



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 05:09 am BST

PDB ID : 6KOZ
Title : Crystal structure of two domain M1 zinc metallopeptidase E323 mutant bound to L-Leucine amino acid
Authors : Agrawal, R.; Kumar, A.; Kumar, A.; Makde, R.D.
Deposited on : 2019-08-13
Resolution : 2.25 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 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 : 2.11
buster-report : 1.1.7 (2018)
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) : 2.11

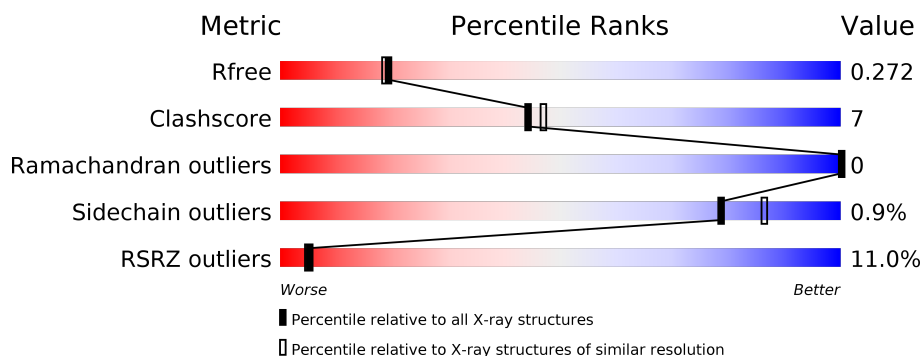
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria 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	474	<div> <div>6%</div> <div> <div></div> <div>76%</div> <div>16%</div> <div>8%</div> </div> </div>
1	B	474	<div> <div>15%</div> <div> <div></div> <div>77%</div> <div>14%</div> <div>9%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7025 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 Zinc metalloprotease, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	433	Total	C	N	O	S	0	1	0
			3289	2092	581	612	4			
1	A	436	Total	C	N	O	S	0	1	0
			3328	2114	591	619	4			

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	MET	-	initiating methionine	UNP Q9RVZ5
B	0	ALA	-	expression tag	UNP Q9RVZ5
B	1	SER	-	expression tag	UNP Q9RVZ5
B	2	TRP	-	expression tag	UNP Q9RVZ5
B	3	SER	-	expression tag	UNP Q9RVZ5
B	4	HIS	-	expression tag	UNP Q9RVZ5
B	5	PRO	-	expression tag	UNP Q9RVZ5
B	6	GLN	-	expression tag	UNP Q9RVZ5
B	7	PHE	-	expression tag	UNP Q9RVZ5
B	8	GLU	-	expression tag	UNP Q9RVZ5
B	9	LYS	-	expression tag	UNP Q9RVZ5
B	10	GLY	-	expression tag	UNP Q9RVZ5
B	11	SER	-	expression tag	UNP Q9RVZ5
B	12	SER	-	expression tag	UNP Q9RVZ5
B	13	HIS	-	expression tag	UNP Q9RVZ5
B	14	HIS	-	expression tag	UNP Q9RVZ5
B	15	HIS	-	expression tag	UNP Q9RVZ5
B	16	HIS	-	expression tag	UNP Q9RVZ5
B	17	HIS	-	expression tag	UNP Q9RVZ5
B	18	HIS	-	expression tag	UNP Q9RVZ5
B	19	SER	-	expression tag	UNP Q9RVZ5
B	20	SER	-	expression tag	UNP Q9RVZ5
B	21	GLY	-	expression tag	UNP Q9RVZ5
B	22	SER	-	expression tag	UNP Q9RVZ5
B	23	GLY	-	expression tag	UNP Q9RVZ5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	24	GLY	-	expression tag	UNP Q9RVZ5
B	25	GLY	-	expression tag	UNP Q9RVZ5
B	26	GLY	-	expression tag	UNP Q9RVZ5
B	27	GLY	-	expression tag	UNP Q9RVZ5
B	28	GLU	-	expression tag	UNP Q9RVZ5
B	29	ASN	-	expression tag	UNP Q9RVZ5
B	30	LEU	-	expression tag	UNP Q9RVZ5
B	31	TYR	-	expression tag	UNP Q9RVZ5
B	32	PHE	-	expression tag	UNP Q9RVZ5
B	33	GLN	-	expression tag	UNP Q9RVZ5
B	34	GLY	-	expression tag	UNP Q9RVZ5
B	35	SER	-	expression tag	UNP Q9RVZ5
B	323	ALA	GLU	engineered mutation	UNP Q9RVZ5
A	-1	MET	-	initiating methionine	UNP Q9RVZ5
A	0	ALA	-	expression tag	UNP Q9RVZ5
A	1	SER	-	expression tag	UNP Q9RVZ5
A	2	TRP	-	expression tag	UNP Q9RVZ5
A	3	SER	-	expression tag	UNP Q9RVZ5
A	4	HIS	-	expression tag	UNP Q9RVZ5
A	5	PRO	-	expression tag	UNP Q9RVZ5
A	6	GLN	-	expression tag	UNP Q9RVZ5
A	7	PHE	-	expression tag	UNP Q9RVZ5
A	8	GLU	-	expression tag	UNP Q9RVZ5
A	9	LYS	-	expression tag	UNP Q9RVZ5
A	10	GLY	-	expression tag	UNP Q9RVZ5
A	11	SER	-	expression tag	UNP Q9RVZ5
A	12	SER	-	expression tag	UNP Q9RVZ5
A	13	HIS	-	expression tag	UNP Q9RVZ5
A	14	HIS	-	expression tag	UNP Q9RVZ5
A	15	HIS	-	expression tag	UNP Q9RVZ5
A	16	HIS	-	expression tag	UNP Q9RVZ5
A	17	HIS	-	expression tag	UNP Q9RVZ5
A	18	HIS	-	expression tag	UNP Q9RVZ5
A	19	SER	-	expression tag	UNP Q9RVZ5
A	20	SER	-	expression tag	UNP Q9RVZ5
A	21	GLY	-	expression tag	UNP Q9RVZ5
A	22	SER	-	expression tag	UNP Q9RVZ5
A	23	GLY	-	expression tag	UNP Q9RVZ5
A	24	GLY	-	expression tag	UNP Q9RVZ5
A	25	GLY	-	expression tag	UNP Q9RVZ5
A	26	GLY	-	expression tag	UNP Q9RVZ5
A	27	GLY	-	expression tag	UNP Q9RVZ5

