



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:20 PM GMT

PDB ID : 1OFQ
Title : CRYSTAL STRUCTURE OF THE TYROSINE-REGULATED 3-DEOXY-D-ARABINO-HEPTULOSONATE-7-PHOSPHATE SYNTHASE FROM SACCCHAROMYCES CEREVISIAE IN COMPLEX WITH MANGANESE(II)
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Deposited on : 2003-04-17
Resolution : 2.70 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

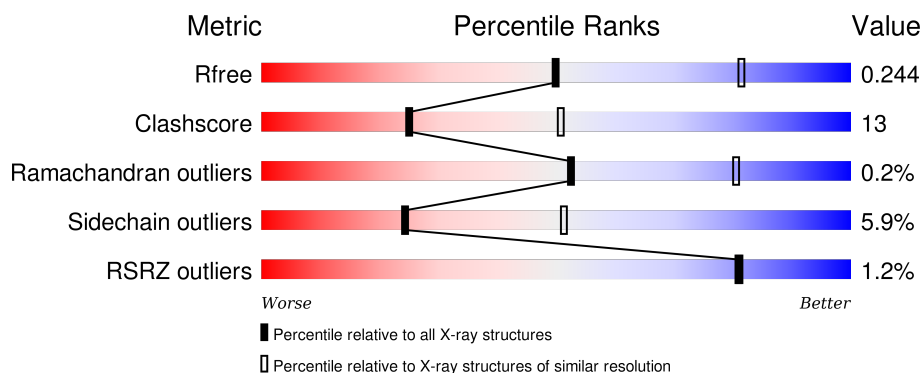
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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	370	<div> <div>%</div> <div> <div></div> <div>70%</div> <div>21%</div> <div>• 7%</div> </div> </div>
1	B	370	<div> <div>2%</div> <div> <div></div> <div>70%</div> <div>20%</div> <div>• 7%</div> </div> </div>
1	C	370	<div> <div>%</div> <div> <div></div> <div>72%</div> <div>19%</div> <div>• 7%</div> </div> </div>
1	D	370	<div> <div>%</div> <div> <div></div> <div>68%</div> <div>22%</div> <div>• 6%</div> </div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 10572 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 PHOSPHO-2-DEHYDRO-3-DEOXYHEPTONATE ALDOLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	343	Total	C	N	O	S	0	0	0
			2562	1595	460	497	10			
1	B	343	Total	C	N	O	S	0	0	0
			2565	1597	457	501	10			
1	C	344	Total	C	N	O	S	0	0	0
			2573	1601	464	498	10			
1	D	346	Total	C	N	O	S	0	0	0
			2583	1609	460	504	10			

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mn	0	0
			1	1		
2	A	1	Total	Mn	0	0
			1	1		
2	D	1	Total	Mn	0	0
			1	1		
2	C	1	Total	Mn	0	0
			1	1		

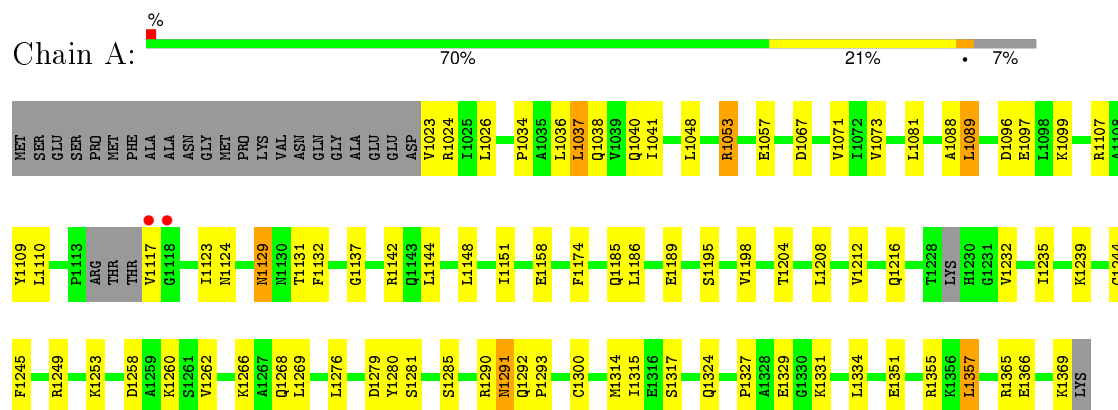
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	76	Total	O	0	0
			76	76		
3	B	57	Total	O	0	0
			57	57		
3	C	65	Total	O	0	0
			65	65		
3	D	87	Total	O	0	0
			87	87		

