Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2023 – 07:17 PM EDT

PDB ID	:	1PU0
Title	:	Structure of Human Cu,Zn Superoxide Dismutase
Authors	:	DiDonato, M.; Craig, L.; Huff, M.E.; Thayer, M.M.; Cardoso, R.M.F.; Kassmann, C.J.; Lo, T.P.; Bruns, C.K.; Powers, E.T.; Kelly, J.W.; Getzoff, E.D.; Tainer, J.A.
Deposited on	:	2003-06-23
Resolution	:	1.70 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
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:

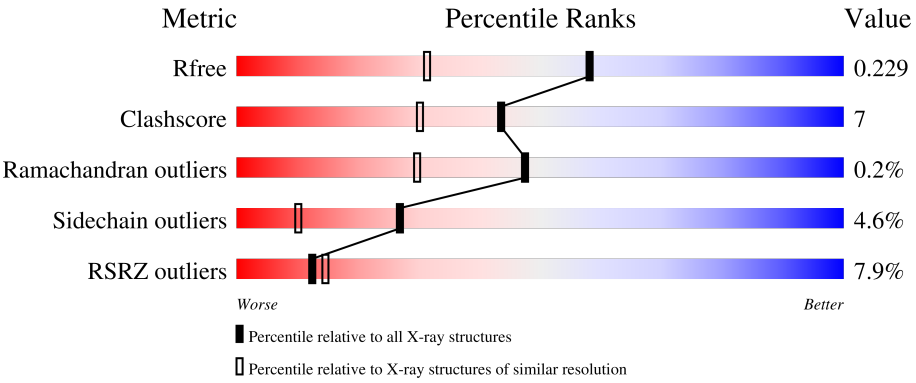
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	130704	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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 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 <=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	153	4% 92% 8%
1	B	153	4% 90% 7%
1	C	153	14% 73% 22% 5%
1	D	153	4% 89% 10%
1	E	153	3% 92% 7%

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Mol	Chain	Length	Quality of chain
1	F	153	6% 90% 9%
1	G	153	13% 76% 20%
1	H	153	5% 89% 10%
1	I	153	16% 83% 14%
1	J	153	10% 87% 12%

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
4	SO4	J	302	-	-	-	X

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 12222 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 Superoxide dismutase [Cu-Zn].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	153	Total	C	N	O	S	0	2	0
			1114	681	203	224	6			
1	B	153	Total	C	N	O	S	0	2	0
			1114	681	203	224	6			
1	C	153	Total	C	N	O	S	0	2	0
			1114	681	203	224	6			
1	D	153	Total	C	N	O	S	0	2	0
			1114	681	203	224	6			
1	E	153	Total	C	N	O	S	0	2	0
			1114	681	203	224	6			
1	F	153	Total	C	N	O	S	0	2	0
			1114	681	203	224	6			
1	G	153	Total	C	N	O	S	0	2	0
			1114	681	203	224	6			
1	H	153	Total	C	N	O	S	0	2	0
			1114	681	203	224	6			
1	I	153	Total	C	N	O	S	0	2	0
			1114	681	203	224	6			
1	J	153	Total	C	N	O	S	0	2	0
			1114	681	203	224	6			

- Molecule 2 is COPPER (I) ION (three-letter code: CU1) (formula: Cu).

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

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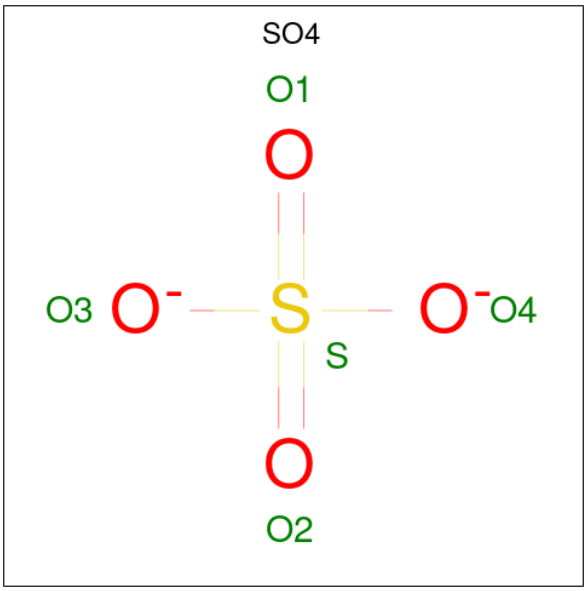
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	E	1	Total Cu 1 1	0	0
2	F	1	Total Cu 1 1	0	0
2	G	1	Total Cu 1 1	0	0
2	H	1	Total Cu 1 1	0	0
2	I	1	Total Cu 1 1	0	0
2	J	1	Total Cu 1 1	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Zn 1 1	0	0
3	B	1	Total Zn 1 1	0	0
3	C	1	Total Zn 1 1	0	0
3	D	1	Total Zn 1 1	0	0
3	E	1	Total Zn 1 1	0	0
3	F	1	Total Zn 1 1	0	0
3	G	1	Total Zn 1 1	0	0
3	H	1	Total Zn 1 1	0	0
3	I	1	Total Zn 1 1	0	0
3	J	1	Total Zn 1 1	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	131	Total O 131 131	0	0
5	B	141	Total O 141 141	0	0
5	C	57	Total O 57 57	0	0
5	D	133	Total O 133 133	0	0
5	E	133	Total O 133 133	0	0
5	F	126	Total O 126 126	0	0
5	G	55	Total O 55 55	0	0
5	H	107	Total O 107 107	0	0
5	I	60	Total O 60 60	0	0

