



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:50 PM UTC

PDB ID : 3RE7 / pdb_00003re7
Title : Copper (II) loaded Bullfrog Ferritin M chain
Authors : Bertini, I.; Lalli, D.; Mangani, S.; Pozzi, C.; Rosa, C.; Turano, P.
Deposited on : 2011-04-02
Resolution : 2.82 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:

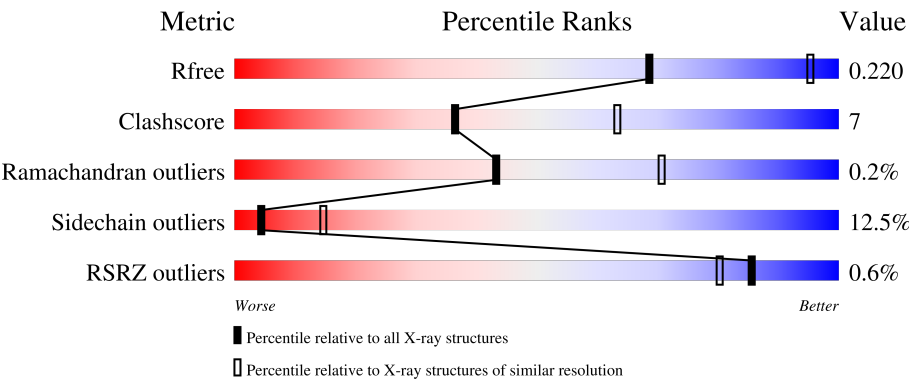
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.82 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.



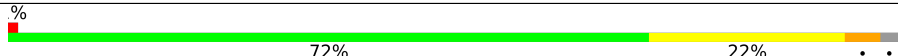
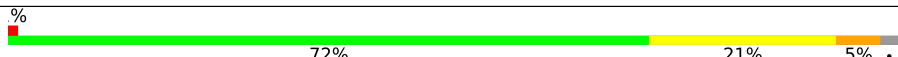
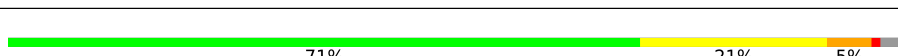
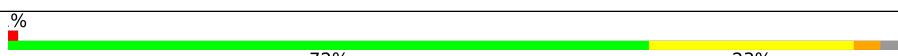
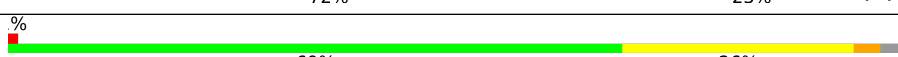
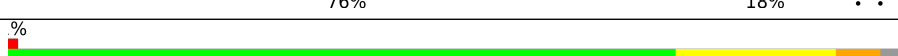
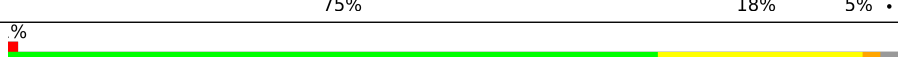
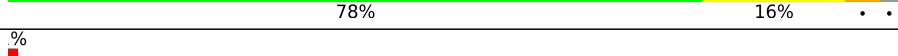
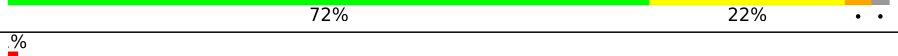
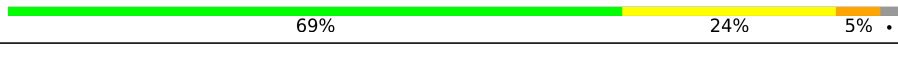
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4591 (2.84-2.80)
Clashscore	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
Sidechain outliers	187428	4918 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

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	176	75%22%..
1	B	176	% 73%22%..
1	C	176	% 72%21%5%.
1	D	176	% 76%19%...
1	E	176	% 68%25%5%.

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Mol	Chain	Length	Quality of chain	
1	F	176		
1	G	176		
1	H	176		
1	I	176		
1	J	176		
1	K	176		
1	L	176		
1	M	176		
1	N	176		
1	O	176		
1	P	176		
1	Q	176		
1	R	176		
1	S	176		
1	T	176		
1	U	176		
1	V	176		
1	W	176		
1	X	176		

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
2	CU	B	184	-	-	-	X
2	CU	D	191	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 34999 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 Ferritin, middle subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	B	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	C	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	D	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	E	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	F	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	G	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	H	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	I	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	J	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	K	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	L	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	M	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	N	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	O	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	P	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	R	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	S	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	T	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	U	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	V	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	W	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			
1	X	172	Total	C	N	O	S	0	0	0
			1418	892	247	272	7			

- Molecule 2 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	8	Total	Cu	0	0
			8	8		
2	B	8	Total	Cu	0	0
			8	8		
2	C	7	Total	Cu	0	0
			7	7		
2	D	10	Total	Cu	0	0
			10	10		
2	E	8	Total	Cu	0	0
			8	8		
2	F	6	Total	Cu	0	0
			6	6		
2	G	8	Total	Cu	0	0
			8	8		
2	H	7	Total	Cu	0	0
			7	7		
2	I	7	Total	Cu	0	0
			7	7		
2	J	8	Total	Cu	0	0
			8	8		
2	K	6	Total	Cu	0	0
			6	6		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	L	7	Total 7	Cu 7	0	0
2	M	8	Total 8	Cu 8	0	0
2	N	7	Total 7	Cu 7	0	0
2	O	6	Total 6	Cu 6	0	0
2	P	6	Total 6	Cu 6	0	0
2	Q	6	Total 6	Cu 6	0	0
2	R	7	Total 7	Cu 7	0	0
2	S	5	Total 5	Cu 5	0	0
2	T	4	Total 4	Cu 4	0	0
2	U	6	Total 6	Cu 6	0	0
2	V	5	Total 5	Cu 5	0	0
2	W	6	Total 6	Cu 6	0	0
2	X	6	Total 6	Cu 6	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	42	Total 42	O 42	0	0
3	B	22	Total 22	O 22	0	0
3	C	34	Total 34	O 34	0	0
3	D	31	Total 31	O 31	0	0
3	E	22	Total 22	O 22	0	0
3	F	28	Total 28	O 28	0	0

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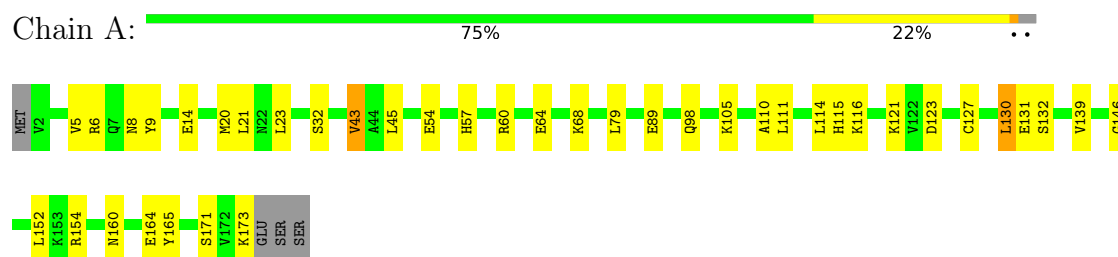
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	31	Total 31	O 31	0	0
3	H	35	Total 35	O 35	0	0
3	I	24	Total 24	O 24	0	0
3	J	44	Total 44	O 44	0	0
3	K	48	Total 48	O 48	0	0
3	L	19	Total 19	O 19	0	0
3	M	30	Total 30	O 30	0	0
3	N	41	Total 41	O 41	0	0
3	O	31	Total 31	O 31	0	0
3	P	32	Total 32	O 32	0	0
3	Q	41	Total 41	O 41	0	0
3	R	41	Total 41	O 41	0	0
3	S	35	Total 35	O 35	0	0
3	T	29	Total 29	O 29	0	0
3	U	41	Total 41	O 41	0	0
3	V	20	Total 20	O 20	0	0
3	W	41	Total 41	O 41	0	0
3	X	43	Total 43	O 43	0	0

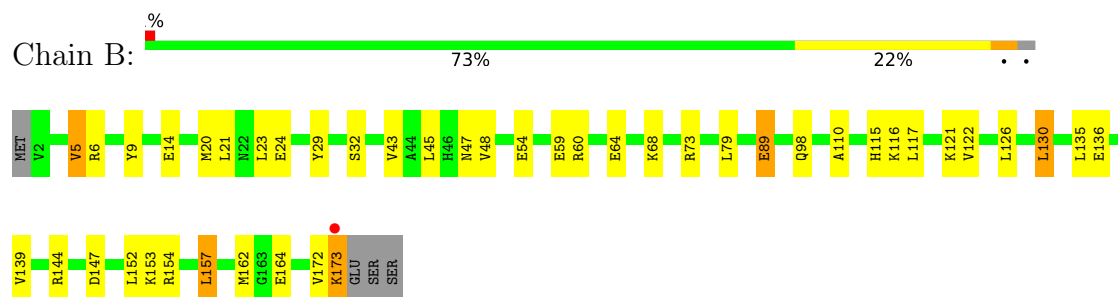
3 Residue-property plots

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.

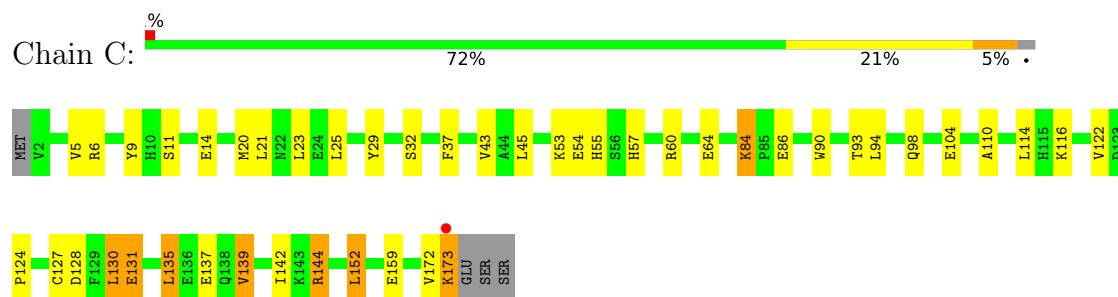
- Molecule 1: Ferritin, middle subunit



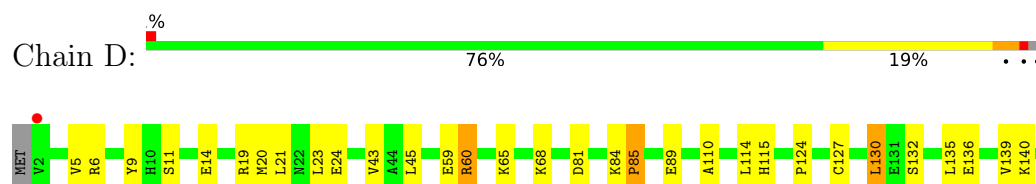
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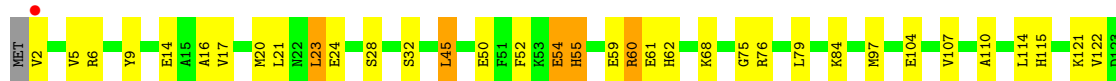


- Molecule 1: Ferritin, middle subunit

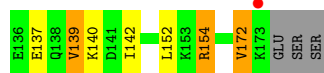
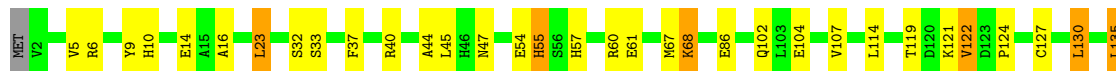
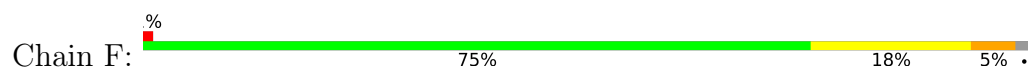




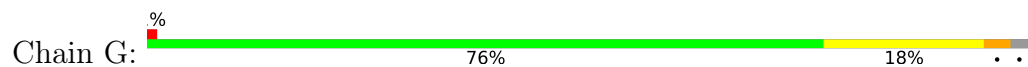
- Molecule 1: Ferritin, middle subunit



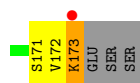
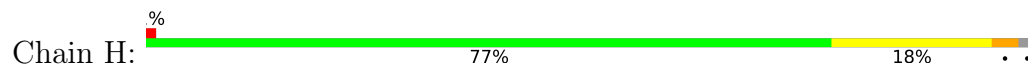
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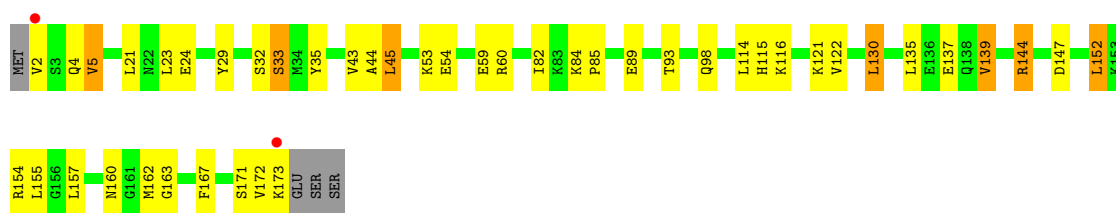