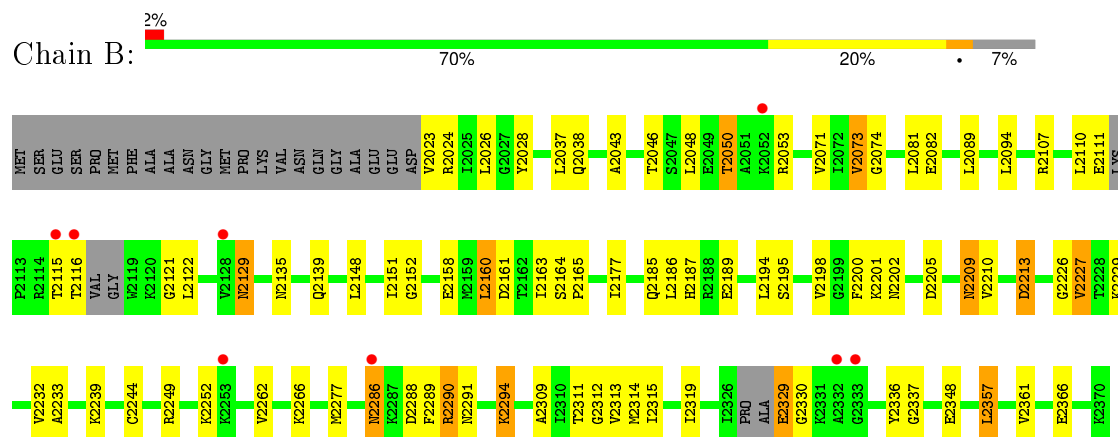
3 Residue-property plots

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.

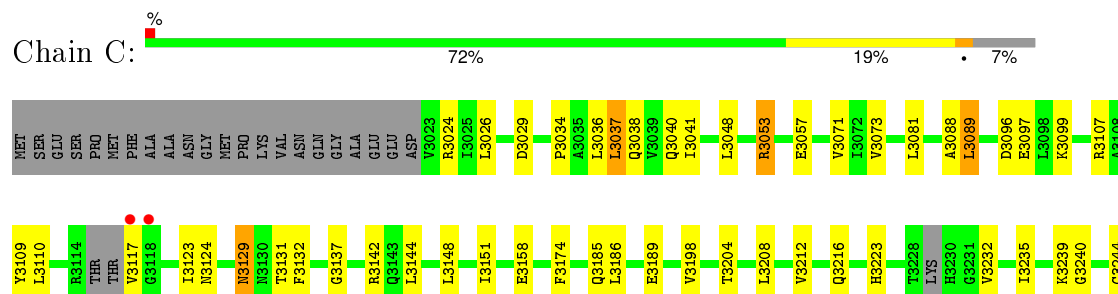
• Molecule 1: PHOSPHO-2-DEHYDRO-3-DEOXYHEPTONATE ALDOLASE

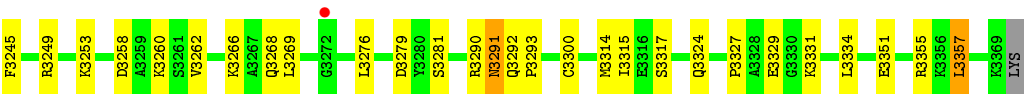


• Molecule 1: PHOSPHO-2-DEHYDRO-3-DEOXYHEPTONATE ALDOLASE

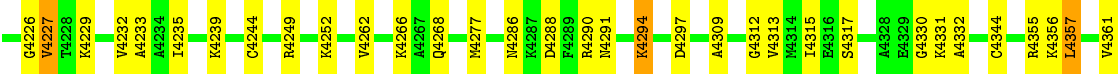
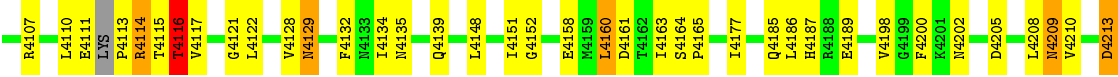
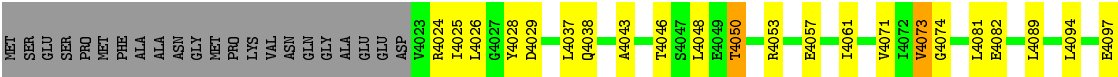


• Molecule 1: PHOSPHO-2-DEHYDRO-3-DEOXYHEPTONATE ALDOLASE





● Molecule 1: PHOSPHO-2-DEHYDRO-3-DEOXYHEPTONATE ALDOLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	105.73Å 105.73Å 268.78Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.98 – 2.70 19.98 – 2.40	Depositor EDS
% Data completeness (in resolution range)	96.6 (19.98-2.70) 95.2 (19.98-2.40)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	11.99 (at 2.41Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.221 , 0.245 0.220 , 0.244	Depositor DCC
R_{free} test set	1375 reflections (2.93%)	DCC
Wilson B-factor (Å ²)	46.1	Xtriage
Anisotropy	0.242	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 31.4	EDS
Estimated twinning fraction	0.019 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 68894 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10572	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% 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: MN