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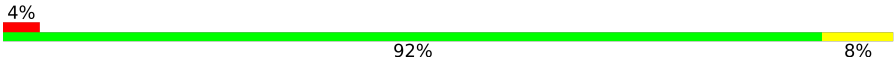
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	J	104	Total 104	O 104	0	0

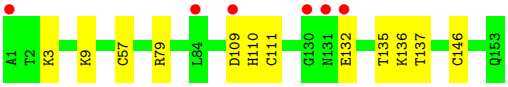
3 Residue-property plots (i)

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: Superoxide dismutase [Cu-Zn]

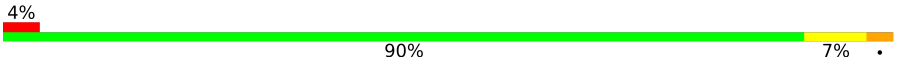
Chain A:

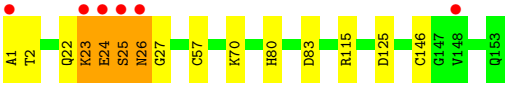




• Molecule 1: Superoxide dismutase [Cu-Zn]

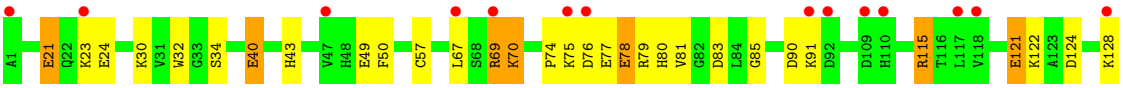
Chain B:





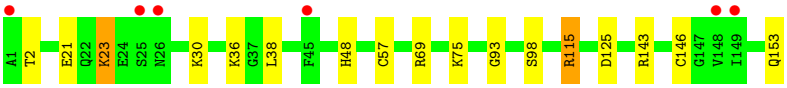
• Molecule 1: Superoxide dismutase [Cu-Zn]

Chain C:



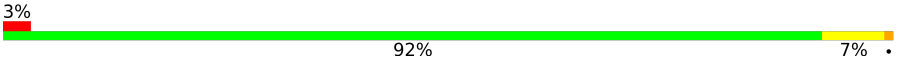
• Molecule 1: Superoxide dismutase [Cu-Zn]

Chain D:



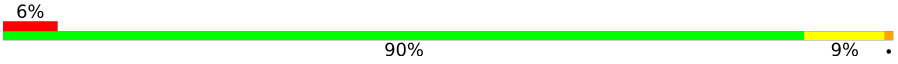
• Molecule 1: Superoxide dismutase [Cu-Zn]

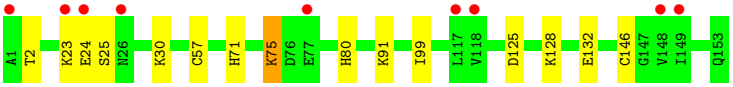
Chain E:



• Molecule 1: Superoxide dismutase [Cu-Zn]

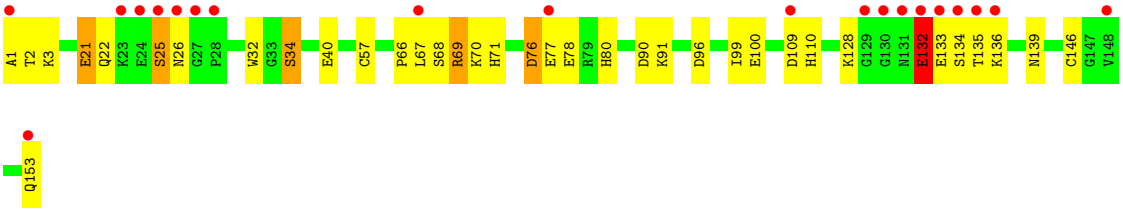
Chain F:





• Molecule 1: Superoxide dismutase [Cu-Zn]