- Molecule 1: Ferritin, middle subunit

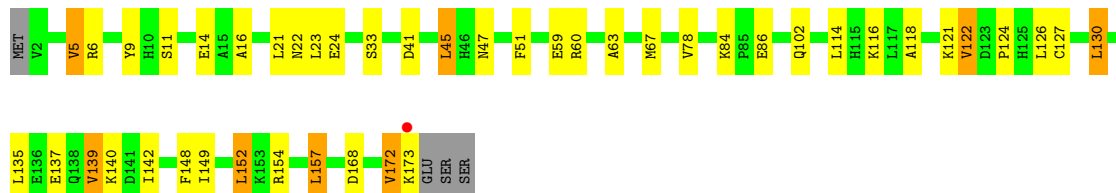


- Molecule 1: Ferritin, middle subunit

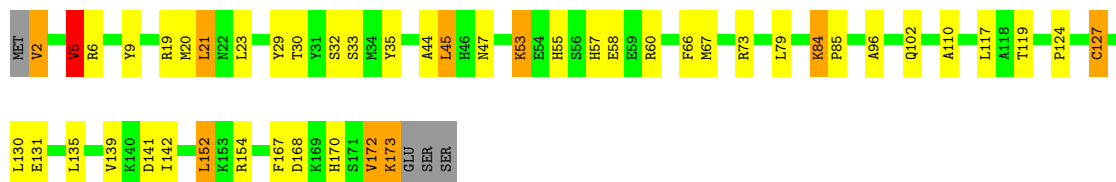




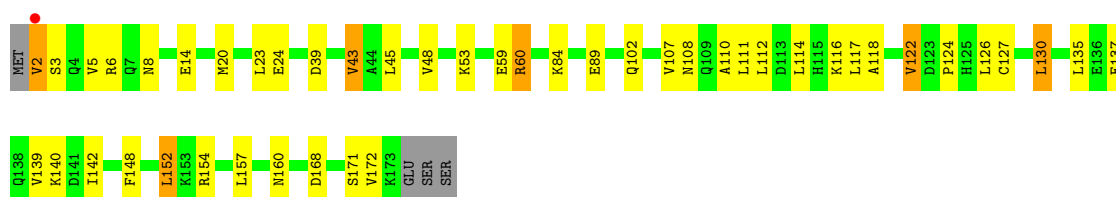
● Molecule 1: Ferritin, middle subunit



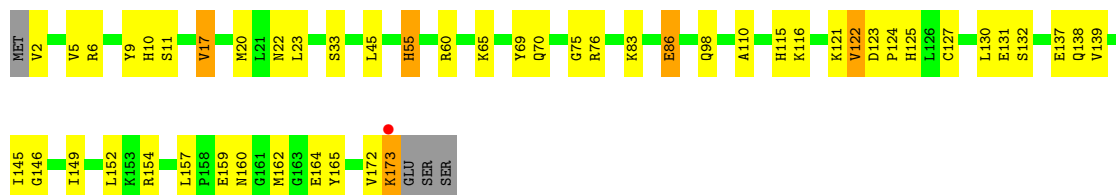
● Molecule 1: Ferritin, middle subunit



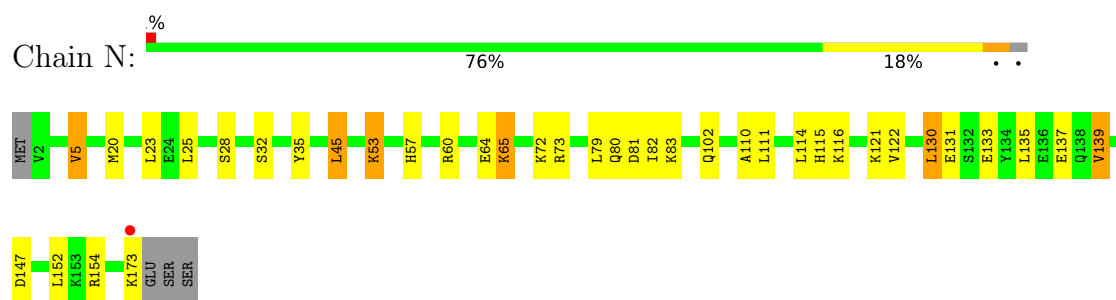
● Molecule 1: Ferritin, middle subunit



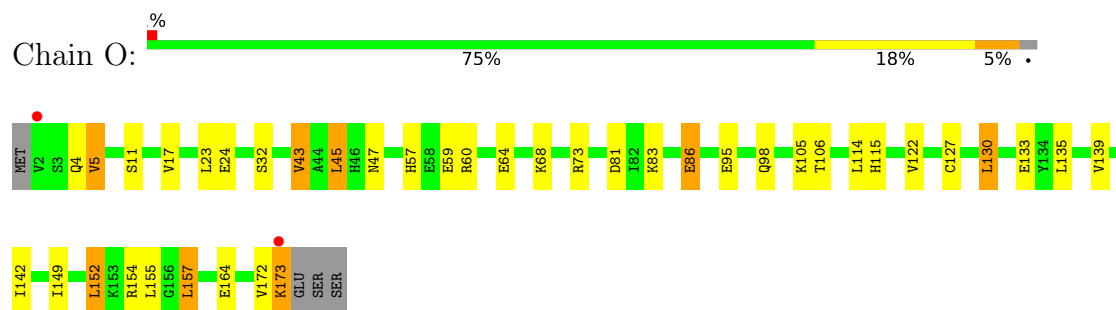
● Molecule 1: Ferritin, middle subunit



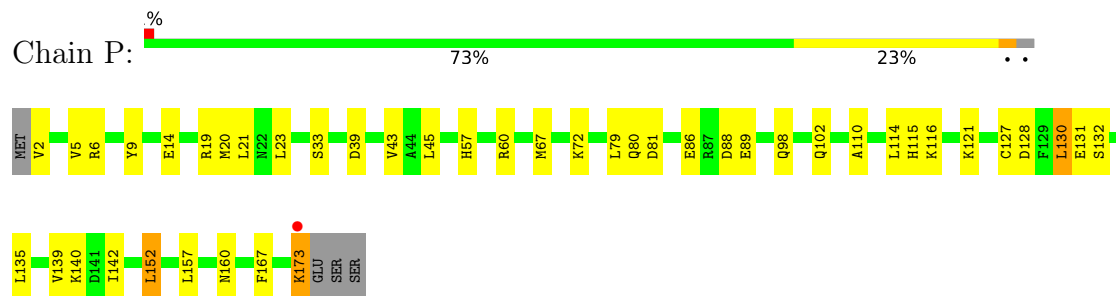
● Molecule 1: Ferritin, middle subunit



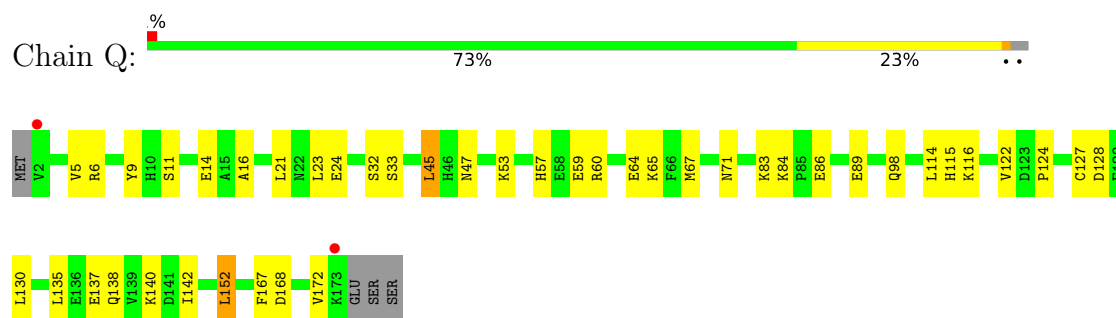
- Molecule 1: Ferritin, middle subunit



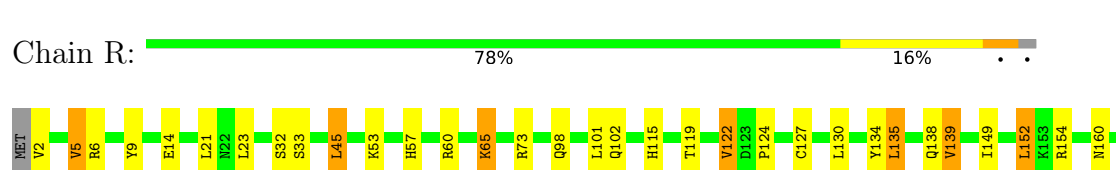
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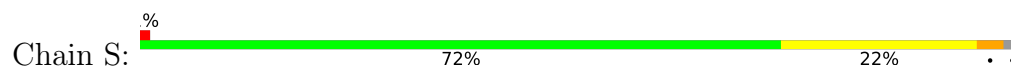


- Molecule 1: Ferritin, middle subunit

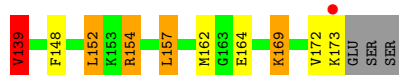




- Molecule 1: Ferritin, middle subunit



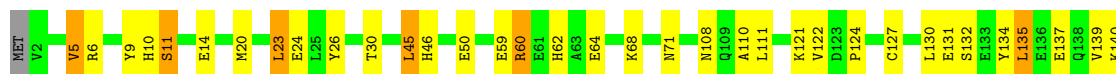
- Molecule 1: Ferritin, middle subunit



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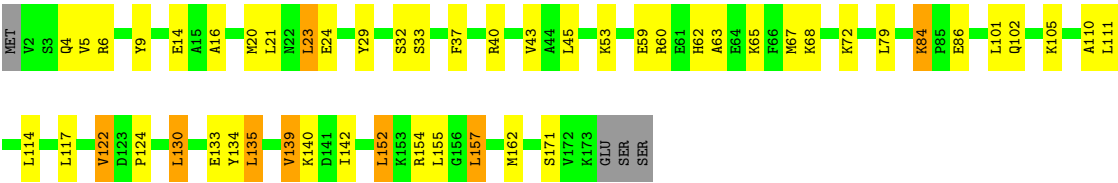


- Molecule 1: Ferritin, middle subunit

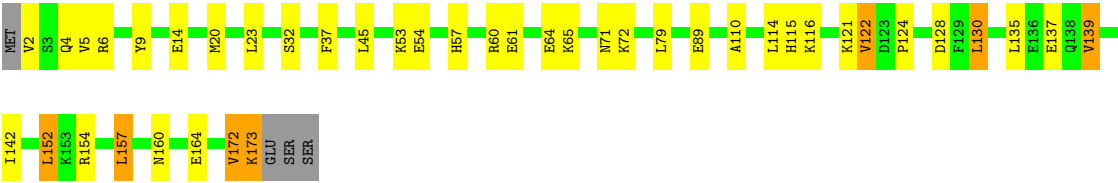


- Molecule 1: Ferritin, middle subunit





● Molecule 1: Ferritin, middle subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	210.21Å 210.21Å 322.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.50 – 2.82 48.50 – 2.82	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.50-2.82) 99.9 (48.50-2.82)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.05 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5	Depositor
R, R_{free}	0.170 , 0.224 0.170 , 0.220	Depositor DCC
R_{free} test set	9961 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	42.4	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 25.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.029 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	34999	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 3.78% 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: CU

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	1.12	0/1446	1.22	4/1944 (0.2%)
1	B	1.04	0/1446	1.16	2/1944 (0.1%)
1	C	1.12	0/1446	1.18	4/1944 (0.2%)
1	D	1.17	1/1446 (0.1%)	1.19	2/1944 (0.1%)
1	E	1.07	0/1446	1.15	3/1944 (0.2%)
1	F	1.09	0/1446	1.15	4/1944 (0.2%)
1	G	1.15	1/1446 (0.1%)	1.18	2/1944 (0.1%)
1	H	1.13	0/1446	1.18	2/1944 (0.1%)
1	I	1.08	1/1446 (0.1%)	1.21	6/1944 (0.3%)
1	J	1.23	2/1446 (0.1%)	1.19	4/1944 (0.2%)
1	K	1.14	2/1446 (0.1%)	1.24	7/1944 (0.4%)
1	L	1.11	0/1446	1.14	3/1944 (0.2%)
1	M	1.14	1/1446 (0.1%)	1.16	7/1944 (0.4%)
1	N	1.15	0/1446	1.21	1/1944 (0.1%)
1	O	1.19	0/1446	1.17	2/1944 (0.1%)
1	P	1.10	0/1446	1.18	3/1944 (0.2%)
1	Q	1.15	0/1446	1.16	1/1944 (0.1%)
1	R	1.14	1/1446 (0.1%)	1.16	3/1944 (0.2%)
1	S	1.12	0/1446	1.17	1/1944 (0.1%)
1	T	1.11	0/1446	1.20	6/1944 (0.3%)
1	U	1.14	0/1446	1.17	3/1944 (0.2%)
1	V	1.10	1/1446 (0.1%)	1.22	3/1944 (0.2%)
1	W	1.15	1/1446 (0.1%)	1.22	4/1944 (0.2%)
1	X	1.20	0/1446	1.19	3/1944 (0.2%)
All	All	1.13	11/34704 (0.0%)	1.18	80/46656 (0.2%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	V	142	ILE	CA-CB	8.00	1.64	1.54
1	J	78	VAL	CA-CB	6.46	1.62	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	139	VAL	CA-CB	5.60	1.62	1.54
1	M	55	HIS	C-O	-5.51	1.17	1.24
1	W	43	VAL	CA-CB	-5.48	1.50	1.55
1	R	139	VAL	CA-CB	5.32	1.61	1.54
1	K	96	ALA	CA-CB	-5.27	1.44	1.53
1	D	85	PRO	CA-C	5.15	1.58	1.52
1	K	141	ASP	CB-CG	5.11	1.64	1.52
1	I	82	ILE	CA-CB	5.05	1.59	1.53
1	J	118	ALA	C-O	-5.04	1.18	1.24

