550	HIS	-	expression tag	UNP Q9P903
B	551	HIS	-	expression tag	UNP Q9P903
B	552	HIS	-	expression tag	UNP Q9P903
B	553	HIS	-	expression tag	UNP Q9P903
B	554	HIS	-	expression tag	UNP Q9P903
B	555	HIS	-	expression tag	UNP Q9P903
B	556	HIS	-	expression tag	UNP Q9P903
B	557	HIS	-	expression tag	UNP Q9P903
B	558	SER	-	expression tag	UNP Q9P903
B	559	GLY	-	expression tag	UNP Q9P903
B	560	ASP	-	expression tag	UNP Q9P903
C	167	KCX	LYS	modified residue	UNP Q9P903
C	543	PRO	-	expression tag	UNP Q9P903
C	544	GLY	-	expression tag	UNP Q9P903
C	545	ASP	-	expression tag	UNP Q9P903
C	546	ASP	-	expression tag	UNP Q9P903
C	547	ASP	-	expression tag	UNP Q9P903
C	548	ASP	-	expression tag	UNP Q9P903
C	549	LYS	-	expression tag	UNP Q9P903
C	550	HIS	-	expression tag	UNP Q9P903
C	551	HIS	-	expression tag	UNP Q9P903
C	552	HIS	-	expression tag	UNP Q9P903
C	553	HIS	-	expression tag	UNP Q9P903
C	554	HIS	-	expression tag	UNP Q9P903
C	555	HIS	-	expression tag	UNP Q9P903
C	556	HIS	-	expression tag	UNP Q9P903
C	557	HIS	-	expression tag	UNP Q9P903
C	558	SER	-	expression tag	UNP Q9P903
C	559	GLY	-	expression tag	UNP Q9P903
C	560	ASP	-	expression tag	UNP Q9P903
D	167	KCX	LYS	modified residue	UNP Q9P903
D	543	PRO	-	expression tag	UNP Q9P903
D	544	GLY	-	expression tag	UNP Q9P903
D	545	ASP	-	expression tag	UNP Q9P903
D	546	ASP	-	expression tag	UNP Q9P903
D	547	ASP	-	expression tag	UNP Q9P903

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Chain	Residue	Modelled	Actual	Comment	Reference
D	548	ASP	-	expression tag	UNP Q9P903
D	549	LYS	-	expression tag	UNP Q9P903
D	550	HIS	-	expression tag	UNP Q9P903
D	551	HIS	-	expression tag	UNP Q9P903
D	552	HIS	-	expression tag	UNP Q9P903
D	553	HIS	-	expression tag	UNP Q9P903
D	554	HIS	-	expression tag	UNP Q9P903
D	555	HIS	-	expression tag	UNP Q9P903
D	556	HIS	-	expression tag	UNP Q9P903
D	557	HIS	-	expression tag	UNP Q9P903
D	558	SER	-	expression tag	UNP Q9P903
D	559	GLY	-	expression tag	UNP Q9P903
D	560	ASP	-	expression tag	UNP Q9P903

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total 2	Zn 2	0	0
2	B	2	Total 2	Zn 2	0	0
2	C	2	Total 2	Zn 2	0	0
2	D	2	Total 2	Zn 2	0	0

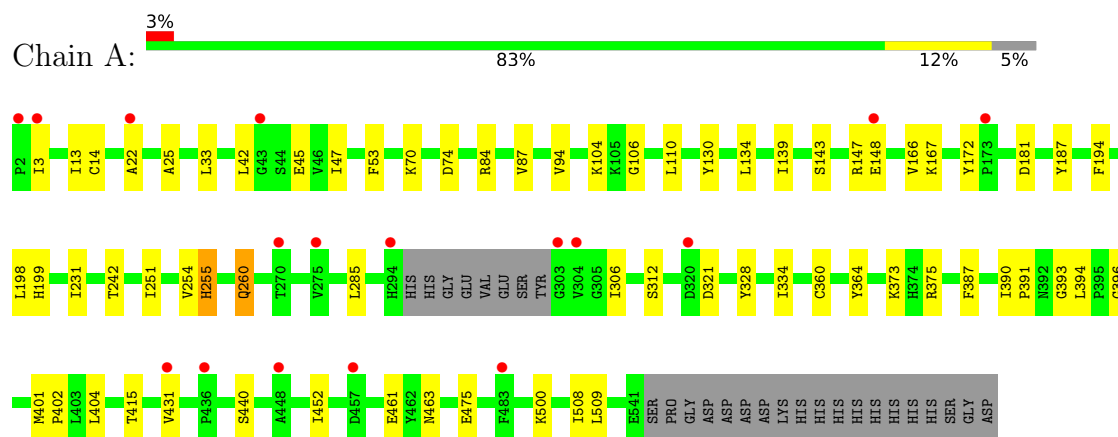
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	189	Total 189	O 189	0	0
3	B	188	Total 188	O 188	0	0
3	C	171	Total 171	O 171	0	0
3	D	133	Total 133	O 133	0	0