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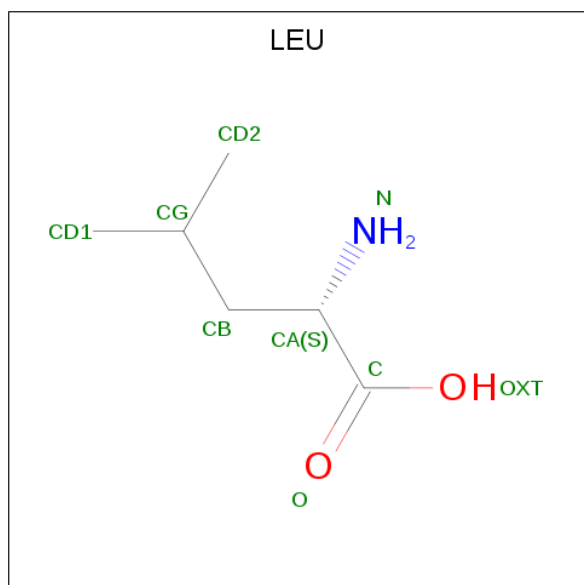
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Chain	Residue	Modelled	Actual	Comment	Reference
A	28	GLU	-	expression tag	UNP Q9RVZ5
A	29	ASN	-	expression tag	UNP Q9RVZ5
A	30	LEU	-	expression tag	UNP Q9RVZ5
A	31	TYR	-	expression tag	UNP Q9RVZ5
A	32	PHE	-	expression tag	UNP Q9RVZ5
A	33	GLN	-	expression tag	UNP Q9RVZ5
A	34	GLY	-	expression tag	UNP Q9RVZ5
A	35	SER	-	expression tag	UNP Q9RVZ5
A	323	ALA	GLU	engineered mutation	UNP Q9RVZ5

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by author).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Zn 1 1	0	0
2	A	1	Total Zn 1 1	0	0

- Molecule 3 is LEUCINE (three-letter code: LEU) (formula: C₆H₁₃NO₂) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C N O 9 6 1 2	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			9	6	1	2		

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Na	0	0
			1	1		
4	A	1	Total	Na	0	0
			1	1		