Chain G:

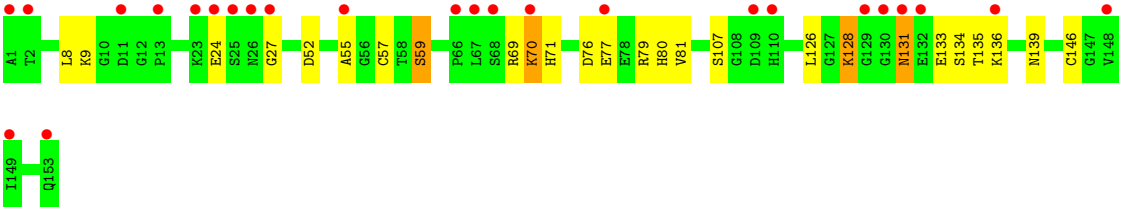


• Molecule 1: Superoxide dismutase [Cu-Zn]

Chain H:

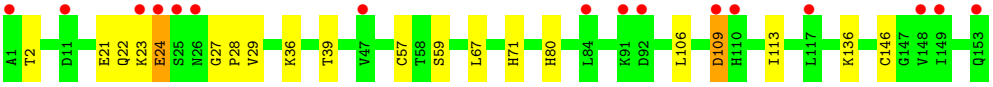
• Molecule 1: Superoxide dismutase [Cu-Zn]

Chain I:



• Molecule 1: Superoxide dismutase [Cu-Zn]

Chain J:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	203.58Å 165.68Å 144.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.70 27.61 – 1.70	Depositor EDS
% Data completeness (in resolution range)	93.6 (30.00-1.70) 97.6 (27.61-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	3.99 (at 1.71Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.210 , 0.243 0.196 , 0.229	Depositor DCC
R_{free} test set	13096 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	23.8	Xtriage
Anisotropy	0.137	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 55.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	12222	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.59% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: ZN, SO4, CU1

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	0.44	0/1140	1.03	1/1536 (0.1%)
1	B	0.45	0/1140	1.03	3/1536 (0.2%)
1	C	0.42	0/1140	1.13	4/1536 (0.3%)
1	D	0.43	0/1140	1.10	5/1536 (0.3%)
1	E	0.43	0/1140	1.11	3/1536 (0.2%)
1	F	0.44	0/1140	1.02	1/1536 (0.1%)
1	G	0.38	0/1140	1.03	0/1536
1	H	0.40	0/1140	1.12	4/1536 (0.3%)
1	I	0.35	0/1140	0.97	0/1536
1	J	0.38	0/1140	0.99	0/1536
All	All	0.41	0/11400	1.05	21/15360 (0.1%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	H	79	ARG	NE-CZ-NH1	-10.11	115.24	120.30
1	B	115	ARG	NE-CZ-NH2	9.96	125.28	120.30
1	E	115	ARG	NE-CZ-NH2	8.39	124.50	120.30
1	D	115	ARG	NE-CZ-NH2	7.42	124.01	120.30
1	H	143	ARG	NE-CZ-NH2	7.21	123.91	120.30
1	E	115	ARG	NE-CZ-NH1	-7.08	116.76	120.30
1	D	115	ARG	NE-CZ-NH1	-6.73	116.93	120.30
1	H	79	ARG	NE-CZ-NH2	6.46	123.53	120.30
1	D	125	ASP	CB-CG-OD2	6.17	123.85	118.30
1	E	115	ARG	CD-NE-CZ	5.90	131.86	123.60
1	C	115	ARG	NE-CZ-NH2	-5.83	117.39	120.30
1	C	79	ARG	NE-CZ-NH2	5.83	123.21	120.30
1	B	115	ARG	NE-CZ-NH1	-5.75	117.43	120.30
1	D	69	ARG	NE-CZ-NH2	5.55	123.08	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	143	ARG	NE-CZ-NH1	-5.50	117.55	120.30
1	C	124	ASP	O-C-N	5.27	131.13	122.70
1	B	125	ASP	CB-CG-OD2	5.24	123.01	118.30
1	A	79	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	F	125	ASP	CB-CG-OD2	5.13	122.92	118.30
1	H	143	ARG	NE-CZ-NH1	-5.05	117.77	120.30
1	C	43	HIS	O-C-N	5.01	131.71	123.20

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	1114	0	1079	6	0
1	B	1114	0	1079	8	0
1	C	1114	0	1079	28	0
1	D	1114	0	1079	13	0
1	E	1114	0	1079	11	0
1	F	1114	0	1079	9	0
1	G	1114	0	1079	25	0
1	H	1114	0	1079	10	0
1	I	1114	0	1079	21	0
1	J	1114	0	1079	13	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
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
3	I	1	0	0	0	0
3	J	1	0	0	0	0
4	B	5	0	0	0	0
4	J	10	0	0	1	0
5	A	131	0	0	2	0
5	B	141	0	0	2	0
5	C	57	0	0	1	0
5	D	133	0	0	4	0
5	E	133	0	0	4	0
5	F	126	0	0	0	0
5	G	55	0	0	2	0
5	H	107	0	0	0	0
5	I	60	0	0	2	0
5	J	104	0	0	1	0
All	All	12222	0	10790	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 7.