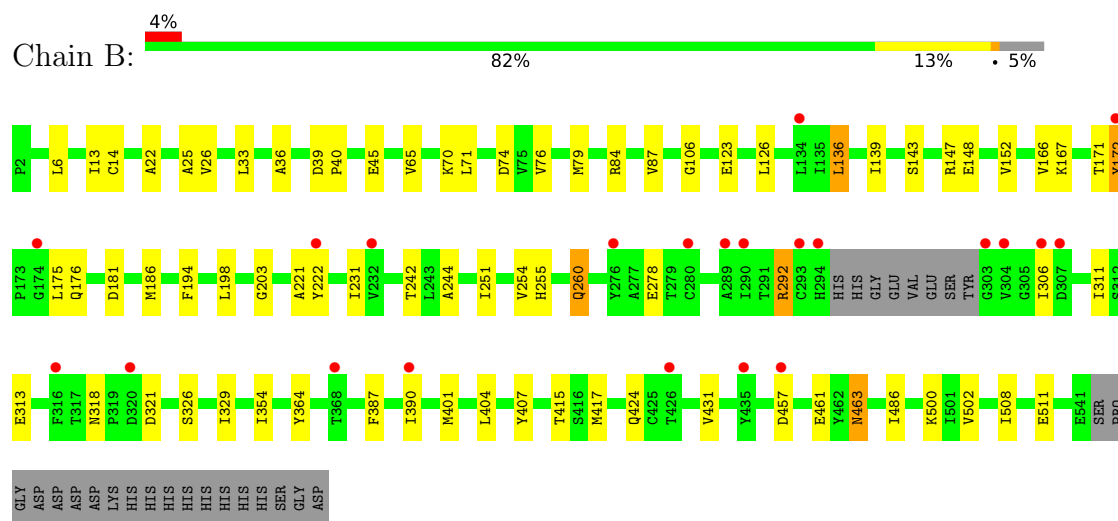
3 Residue-property plots

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.