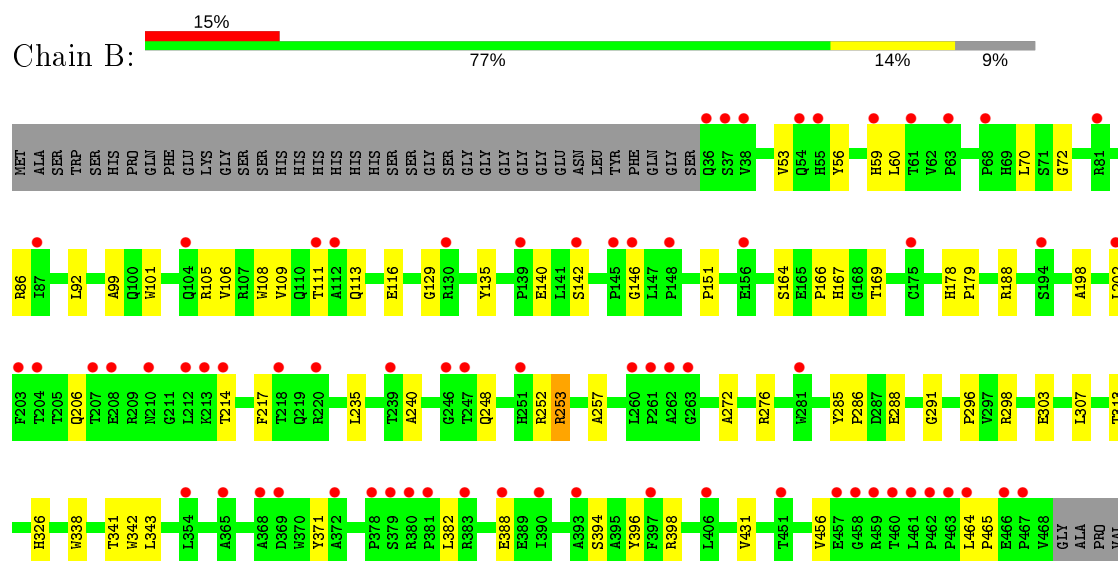
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	165	Total	O	0	4
			169	169		
5	A	213	Total	O	0	4
			217	217		

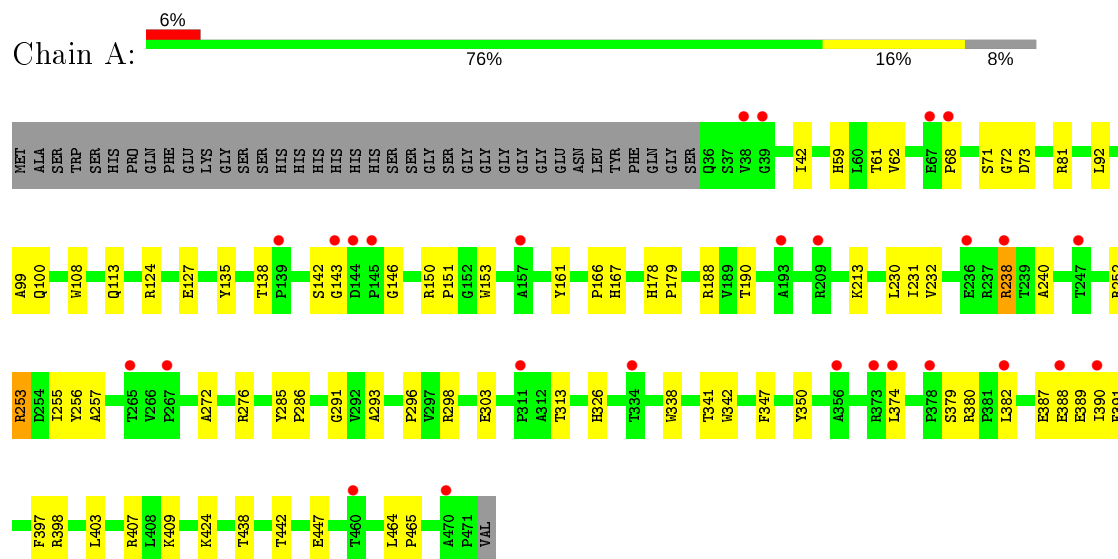
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Zinc metalloprotease, putative



- Molecule 1: Zinc metalloprotease, putative



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	51.21Å 57.42Å 69.01Å 89.97° 82.76° 67.82°	Depositor
Resolution (Å)	26.55 – 2.25 38.92 – 2.25	Depositor EDS
% Data completeness (in resolution range)	98.1 (26.55-2.25) 98.2 (38.92-2.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.09 (at 2.24Å)	Xtriage
Refinement program	PHENIX 1.10.1 _2155	Depositor
R, R_{free}	0.218 , 0.272 0.218 , 0.272	Depositor DCC
R_{free} test set	1653 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	24.1	Xtriage
Anisotropy	0.796	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 45.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7025	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.21% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.25	0/3420	0.48	0/4687
1	B	0.25	0/3380	0.46	0/4636
All	All	0.25	0/6800	0.47	0/9323

There are no bond length outliers.