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:B:57[B]:CYS:HG	1:B:146[B]:CYS:HG	0.88	0.85
1:G:57[B]:CYS:HG	1:G:146[B]:CYS:HG	0.85	0.83
1:H:57[B]:CYS:HG	1:H:146[B]:CYS:HG	0.95	0.81
1:J:57[B]:CYS:HG	1:J:146[B]:CYS:HG	0.88	0.81
1:J:24:GLU:OE1	1:J:28:PRO:HD2	1.84	0.78
1:A:57[B]:CYS:HG	1:A:146[B]:CYS:HG	0.89	0.75
1:D:75:LYS:HE3	5:D:519:HOH:O	1.85	0.75
1:C:121:GLU:HB2	1:C:144:LEU:HD21	1.70	0.72
1:C:70:LYS:HE3	1:C:135:THR:HG21	1.70	0.71
1:I:131:ASN:ND2	1:I:134:SER:H	1.89	0.71
1:A:136:LYS:HD3	1:A:137:THR:HG23	1.74	0.69
1:C:40:GLU:OE2	1:C:91:LYS:HA	1.93	0.68
1:H:108:GLY:O	1:H:111:CYS:HB3	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:131:ASN:HD21	1:I:139:ASN:HD21	1.44	0.66
5:B:563:HOH:O	1:F:75:LYS:HE3	1.94	0.66
1:I:81:VAL:HG23	5:I:1359:HOH:O	1.96	0.65
1:D:2:THR:HG22	5:D:1488:HOH:O	1.97	0.64
1:F:23:LYS:HE3	1:F:30:LYS:HE3	1.79	0.64
1:C:67:LEU:HD22	1:C:69:ARG:NH1	2.11	0.63
1:G:3:LYS:HG2	1:G:21:GLU:HG2	1.80	0.62
1:C:40:GLU:HB2	1:C:90:ASP:O	2.01	0.61
1:J:21:GLU:OE1	1:J:23:LYS:HE3	2.01	0.61
1:E:23:LYS:HZ2	1:E:30:LYS:HD3	1.64	0.61
1:J:24:GLU:OE1	1:J:27:GLY:HA3	2.02	0.60
1:C:57[B]:CYS:HG	1:C:146[B]:CYS:CB	2.15	0.59
1:H:30:LYS:HD3	1:H:100:GLU:HG3	1.83	0.59
1:C:23:LYS:HE2	1:C:24:GLU:HG2	1.85	0.59
1:G:153:GLN:HB2	1:H:50:PHE:CZ	2.38	0.58
1:G:76:ASP:OD1	1:G:128:LYS:HE2	2.04	0.57
1:A:57[B]:CYS:HG	1:A:146[B]:CYS:CB	2.18	0.57
1:H:57[B]:CYS:HG	1:H:146[B]:CYS:CB	2.18	0.57
1:E:100:GLU:HG2	5:E:814:HOH:O	2.05	0.56
1:I:24:GLU:HB2	1:I:27:GLY:HA3	1.87	0.56
1:D:23:LYS:HD3	1:D:23:LYS:N	2.20	0.56
1:D:36:LYS:HE3	5:D:1367:HOH:O	2.06	0.56
1:B:1:ALA:HB3	1:B:22:GLN:O	2.05	0.56
1:E:57[B]:CYS:HG	1:E:146[B]:CYS:CB	2.18	0.56
1:E:21:GLU:OE2	1:E:23:LYS:HE2	2.07	0.55
1:D:23:LYS:HE2	5:D:1281:HOH:O	2.05	0.55
1:I:131:ASN:O	1:I:135:THR:HG23	2.07	0.55
1:I:57[B]:CYS:HG	1:I:146[B]:CYS:HG	0.87	0.54
1:A:9:LYS:HG2	5:A:1378:HOH:O	2.07	0.53
1:E:23:LYS:HD2	5:E:698:HOH:O	2.09	0.52
1:C:49:GLU:O	1:C:115:ARG:HD3	2.10	0.52
1:G:133:GLU:HA	1:G:133:GLU:OE1	2.09	0.52
1:C:70:LYS:HD2	1:C:135:THR:HG22	1.92	0.52
1:C:132:GLU:HG3	1:C:133:GLU:N	2.24	0.52
1:G:57[B]:CYS:HG	1:G:146[B]:CYS:CB	2.22	0.52
1:G:69:ARG:CZ	1:G:78:GLU:HB2	2.40	0.51