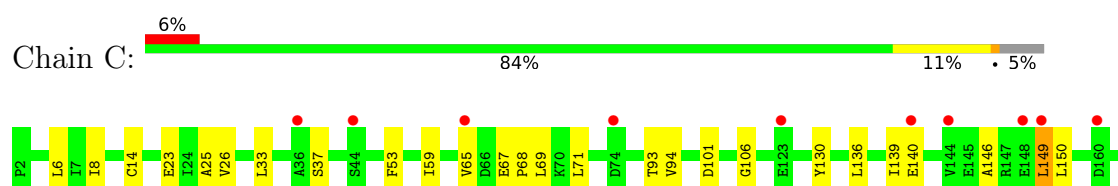
• Molecule 1: dihydropyrimidinase

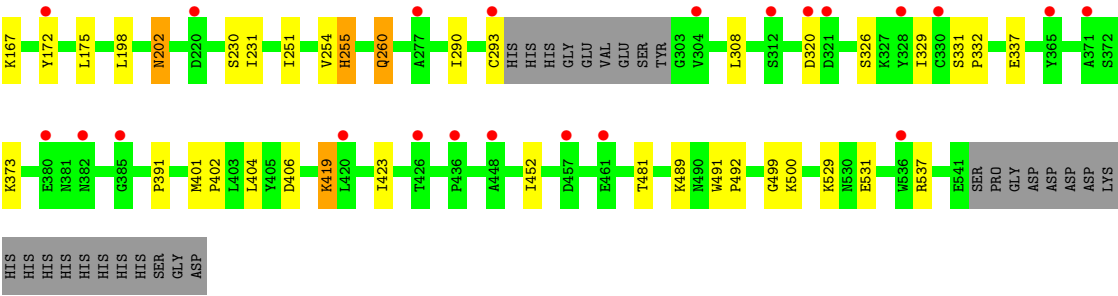


• Molecule 1: dihydropyrimidinase

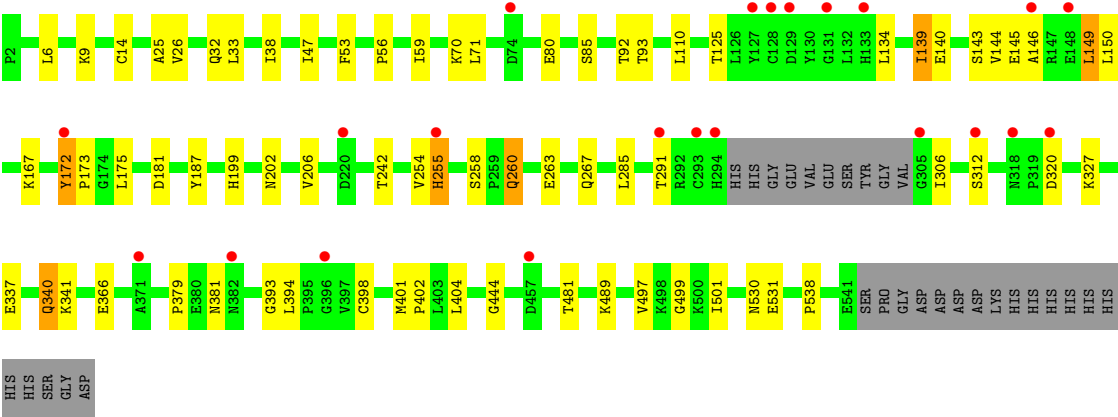
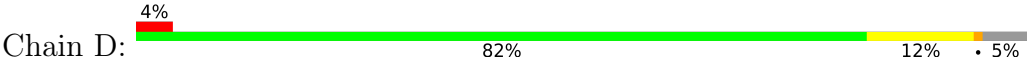


• Molecule 1: dihydropyrimidinase





• Molecule 1: dihydropyrimidinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	91.43Å 73.32Å 162.18Å 90.00° 91.59° 90.00°	Depositor
Resolution (Å)	43.85 – 2.40 43.85 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.0 (43.85-2.40) 99.0 (43.85-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.10 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.217 , 0.274 0.222 , 0.216	Depositor DCC
R_{free} test set	4200 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	35.0	Xtriage
Anisotropy	0.598	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.126 for h,-k,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	17311	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.51% 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: KCX, ZN

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	0.60	0/4252	0.86	2/5771 (0.0%)
1	B	0.59	0/4260	0.85	2/5782 (0.0%)
1	C	0.60	0/4222	0.87	2/5729 (0.0%)
1	D	0.58	0/4234	0.85	0/5745
All	All	0.59	0/16968	0.86	6/23027 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	106	GLY	CA-C-N	5.98	126.41	119.47
1	C	106	GLY	C-N-CA	5.98	126.41	119.47
1	A	106	GLY	CA-C-N	5.93	125.80	119.28
1	A	106	GLY	C-N-CA	5.93	125.80	119.28
1	B	106	GLY	CA-C-N	5.15	124.94	119.28
1	B	106	GLY	C-N-CA	5.15	124.94	119.28

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	4164	0	4093	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	4169	0	4104	42	0
1	C	4142	0	4072	32	0
1	D	4147	0	4077	37	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	189	0	0	2	0
3	B	188	0	0	0	0
3	C	171	0	0	3	0
3	D	133	0	0	3	0
All	All	17311	0	16346	138	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 4.