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	0.34	0/2599	0.78	6/3521 (0.2%)
1	B	0.35	0/2600	0.60	0/3520
1	C	0.35	0/2610	0.74	6/3535 (0.2%)
1	D	0.34	0/2621	0.60	0/3553
All	All	0.34	0/10430	0.69	12/14129 (0.1%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1290	ARG	NE-CZ-NH1	17.03	128.82	120.30
1	A	1290	ARG	NE-CZ-NH2	-14.31	113.14	120.30
1	C	3053	ARG	NE-CZ-NH2	-14.12	113.24	120.30
1	C	3053	ARG	NE-CZ-NH1	13.70	127.15	120.30
1	A	1053	ARG	NE-CZ-NH1	-13.62	113.49	120.30
1	C	3290	ARG	NE-CZ-NH2	12.74	126.67	120.30
1	A	1053	ARG	NE-CZ-NH2	12.51	126.56	120.30
1	C	3290	ARG	NE-CZ-NH1	-9.96	115.32	120.30
1	A	1290	ARG	CD-NE-CZ	8.21	135.09	123.60
1	C	3290	ARG	CD-NE-CZ	7.16	133.62	123.60
1	C	3053	ARG	CD-NE-CZ	6.98	133.37	123.60
1	A	1053	ARG	CD-NE-CZ	6.25	132.35	123.60