1:C:69:ARG:HB3	1:C:78:GLU:OE1	2.10	0.51
1:C:21:GLU:HB2	1:C:32:TRP:CZ3	2.45	0.51
1:D:57[B]:CYS:HG	1:D:146[B]:CYS:CB	2.22	0.51
1:G:1:ALA:HA	1:G:22:GLN:O	2.10	0.51
1:C:81:VAL:HG23	5:C:1347:HOH:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:22:GLN:HB2	1:J:29:VAL:HG22	1.93	0.50
1:E:23:LYS:HE3	5:E:698:HOH:O	2.11	0.50
1:C:70:LYS:HB3	1:C:135:THR:HG22	1.92	0.50
1:C:74:PRO:HD3	1:C:85:GLY:CA	2.42	0.50
1:I:131:ASN:ND2	1:I:134:SER:HB2	2.27	0.50
1:I:80:HIS:HE1	1:I:136:LYS:O	1.94	0.50
1:I:57[B]:CYS:HG	1:I:146[B]:CYS:CB	2.21	0.49
1:J:57[B]:CYS:HG	1:J:146[B]:CYS:CB	2.23	0.49
1:G:40:GLU:HG3	1:G:90:ASP:C	2.33	0.49
1:I:126:LEU:O	1:I:128:LYS:HD2	2.12	0.49
1:G:69:ARG:NH1	1:G:78:GLU:HB2	2.27	0.49
1:C:131:ASN:OD1	1:C:134:SER:HB2	2.13	0.49
1:E:3:LYS:HE2	1:E:21:GLU:OE1	2.13	0.49
1:A:132:GLU:OE2	1:I:131:ASN:HB3	2.13	0.48
1:H:105:SER:O	1:H:111:CYS:HA	2.14	0.48
1:D:98:SER:HB2	1:F:132:GLU:HG2	1.94	0.47
1:I:70:LYS:HE3	1:I:70:LYS:HA	1.96	0.47
1:I:55:ALA:HB3	1:I:59:SER:OG	2.14	0.47
1:F:24:GLU:HA	1:F:24:GLU:OE2	2.14	0.47
1:J:136:LYS:HE3	5:J:1040:HOH:O	2.13	0.47
1:G:21:GLU:HB2	1:G:32:TRP:CZ3	2.50	0.46
1:C:21:GLU:HB2	1:C:32:TRP:HZ3	1.79	0.46
1:G:132:GLU:HG2	1:G:132:GLU:O	2.15	0.46
1:I:131:ASN:HD22	1:I:134:SER:H	1.62	0.46
1:C:129:GLY:HA3	1:C:134:SER:HB2	1.96	0.46
1:G:110:HIS:HB2	5:G:1404:HOH:O	2.16	0.46
1:C:50:PHE:CZ	1:D:153:GLN:HB2	2.50	0.46
1:G:99:ILE:HD11	5:G:1235:HOH:O	2.16	0.46
1:G:69:ARG:NH2	1:G:78:GLU:OE1	2.50	0.45
1:I:8:LEU:O	1:I:9:LYS:HG3	2.16	0.45
1:B:57[B]:CYS:CB	1:B:146[B]:CYS:HG	2.27	0.45
1:G:70:LYS:HD2	1:G:135:THR:HB	1.98	0.45
1:I:69:ARG:HG3	1:I:69:ARG:HH21	1.82	0.45
1:G:3:LYS:HG2	1:G:21:GLU:CG	2.47	0.45
1:B:24:GLU:OE1	1:B:27:GLY:N	2.50	0.45
1:B:25:SER:OG	1:B:26:ASN:ND2	2.50	0.45
1:J:21:GLU:CD	1:J:23:LYS:HZ1	2.20	0.45
1:J:109:ASP:N	1:J:109:ASP:OD1	2.50	0.45
1:C:76:ASP:OD1	1:C:128:LYS:NZ	2.50	0.45
1:C:76:ASP:OD2	1:C:76:ASP:N	2.50	0.45
1:C:122:LYS:NZ	1:C:139:ASN:O	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:109:ASP:N	5:E:806:HOH:O	2.50	0.44
1:G:71:HIS:HB2	1:G:80:HIS:CE1	2.52	0.44
1:G:132:GLU:HG2	1:G:136:LYS:NZ	2.32	0.44
1:G:134:SER:HB2	1:G:139:ASN:ND2	2.33	0.44
1:C:121:GLU:HA	1:C:144:LEU:HD11	1.99	0.44