All (138) 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:260:GLN:HE21	1:C:260:GLN:H	1.14	0.92
1:C:140:GLU:O	1:C:146:ALA:HB2	1.78	0.83
1:D:140:GLU:O	1:D:146:ALA:HB2	1.77	0.83
1:A:198:LEU:HD12	1:A:251:ILE:HD11	1.60	0.83
1:B:198:LEU:HD12	1:B:251[A]:ILE:HD11	1.63	0.81
1:C:198:LEU:HD12	1:C:251:ILE:HD11	1.69	0.73
1:D:337:GLU:HA	1:D:340:GLN:HE22	1.54	0.73
1:A:231:ILE:HG12	1:B:242:THR:HG23	1.71	0.72
1:A:84:ARG:NH1	1:A:475:GLU:OE1	2.24	0.69
1:B:401:MET:HE3	1:B:424:GLN:HB3	1.76	0.68
1:D:71:LEU:HD11	1:D:175:LEU:HD21	1.75	0.67
1:C:140:GLU:O	1:C:146:ALA:CB	2.43	0.66
1:B:171:THR:O	1:B:172[A]:TYR:HB2	1.97	0.64
1:B:326:SER:HA	1:B:329:ILE:HG12	1.80	0.64
1:B:70:LYS:HD2	1:B:74:ASP:OD1	1.99	0.62
1:D:140:GLU:O	1:D:146:ALA:CB	2.47	0.60
1:B:14[B]:CYS:SG	1:B:417:MET:HB3	2.44	0.57
1:A:242:THR:HG23	1:B:231:ILE:HG12	1.86	0.57
1:A:14[A]:CYS:SG	1:A:53:PHE:CD2	2.98	0.57
1:D:143:SER:C	1:D:145:GLU:H	2.13	0.56
1:B:306:ILE:HG22	1:B:407:TYR:CZ	2.41	0.55
1:B:186:MET:HE2	1:B:251[B]:ILE:HD11	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:230:SER:HB3	1:C:260:GLN:HE22	1.71	0.55
1:A:401:MET:HB2	1:A:402:PRO:HD3	1.89	0.55
1:D:199:HIS:ND1	1:D:255:HIS:HE1	2.04	0.54
1:D:56:PRO:HG3	3:D:4642:HOH:O	2.08	0.54
1:D:149:LEU:HB2	3:D:4716:HOH:O	2.08	0.54
1:B:198:LEU:HD12	1:B:251[A]:ILE:CD1	2.35	0.53
1:C:149:LEU:HB3	3:C:3734:HOH:O	2.08	0.53
1:B:25:ALA:HB3	1:B:33:LEU:HB3	1.91	0.53
1:D:145:GLU:HA	1:D:145:GLU:OE1	2.07	0.53
1:B:313:GLU:HA	1:B:318:ASN:ND2	2.24	0.52
1:C:59:ILE:HG12	1:C:93:THR:HB	1.91	0.52
1:D:110:LEU:HB3	1:D:134:LEU:HD21	1.92	0.52
1:B:254:VAL:HG23	1:B:255:HIS:CE1	2.45	0.52
1:B:260:GLN:NE2	1:B:260:GLN:H	2.08	0.51
1:C:260:GLN:HE21	1:C:260:GLN:N	1.96	0.51
1:C:202:ASN:C	1:C:202:ASN:HD22	2.20	0.50
1:B:13:ILE:HD12	1:B:22:ALA:HB3	1.94	0.49
1:A:387:PHE:HA	1:A:390:ILE:HD12	1.94	0.49
1:A:440:SER:HB3	3:C:3658:HOH:O	2.10	0.49
1:C:406:ASP:OD1	1:C:489:LYS:NZ	2.34	0.49
1:C:231:ILE:HG12	1:D:242:THR:HG23	1.93	0.49
1:D:146:ALA:O	1:D:150:LEU:HB2	2.12	0.49
1:C:401:MET:HB2	1:C:402:PRO:HD3	1.95	0.48
1:C:6:LEU:HB3	1:C:26:VAL:HB	1.96	0.48
1:A:147:ARG:HD2	1:A:187:TYR:CE2	2.48	0.48