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	2562	0	2546	62	0
1	B	2565	0	2544	79	0
1	C	2573	0	2559	60	0
1	D	2583	0	2568	85	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	76	0	0	2	0
3	B	57	0	0	9	0
3	C	65	0	0	1	0
3	D	87	0	0	3	0
All	All	10572	0	10217	260	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 (260) 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:2160:LEU:H	1:B:2160:LEU:HD12	1.33	0.94
1:D:4160:LEU:H	1:D:4160:LEU:HD12	1.34	0.93
1:D:4071:VAL:HG12	1:D:4073:VAL:HG22	1.51	0.93
1:B:2071:VAL:HG12	1:B:2073:VAL:HG22	1.51	0.92
1:B:2209:ASN:HD22	1:B:2210:VAL:N	1.70	0.90
1:D:4209:ASN:HD22	1:D:4210:VAL:N	1.71	0.88
1:D:4115:THR:O	1:D:4116:THR:HB	1.76	0.85
1:D:4116:THR:HG22	1:D:4117:VAL:HG23	1.60	0.84
1:C:3071:VAL:HG12	1:C:3073:VAL:HG22	1.61	0.82
1:B:2038:GLN:HE22	1:B:2229:LYS:NZ	1.77	0.82
1:C:3208:LEU:H	1:C:3268:GLN:NE2	1.78	0.82
1:D:4038:GLN:HE22	1:D:4229:LYS:NZ	1.78	0.81
1:A:1071:VAL:HG12	1:A:1073:VAL:HG22	1.62	0.81
1:A:1208:LEU:H	1:A:1268:GLN:NE2	1.78	0.80
1:C:3239:LYS:HB3	1:D:4026:LEU:HD11	1.66	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3037:LEU:HD12	1:C:3142:ARG:HD3	1.69	0.75
1:B:2319:ILE:HB	1:B:2337:GLY:HA3	1.67	0.74
1:C:3291:ASN:HD22	1:C:3291:ASN:N	1.87	0.73
1:A:1037:LEU:HD12	1:A:1142:ARG:HD3	1.70	0.73
1:C:3239:LYS:HB3	1:D:4026:LEU:CD1	2.19	0.72
1:A:1291:ASN:N	1:A:1291:ASN:HD22	1.87	0.71
1:B:2073:VAL:HG13	1:B:2315:ILE:HB	1.73	0.71
1:C:3185:GLN:O	1:C:3189:GLU:HG3	1.91	0.69
1:A:1026:LEU:HD21	1:B:2239:LYS:HB3	1.75	0.69
1:D:4073:VAL:HG13	1:D:4315:ILE:HB	1.74	0.69
1:D:4097:GLU:OE1	1:D:4355:ARG:NH1	2.25	0.69
1:C:3253:LYS:HG2	3:C:2046:HOH:O	1.92	0.69
1:A:1185:GLN:O	1:A:1189:GLU:HG3	1.93	0.68
1:D:4163:ILE:HG12	1:D:4227:VAL:HG21	1.75	0.67
1:D:4160:LEU:N	1:D:4160:LEU:HD12	2.08	0.67
1:B:2160:LEU:N	1:B:2160:LEU:HD12	2.08	0.67
1:D:4038:GLN:HE22	1:D:4229:LYS:HZ2	1.42	0.67
1:B:2329:GLU:N	1:B:2329:GLU:OE1	2.29	0.66
1:B:2163:ILE:HG12	1:B:2227:VAL:HG21	1.76	0.66
1:C:3097:GLU:OE1	1:C:3355:ARG:NH1	2.28	0.66
1:A:1097:GLU:OE1	1:A:1355:ARG:NH1	2.29	0.66
1:B:2038:GLN:HE22	1:B:2229:LYS:HZ2	1.42	0.65
1:C:3174:PHE:HE2	1:C:3245:PHE:HZ	1.44	0.65
1:A:1081:LEU:HD22	1:A:1144:LEU:HD13	1.78	0.64
1:B:2111:GLU:HG2	1:B:2121:GLY:HA3	1.78	0.64
1:A:1174:PHE:HE2	1:A:1245:PHE:HZ	1.44	0.63
1:D:4111:GLU:HG2	1:D:4121:GLY:HA3	1.80	0.63
1:D:4111:GLU:C	1:D:4113:PRO:HD3	2.19	0.63
1:A:1262:VAL:O	1:A:1266:LYS:HG3	1.99	0.63
1:B:2262:VAL:O	1:B:2266:LYS:HG3	1.98	0.63
1:D:4050:THR:HG23	1:D:4053:ARG:HH12	1.63	0.63
1:C:3081:LEU:HD22	1:C:3144:LEU:HD13	1.81	0.63
1:B:2185:GLN:O	1:B:2189:GLU:HG3	1.99	0.63
1:B:2110:LEU:HD22	1:B:2110:LEU:N	2.15	0.62
1:C:3262:VAL:O	1:C:3266:LYS:HG3	1.98	0.62
1:A:1124:ASN:HD22	1:A:1132:PHE:HE2	1.48	0.62
1:C:3186:LEU:HD21	1:D:4161:ASP:HB3	1.80	0.62
1:B:2050:THR:HG23	1:B:2053:ARG:HH12	1.64	0.62
1:D:4110:LEU:N	1:D:4110:LEU:HD22	2.15	0.61
1:A:1208:LEU:HB2	1:A:1268:GLN:HE21	1.66	0.61
1:B:2160:LEU:CD1	1:B:2160:LEU:H	2.09	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:4160:LEU:H	1:D:4160:LEU:CD1	2.10	0.60
1:C:3208:LEU:HB2	1:C:3268:GLN:HE21	1.66	0.60
1:C:3124:ASN:HD22	1:C:3132:PHE:HE2	1.49	0.60
1:A:1034:PRO:O	1:A:1038:GLN:HG3	2.00	0.60
1:D:4037:LEU:HD11	1:D:4139:GLN:HG2	1.83	0.60
1:D:4262:VAL:O	1:D:4266:LYS:HG3	2.01	0.60
1:D:4177:ILE:HG13	1:D:4200:PHE:CE2	2.37	0.60
1:C:3208:LEU:H	1:C:3268:GLN:HE21	1.50	0.60
1:B:2037:LEU:HD11	1:B:2139:GLN:HG2	1.84	0.59
1:A:1081:LEU:CD2	1:A:1144:LEU:HD13	2.32	0.59
1:B:2110:LEU:HD23	3:B:3021:HOH:O	2.03	0.59
1:C:3071:VAL:HG12	1:C:3073:VAL:CG2	2.32	0.59
1:D:4185:GLN:O	1:D:4189:GLU:HG3	2.02	0.59
1:B:2357:LEU:O	1:B:2361:VAL:HG23	2.02	0.59
1:B:2177:ILE:HG13	1:B:2200:PHE:CE2	2.37	0.59
1:A:1208:LEU:H	1:A:1268:GLN:HE21	1.50	0.59
1:C:3331:LYS:HA	1:C:3334:LEU:HD23	1.85	0.59
1:C:3355:ARG:HG2	1:C:3355:ARG:HH11	1.68	0.59
1:B:2071:VAL:CG1	1:B:2073:VAL:HG22	2.29	0.58
1:A:1067:ASP:OD2	1:A:1365:ARG:HD3	2.04	0.58
1:C:3269:LEU:HD11	1:C:3276:LEU:HD21	1.86	0.58
1:D:4073:VAL:CG1	1:D:4315:ILE:HB	2.34	0.58
1:C:3081:LEU:CD2	1:C:3144:LEU:HD13	2.33	0.58
1:B:2073:VAL:CG1	1:B:2315:ILE:HB	2.34	0.57
1:D:4071:VAL:CG1	1:D:4073:VAL:HG22	2.30	0.57
1:A:1266:LYS:HE2	3:A:2051:HOH:O	2.03	0.57