1:J:71:HIS:HB2	1:J:80:HIS:CE1	2.52	0.44
1:C:75:LYS:HB2	1:H:153:GLN:OE1	2.17	0.44
1:J:67:LEU:N	1:J:67:LEU:HD23	2.33	0.44
1:D:38:LEU:O	1:D:93:GLY:HA2	2.17	0.44
1:G:134:SER:HB2	1:G:139:ASN:HD21	1.83	0.43
1:D:23:LYS:HZ1	1:D:30:LYS:HD2	1.83	0.43
1:E:69:ARG:HD3	1:E:78:GLU:OE2	2.18	0.43
1:I:52:ASP:O	1:I:59:SER:HB2	2.18	0.43
1:G:66:PRO:C	1:G:67:LEU:HD23	2.39	0.43
1:H:71:HIS:HB2	1:H:80:HIS:CE1	2.54	0.43
1:E:23:LYS:NZ	1:E:30:LYS:HD3	2.33	0.43
1:I:76:ASP:O	1:I:79:ARG:HG2	2.19	0.43
1:D:75:LYS:HE2	1:F:128:LYS:HG2	1.99	0.43
1:C:80:HIS:HB2	1:C:83:ASP:CG	2.39	0.43
1:E:21:GLU:OE2	1:E:30:LYS:NZ	2.52	0.43
1:H:26:ASN:OD1	1:H:26:ASN:N	2.50	0.43
1:F:57[B]:CYS:HG	1:F:146[B]:CYS:CB	2.28	0.42
1:C:132:GLU:HA	1:C:135:THR:HG1	1.85	0.42
1:J:39:THR:HG22	4:J:302:SO4:O1	2.19	0.42
1:G:34:SER:HB3	1:G:96:ASP:OD1	2.20	0.42
1:A:110:HIS:HD2	5:A:1311:HOH:O	2.03	0.42
1:B:23:LYS:HG2	5:B:719:HOH:O	2.19	0.42
1:G:132:GLU:HG2	1:G:136:LYS:HZ2	1.85	0.41
1:I:133:GLU:HA	1:I:136:LYS:HE3	2.01	0.41
1:D:21:GLU:CD	1:D:23:LYS:HZ2	2.22	0.41
1:J:106:LEU:HD22	1:J:113:ILE:HD11	2.03	0.41
1:C:74:PRO:HD3	1:C:85:GLY:HA3	2.03	0.41
1:C:129:GLY:N	1:C:134:SER:OG	2.51	0.41
1:F:99:ILE:HD13	1:F:99:ILE:HG21	1.89	0.41
1:B:24:GLU:OE1	1:B:26:ASN:N	2.49	0.41
1:H:22:GLN:HG2	1:H:24:GLU:O	2.21	0.41
1:G:2:THR:HG22	1:G:3:LYS:N	2.36	0.41
1:B:80:HIS:HB2	1:B:83:ASP:CG	2.42	0.40
1:I:27:GLY:HA3	5:I:1418:HOH:O	2.20	0.40
1:F:71:HIS:HB2	1:F:80:HIS:CE1	2.56	0.40
1:I:71:HIS:HB2	1:I:80:HIS:CE1	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:48:HIS:O	1:D:115:ARG:HB3	2.22	0.40
1:F:23:LYS:CE	1:F:30:LYS:HE3	2.49	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	153/153 (100%)	153 (100%)	0	0	100 100
1	B	153/153 (100%)	152 (99%)	0	1 (1%)	22 8
1	C	153/153 (100%)	147 (96%)	6 (4%)	0	100 100
1	D	153/153 (100%)	152 (99%)	1 (1%)	0	100 100
1	E	153/153 (100%)	151 (99%)	2 (1%)	0	100 100
1	F	153/153 (100%)	153 (100%)	0	0	100 100
1	G	153/153 (100%)	147 (96%)	4 (3%)	2 (1%)	12 2
1	H	153/153 (100%)	152 (99%)	1 (1%)	0	100 100
1	I	153/153 (100%)	150 (98%)	3 (2%)	0	100 100
1	J	153/153 (100%)	150 (98%)	3 (2%)	0	100 100
All	All	1530/1530 (100%)	1507 (98%)	20 (1%)	3 (0%)	47 30