1:C:71:LEU:HD11	1:C:175:LEU:HD21	1.94	0.48
1:D:327:LYS:O	1:D:394:LEU:HD22	2.14	0.48
1:A:321:ASP:OD2	1:A:364:TYR:OH	2.25	0.48
1:B:401:MET:HA	1:B:424:GLN:HE22	1.79	0.47
1:B:39:ASP:OD2	1:B:40:PRO:HD2	2.14	0.47
1:C:260:GLN:H	1:C:260:GLN:NE2	1.95	0.47
1:D:47:ILE:HD11	1:D:501:ILE:HD11	1.96	0.47
1:C:290:ILE:HD13	1:C:308:LEU:HD22	1.96	0.47
1:C:419:LYS:O	1:C:423:ILE:HG12	2.14	0.47
1:D:139:ILE:HG13	1:D:181:ASP:HB3	1.95	0.47
1:B:387:PHE:HA	1:B:390:ILE:HD12	1.96	0.47
1:D:9:LYS:NZ	1:D:38:ILE:O	2.46	0.47
1:D:80:GLU:HG3	1:D:125:THR:O	2.14	0.47
1:D:379:PRO:C	1:D:381:ASN:H	2.22	0.47
1:D:14[A]:CYS:SG	1:D:53:PHE:CD2	3.08	0.47
1:B:84:ARG:NH1	1:B:508:ILE:HG21	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:401:MET:HB2	1:D:402:PRO:HD3	1.97	0.47
1:D:59:ILE:HG12	1:D:93:THR:HB	1.96	0.46
1:B:87:VAL:HG11	1:B:502:VAL:HG12	1.98	0.46
1:D:187:TYR:CE1	1:D:538:PRO:HB3	2.50	0.46
1:C:69:LEU:HD11	1:C:101:ASP:HB2	1.98	0.46
1:A:199:HIS:ND1	1:A:255:HIS:HE1	2.13	0.46
1:C:537:ARG:HB2	3:C:3708:HOH:O	2.16	0.46
1:A:166:VAL:HG23	1:A:194:PHE:HD2	1.81	0.45
1:A:260:GLN:NE2	1:A:260:GLN:H	2.14	0.45
1:B:244:ALA:HB2	1:B:251[B]:ILE:HG13	1.97	0.45
1:B:306:ILE:HD11	1:B:311:ILE:HD11	1.99	0.45
1:A:387:PHE:HA	1:A:390:ILE:CD1	2.47	0.44
1:B:166:VAL:HG23	1:B:194:PHE:HD1	1.81	0.44
1:D:393:GLY:O	1:D:394:LEU:HD23	2.16	0.44
1:C:8:ILE:HD13	1:C:452:ILE:HD11	2.00	0.44
1:D:530:ASN:ND2	3:D:4671:HOH:O	2.50	0.44
1:A:74:ASP:O	1:A:375:ARG:NH2	2.51	0.44
1:A:139:ILE:HG13	1:A:181:ASP:HB3	2.00	0.44
1:B:6:LEU:HB3	1:B:26:VAL:HB	1.99	0.44
1:B:260:GLN:H	1:B:260:GLN:HE21	1.64	0.44
1:C:331:SER:HA	1:C:332:PRO:C	2.41	0.44
1:B:401:MET:HA	1:B:424:GLN:NE2	2.33	0.44
1:D:202:ASN:O	1:D:206:VAL:HG23	2.18	0.44
1:A:508:ILE:HG22	3:A:1661:HOH:O	2.16	0.44
1:A:509:LEU:HD23	3:A:1790:HOH:O	2.17	0.44
1:A:110:LEU:HB3	1:A:134:LEU:HD21	2.00	0.44
1:B:136:LEU:HD22	1:B:166:VAL:HG11	2.00	0.44
1:B:221:ALA:HB1	1:B:329:ILE:HD11	2.00	0.44
1:B:401:MET:HE3	1:B:424:GLN:CB	2.45	0.43
1:B:79:MET:HG3	1:B:126:LEU:HD22	1.99	0.43
1:B:321:ASP:OD2	1:B:364:TYR:OH	2.31	0.43
1:C:25:ALA:HB3	1:C:33:LEU:HB3	2.00	0.43
1:D:85:SER:OG	1:D:398:CYS:HB2	2.18	0.43