All (80) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	68	LYS	N-CA-C	-7.59	103.00	111.28
1	I	84	LYS	CA-C-N	-7.55	112.96	120.21
1	I	84	LYS	C-N-CA	-7.55	112.96	120.21
1	K	5	VAL	N-CA-C	-7.20	106.48	113.53
1	T	48	VAL	N-CA-C	-7.02	103.63	110.72
1	A	43	VAL	N-CA-C	-6.91	106.16	111.62
1	L	43	VAL	N-CA-C	-6.91	106.16	111.62
1	N	139	VAL	N-CA-CB	6.75	119.72	110.54
1	W	84	LYS	CA-C-N	-6.58	113.89	120.21
1	W	84	LYS	C-N-CA	-6.58	113.89	120.21
1	M	33	SER	N-CA-C	-6.55	104.22	111.36
1	W	86	GLU	N-CA-C	6.55	121.10	113.18
1	K	172	VAL	CB-CA-C	6.43	119.72	110.33
1	F	104	GLU	N-CA-C	-6.40	104.39	111.36
1	A	160	ASN	N-CA-C	6.19	118.38	108.41
1	F	68	LYS	N-CA-C	-6.14	104.59	111.28
1	X	139	VAL	N-CA-CB	6.13	119.78	110.58
1	V	132	SER	N-CA-C	6.11	118.02	111.36
1	W	139	VAL	N-CA-CB	5.99	119.56	110.58
1	B	89	GLU	N-CA-C	-5.98	101.14	110.42
1	D	132	SER	N-CA-C	5.95	117.77	111.28
1	J	139	VAL	N-CA-CB	5.92	118.60	110.54
1	T	43	VAL	N-CA-C	-5.91	106.48	111.56
1	K	84	LYS	CA-C-N	-5.90	114.68	120.52
1	K	84	LYS	C-N-CA	-5.90	114.68	120.52
1	H	65	LYS	N-CA-C	5.89	117.70	111.28
1	L	160	ASN	N-CA-C	5.88	118.68	107.75
1	I	44	ALA	N-CA-C	5.87	119.45	111.17
1	V	142	ILE	N-CA-CB	5.85	118.49	110.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	P	132	SER	N-CA-C	5.76	118.78	111.69
1	X	71	ASN	N-CA-C	5.76	117.64	111.36
1	K	2	VAL	CB-CA-C	-5.76	100.46	111.40
1	F	55	HIS	N-CA-C	5.76	117.55	111.28
1	M	146	GLY	N-CA-C	-5.67	105.92	112.50
1	Q	138	GLN	N-CA-C	5.66	117.53	111.36
1	I	160	ASN	N-CA-C	5.65	117.29	109.15
1	O	133	GLU	N-CA-C	5.62	120.14	113.28
1	M	160	ASN	N-CA-C	5.59	117.83	107.99
1	H	57	HIS	N-CA-C	-5.58	105.28	111.36
1	P	88	ASP	N-CA-C	-5.58	107.02	113.88
1	I	139	VAL	N-CA-CB	5.55	118.09	110.54
1	C	139	VAL	N-CA-CB	5.52	120.09	110.65
1	M	17	VAL	N-CA-C	-5.51	105.35	110.53
1	O	43	VAL	N-CA-C	-5.48	107.29	111.62
1	U	139	VAL	N-CA-CB	5.47	118.78	110.58
1	C	86	GLU	N-CA-C	5.44	119.02	112.38
1	K	170	HIS	N-CA-C	5.43	120.56	113.88
1	M	138	GLN	N-CA-C	5.39	117.58	111.11
1	X	160	ASN	N-CA-C	5.39	117.48	107.99
1	I	43	VAL	O-C-N	5.38	126.12	121.98
1	R	138	GLN	N-CA-C	5.32	116.77	111.07
1	K	44	ALA	N-CA-C	5.29	118.92	111.52
1	U	30	THR	N-CA-C	-5.25	105.64	111.36
1	T	84	LYS	CA-C-N	-5.24	115.33	120.52
1	T	84	LYS	C-N-CA	-5.24	115.33	120.52
1	S	33	SER	N-CA-C	-5.23	105.66	111.36
1	M	123	ASP	N-CA-C	-5.23	100.19	108.82
1	T	71	ASN	N-CA-C	5.22	117.05	111.36
1	E	54	GLU	N-CA-C	-5.20	105.51	111.07
1	F	44	ALA	N-CA-C	5.19	118.40	111.24
1	A	164	GLU	N-CA-C	-5.18	105.81	111.82
1	E	50	GLU	N-CA-C	-5.18	105.81	111.82
1	M	22	ASN	N-CA-C	5.18	117.59	111.33
1	V	144	ARG	N-CA-C	-5.17	105.54	111.07
1	C	84	LYS	CA-C-N	-5.17	115.25	120.21
1	C	84	LYS	C-N-CA	-5.17	115.25	120.21
1	B	48	VAL	N-CA-C	-5.15	105.52	110.72
1	U	160	ASN	N-CA-C	5.13	117.66	108.17
1	D	144	ARG	NE-CZ-NH2	5.11	123.80	119.20
1	J	172	VAL	CB-CA-C	5.09	118.20	110.62
1	T	139	VAL	N-CA-CB	5.08	119.96	110.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	146	GLY	N-CA-C	-5.05	106.64	112.50
1	J	22	ASN	N-CA-C	5.05	116.78	111.28
1	G	160	ASN	N-CA-C	5.04	117.12	107.75
1	L	89	GLU	N-CA-C	-5.03	100.40	108.55
1	P	160	ASN	N-CA-C	5.03	116.50	108.41
1	J	126	LEU	N-CA-C	-5.02	105.89	111.36
1	R	160	ASN	N-CA-C	5.02	117.42	107.98
1	E	62	HIS	N-CA-C	5.01	117.39	111.33
1	R	152	LEU	N-CA-C	-5.00	106.02	111.82