1:C:3026:LEU:HD21	1:D:4239:LYS:HB3	1.87	0.57
1:A:1269:LEU:HD11	1:A:1276:LEU:HD21	1.86	0.57
1:B:2232:VAL:HG21	1:D:4232:VAL:HG21	1.86	0.57
1:D:4291:ASN:HD22	1:D:4294:LYS:HD3	1.69	0.56
1:C:3073:VAL:HG13	1:C:3315:ILE:HB	1.88	0.56
1:C:3024:ARG:HA	1:D:4239:LYS:HG2	1.88	0.56
1:A:1331:LYS:HA	1:A:1334:LEU:HD23	1.86	0.56
1:A:1026:LEU:CD2	1:B:2239:LYS:HB3	2.34	0.56
1:C:3235:ILE:HB	1:D:4029:ASP:HB2	1.87	0.56
1:A:1355:ARG:HG2	1:A:1355:ARG:HH11	1.69	0.56
1:B:2201:LYS:NZ	3:B:3031:HOH:O	2.39	0.56
1:A:1239:LYS:HB3	1:B:2026:LEU:CD1	2.35	0.56
1:D:4288:ASP:OD1	1:D:4290:ARG:HG3	2.05	0.56
1:B:2209:ASN:HD22	1:B:2210:VAL:H	1.51	0.56
1:C:3034:PRO:O	1:C:3038:GLN:HG3	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2291:ASN:HD22	1:B:2294:LYS:HD3	1.72	0.55
1:D:4266:LYS:HE2	3:D:2066:HOH:O	2.06	0.55
1:D:4369:LYS:HE2	1:D:4369:LYS:HA	1.88	0.55
1:A:1071:VAL:HG12	1:A:1073:VAL:CG2	2.34	0.54
1:C:3239:LYS:HG2	1:D:4024:ARG:HA	1.88	0.54
1:D:4074:GLY:HA3	1:D:4107:ARG:HB2	1.90	0.54
1:C:3327:PRO:HB2	1:C:3329:GLU:OE1	2.07	0.54
1:C:3300:CYS:SG	1:C:3357:LEU:HA	2.48	0.54
1:C:3129:ASN:HD22	1:C:3129:ASN:C	2.10	0.54
1:A:1129:ASN:C	1:A:1129:ASN:HD22	2.12	0.53
1:A:1327:PRO:HB2	1:A:1329:GLU:OE1	2.08	0.53
1:A:1073:VAL:HG13	1:A:1315:ILE:HB	1.89	0.53
1:B:2209:ASN:C	1:B:2209:ASN:HD22	2.12	0.53
1:A:1097:GLU:CD	1:A:1355:ARG:HH12	2.12	0.53
1:B:2074:GLY:HA3	1:B:2107:ARG:HB2	1.90	0.53
1:A:1239:LYS:HB3	1:B:2026:LEU:HD13	1.91	0.53
1:A:1300:CYS:SG	1:A:1357:LEU:HA	2.49	0.53
1:B:2288:ASP:OD1	1:B:2290:ARG:HG3	2.08	0.53
1:A:1024:ARG:HD2	1:B:2194:LEU:O	2.10	0.52
1:A:1186:LEU:HD21	1:B:2161:ASP:HB3	1.91	0.52
1:D:4135:ASN:O	1:D:4139:GLN:HG3	2.08	0.52
1:B:2314:MET:HB2	3:B:3045:HOH:O	2.08	0.52
1:B:2209:ASN:O	1:B:2213:ASP:HB2	2.09	0.52
1:C:3036:LEU:O	1:C:3040:GLN:HG3	2.09	0.52
1:C:3258:ASP:OD2	1:C:3260:LYS:HB3	2.10	0.51
1:D:4209:ASN:O	1:D:4213:ASP:HB2	2.09	0.51
1:B:2348:GLU:HG2	3:B:3054:HOH:O	2.09	0.51
1:C:3097:GLU:CD	1:C:3355:ARG:HH12	2.13	0.51
1:A:1096:ASP:HA	1:A:1099:LYS:HG3	1.92	0.51
1:B:2135:ASN:O	1:B:2139:GLN:HG3	2.10	0.51
1:D:4209:ASN:HD22	1:D:4210:VAL:H	1.51	0.51
1:D:4050:THR:HG21	1:D:4152:GLY:C	2.31	0.51
1:B:2038:GLN:HE22	1:B:2229:LYS:HZ3	1.55	0.51
1:C:3096:ASP:HA	1:C:3099:LYS:HG3	1.91	0.51
1:D:4050:THR:HA	1:D:4053:ARG:NH1	2.26	0.51
1:B:2050:THR:HA	1:B:2053:ARG:NH1	2.26	0.51
1:B:2038:GLN:NE2	1:B:2229:LYS:NZ	2.55	0.51
1:B:2314:MET:HE3	3:B:3045:HOH:O	2.11	0.50
1:A:1036:LEU:O	1:A:1040:GLN:HG3	2.11	0.50
1:C:3089:LEU:HD13	1:C:3151:ILE:HD13	1.92	0.50
1:D:4209:ASN:HD22	1:D:4209:ASN:C	2.14	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:4038:GLN:NE2	1:D:4229:LYS:NZ	2.56	0.50
1:A:1258:ASP:OD2	1:A:1260:LYS:HB3	2.10	0.50
1:B:2160:LEU:HD11	3:B:3021:HOH:O	2.12	0.50
1:C:3291:ASN:N	1:C:3291:ASN:ND2	2.59	0.50
1:B:2266:LYS:HD2	1:B:2309:ALA:CB	2.42	0.50
1:C:3235:ILE:HD13	1:D:4134:ILE:HB	1.94	0.50
1:D:4198:VAL:O	1:D:4244:CYS:HA	2.12	0.50
1:A:1253:LYS:HG2	3:A:2047:HOH:O	2.11	0.50
1:A:1195:SER:OG	1:B:2023:VAL:HG22	2.11	0.50
1:A:1351:GLU:O	1:A:1355:ARG:HG3	2.12	0.50
1:D:4205:ASP:HA	1:D:4252:LYS:HB2	1.94	0.49
1:B:2311:THR:HA	3:B:3008:HOH:O	2.12	0.49
1:A:1292:GLN:N	1:A:1293:PRO:HD2	2.27	0.49
1:A:1089:LEU:HD13	1:A:1151:ILE:HD13	1.93	0.49
1:A:1208:LEU:H	1:A:1268:GLN:HE22	1.60	0.49
1:B:2050:THR:HG21	1:B:2152:GLY:C	2.33	0.49
1:B:2129:ASN:C	1:B:2129:ASN:HD22	2.16	0.49
1:A:1198:VAL:O	1:A:1244:CYS:HA	2.12	0.48
1:B:2050:THR:HA	1:B:2053:ARG:HH12	1.79	0.48
1:C:3081:LEU:CD2	1:C:3144:LEU:HB2	2.43	0.48
1:C:3088:ALA:HB2	1:C:3148:LEU:HD22	1.94	0.48
1:D:4160:LEU:HD11	3:D:2032:HOH:O	2.12	0.48
1:A:1088:ALA:HB2	1:A:1148:LEU:HD22	1.96	0.48
1:B:2205:ASP:HA	1:B:2252:LYS:HB2	1.95	0.48
1:C:3351:GLU:O	1:C:3355:ARG:HG3	2.13	0.48
1:A:1081:LEU:CD2	1:A:1144:LEU:HB2	2.43	0.48
1:D:4129:ASN:HD22	1:D:4129:ASN:C	2.15	0.48
1:C:3037:LEU:HD12	1:C:3142:ARG:CD	2.43	0.48
1:C:3292:GLN:N	1:C:3293:PRO:HD2	2.28	0.48
1:A:1129:ASN:ND2	1:A:1131:THR:H	2.12	0.47
1:D:4110:LEU:HD23	3:D:2032:HOH:O	2.14	0.47
1:D:4038:GLN:HE22	1:D:4229:LYS:HZ3	1.57	0.47
1:B:2198:VAL:O	1:B:2244:CYS:HA	2.15	0.47
1:D:4050:THR:HA	1:D:4053:ARG:HH12	1.78	0.47
1:D:4266:LYS:HD2	1:D:4309:ALA:CB	2.44	0.47
1:B:2289:PHE:HA	3:B:3047:HOH:O	2.13	0.47