1:C:230:SER:HB3	1:C:260:GLN:NE2	2.33	0.43
1:B:71:LEU:HD11	1:B:175:LEU:HD21	2.00	0.43
1:B:306:ILE:HD12	1:B:486:ILE:HG21	2.01	0.43
1:C:254:VAL:HG23	1:C:255:HIS:CE1	2.53	0.43
1:C:491:TRP:CG	1:C:492:PRO:HD2	2.54	0.43
1:B:143:SER:O	1:B:147:ARG:CG	2.66	0.43
1:B:139:ILE:HG13	1:B:181:ASP:HB3	2.00	0.43
1:C:14:CYS:SG	1:C:53:PHE:CD2	3.12	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:94:VAL:O	1:C:130:TYR:HA	2.19	0.43
1:A:147:ARG:HD2	1:A:187:TYR:CD2	2.54	0.43
1:A:328:TYR:HA	1:A:394:LEU:HD21	2.00	0.43
1:B:387:PHE:HA	1:B:390:ILE:CD1	2.49	0.43
1:B:463:ASN:OD1	1:B:463:ASN:N	2.52	0.43
1:D:258:SER:OG	1:D:260:GLN:HG2	2.19	0.42
1:C:146:ALA:O	1:C:150:LEU:HB2	2.19	0.42
1:A:3:ILE:HG22	1:A:42:LEU:O	2.19	0.42
1:D:6:LEU:HB3	1:D:26:VAL:HB	2.02	0.42
1:A:285:LEU:HB3	1:A:306:ILE:HD12	2.00	0.42
1:D:285:LEU:HD22	1:D:306:ILE:HG13	2.01	0.42
1:B:36:ALA:HB2	1:D:444:GLY:HA3	2.01	0.41
1:B:222:TYR:CD1	1:B:292:ARG:HG2	2.55	0.41
1:D:92:THR:HG22	1:D:497:VAL:HG22	2.02	0.41
1:A:94:VAL:O	1:A:130:TYR:HA	2.20	0.41
1:A:143:SER:O	1:A:147:ARG:HG3	2.19	0.41
1:A:360:CYS:HB3	1:A:393:GLY:HA3	2.02	0.41
1:B:176:GLN:OE1	1:B:203:GLY:HA3	2.19	0.41
1:D:26:VAL:HG11	1:D:499:GLY:HA2	2.01	0.41
1:A:47:ILE:HG21	1:A:452:ILE:HD13	2.02	0.41
1:A:373:LYS:HD3	1:A:391:PRO:HG2	2.03	0.41
1:C:67:GLU:HB2	1:C:68:PRO:HD2	2.01	0.41
1:D:254:VAL:HG23	1:D:255:HIS:CE1	2.56	0.41
1:A:25:ALA:HB3	1:A:33:LEU:HB3	2.02	0.41
1:D:25:ALA:HB3	1:D:33:LEU:HB3	2.02	0.41
1:D:172:TYR:HA	1:D:173:PRO:HD3	1.95	0.41
1:A:13:ILE:HD12	1:A:22:ALA:HB3	2.02	0.41
1:C:26:VAL:HG11	1:C:499:GLY:HA2	2.02	0.41
1:B:278:GLU:HG3	1:B:354:ILE:HG13	2.02	0.41
1:C:373:LYS:HA	1:C:391:PRO:HG2	2.03	0.40
1:D:263:GLU:O	1:D:267:GLN:HG3	2.22	0.40
1:D:285:LEU:HB3	1:D:306:ILE:HD12	2.03	0.40
1:A:254:VAL:HG23	1:A:255:HIS:CE1	2.56	0.40
1:C:326:SER:HA	1:C:329:ILE:HG12	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	529/559 (95%)	510 (96%)	15 (3%)	4 (1%)	16	25
1	B	530/559 (95%)	512 (97%)	16 (3%)	2 (0%)	30	43
1	C	526/559 (94%)	506 (96%)	18 (3%)	2 (0%)	30	43
1	D	527/559 (94%)	505 (96%)	19 (4%)	3 (1%)	21	32
All	All	2112/2236 (94%)	2033 (96%)	68 (3%)	11 (0%)	30	37