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	1418	0	1369	27	0
1	B	1418	0	1369	25	0
1	C	1418	0	1369	29	0
1	D	1418	0	1370	24	0
1	E	1418	0	1369	31	0
1	F	1418	0	1369	21	0
1	G	1418	0	1369	27	0
1	H	1418	0	1369	20	0
1	I	1418	0	1369	24	0
1	J	1418	0	1369	22	0
1	K	1418	0	1369	33	0
1	L	1418	0	1369	24	0
1	M	1418	0	1369	26	0
1	N	1418	0	1369	21	0
1	O	1418	0	1369	25	0
1	P	1418	0	1369	24	0
1	Q	1418	0	1369	20	0
1	R	1418	0	1369	20	0
1	S	1418	0	1369	32	0
1	T	1418	0	1369	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	U	1418	0	1369	25	0
1	V	1418	0	1369	28	0
1	W	1418	0	1369	35	0
1	X	1418	0	1370	24	0
2	A	8	0	0	0	0
2	B	8	0	0	0	0
2	C	7	0	0	0	0
2	D	10	0	0	3	0
2	E	8	0	0	0	0
2	F	6	0	0	0	0
2	G	8	0	0	0	0
2	H	7	0	0	0	0
2	I	7	0	0	0	0
2	J	8	0	0	0	0
2	K	6	0	0	0	0
2	L	7	0	0	0	0
2	M	8	0	0	0	0
2	N	7	0	0	0	0
2	O	6	0	0	0	0
2	P	6	0	0	0	0
2	Q	6	0	0	0	0
2	R	7	0	0	0	0
2	S	5	0	0	0	0
2	T	4	0	0	0	0
2	U	6	0	0	0	0
2	V	5	0	0	0	0
2	W	6	0	0	0	0
2	X	6	0	0	0	0
3	A	42	0	0	3	0
3	B	22	0	0	1	0
3	C	34	0	0	1	0
3	D	31	0	0	5	0
3	E	22	0	0	2	0
3	F	28	0	0	2	0
3	G	31	0	0	1	0
3	H	35	0	0	3	0
3	I	24	0	0	3	0
3	J	44	0	0	1	0
3	K	48	0	0	0	0
3	L	19	0	0	1	0
3	M	30	0	0	2	0
3	N	41	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	O	31	0	0	5	0
3	P	32	0	0	2	0
3	Q	41	0	0	3	0
3	R	41	0	0	0	0
3	S	35	0	0	4	0
3	T	29	0	0	1	0
3	U	41	0	0	3	0
3	V	20	0	0	2	0
3	W	41	0	0	6	0
3	X	43	0	0	1	0
All	All	34999	0	32858	482	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 (482) 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:20:MET:HE1	1:B:110:ALA:HB1	1.29	1.11
1:D:144:ARG:HD2	3:D:663:HOH:O	1.52	1.09
1:I:144:ARG:HD2	3:I:274:HOH:O	1.58	1.03
1:P:173:LYS:H	1:P:173:LYS:HD3	1.22	1.01
1:A:57:HIS:HB3	3:A:673:HOH:O	1.61	0.99
1:A:60:ARG:HD3	1:G:60:ARG:NE	1.76	0.99
1:M:20:MET:HE1	1:M:110:ALA:HB1	1.50	0.94
1:C:60:ARG:HD3	1:O:60:ARG:HD2	1.51	0.92
1:D:127:CYS:HG	2:D:191:CU:CU	0.82	0.91
1:M:20:MET:HE2	1:M:20:MET:HA	1.54	0.89
1:S:118:ALA:HB2	1:S:126:LEU:HD23	1.54	0.89
1:B:20:MET:HE1	1:B:110:ALA:CB	2.05	0.87
1:T:6:ARG:NH2	1:T:14:GLU:OE1	2.08	0.86
1:P:173:LYS:H	1:P:173:LYS:CD	1.87	0.86
1:V:148:PHE:O	1:V:152:LEU:HD22	1.76	0.84
1:C:142:ILE:HG22	1:S:5:VAL:HG22	1.57	0.83
1:K:60:ARG:HD3	1:W:60:ARG:HD3	1.58	0.83
1:D:20:MET:HE1	1:D:110:ALA:HB1	1.60	0.82
1:M:20:MET:HE1	1:M:110:ALA:CB	2.09	0.82
1:X:6:ARG:NH1	1:X:9:TYR:O	2.13	0.82
1:A:20:MET:HE1	1:A:110:ALA:C	2.04	0.81
1:C:6:ARG:NH2	1:C:14:GLU:OE1	2.13	0.81
1:W:20:MET:HE3	1:W:111:LEU:HD23	1.61	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:W:6:ARG:NH1	1:W:9:TYR:O	2.14	0.80
1:D:84:LYS:NZ	3:D:193:HOH:O	2.10	0.80
1:P:6:ARG:NH1	1:P:9:TYR:O	2.14	0.80
1:C:131:GLU:HG2	1:S:128:ASP:HB2	1.64	0.78
1:X:20:MET:HE1	1:X:110:ALA:C	2.08	0.78
1:J:142:ILE:HG22	1:N:5:VAL:HG22	1.66	0.78
1:D:154:ARG:NH2	1:K:45:LEU:HD13	2.00	0.77
1:A:60:ARG:HD3	1:G:60:ARG:HE	1.46	0.77
1:L:122:VAL:HG12	1:L:122:VAL:O	1.84	0.76
1:L:6:ARG:NH2	1:L:14:GLU:OE1	2.16	0.76
1:L:20:MET:HE1	1:L:110:ALA:C	2.09	0.76
1:Q:6:ARG:NH2	1:Q:14:GLU:OE1	2.17	0.76
1:R:5:VAL:HG22	1:S:142:ILE:HG22	1.68	0.75
1:D:6:ARG:NH2	1:D:14:GLU:OE1	2.19	0.75
1:L:20:MET:HE3	1:L:111:LEU:HD23	1.69	0.75
1:W:20:MET:HE1	1:W:110:ALA:C	2.12	0.75
1:A:6:ARG:NH2	1:A:14:GLU:OE1	2.17	0.75
1:Q:6:ARG:NH1	1:Q:9:TYR:O	2.19	0.75
1:J:6:ARG:NH1	1:J:9:TYR:O	2.19	0.75
1:K:142:ILE:HG22	1:U:5:VAL:HG22	1.69	0.75
1:D:115:HIS:CE1	1:W:124:PRO:HB3	2.22	0.74
1:S:6:ARG:NH2	1:S:14:GLU:OE1	2.16	0.74
1:U:6:ARG:NH1	1:U:9:TYR:O	2.20	0.74
1:P:128:ASP:HB2	1:U:131:GLU:HG2	1.68	0.74
1:R:6:ARG:NH2	1:R:14:GLU:OE1	2.18	0.74
1:S:57:HIS:HB3	3:S:259:HOH:O	1.87	0.73
1:T:57:HIS:HB3	3:T:381:HOH:O	1.86	0.73
1:S:152:LEU:HD12	1:S:157:LEU:HD13	1.69	0.73
1:K:60:ARG:CD	1:W:60:ARG:HD3	2.18	0.73
1:B:60:ARG:CD	1:F:60:ARG:HD2	2.19	0.73
1:X:6:ARG:NH2	1:X:14:GLU:OE1	2.20	0.72
1:U:58:GLU:HG2	3:U:187:HOH:O	1.89	0.72
1:W:6:ARG:NH2	1:W:14:GLU:OE1	2.22	0.72
1:P:6:ARG:NH2	1:P:14:GLU:OE1	2.22	0.72
1:U:6:ARG:NH2	1:U:14:GLU:OE1	2.23	0.71
1:H:5:VAL:HG22	1:X:142:ILE:HG22	1.71	0.71
1:D:127:CYS:SG	2:D:191:CU:CU	1.81	0.71
1:S:11:SER:HB2	3:S:250:HOH:O	1.91	0.71
1:D:85:PRO:HG2	3:D:403:HOH:O	1.91	0.70
1:P:173:LYS:HD3	1:P:173:LYS:N	2.03	0.70
1:Q:57:HIS:HB3	3:Q:234:HOH:O	1.90	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:20:MET:HE1	1:N:110:ALA:C	2.16	0.70
1:X:173:LYS:H	1:X:173:LYS:HE3	1.56	0.70
3:N:178:HOH:O	1:R:57:HIS:HB3	1.91	0.70
1:V:144:ARG:HD3	3:V:672:HOH:O	1.92	0.70
1:B:173:LYS:HE2	1:B:173:LYS:H	1.56	0.70
1:J:47:ASN:HB2	1:J:168:ASP:OD1	1.91	0.70
1:G:6:ARG:NH2	1:G:14:GLU:OE1	2.19	0.69
1:F:122:VAL:HG13	3:F:627:HOH:O	1.92	0.69
1:V:152:LEU:HD12	1:V:157:LEU:HD13	1.73	0.69
1:N:154:ARG:NH2	1:U:43:VAL:O	2.24	0.69
1:P:19:ARG:HD3	3:P:346:HOH:O	1.92	0.68
1:P:60:ARG:NE	1:T:60:ARG:HD3	2.07	0.68
1:I:60:ARG:HD2	3:I:179:HOH:O	1.93	0.68
1:O:4:GLN:HG2	3:O:187:HOH:O	1.94	0.68
1:G:148:PHE:O	1:G:152:LEU:HD22	1.92	0.68
1:J:51:PHE:HB2	1:J:172:VAL:HG21	1.76	0.68
1:P:20:MET:HE1	1:P:110:ALA:C	2.20	0.67
1:A:123:ASP:OD2	3:A:746:HOH:O	2.12	0.67
1:Q:60:ARG:HD3	1:U:60:ARG:HD2	1.76	0.67
1:D:127:CYS:SG	2:D:192:CU:CU	1.85	0.67
1:J:152:LEU:HD12	1:J:157:LEU:HD13	1.77	0.66
1:J:5:VAL:HG22	1:Q:142:ILE:HG22	1.77	0.66
1:V:6:ARG:NH2	1:V:14:GLU:OE1	2.19	0.66
1:A:131:GLU:HG2	1:X:128:ASP:HB2	1.77	0.66
1:C:55:HIS:NE2	1:C:104:GLU:OE2	2.28	0.66
1:T:43:VAL:O	1:U:154:ARG:NH2	2.22	0.66
1:C:6:ARG:NH1	1:C:9:TYR:O	2.29	0.66
1:H:57:HIS:HB3	3:H:648:HOH:O	1.96	0.65
1:V:137:GLU:HG3	3:V:557:HOH:O	1.96	0.65
1:W:24:GLU:OE1	1:W:59:GLU:OE1	2.15	0.65
1:C:60:ARG:CD	1:O:60:ARG:HD2	2.27	0.65
1:F:23:LEU:HD13	1:F:107:VAL:HG22	1.79	0.65
1:D:172:VAL:O	1:D:173:LYS:HG3	1.97	0.65
1:D:6:ARG:NH1	1:D:9:TYR:O	2.30	0.65
1:N:64:GLU:OE1	1:R:57:HIS:CE1	2.50	0.64
1:E:154:ARG:NH2	1:L:43:VAL:O	2.29	0.64
1:K:20:MET:HE1	1:K:110:ALA:HB1	1.80	0.64
1:A:57:HIS:CE1	1:G:64:GLU:OE2	2.50	0.64
1:B:47:ASN:HB3	1:B:172:VAL:HG23	1.78	0.64
1:G:144:ARG:HG2	1:G:144:ARG:HH11	1.62	0.64
1:J:6:ARG:NH2	1:J:14:GLU:OE1	2.26	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:20:MET:HE1	1:C:110:ALA:HB1	1.79	0.63
1:O:45:LEU:HD13	1:R:154:ARG:NH2	2.14	0.63
1:G:124:PRO:HB3	1:O:115:HIS:CE1	2.33	0.63
1:J:24:GLU:OE1	1:J:59:GLU:OE1	2.16	0.63
1:U:144:ARG:HD3	3:U:200:HOH:O	1.98	0.63
1:F:57:HIS:HB3	3:F:379:HOH:O	1.98	0.62
1:C:142:ILE:HG22	1:S:5:VAL:CG2	2.27	0.62
1:L:122:VAL:O	1:L:122:VAL:CG1	2.48	0.62
1:E:6:ARG:NH1	1:E:9:TYR:O	2.32	0.62
1:F:6:ARG:NH2	1:F:14:GLU:OE1	2.27	0.62
1:E:20:MET:HE1	1:E:110:ALA:C	2.24	0.62
1:T:169:LYS:HG2	1:U:170:HIS:HB3	1.81	0.62
1:I:21:LEU:C	1:I:21:LEU:HD23	2.25	0.62
1:B:5:VAL:HG22	1:V:142:ILE:HG13	1.82	0.61
1:M:60:ARG:HD2	1:S:60:ARG:CZ	2.31	0.61
1:U:23:LEU:HD13	1:U:107:VAL:HG22	1.82	0.61
1:S:43:VAL:O	1:T:154:ARG:NH2	2.27	0.61
1:K:60:ARG:HD3	1:W:60:ARG:CD	2.29	0.61
1:E:60:ARG:HD2	1:I:60:ARG:CD	2.31	0.61
1:G:144:ARG:HG2	1:G:144:ARG:NH1	2.15	0.61
1:X:114:LEU:HG	1:X:130:LEU:HD21	1.83	0.60
1:A:20:MET:HE1	1:A:111:LEU:N	2.17	0.60
1:M:124:PRO:HA	1:M:127:CYS:HB3	1.84	0.60
1:E:6:ARG:NH2	1:E:14:GLU:OE1	2.33	0.60
1:E:157:LEU:HD21	1:E:164:GLU:HG3	1.83	0.60
1:S:11:SER:CB	3:S:250:HOH:O	2.49	0.60
1:C:114:LEU:HG	1:C:130:LEU:HD21	1.83	0.59
1:E:162:MET:SD	1:W:162:MET:HE2	2.42	0.59
1:M:83:LYS:HE3	3:M:341:HOH:O	2.02	0.59
1:S:157:LEU:HD21	1:S:164:GLU:HG3	1.83	0.59
1:T:157:LEU:HD21	1:T:164:GLU:HG3	1.83	0.59
1:D:43:VAL:O	1:X:154:ARG:NH2	2.32	0.59
1:A:60:ARG:CD	1:G:60:ARG:HE	2.15	0.59
1:E:60:ARG:HD2	1:I:60:ARG:HD3	1.85	0.59
1:A:6:ARG:NH1	1:A:9:TYR:O	2.34	0.59
1:B:21:LEU:C	1:B:21:LEU:HD23	2.27	0.59
1:J:154:ARG:NH2	1:R:45:LEU:HD13	2.17	0.59
1:Q:152:LEU:HD13	1:Q:167:PHE:CD1	2.38	0.59
1:A:64:GLU:OE1	1:G:57:HIS:HE1	1.85	0.58
1:I:29:TYR:O	1:I:32:SER:HB3	2.03	0.58
1:C:55:HIS:CE1	1:C:104:GLU:OE2	2.57	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:W:20:MET:HE1	1:W:111:LEU:N	2.19	0.58
1:L:154:ARG:NH2	1:P:43:VAL:O	2.36	0.58
1:N:60:ARG:HD2	1:R:60:ARG:HD3	1.85	0.58
3:E:708:HOH:O	1:I:4:GLN:HG2	2.03	0.58
1:O:43:VAL:O	1:R:154:ARG:NH2	2.24	0.58
1:T:45:LEU:HD13	1:U:154:ARG:NH2	2.18	0.58
1:B:43:VAL:O	1:G:154:ARG:NH2	2.30	0.58
1:E:61:GLU:HG3	3:E:532:HOH:O	2.03	0.58
1:A:20:MET:HE3	1:A:111:LEU:HA	1.86	0.58
1:H:173:LYS:HB3	3:H:635:HOH:O	2.03	0.58
1:V:20:MET:HE1	1:V:110:ALA:C	2.29	0.58
1:W:62:HIS:HB3	1:W:134:TYR:HE2	1.68	0.57
1:S:6:ARG:NH1	1:S:9:TYR:O	2.36	0.57
1:L:142:ILE:HG22	1:T:5:VAL:HG22	1.85	0.57
1:O:98:GLN:HG2	1:O:149:ILE:HD13	1.86	0.57
1:A:114:LEU:HG	1:A:130:LEU:HD21	1.86	0.57
1:C:21:LEU:HD23	1:C:21:LEU:O	2.03	0.57
1:D:152:LEU:HD13	1:D:167:PHE:CD1	2.39	0.57
1:C:43:VAL:O	1:F:154:ARG:NH2	2.28	0.57
1:E:122:VAL:HG12	1:E:122:VAL:O	2.03	0.57
1:W:29:TYR:O	1:W:32:SER:HB3	2.05	0.57
1:K:154:ARG:NH2	1:Q:45:LEU:HD13	2.20	0.57
1:B:126:LEU:O	1:B:130:LEU:HD22	2.05	0.56
1:F:142:ILE:HG22	1:O:5:VAL:HG22	1.87	0.56
1:X:152:LEU:HD12	1:X:157:LEU:HD13	1.87	0.56
1:A:20:MET:HE1	1:A:110:ALA:CB	2.35	0.56
1:R:65:LYS:HB3	1:R:134:TYR:OH	2.05	0.56
1:N:45:LEU:HD13	1:S:154:ARG:NH2	2.21	0.56
1:C:21:LEU:HD23	1:C:21:LEU:C	2.31	0.55
1:K:84:LYS:HE3	3:W:187:HOH:O	2.06	0.55
1:M:20:MET:CE	1:M:110:ALA:HB1	2.32	0.55
3:D:189:HOH:O	1:X:154:ARG:HG3	2.05	0.55
1:S:80:GLN:HA	1:S:80:GLN:NE2	2.21	0.55
1:E:45:LEU:HD13	1:W:154:ARG:NH2	2.22	0.55
1:E:162:MET:SD	1:W:162:MET:CE	2.95	0.55
1:S:84:LYS:HE2	3:S:180:HOH:O	2.06	0.55
1:B:60:ARG:HD2	3:B:179:HOH:O	2.07	0.54
1:D:60:ARG:HD3	1:H:60:ARG:HD3	1.90	0.54
1:L:2:VAL:HG22	1:L:3:SER:H	1.72	0.54
1:B:115:HIS:CE1	1:E:124:PRO:HB3	2.42	0.54
1:E:20:MET:HE1	1:E:110:ALA:HB1	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:60:ARG:HD2	1:X:60:ARG:CZ	2.38	0.54
1:V:20:MET:HE1	1:V:110:ALA:CB	2.38	0.54
1:X:157:LEU:HD21	1:X:164:GLU:HG3	1.89	0.54
1:L:48:VAL:HG23	1:L:168:ASP:OD1	2.08	0.54
1:P:21:LEU:C	1:P:21:LEU:HD23	2.33	0.54
1:X:173:LYS:HE3	1:X:173:LYS:N	2.23	0.54
1:O:173:LYS:HE3	1:O:173:LYS:HA	1.89	0.54
1:Q:64:GLU:OE2	1:U:57:HIS:CE1	2.61	0.54
1:V:148:PHE:O	1:V:152:LEU:CD2	2.54	0.53
1:K:152:LEU:HD13	1:K:167:PHE:CD1	2.44	0.53
1:N:147:ASP:OD1	1:U:41:ASP:HA	2.08	0.53
1:Q:21:LEU:HD11	1:Q:67:MET:HG3	1.89	0.53
1:V:24:GLU:OE1	1:V:59:GLU:OE1	2.27	0.53
1:H:6:ARG:NH1	1:H:9:TYR:O	2.39	0.53
1:G:5:VAL:HG22	1:O:142:ILE:HG22	1.90	0.53
1:A:57:HIS:HE1	1:G:64:GLU:OE2	1.89	0.53
1:K:21:LEU:HD23	1:K:21:LEU:C	2.34	0.53
1:T:148:PHE:O	1:T:152:LEU:HD23	2.10	0.52
1:E:142:ILE:HG22	1:V:5:VAL:HG22	1.91	0.52
1:V:20:MET:HE1	1:V:110:ALA:HB1	1.90	0.52
1:C:20:MET:HE1	1:C:110:ALA:CB	2.38	0.52
1:D:157:LEU:HD21	1:D:164:GLU:HG3	1.92	0.52
1:E:23:LEU:HD13	1:E:107:VAL:HG22	1.92	0.52
1:J:60:ARG:HD2	1:X:60:ARG:NE	2.24	0.52
1:G:6:ARG:NH1	1:G:9:TYR:O	2.42	0.52
1:T:148:PHE:O	1:T:152:LEU:CD2	2.57	0.52
1:S:20:MET:HE1	1:S:110:ALA:HB3	1.91	0.52
1:U:21:LEU:C	1:U:21:LEU:HD23	2.35	0.52
1:K:173:LYS:C	1:K:173:LYS:HE3	2.35	0.52
1:G:43:VAL:O	1:H:154:ARG:NH2	2.36	0.52
1:H:6:ARG:NH2	1:H:14:GLU:OE1	2.29	0.52
1:A:20:MET:CE	1:A:110:ALA:C	2.80	0.51
1:Q:47:ASN:HB2	1:Q:168:ASP:OD1	2.10	0.51
1:P:57:HIS:HB3	3:P:393:HOH:O	2.10	0.51
1:T:6:ARG:NH1	1:T:9:TYR:O	2.37	0.51
1:L:114:LEU:HG	1:L:130:LEU:HD21	1.91	0.51
1:C:64:GLU:OE2	1:O:57:HIS:CE1	2.63	0.51
1:V:108:ASN:O	1:V:111:LEU:HB2	2.11	0.51
1:E:79:LEU:CD1	1:I:33:SER:HB2	2.40	0.51
1:D:19:ARG:HB3	3:D:615:HOH:O	2.10	0.51
1:C:94:LEU:O	1:C:98:GLN:HG3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:24:GLU:OE1	1:O:59:GLU:OE1	2.29	0.51
1:A:43:VAL:O	1:O:154:ARG:NH2	2.39	0.51
1:L:39:ASP:HB3	1:V:71:ASN:ND2	2.26	0.51