1:C:3208:LEU:H	1:C:3268:GLN:HE22	1.60	0.47
1:C:3129:ASN:ND2	1:C:3131:THR:H	2.12	0.47
1:A:1186:LEU:HD12	1:B:2187:HIS:HE1	1.79	0.47
1:A:1235:ILE:O	1:B:2028:TYR:HA	2.15	0.47
1:A:1037:LEU:HD11	1:A:1041:ILE:HD12	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3198:VAL:O	1:C:3244:CYS:HA	2.14	0.46
1:C:3223:HIS:CD2	1:D:4132:PHE:HB3	2.51	0.46
1:D:4113:PRO:O	1:D:4114:ARG:C	2.54	0.46
1:A:1258:ASP:O	1:A:1262:VAL:HG23	2.16	0.46
1:D:4313:VAL:HG21	1:D:4357:LEU:HD11	1.98	0.46
1:B:2313:VAL:HG21	1:B:2357:LEU:HD11	1.98	0.46
1:B:2232:VAL:HG21	1:D:4232:VAL:CG2	2.46	0.46
1:B:2330:GLY:HA3	3:B:3018:HOH:O	2.16	0.46
1:C:3029:ASP:HB2	1:D:4235:ILE:HB	1.98	0.46
1:D:4288:ASP:OD1	1:D:4290:ARG:CG	2.64	0.45
1:D:4046:THR:O	1:D:4050:THR:OG1	2.33	0.45
1:D:4128:VAL:HB	1:D:4331:LYS:HG3	1.99	0.45
1:A:1109:TYR:CE2	1:A:1158:GLU:HB2	2.51	0.45
1:B:2348:GLU:H	1:B:2348:GLU:CD	2.19	0.45
1:B:2082:GLU:CD	1:B:2082:GLU:H	2.20	0.45
1:A:1053:ARG:O	1:A:1057:GLU:HG3	2.16	0.45
1:D:4369:LYS:HA	1:D:4369:LYS:CE	2.46	0.45
1:D:4128:VAL:HB	1:D:4331:LYS:CG	2.47	0.45
1:B:2043:ALA:HB3	1:B:2048:LEU:HD11	1.99	0.44
1:D:4053:ARG:NH1	1:D:4053:ARG:HB3	2.33	0.44
1:D:4291:ASN:HA	1:D:4294:LYS:HD3	2.00	0.44
1:C:3037:LEU:HD11	1:C:3041:ILE:HD12	2.00	0.44
1:A:1291:ASN:N	1:A:1291:ASN:ND2	2.59	0.44
1:B:2115:THR:O	1:B:2116:THR:C	2.55	0.44
1:A:1023:VAL:HG12	1:B:2195:SER:OG	2.18	0.44
1:C:3053:ARG:O	1:C:3057:GLU:HG3	2.17	0.44
1:A:1123:ILE:O	1:A:1137:GLY:HA3	2.17	0.44
1:D:4050:THR:CG2	1:D:4053:ARG:HH12	2.29	0.44
1:C:3279:ASP:HA	1:C:3314:MET:HB2	1.98	0.44
1:B:2286:ASN:HA	1:B:2286:ASN:HD22	1.66	0.44
1:C:3239:LYS:HD3	1:D:4024:ARG:HG2	2.00	0.44
1:D:4291:ASN:HD22	1:D:4291:ASN:HA	1.66	0.43
1:D:4082:GLU:CD	1:D:4082:GLU:H	2.19	0.43
1:B:2158:GLU:O	1:B:2160:LEU:HD12	2.18	0.43
1:C:3258:ASP:O	1:C:3262:VAL:HG23	2.17	0.43
1:C:3240:GLY:O	1:D:4024:ARG:HD3	2.19	0.43
1:B:2046:THR:O	1:B:2050:THR:OG1	2.36	0.43
1:D:4074:GLY:CA	1:D:4107:ARG:HB2	2.49	0.43
1:B:2288:ASP:OD1	1:B:2290:ARG:CG	2.67	0.43
1:A:1366:GLU:HA	1:A:1369:LYS:HG2	2.01	0.43
1:B:2110:LEU:HD12	1:B:2122:LEU:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:4110:LEU:HD12	1:D:4122:LEU:HB3	1.99	0.43
1:B:2053:ARG:NH1	1:B:2053:ARG:HB3	2.33	0.43
1:A:1212:VAL:O	1:A:1216:GLN:HG3	2.19	0.43
1:A:1037:LEU:HD12	1:A:1142:ARG:CD	2.43	0.43
1:C:3186:LEU:HD12	1:D:4187:HIS:HE1	1.84	0.43
1:B:2291:ASN:HA	1:B:2294:LYS:HD3	2.00	0.42
1:C:3123:ILE:O	1:C:3137:GLY:HA3	2.18	0.42
1:C:3129:ASN:ND2	1:C:3129:ASN:C	2.73	0.42
1:A:1279:ASP:HA	1:A:1314:MET:HB2	1.99	0.42
1:D:4129:ASN:ND2	1:D:4129:ASN:C	2.72	0.42
1:B:2148:LEU:O	1:B:2151:ILE:HG12	2.20	0.42
1:D:4317:SER:HA	1:D:4344:CYS:HB3	2.01	0.42
1:A:1324:GLN:CD	1:A:1334:LEU:HD12	2.40	0.42
1:C:3109:TYR:CE2	1:C:3158:GLU:HB2	2.54	0.42
1:B:2129:ASN:C	1:B:2129:ASN:ND2	2.73	0.42
1:D:4148:LEU:O	1:D:4151:ILE:HG12	2.20	0.42
1:D:4297:ASP:OD1	1:D:4356:LYS:HE3	2.19	0.41
1:D:4043:ALA:HB3	1:D:4048:LEU:HD11	2.03	0.41
1:C:3324:GLN:CD	1:C:3334:LEU:HD12	2.41	0.41
1:A:1239:LYS:HG2	1:B:2024:ARG:HA	2.02	0.41
1:B:2164:SER:OG	1:B:2165:PRO:HD3	2.20	0.41
1:A:1280:TYR:O	1:A:1285:SER:HB3	2.20	0.41
1:B:2074:GLY:CA	1:B:2107:ARG:HB2	2.49	0.41
1:A:1148:LEU:O	1:A:1151:ILE:HG12	2.21	0.41
1:D:4164:SER:OG	1:D:4165:PRO:HD3	2.21	0.41
1:D:4208:LEU:HB2	1:D:4268:GLN:HE22	1.86	0.41
1:B:2277:MET:HA	1:B:2312:GLY:O	2.21	0.41
1:D:4226:GLY:O	1:D:4233:ALA:HA	2.21	0.41
1:D:4277:MET:HA	1:D:4312:GLY:O	2.21	0.41
1:C:3081:LEU:HD21	1:C:3144:LEU:HB2	2.03	0.41
1:D:4357:LEU:O	1:D:4361:VAL:HG23	2.21	0.41
1:B:2115:THR:HG22	1:B:2115:THR:O	2.21	0.41
1:C:3212:VAL:O	1:C:3216:GLN:HG3	2.21	0.41
1:A:1129:ASN:C	1:A:1129:ASN:ND2	2.74	0.40
1:B:2226:GLY:O	1:B:2233:ALA:HA	2.21	0.40
1:D:4025:ILE:HG21	1:D:4028:TYR:CZ	2.56	0.40
1:D:4330:GLY:C	1:D:4332:ALA:N	2.75	0.40
1:D:4057:GLU:O	1:D:4061:ILE:HG13	2.21	0.40
1:D:4158:GLU:O	1:D:4160:LEU:HD12	2.22	0.40
1:B:2110:LEU:CD2	1:B:2110:LEU:N	2.83	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	337/370 (91%)	323 (96%)	14 (4%)	0	100	100
1	B	335/370 (90%)	318 (95%)	16 (5%)	1 (0%)	46	75
1	C	338/370 (91%)	324 (96%)	14 (4%)	0	100	100
1	D	342/370 (92%)	329 (96%)	11 (3%)	2 (1%)	30	59
All	All	1352/1480 (91%)	1294 (96%)	55 (4%)	3 (0%)	52	80