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	25	SER
1	G	25	SER
1	G	132	GLU

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	120/118 (102%)	116 (97%)	4 (3%)	38 19
1	B	120/118 (102%)	115 (96%)	5 (4%)	30 12
1	C	120/118 (102%)	109 (91%)	11 (9%)	9 2
1	D	120/118 (102%)	119 (99%)	1 (1%)	81 74
1	E	120/118 (102%)	116 (97%)	4 (3%)	38 19
1	F	120/118 (102%)	116 (97%)	4 (3%)	38 19
1	G	120/118 (102%)	108 (90%)	12 (10%)	7 1
1	H	120/118 (102%)	118 (98%)	2 (2%)	60 46
1	I	120/118 (102%)	114 (95%)	6 (5%)	24 8
1	J	120/118 (102%)	115 (96%)	5 (4%)	30 12
All	All	1200/1180 (102%)	1146 (96%)	54 (4%)	27 10

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	109	ASP
1	A	111	CYS
1	A	135	THR
1	B	2	THR
1	B	23	LYS
1	B	24	GLU
1	B	26	ASN
1	B	70	LYS
1	C	21	GLU
1	C	30	LYS
1	C	34	SER
1	C	40	GLU
1	C	69	ARG
1	C	70	LYS
1	C	77	GLU

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Mol	Chain	Res	Type
1	C	78	GLU
1	C	121	GLU
1	C	135	THR
1	C	136	LYS
1	D	23	LYS
1	E	9	LYS
1	E	23	LYS
1	E	30	LYS
1	E	34	SER
1	F	2	THR
1	F	25	SER
1	F	75	LYS
1	F	91	LYS
1	G	21	GLU
1	G	25	SER
1	G	26	ASN
1	G	34	SER
1	G	68	SER
1	G	69	ARG
1	G	76	ASP
1	G	77	GLU
1	G	91	LYS
1	G	100	GLU
1	G	109	ASP
1	G	132	GLU
1	H	3	LYS
1	H	111	CYS
1	I	59	SER
1	I	70	LYS
1	I	77	GLU
1	I	107	SER
1	I	128	LYS
1	I	131	ASN
1	J	2	THR
1	J	24	GLU
1	J	36	LYS
1	J	59	SER
1	J	109	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	26	ASN
1	F	153	GLN
1	G	53	ASN
1	G	139	ASN
1	H	19	ASN
1	H	53	ASN
1	I	131	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 23 ligands modelled in this entry, 20 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	SO4	J	302	-	4,4,4	0.27	0	6,6,6	0.08	0
4	SO4	J	301	-	4,4,4	0.29	0	6,6,6	0.12	0
4	SO4	B	300	-	4,4,4	0.33	0	6,6,6	0.34	0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	J	302	SO4	1	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, 95^{th} 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	153/153 (100%)	0.22	6 (3%) 39 44	20, 28, 46, 59	0
1	B	153/153 (100%)	0.05	6 (3%) 39 44	18, 24, 49, 71	0
1	C	153/153 (100%)	0.99	21 (13%) 3 3	23, 39, 70, 89	0
1	D	153/153 (100%)	0.17	6 (3%) 39 44	19, 25, 48, 64	0
1	E	153/153 (100%)	0.14	5 (3%) 46 51	19, 25, 44, 64	0
1	F	153/153 (100%)	0.23	9 (5%) 22 24	19, 25, 50, 76	0
1	G	153/153 (100%)	0.80	20 (13%) 3 3	28, 39, 75, 81	0
1	H	153/153 (100%)	0.15	7 (4%) 32 36	22, 29, 54, 86	0
1	I	153/153 (100%)	0.72	25 (16%) 1 1	24, 37, 71, 95	0
1	J	153/153 (100%)	0.45	16 (10%) 6 7	21, 31, 56, 88	0
All	All	1530/1530 (100%)	0.39	121 (7%) 12 14	18, 30, 60, 95	0

All (121) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	1	ALA	7.7
1	C	1	ALA	7.6
1	J	1	ALA	7.0
1	C	129	GLY	6.5
1	J	26	ASN	5.8
1	I	130	GLY	5.7
1	G	131	ASN	5.4
1	C	109	ASP	5.4
1	G	132	GLU	5.1
1	F	26	ASN	5.1
1	F	1	ALA	4.8
1	B	1	ALA	4.6
1	J	25	SER	4.6