1:F:47:ASN:HB3	1:F:172:VAL:HG12	1.93	0.50
1:D:124:PRO:HB3	1:I:115:HIS:CE1	2.47	0.50
1:F:6:ARG:NH1	1:F:9:TYR:O	2.39	0.50
1:P:114:LEU:HG	1:P:130:LEU:HD21	1.93	0.50
1:E:55:HIS:NE2	1:E:104:GLU:OE2	2.43	0.50
1:Q:71:ASN:ND2	1:U:39:ASP:HB3	2.26	0.50
1:V:60:ARG:HG3	1:V:60:ARG:HH11	1.76	0.50
1:M:115:HIS:NE2	1:M:131:GLU:OE2	2.38	0.50
1:A:60:ARG:CD	1:G:60:ARG:NE	2.62	0.50
1:C:93:THR:HG22	1:C:152:LEU:HD21	1.94	0.50
1:K:60:ARG:NH1	1:W:60:ARG:HD3	2.26	0.50
1:V:142:ILE:HG12	1:V:143:LYS:N	2.27	0.50
1:K:20:MET:HE1	1:K:110:ALA:CB	2.42	0.50
1:M:98:GLN:HG2	1:M:149:ILE:HD13	1.93	0.50
3:J:180:HOH:O	1:X:57:HIS:HB3	2.11	0.49
1:K:53:LYS:HD2	1:K:57:HIS:CE1	2.46	0.49
1:L:24:GLU:OE1	1:L:59:GLU:OE1	2.30	0.49
1:W:102:GLN:HB2	3:W:683:HOH:O	2.12	0.49
1:B:60:ARG:HD3	1:F:60:ARG:HD2	1.93	0.49
1:N:25:LEU:HB3	1:N:82:ILE:HD13	1.93	0.49
1:H:173:LYS:H	1:H:173:LYS:NZ	2.10	0.49
1:O:157:LEU:HD21	1:O:164:GLU:HG3	1.95	0.49
1:R:6:ARG:NH1	1:R:9:TYR:O	2.43	0.49
1:E:126:LEU:O	1:E:130:LEU:HD22	2.12	0.49
1:F:124:PRO:HB3	1:G:115:HIS:CE1	2.47	0.49
1:E:115:HIS:CE1	1:V:124:PRO:HG3	2.47	0.49
1:K:60:ARG:HD2	3:W:188:HOH:O	2.11	0.49
1:S:118:ALA:HB2	1:S:126:LEU:CD2	2.37	0.49
1:A:79:LEU:HD12	1:G:33:SER:HB2	1.95	0.49
1:H:45:LEU:HD13	1:I:154:ARG:NH2	2.27	0.49
1:D:24:GLU:OE1	1:D:59:GLU:OE1	2.30	0.49
1:N:65:LYS:HE2	1:N:133:GLU:OE1	2.13	0.49
1:T:114:LEU:HG	1:T:130:LEU:HD21	1.94	0.49
1:C:124:PRO:HB3	1:R:115:HIS:CE1	2.48	0.48
1:F:37:PHE:O	1:F:40:ARG:HB2	2.12	0.48
1:I:93:THR:HG22	1:I:152:LEU:HD21	1.95	0.48
1:K:57:HIS:HB3	3:W:772:HOH:O	2.12	0.48
1:N:60:ARG:CD	1:R:60:ARG:HD3	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:114:LEU:HG	1:I:130:LEU:HD21	1.95	0.48
1:N:114:LEU:HG	1:N:130:LEU:HD21	1.94	0.48
1:X:60:ARG:NH2	3:X:537:HOH:O	2.46	0.48
1:L:6:ARG:HD3	3:L:465:HOH:O	2.13	0.48
1:M:20:MET:HE2	1:M:20:MET:CA	2.29	0.48
1:B:157:LEU:HD11	1:B:164:GLU:HG3	1.95	0.48
1:H:86:GLU:HB2	3:H:189:HOH:O	2.11	0.48
1:N:35:TYR:CE1	1:N:53:LYS:HD3	2.49	0.48
1:O:95:GLU:HG3	3:O:383:HOH:O	2.13	0.48
1:Q:60:ARG:HD2	3:Q:568:HOH:O	2.13	0.48
1:S:45:LEU:HD13	1:T:154:ARG:NH2	2.29	0.48
1:J:24:GLU:HB2	1:J:63:ALA:HB2	1.96	0.48
1:L:53:LYS:HE3	1:V:64:GLU:OE2	2.14	0.48
1:V:62:HIS:HB3	1:V:134:TYR:HE2	1.79	0.48
1:K:29:TYR:O	1:K:32:SER:HB2	2.14	0.47
1:S:69:TYR:HE1	1:S:125:HIS:CE1	2.32	0.47
1:D:114:LEU:HG	1:D:130:LEU:HD21	1.96	0.47
1:O:152:LEU:HD12	1:O:157:LEU:HD13	1.96	0.47
1:L:60:ARG:CZ	1:V:60:ARG:HD3	2.44	0.47
1:N:57:HIS:HB3	3:N:734:HOH:O	2.13	0.47
1:W:65:LYS:HE2	1:W:133:GLU:OE1	2.14	0.47
1:A:154:ARG:NH2	1:J:45:LEU:HD13	2.28	0.47
1:M:60:ARG:HD2	1:S:60:ARG:NH1	2.29	0.47
1:I:85:PRO:HG2	3:I:186:HOH:O	2.15	0.47
1:S:69:TYR:CE1	1:S:125:HIS:CE1	3.02	0.47
1:T:127:CYS:SG	1:T:131:GLU:OE2	2.73	0.47
1:W:152:LEU:HD12	1:W:157:LEU:HD13	1.97	0.47
3:A:179:HOH:O	1:X:122:VAL:HG12	2.15	0.47
1:G:160:ASN:C	1:G:160:ASN:OD1	2.55	0.47
1:Q:16:ALA:HB1	1:Q:114:LEU:HD13	1.96	0.47
1:S:118:ALA:CB	1:S:126:LEU:HD23	2.35	0.47
1:L:148:PHE:O	1:L:152:LEU:HD22	2.14	0.47
1:M:157:LEU:HD11	1:M:164:GLU:HG3	1.96	0.47
1:F:135:LEU:O	1:F:139:VAL:HG13	2.15	0.47
1:G:27:ALA:HA	1:G:103:LEU:HD21	1.96	0.47
1:I:24:GLU:OE1	1:I:59:GLU:OE1	2.33	0.46
1:K:6:ARG:NH1	1:K:9:TYR:O	2.38	0.46
1:R:101:LEU:HD22	1:R:149:ILE:HD12	1.96	0.46
1:G:2:VAL:N	3:G:190:HOH:O	2.48	0.46
1:J:67:MET:HE1	1:X:32:SER:OG	2.15	0.46
1:E:165:TYR:CD2	1:W:155:LEU:HD21	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:W:16:ALA:HB1	1:W:114:LEU:HD13	1.97	0.46
1:K:79:LEU:HD12	1:W:33:SER:HB2	1.96	0.46
1:K:124:PRO:HB3	1:P:115:HIS:CE1	2.51	0.46
3:Q:178:HOH:O	1:U:4:GLN:HG2	2.15	0.46
1:S:46:HIS:O	1:S:49:ALA:HB3	2.16	0.46
1:B:60:ARG:NE	1:F:60:ARG:HD2	2.30	0.46
1:O:57:HIS:HB3	3:O:188:HOH:O	2.16	0.46
1:B:6:ARG:NH1	1:B:9:TYR:O	2.43	0.46
1:B:79:LEU:HD12	1:F:33:SER:HB2	1.96	0.46
1:I:155:LEU:HD22	1:I:163:GLY:HA2	1.98	0.46
1:M:10:HIS:CD2	1:M:121:LYS:HE3	2.51	0.46
1:R:135:LEU:HD12	1:R:135:LEU:HA	1.84	0.46
1:G:140:LYS:HB2	1:G:140:LYS:HE3	1.69	0.45
1:N:20:MET:HE3	1:N:111:LEU:HD23	1.97	0.45
1:O:114:LEU:HG	1:O:130:LEU:HD21	1.98	0.45
1:U:57:HIS:HB3	3:U:287:HOH:O	2.15	0.45
1:U:114:LEU:HD23	1:U:130:LEU:HD11	1.98	0.45
1:W:37:PHE:O	1:W:40:ARG:HB2	2.15	0.45
1:N:131:GLU:HG2	1:Q:128:ASP:HB2	1.98	0.45
1:Q:33:SER:HB2	1:U:79:LEU:HD12	1.98	0.45
1:G:148:PHE:O	1:G:152:LEU:CD2	2.62	0.45
1:D:20:MET:HE1	1:D:110:ALA:CB	2.40	0.45
1:H:41:ASP:HA	1:I:147:ASP:OD1	2.17	0.45
1:J:154:ARG:NH2	1:R:45:LEU:CD1	2.79	0.45
1:K:131:GLU:HB3	1:U:128:ASP:HB2	1.99	0.45
1:V:127:CYS:SG	1:V:131:GLU:OE2	2.74	0.45
1:M:75:GLY:O	1:M:76:ARG:NH1	2.47	0.45
1:T:42:ASP:HA	1:U:150:THR:OG1	2.17	0.45
1:D:60:ARG:NH1	1:H:60:ARG:HD2	2.31	0.45
1:I:21:LEU:HD23	1:I:21:LEU:O	2.16	0.45
1:L:20:MET:HE1	1:L:111:LEU:N	2.31	0.45
3:N:177:HOH:O	1:R:60:ARG:HD2	2.16	0.45
1:Q:21:LEU:C	1:Q:21:LEU:HD23	2.41	0.45
1:C:20:MET:HE2	1:C:20:MET:HB2	1.69	0.45
1:I:152:LEU:HD13	1:I:167:PHE:CD1	2.52	0.45
1:L:124:PRO:HB3	1:M:115:HIS:CE1	2.52	0.45
1:P:21:LEU:HD11	1:P:67:MET:HG2	1.98	0.45
1:E:115:HIS:ND1	1:V:124:PRO:HG3	2.32	0.45
1:J:16:ALA:HB1	1:J:114:LEU:HD13	1.99	0.45
1:F:10:HIS:CG	1:F:121:LYS:HD2	2.52	0.44
1:L:118:ALA:HB2	1:L:126:LEU:HD23	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:173:LYS:CD	1:P:173:LYS:N	2.67	0.44
1:R:73:ARG:HD2	1:R:73:ARG:HA	1.77	0.44
1:K:60:ARG:CZ	1:W:60:ARG:HD3	2.48	0.44
1:N:79:LEU:HD12	1:R:33:SER:HB2	1.99	0.44
1:T:80:GLN:O	1:T:81:ASP:C	2.59	0.44
1:C:21:LEU:C	1:C:21:LEU:CD2	2.90	0.44
1:C:144:ARG:HD2	3:C:177:HOH:O	2.16	0.44
1:E:24:GLU:OE1	1:E:59:GLU:OE1	2.34	0.44
1:F:16:ALA:HB1	1:F:114:LEU:HD13	1.99	0.44
1:W:114:LEU:HG	1:W:130:LEU:HD21	1.99	0.44
1:A:60:ARG:HD3	1:G:60:ARG:CZ	2.44	0.44
1:H:69:TYR:CE2	1:H:126:LEU:HD13	2.52	0.44
1:P:79:LEU:HD13	1:T:29:TYR:CE1	2.52	0.44
1:L:112:LEU:HD23	1:L:112:LEU:HA	1.84	0.44
1:T:75:GLY:O	1:T:76:ARG:NH1	2.51	0.44
1:A:64:GLU:OE1	1:G:57:HIS:CE1	2.69	0.44
1:I:5:VAL:HG22	1:W:142:ILE:HG22	1.98	0.44
1:C:25:LEU:HD23	1:C:25:LEU:HA	1.85	0.44
1:C:173:LYS:HB3	1:C:173:LYS:HZ2	1.82	0.44
1:Q:140:LYS:HB2	1:Q:140:LYS:HE3	1.72	0.44
1:S:114:LEU:HG	1:S:130:LEU:HD21	1.98	0.44
1:V:23:LEU:O	1:V:26:TYR:HB3	2.17	0.44
1:B:162:MET:HE3	1:I:162:MET:SD	2.58	0.44
1:D:60:ARG:HD3	1:H:60:ARG:CD	2.48	0.44
1:N:20:MET:HE1	1:N:110:ALA:CB	2.47	0.44
1:B:154:ARG:NH2	1:I:45:LEU:HD13	2.33	0.43
1:C:60:ARG:HD3	1:O:60:ARG:CD	2.36	0.43
1:V:46:HIS:CE1	1:V:50:GLU:OE2	2.71	0.43
1:C:116:LYS:HD3	1:S:122:VAL:HG11	2.00	0.43
1:E:79:LEU:HD12	1:I:33:SER:HB2	2.00	0.43
1:H:124:PRO:HB3	1:X:115:HIS:CE1	2.53	0.43
1:J:124:PRO:HB3	1:Q:115:HIS:CE1	2.53	0.43
1:C:57:HIS:HE1	1:O:64:GLU:OE1	2.00	0.43
1:M:20:MET:HE1	1:M:110:ALA:HB3	1.96	0.43
1:K:35:TYR:CD1	1:K:53:LYS:HB2	2.53	0.43
1:M:6:ARG:NH1	1:M:9:TYR:O	2.38	0.43
1:N:115:HIS:CE1	1:Q:124:PRO:HB3	2.54	0.43
1:O:47:ASN:HB3	1:O:172:VAL:HG23	1.99	0.43
1:P:33:SER:HB2	1:T:79:LEU:HD12	2.01	0.43
1:B:21:LEU:C	1:B:21:LEU:CD2	2.92	0.43
1:M:125:HIS:CD2	1:T:139:VAL:HG13	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:145:ILE:O	1:M:149:ILE:HG13	2.17	0.43
1:I:21:LEU:C	1:I:21:LEU:CD2	2.92	0.43
1:E:16:ALA:HB1	1:E:114:LEU:HD13	2.00	0.43
1:H:73:ARG:HD2	1:H:73:ARG:HA	1.50	0.43
1:T:162:MET:HE3	1:T:162:MET:HB2	1.82	0.43
1:D:136:GLU:OE2	1:W:72:LYS:NZ	2.52	0.43
1:R:124:PRO:HB3	1:S:115:HIS:CE1	2.54	0.43
1:N:80:GLN:O	1:N:81:ASP:C	2.61	0.43
1:U:73:ARG:HA	1:U:73:ARG:HD2	1.59	0.43
1:B:24:GLU:OE1	1:B:59:GLU:OE1	2.37	0.43
1:B:32:SER:HB3	1:F:67:MET:HE1	2.00	0.43
1:K:5:VAL:HG22	1:P:142:ILE:HG22	2.00	0.43
1:K:127:CYS:SG	1:P:131:GLU:OE1	2.77	0.43
1:U:20:MET:O	1:U:24:GLU:HG2	2.19	0.43
1:W:21:LEU:HD11	1:W:67:MET:HG2	2.00	0.43
1:C:37:PHE:CD2	1:C:90:TRP:HB2	2.54	0.42
1:E:20:MET:HE1	1:E:110:ALA:O	2.18	0.42
1:F:130:LEU:HD12	1:F:130:LEU:HA	1.86	0.42
1:G:130:LEU:HD12	1:G:130:LEU:HA	1.88	0.42
1:M:125:HIS:CD2	1:T:139:VAL:CG1	3.02	0.42
1:N:73:ARG:HA	1:N:73:ARG:HD2	1.76	0.42
1:O:86:GLU:HG2	3:O:709:HOH:O	2.18	0.42
1:E:168:ASP:OD2	1:E:169:LYS:HD3	2.18	0.42
1:H:173:LYS:H	1:H:173:LYS:HZ3	1.67	0.42
1:S:160:ASN:OD1	1:S:160:ASN:C	2.62	0.42
1:A:165:TYR:CD2	1:O:155:LEU:HD21	2.54	0.42
1:P:152:LEU:HD13	1:P:167:PHE:CD1	2.53	0.42
1:A:21:LEU:HD23	1:A:21:LEU:C	2.45	0.42
1:B:153:LYS:O	1:B:154:ARG:C	2.63	0.42
1:J:33:SER:HB2	1:X:79:LEU:HD12	2.01	0.42
1:L:107:VAL:O	1:L:108:ASN:C	2.59	0.42
1:O:60:ARG:NH1	3:O:178:HOH:O	2.52	0.42
1:F:16:ALA:CB	1:F:114:LEU:HD13	2.50	0.42
1:P:60:ARG:CD	1:T:60:ARG:HD3	2.49	0.42
1:F:23:LEU:CD1	1:F:107:VAL:HG22	2.48	0.42
1:C:135:LEU:HD21	1:S:124:PRO:HB2	2.01	0.42
1:W:4:GLN:HB2	3:W:363:HOH:O	2.19	0.42
1:A:115:HIS:CE1	1:X:124:PRO:HB3	2.55	0.42
1:E:52:PHE:CZ	1:E:97:MET:HE2	2.55	0.42
1:K:30:THR:OG1	1:K:85:PRO:HB3	2.19	0.42
1:L:8:ASN:CG	1:L:8:ASN:O	2.63	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:86:GLU:HB2	3:M:186:HOH:O	2.19	0.42
1:E:20:MET:HE1	1:E:110:ALA:CB	2.50	0.41
1:E:124:PRO:HA	1:E:127:CYS:HB3	2.01	0.41
1:B:9:TYR:CE2	1:B:14:GLU:HB2	2.55	0.41
1:B:144:ARG:O	1:B:147:ASP:HB2	2.20	0.41
1:H:25:LEU:HD23	1:H:25:LEU:HA	1.83	0.41
1:J:41:ASP:OD2	1:X:4:GLN:HG2	2.20	0.41
1:P:80:GLN:O	1:P:81:ASP:C	2.62	0.41
1:I:35:TYR:CD1	1:I:53:LYS:HB2	2.55	0.41
1:K:33:SER:HB2	1:W:79:LEU:HD12	2.02	0.41
1:M:154:ARG:NH2	1:V:45:LEU:HD13	2.34	0.41
1:M:162:MET:O	1:M:165:TYR:HB3	2.19	0.41
1:W:101:LEU:O	1:W:105:LYS:HG3	2.19	0.41
1:K:84:LYS:HD2	3:W:518:HOH:O	2.20	0.41
1:N:60:ARG:NE	1:R:60:ARG:HH11	2.19	0.41
1:O:73:ARG:HA	1:O:73:ARG:HD2	1.69	0.41
1:P:39:ASP:HB3	1:T:71:ASN:ND2	2.35	0.41
1:H:145:ILE:O	1:H:146:GLY:C	2.64	0.41
1:J:51:PHE:CD2	1:J:51:PHE:C	2.97	0.41
1:J:148:PHE:O	1:J:149:ILE:C	2.62	0.41
1:M:173:LYS:H	1:M:173:LYS:NZ	2.18	0.41
1:X:37:PHE:CE1	1:X:89:GLU:HB2	2.55	0.41
1:E:75:GLY:O	1:E:76:ARG:NH1	2.53	0.41
1:H:47:ASN:O	1:H:172:VAL:HG11	2.21	0.41
1:J:114:LEU:HG	1:J:130:LEU:HD21	2.02	0.41
1:K:21:LEU:HD11	1:K:67:MET:HG2	2.03	0.41
1:K:73:ARG:HD2	1:K:73:ARG:HA	1.89	0.41
1:L:24:GLU:OE1	1:L:24:GLU:HA	2.21	0.41
1:M:69:TYR:O	1:M:70:GLN:C	2.63	0.41
1:V:10:HIS:CD2	1:V:121:LYS:HE3	2.54	0.41
1:K:47:ASN:HB3	1:K:172:VAL:HG13	2.02	0.41
1:W:23:LEU:HD23	1:W:23:LEU:HA	1.94	0.41
1:W:135:LEU:HD12	1:W:135:LEU:HA	1.90	0.41
1:M:60:ARG:NE	1:S:60:ARG:HD2	2.35	0.40
1:Q:24:GLU:OE1	1:Q:59:GLU:OE1	2.38	0.40
1:S:69:TYR:CE1	1:S:125:HIS:HE1	2.39	0.40
1:T:172:VAL:O	1:T:173:LYS:C	2.64	0.40
1:T:27:ALA:O	1:T:28:SER:C	2.64	0.40
1:X:172:VAL:O	1:X:173:LYS:C	2.64	0.40
1:K:60:ARG:NH1	1:W:60:ARG:CD	2.85	0.40
1:V:135:LEU:HD12	1:V:135:LEU:HA	1.92	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:W:62:HIS:O	1:W:63:ALA:C	2.62	0.40
1:A:8:ASN:O	1:A:8:ASN:CG	2.64	0.40
1:G:21:LEU:HD11	1:G:67:MET:HG2	2.03	0.40
1:K:20:MET:HG3	1:K:66:PHE:CE2	2.57	0.40
1:T:24:GLU:OE1	1:T:59:GLU:OE1	2.39	0.40
1:U:101:LEU:O	1:U:105:LYS:HG3	2.21	0.40
1:B:73:ARG:HA	1:B:73:ARG:HD2	1.66	0.40
1:V:6:ARG:NH1	1:V:9:TYR:O	2.48	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	170/176 (97%)	166 (98%)	4 (2%)	0	100	100
1	B	170/176 (97%)	164 (96%)	6 (4%)	0	100	100
1	C	170/176 (97%)	165 (97%)	5 (3%)	0	100	100
1	D	170/176 (97%)	159 (94%)	10 (6%)	1 (1%)	21	48
1	E	170/176 (97%)	164 (96%)	6 (4%)	0	100	100
1	F	170/176 (97%)	166 (98%)	4 (2%)	0	100	100
1	G	170/176 (97%)	164 (96%)	6 (4%)	0	100	100
1	H	170/176 (97%)	166 (98%)	4 (2%)	0	100	100
1	I	170/176 (97%)	164 (96%)	6 (4%)	0	100	100
1	J	170/176 (97%)	166 (98%)	3 (2%)	1 (1%)	21	48
1	K	170/176 (97%)	165 (97%)	5 (3%)	0	100	100
1	L	170/176 (97%)	162 (95%)	8 (5%)	0	100	100
1	M	170/176 (97%)	159 (94%)	10 (6%)	1 (1%)	21	48