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	2336	TYR
1	D	4116	THR
1	D	4114	ARG

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	270/298 (91%)	256 (95%)	14 (5%)	29	58
1	B	271/298 (91%)	252 (93%)	19 (7%)	19	42
1	C	271/298 (91%)	257 (95%)	14 (5%)	29	58
1	D	273/298 (92%)	256 (94%)	17 (6%)	23	49
All	All	1085/1192 (91%)	1021 (94%)	64 (6%)	24	51

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1037	LEU
1	A	1048	LEU
1	A	1089	LEU
1	A	1107	ARG
1	A	1110	LEU
1	A	1117	VAL
1	A	1129	ASN
1	A	1204	THR
1	A	1232	VAL
1	A	1249	ARG
1	A	1281	SER
1	A	1291	ASN
1	A	1317	SER
1	A	1357	LEU
1	B	2050	THR
1	B	2073	VAL
1	B	2081	LEU
1	B	2089	LEU
1	B	2094	LEU
1	B	2129	ASN
1	B	2160	LEU
1	B	2186	LEU
1	B	2202	ASN
1	B	2209	ASN
1	B	2213	ASP
1	B	2227	VAL
1	B	2249	ARG
1	B	2286	ASN
1	B	2290	ARG
1	B	2294	LYS
1	B	2329	GLU
1	B	2357	LEU
1	B	2366	GLU
1	C	3037	LEU
1	C	3048	LEU
1	C	3089	LEU
1	C	3107	ARG
1	C	3110	LEU
1	C	3117	VAL
1	C	3129	ASN
1	C	3204	THR
1	C	3232	VAL
1	C	3249	ARG