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	144	VAL
1	A	172[A]	TYR
1	A	172[B]	TYR
1	A	255	HIS
1	B	172[A]	TYR
1	B	172[B]	TYR
1	C	172	TYR
1	D	172	TYR
1	D	255	HIS
1	C	255	HIS
1	A	396	GLY

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	457/479 (95%)	443 (97%)	14 (3%)	35	57
1	B	458/479 (96%)	441 (96%)	17 (4%)	30	51
1	C	454/479 (95%)	437 (96%)	17 (4%)	30	51
1	D	456/479 (95%)	441 (97%)	15 (3%)	33	55
All	All	1825/1916 (95%)	1762 (96%)	63 (4%)	32	53

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	GLU
1	A	70	LYS
1	A	87	VAL
1	A	104	LYS
1	A	148	GLU
1	A	260	GLN
1	A	312	SER
1	A	334	ILE
1	A	404	LEU
1	A	415	THR
1	A	431	VAL
1	A	461	GLU
1	A	463	ASN
1	A	500	LYS
1	B	45	GLU
1	B	65	VAL
1	B	76	VAL
1	B	123	GLU
1	B	136	LEU
1	B	148	GLU
1	B	152	VAL
1	B	260	GLN
1	B	292	ARG
1	B	404	LEU
1	B	415	THR
1	B	431	VAL
1	B	457	ASP
1	B	461	GLU
1	B	463	ASN
1	B	500	LYS
1	B	511	GLU
1	C	23	GLU
1	C	37	SER

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Mol	Chain	Res	Type
1	C	65	VAL
1	C	136	LEU
1	C	139	ILE
1	C	149	LEU
1	C	202	ASN
1	C	260	GLN
1	C	293	CYS
1	C	320	ASP
1	C	337	GLU
1	C	404	LEU
1	C	419	LYS
1	C	481	THR
1	C	500	LYS
1	C	529	LYS
1	C	531	GLU
1	D	32	GLN
1	D	70	LYS
1	D	139	ILE
1	D	149	LEU
1	D	260	GLN
1	D	291	THR
1	D	312	SER
1	D	320	ASP
1	D	340	GLN
1	D	341	LYS
1	D	366	GLU
1	D	404	LEU
1	D	481	THR
1	D	489	LYS
1	D	531	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	GLN
1	A	153	GLN
1	A	216	GLN
1	A	260	GLN
1	A	267	GLN
1	A	294	HIS
1	A	424	GLN
1	B	192	ASN

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Mol	Chain	Res	Type
1	B	260	GLN
1	B	318	ASN
1	B	392	ASN
1	B	424	GLN
1	C	27	ASN
1	C	202	ASN
1	C	238	ASN
1	C	260	GLN
1	C	374	HIS
1	C	437	GLN
1	D	267	GLN
1	D	340	GLN
1	D	530	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	KCX	A	167	1,2	10,11,12	0.96	0	6,12,14	1.37	1 (16%)
1	KCX	D	167	1,2	10,11,12	0.93	0	6,12,14	1.68	1 (16%)
1	KCX	B	167	1,2	10,11,12	0.96	0	6,12,14	1.76	2 (33%)
1	KCX	C	167	1,2	10,11,12	0.87	0	6,12,14	1.22	1 (16%)

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
1	KCX	A	167	1,2	-	1/9/10/12	-
1	KCX	D	167	1,2	-	0/9/10/12	-
1	KCX	B	167	1,2	-	1/9/10/12	-
1	KCX	C	167	1,2	-	1/9/10/12	-

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	D	167	KCX	OQ1-CX-NZ	-3.90	118.99	124.92
1	B	167	KCX	OQ1-CX-NZ	-3.63	119.40	124.92
1	A	167	KCX	OQ1-CX-NZ	-3.21	120.04	124.92
1	C	167	KCX	OQ1-CX-NZ	-2.87	120.56	124.92
1	B	167	KCX	CE-NZ-CX	-2.11	118.40	121.98

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	167	KCX	C-CA-CB-CG
1	C	167	KCX	C-CA-CB-CG
1	A	167	KCX	C-CA-CB-CG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	531/559 (94%)	0.75	17 (3%)	50 46	21, 36, 38, 45	2 (0%)
1	B	531/559 (94%)	0.71	22 (4%)	41 37	22, 36, 38, 42	3 (0%)
1	C	530/559 (94%)	0.87	32 (6%)	27 24	31, 36, 38, 42	0
1	D	529/559 (94%)	0.80	22 (4%)	40 36	24, 36, 38, 42	2 (0%)
All	All	2121/2236 (94%)	0.78	93 (4%)	39 34	21, 36, 38, 45	7 (0%)