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	N	170/176 (97%)	164 (96%)	6 (4%)	0	100	100
1	O	170/176 (97%)	164 (96%)	5 (3%)	1 (1%)	21	48
1	P	170/176 (97%)	163 (96%)	7 (4%)	0	100	100
1	Q	170/176 (97%)	163 (96%)	7 (4%)	0	100	100
1	R	170/176 (97%)	163 (96%)	6 (4%)	1 (1%)	21	48
1	S	170/176 (97%)	167 (98%)	3 (2%)	0	100	100
1	T	170/176 (97%)	164 (96%)	6 (4%)	0	100	100
1	U	170/176 (97%)	164 (96%)	6 (4%)	0	100	100
1	V	170/176 (97%)	163 (96%)	5 (3%)	2 (1%)	10	31
1	W	170/176 (97%)	166 (98%)	3 (2%)	1 (1%)	21	48
1	X	170/176 (97%)	166 (98%)	4 (2%)	0	100	100
All	All	4080/4224 (97%)	3937 (96%)	135 (3%)	8 (0%)	43	71

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	M	122	VAL
1	D	81	ASP
1	J	122	VAL
1	V	11	SER
1	O	81	ASP
1	V	122	VAL
1	R	122	VAL
1	W	122	VAL

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	153/157 (98%)	135 (88%)	18 (12%)	5	16
1	B	153/157 (98%)	133 (87%)	20 (13%)	4	13

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	153/157 (98%)	131 (86%)	22 (14%)	3	10
1	D	153/157 (98%)	137 (90%)	16 (10%)	6	21
1	E	153/157 (98%)	131 (86%)	22 (14%)	3	10
1	F	153/157 (98%)	132 (86%)	21 (14%)	3	12
1	G	153/157 (98%)	135 (88%)	18 (12%)	5	16
1	H	153/157 (98%)	134 (88%)	19 (12%)	4	15
1	I	153/157 (98%)	132 (86%)	21 (14%)	3	12
1	J	153/157 (98%)	133 (87%)	20 (13%)	4	13
1	K	153/157 (98%)	134 (88%)	19 (12%)	4	15
1	L	153/157 (98%)	133 (87%)	20 (13%)	4	13
1	M	153/157 (98%)	134 (88%)	19 (12%)	4	15
1	N	153/157 (98%)	134 (88%)	19 (12%)	4	15
1	O	153/157 (98%)	134 (88%)	19 (12%)	4	15
1	P	153/157 (98%)	134 (88%)	19 (12%)	4	15
1	Q	153/157 (98%)	133 (87%)	20 (13%)	4	13
1	R	153/157 (98%)	133 (87%)	20 (13%)	4	13
1	S	153/157 (98%)	134 (88%)	19 (12%)	4	15
1	T	153/157 (98%)	136 (89%)	17 (11%)	6	19
1	U	153/157 (98%)	133 (87%)	20 (13%)	4	13
1	V	153/157 (98%)	138 (90%)	15 (10%)	7	24
1	W	153/157 (98%)	138 (90%)	15 (10%)	7	24
1	X	153/157 (98%)	132 (86%)	21 (14%)	3	12
All	All	3672/3768 (98%)	3213 (88%)	459 (12%)	4	15

All (459) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	VAL
1	A	23	LEU
1	A	32	SER
1	A	45	LEU
1	A	54	GLU
1	A	68	LYS
1	A	89	GLU

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Mol	Chain	Res	Type
1	A	98	GLN
1	A	105	LYS
1	A	116	LYS
1	A	121	LYS
1	A	127	CYS
1	A	130	LEU
1	A	132	SER
1	A	139	VAL
1	A	152	LEU
1	A	171	SER
1	A	173	LYS
1	B	5	VAL
1	B	23	LEU
1	B	29	TYR
1	B	45	LEU
1	B	54	GLU
1	B	64	GLU
1	B	68	LYS
1	B	89	GLU
1	B	98	GLN
1	B	116	LYS
1	B	117	LEU
1	B	121	LYS
1	B	122	VAL
1	B	130	LEU
1	B	135	LEU
1	B	136	GLU
1	B	139	VAL
1	B	152	LEU
1	B	157	LEU
1	B	173	LYS
1	C	5	VAL
1	C	11	SER
1	C	23	LEU
1	C	29	TYR
1	C	32	SER
1	C	45	LEU
1	C	53	LYS
1	C	54	GLU
1	C	84	LYS
1	C	122	VAL
1	C	127	CYS

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Mol	Chain	Res	Type
1	C	128	ASP
1	C	130	LEU
1	C	131	GLU
1	C	135	LEU
1	C	137	GLU
1	C	139	VAL
1	C	144	ARG
1	C	152	LEU
1	C	159	GLU
1	C	172	VAL
1	C	173	LYS
1	D	5	VAL
1	D	11	SER
1	D	21	LEU
1	D	23	LEU
1	D	45	LEU
1	D	60	ARG
1	D	65	LYS
1	D	68	LYS
1	D	89	GLU
1	D	130	LEU
1	D	135	LEU
1	D	139	VAL
1	D	140	LYS
1	D	144	ARG
1	D	152	LEU
1	D	172	VAL
1	E	2	VAL
1	E	5	VAL
1	E	17	VAL
1	E	21	LEU
1	E	23	LEU
1	E	28	SER
1	E	32	SER
1	E	45	LEU
1	E	54	GLU
1	E	55	HIS
1	E	60	ARG
1	E	68	LYS
1	E	84	LYS
1	E	121	LYS
1	E	130	LEU