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Mol	Chain	Res	Type
1	C	3281	SER
1	C	3291	ASN
1	C	3317	SER
1	C	3357	LEU
1	D	4050	THR
1	D	4073	VAL
1	D	4081	LEU
1	D	4089	LEU
1	D	4094	LEU
1	D	4116	THR
1	D	4129	ASN
1	D	4160	LEU
1	D	4186	LEU
1	D	4202	ASN
1	D	4209	ASN
1	D	4213	ASP
1	D	4227	VAL
1	D	4249	ARG
1	D	4286	ASN
1	D	4294	LYS
1	D	4357	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1124	ASN
1	A	1129	ASN
1	A	1150	ASN
1	A	1268	GLN
1	A	1291	ASN
1	B	2038	GLN
1	B	2129	ASN
1	B	2150	ASN
1	B	2209	ASN
1	B	2220	HIS
1	B	2268	GLN
1	B	2286	ASN
1	B	2291	ASN
1	C	3124	ASN
1	C	3129	ASN
1	C	3150	ASN
1	C	3268	GLN

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Mol	Chain	Res	Type
1	C	3291	ASN
1	C	3308	ASN
1	C	3368	ASN
1	D	4038	GLN
1	D	4129	ASN
1	D	4150	ASN
1	D	4209	ASN
1	D	4220	HIS
1	D	4268	GLN
1	D	4286	ASN
1	D	4291	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](#)

Of 4 ligands modelled in this entry, 4 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

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	343/370 (92%)	-0.45	2 (0%) 90 91	34, 45, 59, 72	0
1	B	343/370 (92%)	-0.34	8 (2%) 64 64	32, 45, 66, 99	0
1	C	344/370 (92%)	-0.46	3 (0%) 85 86	33, 45, 59, 73	0
1	D	346/370 (93%)	-0.40	3 (0%) 85 86	31, 44, 62, 72	0
All	All	1376/1480 (92%)	-0.41	16 (1%) 81 81	31, 45, 62, 99	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2116	THR	5.1
1	D	4115	THR	4.1
1	C	3117	VAL	3.5
1	B	2332	ALA	3.0
1	B	2115	THR	3.0
1	B	2128	VAL	2.8
1	D	4328	ALA	2.7
1	D	4116	THR	2.7
1	A	1117	VAL	2.5
1	B	2286	ASN	2.5
1	B	2253	LYS	2.4
1	A	1118	GLY	2.3
1	C	3272	GLY	2.2
1	B	2333	GLY	2.2
1	C	3118	GLY	2.1
1	B	2052	LYS	2.0

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

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(\AA^2)	Q<0.9
2	MN	C	3000	1/1	0.97	0.06	-1.75	39,39,39,39	0
2	MN	A	1000	1/1	0.99	0.02	-1.75	38,38,38,38	0
2	MN	B	2000	1/1	0.97	0.06	-2.69	52,52,52,52	0
2	MN	D	4000	1/1	0.99	0.04	-	42,42,42,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.