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	146	ALA	8.0
1	C	371	ALA	4.5
1	C	385	GLY	4.1
1	C	304	VAL	4.0
1	C	457	ASP	3.9
1	B	172[A]	TYR	3.9
1	C	380	GLU	3.8
1	A	320	ASP	3.7
1	B	306	ILE	3.4
1	B	304	VAL	3.3
1	B	457	ASP	3.3
1	A	457	ASP	3.2
1	C	328	TYR	3.2
1	C	293	CYS	3.2
1	C	123	GLU	3.1
1	B	174	GLY	3.1
1	D	128	CYS	3.0
1	A	303	GLY	3.0
1	B	280	CYS	3.0
1	D	293	CYS	3.0
1	D	294	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	148	GLU	2.9
1	D	127	TYR	2.8
1	C	426	THR	2.8
1	C	448	ALA	2.8
1	B	290	ILE	2.8
1	A	304	VAL	2.7
1	D	320	ASP	2.7
1	B	303	GLY	2.7
1	B	316	PHE	2.7
1	A	436	PRO	2.7
1	C	160	ASP	2.6
1	C	320	ASP	2.6
1	C	36	ALA	2.6
1	C	65	VAL	2.6
1	B	294	HIS	2.6
1	D	291	THR	2.5
1	C	74	ASP	2.5
1	D	129	ASP	2.5
1	C	330	CYS	2.5
1	C	461	GLU	2.5
1	C	536	TRP	2.5
1	B	307	ASP	2.4
1	C	365	TYR	2.4
1	D	131	GLY	2.4
1	A	270	THR	2.4
1	A	22	ALA	2.4
1	D	148	GLU	2.4
1	C	172	TYR	2.4
1	A	275	VAL	2.4
1	A	173	PRO	2.3
1	C	321	ASP	2.3
1	C	382	ASN	2.3
1	D	371	ALA	2.3
1	B	293	CYS	2.3
1	A	448	ALA	2.3
1	C	277	ALA	2.3
1	D	312	SER	2.3
1	C	148	GLU	2.3
1	D	172	TYR	2.3
1	C	149	LEU	2.2
1	A	431	VAL	2.2
1	D	396	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	3	ILE	2.2
1	B	390	ILE	2.2
1	B	320	ASP	2.2
1	D	220	ASP	2.2
1	D	318	ASN	2.2
1	D	133	HIS	2.2
1	D	255	HIS	2.2
1	B	368	THR	2.2
1	B	289	ALA	2.2
1	B	222	TYR	2.2
1	D	74	ASP	2.2
1	C	44	SER	2.1
1	D	305	GLY	2.1
1	B	232	VAL	2.1
1	D	382	ASN	2.1
1	C	140	GLU	2.1
1	D	457	ASP	2.1
1	A	294	HIS	2.1
1	A	2	PRO	2.1
1	C	144	VAL	2.0
1	C	420	LEU	2.0
1	A	43	GLY	2.0
1	B	426	THR	2.0
1	B	435	TYR	2.0
1	A	483	PHE	2.0
1	B	134	LEU	2.0
1	C	220	ASP	2.0
1	C	436	PRO	2.0
1	B	276	TYR	2.0
1	C	312	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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. 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	B-factors(Å ²)	Q<0.9
1	KCX	D	167	12/13	0.84	0.15	35,36,44,44	3
1	KCX	C	167	12/13	0.89	0.15	35,36,44,44	3
1	KCX	A	167	12/13	0.89	0.13	35,36,44,45	3

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
1	KCX	B	167	12/13	0.91	0.10	35,35,44,44	3

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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. 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	B-factors(\AA^2)	Q<0.9
2	ZN	D	602	1/1	0.95	0.21	46,46,46,46	1
2	ZN	D	601	1/1	0.96	0.12	44,44,44,44	1
2	ZN	B	602	1/1	0.97	0.18	46,46,46,46	1
2	ZN	C	602	1/1	0.97	0.15	45,45,45,45	1
2	ZN	A	602	1/1	0.98	0.16	45,45,45,45	1
2	ZN	B	601	1/1	0.99	0.13	43,43,43,43	1
2	ZN	A	601	1/1	0.99	0.15	45,45,45,45	1
2	ZN	C	601	1/1	0.99	0.11	44,44,44,44	1

6.5 Other polymers [i](#)

There are no such residues in this entry.