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Mol	Chain	Res	Type
1	E	135	LEU
1	E	137	GLU
1	E	139	VAL
1	E	140	LYS
1	E	152	LEU
1	E	157	LEU
1	E	169	LYS
1	F	5	VAL
1	F	23	LEU
1	F	32	SER
1	F	45	LEU
1	F	54	GLU
1	F	55	HIS
1	F	61	GLU
1	F	68	LYS
1	F	86	GLU
1	F	102	GLN
1	F	119	THR
1	F	122	VAL
1	F	127	CYS
1	F	130	LEU
1	F	135	LEU
1	F	137	GLU
1	F	139	VAL
1	F	140	LYS
1	F	152	LEU
1	F	154	ARG
1	F	172	VAL
1	G	5	VAL
1	G	11	SER
1	G	23	LEU
1	G	45	LEU
1	G	54	GLU
1	G	55	HIS
1	G	61	GLU
1	G	68	LYS
1	G	83	LYS
1	G	86	GLU
1	G	88	ASP
1	G	98	GLN
1	G	130	LEU
1	G	135	LEU

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Mol	Chain	Res	Type
1	G	137	GLU
1	G	139	VAL
1	G	152	LEU
1	G	157	LEU
1	H	2	VAL
1	H	3	SER
1	H	5	VAL
1	H	6	ARG
1	H	23	LEU
1	H	45	LEU
1	H	64	GLU
1	H	83	LYS
1	H	89	GLU
1	H	116	LYS
1	H	121	LYS
1	H	122	VAL
1	H	127	CYS
1	H	130	LEU
1	H	135	LEU
1	H	139	VAL
1	H	152	LEU
1	H	171	SER
1	H	173	LYS
1	I	2	VAL
1	I	5	VAL
1	I	23	LEU
1	I	33	SER
1	I	45	LEU
1	I	54	GLU
1	I	89	GLU
1	I	98	GLN
1	I	116	LYS
1	I	121	LYS
1	I	122	VAL
1	I	130	LEU
1	I	135	LEU
1	I	137	GLU
1	I	139	VAL
1	I	144	ARG
1	I	152	LEU
1	I	157	LEU
1	I	171	SER

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Mol	Chain	Res	Type
1	I	172	VAL
1	I	173	LYS
1	J	5	VAL
1	J	11	SER
1	J	21	LEU
1	J	23	LEU
1	J	45	LEU
1	J	84	LYS
1	J	86	GLU
1	J	102	GLN
1	J	116	LYS
1	J	121	LYS
1	J	122	VAL
1	J	127	CYS
1	J	130	LEU
1	J	135	LEU
1	J	137	GLU
1	J	139	VAL
1	J	140	LYS
1	J	152	LEU
1	J	157	LEU
1	J	173	LYS
1	K	2	VAL
1	K	5	VAL
1	K	19	ARG
1	K	21	LEU
1	K	23	LEU
1	K	45	LEU
1	K	53	LYS
1	K	55	HIS
1	K	58	GLU
1	K	102	GLN
1	K	117	LEU
1	K	119	THR
1	K	127	CYS
1	K	130	LEU
1	K	135	LEU
1	K	139	VAL
1	K	152	LEU
1	K	168	ASP
1	K	173	LYS
1	L	2	VAL

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Mol	Chain	Res	Type
1	L	5	VAL
1	L	23	LEU
1	L	45	LEU
1	L	60	ARG
1	L	84	LYS
1	L	102	GLN
1	L	116	LYS
1	L	117	LEU
1	L	122	VAL
1	L	127	CYS
1	L	130	LEU
1	L	135	LEU
1	L	137	GLU
1	L	139	VAL
1	L	140	LYS
1	L	152	LEU
1	L	157	LEU
1	L	171	SER
1	L	172	VAL
1	M	2	VAL
1	M	5	VAL
1	M	11	SER
1	M	17	VAL
1	M	23	LEU
1	M	45	LEU
1	M	55	HIS
1	M	65	LYS
1	M	86	GLU
1	M	116	LYS
1	M	122	VAL
1	M	130	LEU
1	M	132	SER
1	M	137	GLU
1	M	139	VAL
1	M	152	LEU
1	M	159	GLU
1	M	172	VAL
1	M	173	LYS
1	N	5	VAL
1	N	23	LEU
1	N	28	SER
1	N	32	SER

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Mol	Chain	Res	Type
1	N	45	LEU
1	N	53	LYS
1	N	65	LYS
1	N	72	LYS
1	N	83	LYS
1	N	102	GLN
1	N	116	LYS
1	N	121	LYS
1	N	122	VAL
1	N	130	LEU
1	N	135	LEU
1	N	137	GLU
1	N	139	VAL
1	N	152	LEU
1	N	173	LYS
1	O	5	VAL
1	O	11	SER
1	O	17	VAL
1	O	23	LEU
1	O	32	SER
1	O	45	LEU
1	O	68	LYS
1	O	83	LYS
1	O	86	GLU
1	O	105	LYS
1	O	106	THR
1	O	122	VAL
1	O	127	CYS
1	O	130	LEU
1	O	135	LEU
1	O	139	VAL
1	O	152	LEU
1	O	157	LEU
1	O	173	LYS
1	P	2	VAL
1	P	5	VAL
1	P	23	LEU
1	P	45	LEU
1	P	72	LYS
1	P	86	GLU
1	P	89	GLU
1	P	98	GLN

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Mol	Chain	Res	Type
1	P	102	GLN
1	P	116	LYS
1	P	121	LYS
1	P	127	CYS
1	P	130	LEU
1	P	135	LEU
1	P	139	VAL
1	P	140	LYS
1	P	152	LEU
1	P	157	LEU
1	P	173	LYS
1	Q	5	VAL
1	Q	11	SER
1	Q	23	LEU
1	Q	32	SER
1	Q	45	LEU
1	Q	53	LYS
1	Q	65	LYS
1	Q	83	LYS
1	Q	84	LYS
1	Q	86	GLU
1	Q	89	GLU
1	Q	98	GLN
1	Q	116	LYS
1	Q	122	VAL
1	Q	127	CYS
1	Q	130	LEU
1	Q	135	LEU
1	Q	137	GLU
1	Q	152	LEU
1	Q	172	VAL
1	R	2	VAL
1	R	5	VAL
1	R	21	LEU
1	R	23	LEU
1	R	32	SER
1	R	45	LEU
1	R	53	LYS
1	R	65	LYS
1	R	98	GLN
1	R	102	GLN
1	R	119	THR

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Mol	Chain	Res	Type
1	R	122	VAL
1	R	127	CYS
1	R	130	LEU
1	R	135	LEU
1	R	139	VAL
1	R	152	LEU
1	R	169	LYS
1	R	172	VAL
1	R	173	LYS
1	S	2	VAL
1	S	5	VAL
1	S	21	LEU
1	S	23	LEU
1	S	28	SER
1	S	29	TYR
1	S	45	LEU
1	S	68	LYS
1	S	72	LYS
1	S	84	LYS
1	S	98	GLN
1	S	119	THR
1	S	130	LEU
1	S	135	LEU
1	S	139	VAL
1	S	144	ARG
1	S	152	LEU
1	S	157	LEU
1	S	173	LYS
1	T	2	VAL
1	T	5	VAL
1	T	11	SER
1	T	23	LEU
1	T	45	LEU
1	T	65	LYS
1	T	98	GLN
1	T	109	GLN
1	T	121	LYS
1	T	122	VAL
1	T	130	LEU
1	T	135	LEU
1	T	139	VAL
1	T	152	LEU

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Mol	Chain	Res	Type
1	T	154	ARG
1	T	157	LEU
1	T	169	LYS
1	U	2	VAL
1	U	5	VAL
1	U	11	SER
1	U	23	LEU
1	U	45	LEU
1	U	68	LYS
1	U	84	LYS
1	U	86	GLU
1	U	116	LYS
1	U	122	VAL
1	U	127	CYS
1	U	130	LEU
1	U	131	GLU
1	U	135	LEU
1	U	137	GLU
1	U	139	VAL
1	U	144	ARG
1	U	152	LEU
1	U	157	LEU
1	U	173	LYS
1	V	5	VAL
1	V	11	SER
1	V	23	LEU
1	V	30	THR
1	V	45	LEU
1	V	60	ARG
1	V	68	LYS
1	V	130	LEU
1	V	135	LEU
1	V	139	VAL
1	V	140	LYS
1	V	142	ILE
1	V	152	LEU
1	V	157	LEU
1	V	173	LYS
1	W	5	VAL
1	W	23	LEU
1	W	45	LEU
1	W	53	LYS

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Mol	Chain	Res	Type
1	W	68	LYS
1	W	84	LYS
1	W	117	LEU
1	W	122	VAL
1	W	130	LEU
1	W	135	LEU
1	W	139	VAL
1	W	140	LYS
1	W	152	LEU
1	W	157	LEU
1	W	171	SER
1	X	2	VAL
1	X	5	VAL
1	X	23	LEU
1	X	45	LEU
1	X	53	LYS
1	X	54	GLU
1	X	61	GLU
1	X	64	GLU
1	X	65	LYS
1	X	72	LYS
1	X	116	LYS
1	X	121	LYS
1	X	122	VAL
1	X	130	LEU
1	X	135	LEU
1	X	137	GLU
1	X	139	VAL
1	X	152	LEU
1	X	157	LEU
1	X	172	VAL
1	X	173	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (82) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	GLN
1	A	98	GLN
1	B	62	HIS
1	B	70	GLN
1	B	138	GLN
1	C	22	ASN

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Mol	Chain	Res	Type
1	C	57	HIS
1	C	71	ASN
1	C	108	ASN
1	C	138	GLN
1	C	151	ASN
1	E	62	HIS
1	E	71	ASN
1	E	98	GLN
1	E	108	ASN
1	E	138	GLN
1	F	46	HIS
1	F	62	HIS
1	F	98	GLN
1	F	151	ASN
1	G	57	HIS
1	G	98	GLN
1	G	102	GLN
1	G	151	ASN
1	G	170	HIS
1	H	10	HIS
1	H	46	HIS
1	H	80	GLN
1	H	98	GLN
1	H	138	GLN
1	I	22	ASN
1	I	57	HIS
1	I	62	HIS
1	I	98	GLN
1	J	46	HIS
1	J	57	HIS
1	J	138	GLN
1	K	46	HIS
1	K	57	HIS
1	K	80	GLN
1	L	98	GLN
1	L	138	GLN
1	M	4	GLN
1	M	22	ASN
1	M	46	HIS
1	M	138	GLN
1	N	4	GLN
1	N	98	GLN