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Mol	Chain	Res	Type	RSRZ
1	I	132	GLU	4.5
1	D	26	ASN	4.4
1	G	1	ALA	4.4
1	B	26	ASN	4.4
1	C	132	GLU	4.3
1	C	135	THR	4.2
1	I	55	ALA	4.2
1	I	131	ASN	4.1
1	A	109	ASP	4.1
1	J	23	LYS	4.0
1	I	27	GLY	3.9
1	G	23	LYS	3.9
1	I	24	GLU	3.9
1	G	25	SER	3.9
1	C	131	ASN	3.8
1	D	1	ALA	3.8
1	J	109	ASP	3.8
1	G	133	GLU	3.7
1	G	130	GLY	3.7
1	I	109	ASP	3.6
1	C	91	LYS	3.6
1	I	11	ASP	3.6
1	J	92	ASP	3.3
1	E	117	LEU	3.3
1	I	25	SER	3.3
1	J	24	GLU	3.3
1	J	110	HIS	3.3
1	F	148	VAL	3.3
1	G	77	GLU	3.2
1	B	25	SER	3.2
1	A	130	GLY	3.2
1	C	117	LEU	3.1
1	I	129	GLY	3.1
1	F	23	LYS	3.1
1	C	47	VAL	3.1
1	C	67	LEU	3.1
1	C	75	LYS	3.1
1	I	149	ILE	3.1
1	I	66	PRO	3.1
1	C	130	GLY	3.1
1	G	24	GLU	3.0
1	G	136	LYS	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	132	GLU	3.0
1	A	1	ALA	3.0
1	I	136	LYS	3.0
1	D	149	ILE	3.0
1	I	77	GLU	3.0
1	E	149	ILE	2.9
1	J	149	ILE	2.9
1	H	108	GLY	2.9
1	H	1	ALA	2.8
1	G	27	GLY	2.8
1	I	148	VAL	2.8
1	E	1	ALA	2.8
1	C	149	ILE	2.8
1	I	26	ASN	2.8
1	I	153	GLN	2.7
1	C	69	ARG	2.7
1	G	109	ASP	2.7
1	J	153	GLN	2.7
1	J	84	LEU	2.6
1	G	129	GLY	2.6
1	H	111	CYS	2.6
1	I	110	HIS	2.6
1	C	76	ASP	2.6
1	J	11	ASP	2.6
1	G	67	LEU	2.6
1	G	26	ASN	2.5
1	I	67	LEU	2.5
1	H	109	ASP	2.5
1	G	28	PRO	2.5
1	A	131	ASN	2.4
1	G	135	THR	2.4
1	I	68	SER	2.4
1	C	110	HIS	2.4
1	H	26	ASN	2.4
1	C	92	ASP	2.3
1	F	24	GLU	2.3
1	C	23	LYS	2.3
1	H	110	HIS	2.3
1	J	91	LYS	2.3
1	E	153	GLN	2.3
1	I	23	LYS	2.3
1	J	47	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	F	117	LEU	2.3
1	I	70	LYS	2.3
1	B	24	GLU	2.3
1	D	45	PHE	2.2
1	J	117	LEU	2.2
1	F	77	GLU	2.2
1	J	148	VAL	2.2
1	B	23	LYS	2.2
1	H	149	ILE	2.2
1	C	128	LYS	2.1
1	G	153	GLN	2.1
1	B	148	VAL	2.1
1	C	148	VAL	2.1
1	D	148	VAL	2.1
1	D	25	SER	2.1
1	I	13	PRO	2.1
1	E	148	VAL	2.1
1	C	118	VAL	2.1
1	G	148	VAL	2.0
1	A	84	LEU	2.0
1	G	134	SER	2.0
1	F	149	ILE	2.0
1	I	2	THR	2.0
1	F	118	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 monosaccharides 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, 95^{th} 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
4	SO4	J	302	5/5	0.77	0.46	52,105,121,128	0
4	SO4	B	300	5/5	0.92	0.20	43,55,64,85	0
4	SO4	J	301	5/5	0.96	0.25	56,83,87,88	0
2	CU1	C	200	1/1	0.99	0.04	40,40,40,40	0
2	CU1	G	200	1/1	0.99	0.04	35,35,35,35	0
3	ZN	C	201	1/1	0.99	0.04	46,46,46,46	0
2	CU1	B	200	1/1	1.00	0.04	21,21,21,21	0
2	CU1	H	200	1/1	1.00	0.04	25,25,25,25	0
2	CU1	I	200	1/1	1.00	0.02	35,35,35,35	0
2	CU1	J	200	1/1	1.00	0.05	24,24,24,24	0
3	ZN	A	201	1/1	1.00	0.05	25,25,25,25	0
3	ZN	B	201	1/1	1.00	0.05	20,20,20,20	0
2	CU1	A	200	1/1	1.00	0.04	24,24,24,24	0
3	ZN	D	201	1/1	1.00	0.05	20,20,20,20	0
3	ZN	E	201	1/1	1.00	0.05	22,22,22,22	0
3	ZN	F	201	1/1	1.00	0.04	21,21,21,21	0
3	ZN	G	201	1/1	1.00	0.02	44,44,44,44	0
3	ZN	H	201	1/1	1.00	0.05	24,24,24,24	0
3	ZN	I	201	1/1	1.00	0.05	40,40,40,40	0
3	ZN	J	201	1/1	1.00	0.04	23,23,23,23	0
2	CU1	D	200	1/1	1.00	0.04	22,22,22,22	0
2	CU1	E	200	1/1	1.00	0.04	22,22,22,22	0
2	CU1	F	200	1/1	1.00	0.05	21,21,21,21	0

6.5 Other polymers (i)

There are no such residues in this entry.