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Mol	Chain	Res	Type
1	O	46	HIS
1	O	102	GLN
1	O	151	ASN
1	P	46	HIS
1	P	55	HIS
1	P	71	ASN
1	P	98	GLN
1	Q	98	GLN
1	R	4	GLN
1	R	46	HIS
1	R	98	GLN
1	S	4	GLN
1	S	80	GLN
1	S	98	GLN
1	T	46	HIS
1	T	80	GLN
1	T	98	GLN
1	T	151	ASN
1	U	46	HIS
1	U	71	ASN
1	U	98	GLN
1	U	138	GLN
1	U	151	ASN
1	V	46	HIS
1	V	80	GLN
1	V	108	ASN
1	W	57	HIS
1	W	80	GLN
1	W	98	GLN
1	W	108	ASN
1	W	138	GLN
1	X	46	HIS
1	X	57	HIS
1	X	62	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 162 ligands modelled in this entry, 162 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	172/176 (97%)	-0.88	0 100 100	22, 32, 46, 54	0
1	B	172/176 (97%)	-0.69	1 (0%) 85 80	31, 40, 54, 66	0
1	C	172/176 (97%)	-0.82	1 (0%) 85 80	24, 35, 47, 64	0
1	D	172/176 (97%)	-0.87	1 (0%) 85 80	21, 31, 48, 57	0
1	E	172/176 (97%)	-0.69	1 (0%) 85 80	28, 39, 53, 60	0
1	F	172/176 (97%)	-0.65	1 (0%) 85 80	27, 38, 50, 68	0
1	G	172/176 (97%)	-0.77	1 (0%) 85 80	27, 36, 47, 68	0
1	H	172/176 (97%)	-0.83	1 (0%) 85 80	22, 33, 46, 61	0
1	I	172/176 (97%)	-0.72	2 (1%) 76 68	30, 39, 51, 66	0
1	J	172/176 (97%)	-0.93	1 (0%) 85 80	19, 29, 44, 63	0
1	K	172/176 (97%)	-1.01	0 100 100	21, 29, 43, 54	0
1	L	172/176 (97%)	-0.62	1 (0%) 85 80	30, 40, 52, 62	0
1	M	172/176 (97%)	-0.61	1 (0%) 85 80	30, 40, 52, 61	0
1	N	172/176 (97%)	-0.95	1 (0%) 85 80	18, 30, 45, 60	0
1	O	172/176 (97%)	-0.87	2 (1%) 76 68	20, 31, 45, 57	0
1	P	172/176 (97%)	-0.80	1 (0%) 85 80	26, 35, 47, 62	0
1	Q	172/176 (97%)	-0.94	2 (1%) 76 68	21, 30, 43, 55	0
1	R	172/176 (97%)	-0.90	0 100 100	19, 31, 46, 61	0
1	S	172/176 (97%)	-0.70	2 (1%) 76 68	27, 37, 50, 65	0
1	T	172/176 (97%)	-0.67	2 (1%) 76 68	28, 37, 51, 65	0
1	U	172/176 (97%)	-0.88	1 (0%) 85 80	22, 32, 45, 65	0
1	V	172/176 (97%)	-0.58	0 100 100	31, 42, 55, 66	0
1	W	172/176 (97%)	-0.89	0 100 100	23, 32, 45, 56	0
1	X	172/176 (97%)	-0.95	0 100 100	20, 29, 43, 57	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
All	All	4128/4224 (97%)	-0.80	23 (0%) 85 80	18, 35, 50, 68	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	T	173	LYS	3.7
1	S	173	LYS	3.3
1	N	173	LYS	3.2
1	I	173	LYS	3.1
1	F	173	LYS	2.8
1	U	173	LYS	2.8
1	I	2	VAL	2.8
1	Q	2	VAL	2.7
1	M	173	LYS	2.7
1	B	173	LYS	2.6
1	S	2	VAL	2.5
1	L	2	VAL	2.5
1	O	173	LYS	2.4
1	C	173	LYS	2.4
1	G	173	LYS	2.4
1	T	2	VAL	2.4
1	J	173	LYS	2.4
1	O	2	VAL	2.3
1	Q	173	LYS	2.3
1	P	173	LYS	2.2
1	D	2	VAL	2.2
1	E	2	VAL	2.1
1	H	173	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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
2	CU	H	192	1/1	0.58	0.19	69,69,69,69	1
2	CU	R	192	1/1	0.64	0.22	78,78,78,78	1
2	CU	K	177	1/1	0.68	0.39	66,66,66,66	1
2	CU	B	184	1/1	0.70	0.45	75,75,75,75	1
2	CU	U	192	1/1	0.70	0.28	72,72,72,72	1
2	CU	E	192	1/1	0.73	0.14	71,71,71,71	1
2	CU	S	192	1/1	0.74	0.17	74,74,74,74	1
2	CU	D	184	1/1	0.75	0.33	65,65,65,65	1
2	CU	E	184	1/1	0.75	0.38	59,59,59,59	1
2	CU	P	184	1/1	0.77	0.23	63,63,63,63	1
2	CU	V	181	1/1	0.78	0.20	66,66,66,66	1
2	CU	W	184	1/1	0.78	0.26	66,66,66,66	1
2	CU	J	177	1/1	0.80	0.23	65,65,65,65	1
2	CU	L	192	1/1	0.80	0.16	73,73,73,73	1
2	CU	C	185	1/1	0.81	0.19	59,59,59,59	1
2	CU	K	192	1/1	0.81	0.25	56,56,56,56	1
2	CU	N	192	1/1	0.81	0.14	67,67,67,67	1
2	CU	F	184	1/1	0.82	0.25	57,57,57,57	1
2	CU	H	177	1/1	0.83	0.16	44,44,44,44	1
2	CU	R	185	1/1	0.83	0.15	57,57,57,57	1
2	CU	I	184	1/1	0.83	0.34	59,59,59,59	1
2	CU	Q	184	1/1	0.84	0.32	57,57,57,57	1
2	CU	G	184	1/1	0.85	0.16	64,64,64,64	1
2	CU	W	183	1/1	0.86	0.21	61,61,61,61	1
2	CU	T	191	1/1	0.86	0.16	69,69,69,69	1
2	CU	G	192	1/1	0.87	0.14	61,61,61,61	1
2	CU	D	178	1/1	0.87	0.14	61,61,61,61	1
2	CU	X	184	1/1	0.87	0.30	58,58,58,58	1
2	CU	R	184	1/1	0.88	0.14	64,64,64,64	1
2	CU	K	185	1/1	0.88	0.20	55,55,55,55	1
2	CU	P	181	1/1	0.88	0.16	72,72,72,72	1
2	CU	D	177	1/1	0.88	0.27	64,64,64,64	1
2	CU	B	192	1/1	0.89	0.12	64,64,64,64	1
2	CU	E	185	1/1	0.89	0.14	49,49,49,49	1
2	CU	C	184	1/1	0.89	0.28	59,59,59,59	1
2	CU	H	184	1/1	0.89	0.19	62,62,62,62	1
2	CU	E	181	1/1	0.89	0.24	65,65,65,65	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CU	R	181	1/1	0.90	0.18	60,60,60,60	1
2	CU	B	177	1/1	0.90	0.12	70,70,70,70	1
2	CU	M	182	1/1	0.90	0.17	63,63,63,63	1
2	CU	C	192	1/1	0.91	0.09	57,57,57,57	1
2	CU	J	192	1/1	0.91	0.10	60,60,60,60	1
2	CU	M	177	1/1	0.91	0.11	58,58,58,58	1
2	CU	G	177	1/1	0.91	0.10	62,62,62,62	1
2	CU	O	185	1/1	0.91	0.07	53,53,53,53	1
2	CU	A	192	1/1	0.91	0.11	58,58,58,58	1
2	CU	G	181	1/1	0.91	0.20	67,67,67,67	1
2	CU	P	185	1/1	0.91	0.23	57,57,57,57	1
2	CU	G	183	1/1	0.91	0.12	53,53,53,53	1
2	CU	L	184	1/1	0.91	0.16	62,62,62,62	1
2	CU	M	191	1/1	0.92	0.11	67,67,67,67	1
2	CU	U	185	1/1	0.92	0.12	50,50,50,50	1
2	CU	D	185	1/1	0.92	0.13	71,71,71,71	1
2	CU	M	185	1/1	0.92	0.14	62,62,62,62	1
2	CU	F	181	1/1	0.92	0.15	74,74,74,74	1
2	CU	F	183	1/1	0.92	0.15	58,58,58,58	1
2	CU	A	181	1/1	0.93	0.15	65,65,65,65	1
2	CU	S	182	1/1	0.93	0.13	62,62,62,62	1
2	CU	S	191	1/1	0.93	0.10	66,66,66,66	1
2	CU	T	181	1/1	0.93	0.15	64,64,64,64	1
2	CU	P	191	1/1	0.93	0.09	53,53,53,53	1
2	CU	I	192	1/1	0.93	0.09	64,64,64,64	1
2	CU	F	182	1/1	0.93	0.08	56,56,56,56	1
2	CU	L	181	1/1	0.93	0.07	61,61,61,61	1
2	CU	I	182	1/1	0.93	0.14	57,57,57,57	1
2	CU	K	191	1/1	0.93	0.07	46,46,46,46	1
2	CU	L	185	1/1	0.93	0.08	62,62,62,62	1
2	CU	L	182	1/1	0.94	0.09	51,51,51,51	1
2	CU	L	183	1/1	0.94	0.06	63,63,63,63	1
2	CU	I	181	1/1	0.94	0.14	50,50,50,50	1
2	CU	A	184	1/1	0.94	0.09	57,57,57,57	1
2	CU	C	182	1/1	0.94	0.07	53,53,53,53	1
2	CU	E	182	1/1	0.94	0.12	66,66,66,66	1
2	CU	M	183	1/1	0.94	0.11	61,61,61,61	1
2	CU	J	182	1/1	0.94	0.08	62,62,62,62	1
2	CU	D	181	1/1	0.94	0.20	65,65,65,65	1
2	CU	D	183	1/1	0.94	0.14	66,66,66,66	1
2	CU	N	183	1/1	0.94	0.08	63,63,63,63	1
2	CU	C	183	1/1	0.94	0.11	60,60,60,60	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CU	A	183	1/1	0.94	0.09	61,61,61,61	1
2	CU	O	183	1/1	0.94	0.14	59,59,59,59	1
2	CU	V	191	1/1	0.94	0.06	56,56,56,56	1
2	CU	D	191	1/1	0.94	0.09	61,61,61,61	1
2	CU	H	191	1/1	0.94	0.07	49,49,49,49	1
2	CU	X	185	1/1	0.94	0.08	54,54,54,54	1
2	CU	X	182	1/1	0.94	0.07	62,62,62,62	1
2	CU	C	191	1/1	0.94	0.07	61,61,61,61	1
2	CU	J	183	1/1	0.95	0.07	57,57,57,57	1
2	CU	R	191	1/1	0.95	0.07	45,45,45,45	1
2	CU	J	184	1/1	0.95	0.26	55,55,55,55	1
2	CU	O	191	1/1	0.95	0.09	52,52,52,52	1
2	CU	A	182	1/1	0.95	0.13	65,65,65,65	1
2	CU	L	191	1/1	0.95	0.06	62,62,62,62	1
2	CU	H	183	1/1	0.95	0.08	52,52,52,52	1
2	CU	N	181	1/1	0.95	0.13	63,63,63,63	1
2	CU	Q	181	1/1	0.95	0.09	59,59,59,59	1
2	CU	V	185	1/1	0.95	0.09	61,61,61,61	1
2	CU	Q	182	1/1	0.95	0.09	67,67,67,67	1
2	CU	K	181	1/1	0.95	0.10	65,65,65,65	1
2	CU	Q	191	1/1	0.95	0.07	46,46,46,46	1
2	CU	Q	185	1/1	0.95	0.05	60,60,60,60	1
2	CU	N	184	1/1	0.95	0.16	55,55,55,55	1
2	CU	N	191	1/1	0.95	0.12	46,46,46,46	1
2	CU	M	192	1/1	0.95	0.09	69,69,69,69	1
2	CU	X	191	1/1	0.95	0.06	51,51,51,51	1
2	CU	F	191	1/1	0.96	0.06	53,53,53,53	1
2	CU	N	182	1/1	0.96	0.12	62,62,62,62	1
2	CU	I	183	1/1	0.96	0.05	60,60,60,60	1
2	CU	H	181	1/1	0.96	0.09	65,65,65,65	1
2	CU	Q	183	1/1	0.96	0.07	54,54,54,54	1
2	CU	A	185	1/1	0.96	0.06	58,58,58,58	1
2	CU	U	183	1/1	0.96	0.10	56,56,56,56	1
2	CU	D	192	1/1	0.96	0.07	56,56,56,56	1
2	CU	C	181	1/1	0.96	0.11	80,80,80,80	1
2	CU	V	182	1/1	0.96	0.10	62,62,62,62	1
2	CU	V	183	1/1	0.96	0.12	56,56,56,56	1
2	CU	O	182	1/1	0.96	0.14	63,63,63,63	1
2	CU	W	185	1/1	0.96	0.09	56,56,56,56	1
2	CU	M	184	1/1	0.96	0.11	56,56,56,56	1
2	CU	E	191	1/1	0.96	0.04	58,58,58,58	1
2	CU	R	183	1/1	0.96	0.13	58,58,58,58	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CU	O	192	1/1	0.96	0.10	59,59,59,59	1
2	CU	I	185	1/1	0.96	0.25	62,62,62,62	1
2	CU	A	191	1/1	0.96	0.06	41,41,41,41	1
2	CU	K	182	1/1	0.97	0.06	54,54,54,54	1
2	CU	U	181	1/1	0.97	0.08	66,66,66,66	1
2	CU	U	182	1/1	0.97	0.10	68,68,68,68	1
2	CU	G	185	1/1	0.97	0.05	55,55,55,55	1
2	CU	G	191	1/1	0.97	0.05	55,55,55,55	1
2	CU	P	182	1/1	0.97	0.05	60,60,60,60	1
2	CU	R	182	1/1	0.97	0.06	55,55,55,55	1
2	CU	P	183	1/1	0.97	0.04	61,61,61,61	1
2	CU	B	182	1/1	0.97	0.05	63,63,63,63	1
2	CU	M	181	1/1	0.97	0.12	54,54,54,54	1
2	CU	W	181	1/1	0.97	0.06	61,61,61,61	1
2	CU	J	191	1/1	0.97	0.04	52,52,52,52	1
2	CU	B	181	1/1	0.97	0.05	67,67,67,67	1
2	CU	I	191	1/1	0.97	0.04	56,56,56,56	1
2	CU	O	181	1/1	0.97	0.07	59,59,59,59	1
2	CU	J	185	1/1	0.97	0.11	55,55,55,55	1
2	CU	D	182	1/1	0.97	0.11	64,64,64,64	1
2	CU	T	182	1/1	0.98	0.10	58,58,58,58	1
2	CU	T	183	1/1	0.98	0.04	57,57,57,57	1
2	CU	H	182	1/1	0.98	0.06	54,54,54,54	1
2	CU	E	183	1/1	0.98	0.09	61,61,61,61	1
2	CU	J	181	1/1	0.98	0.04	58,58,58,58	1
2	CU	W	182	1/1	0.98	0.07	53,53,53,53	1
2	CU	S	181	1/1	0.98	0.07	71,71,71,71	1
2	CU	G	182	1/1	0.98	0.07	67,67,67,67	1
2	CU	W	191	1/1	0.98	0.04	56,56,56,56	1
2	CU	S	183	1/1	0.98	0.10	59,59,59,59	1
2	CU	X	181	1/1	0.98	0.04	55,55,55,55	1
2	CU	U	191	1/1	0.98	0.06	46,46,46,46	1
2	CU	X	183	1/1	0.98	0.06	54,54,54,54	1
2	CU	B	183	1/1	0.98	0.04	56,56,56,56	1
2	CU	B	191	1/1	0.98	0.05	52,52,52,52	1
2	CU	N	190	1/1	0.99	0.04	32,32,32,32	1
2	CU	D	190	1/1	0.99	0.03	32,32,32,32	1
2	CU	B	190	1/1	0.99	0.03	34,34,34,34	1
2	CU	A	190	1/1	0.99	0.04	37,37,37,37	1
2	CU	F	190	1/1	0.99	0.03	33,33,33,33	1
2	CU	E	190	1/1	1.00	0.03	34,34,34,34	1

6.5 Other polymers [i](#)

There are no such residues in this entry.