There are no bond angle outliers.

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	3328	0	3281	47	0
1	B	3289	0	3227	42	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	9	0	10	0	0
3	B	9	0	10	2	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	217	0	0	3	0
5	B	169	0	0	0	0
All	All	7025	0	6528	87	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 (87) 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:A:146:GLY:HA3	1:A:388:GLU:HG3	1.59	0.84
1:B:146:GLY:HA3	1:B:388:GLU:HG3	1.69	0.75
1:B:92:LEU:HD13	1:B:113:GLN:HG2	1.73	0.69
1:A:92:LEU:HD13	1:A:113:GLN:HG2	1.75	0.69
1:B:111:THR:HG21	1:A:424:LYS:HB3	1.81	0.62
1:A:298:ARG:HG3	1:A:313:THR:HG21	1.82	0.61
1:A:238:ARG:NH2	5:A:609:HOH:O	2.36	0.59
1:A:272:ALA:O	1:A:276:ARG:HG3	2.03	0.58
1:A:382:LEU:HD21	1:A:398:ARG:HG2	1.85	0.58
1:B:298:ARG:HG3	1:B:313:THR:HG21	1.86	0.58
1:B:382:LEU:HD21	1:B:398:ARG:HG2	1.88	0.56
1:B:105:ARG:NH1	1:B:106:VAL:O	2.39	0.56
1:A:81:ARG:NH1	5:A:616:HOH:O	2.39	0.56
1:A:151:PRO:HG3	1:A:166:PRO:HD2	1.88	0.56
1:B:240:ALA:HB3	1:B:252:ARG:HB3	1.89	0.55
1:A:240:ALA:HB3	1:A:252:ARG:HB3	1.89	0.53
1:B:396:TYR:OH	3:B:502:LEU:O	2.20	0.53
1:A:143:GLY:HA2	1:A:150:ARG:HG2	1.91	0.53
1:A:62:VAL:HG11	1:A:232:VAL:HG21	1.91	0.53
1:A:388:GLU:N	1:A:388:GLU:OE1	2.36	0.52
1:A:438:THR:O	1:A:442:THR:HG23	2.10	0.51
1:B:272:ALA:O	1:B:276:ARG:HG3	2.09	0.51
1:B:109:VAL:HG22	1:B:111:THR:HG23	1.93	0.51
1:A:59:HIS:HB3	1:A:73:ASP:HB3	1.93	0.50
1:B:382:LEU:HD12	1:B:456:VAL:HG22	1.93	0.50
1:A:42:ILE:CG2	1:A:387:GLU:HG3	2.42	0.49
1:B:72:GLY:HA3	1:B:135:TYR:CZ	2.47	0.49
1:A:238:ARG:HD3	1:A:256:TYR:CD1	2.47	0.49
1:B:257:ALA:HB3	1:B:296:PRO:HD3	1.93	0.49
1:A:255:ILE:HB	1:A:293:ALA:HA	1.94	0.48
1:A:342:TRP:CD2	1:A:382:LEU:HD13	2.48	0.48
1:A:253:ARG:O	1:A:291:GLY:HA3	2.12	0.48
1:A:380:ARG:NE	1:A:389:GLU:HB3	2.28	0.48
1:A:100:GLN:OE1	5:A:601:HOH:O	2.20	0.48
1:B:303:GLU:OE1	3:B:502:LEU:N	2.47	0.47
1:B:99:ALA:HB3	1:B:108:TRP:CD1	2.50	0.47
1:B:105:ARG:HH11	1:B:106:VAL:H	1.61	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:253:ARG:NH2	1:B:307:LEU:HB2	2.29	0.47
1:A:124:ARG:O	1:A:127:GLU:HG2	2.15	0.47
1:B:151:PRO:HG3	1:B:166:PRO:HD2	1.97	0.47
1:A:303:GLU:HG3	1:A:326:HIS:HB3	1.96	0.46
1:A:99:ALA:HB3	1:A:108:TRP:CD1	2.49	0.46
1:A:257:ALA:HB3	1:A:296:PRO:HD3	1.97	0.46
1:A:142:SER:O	1:A:151:PRO:HD2	2.15	0.46
1:A:153:TRP:HH2	1:A:230:LEU:HD22	1.81	0.46
1:B:253:ARG:O	1:B:291:GLY:HA3	2.15	0.46
1:B:303:GLU:HG3	1:B:326:HIS:HB3	1.98	0.46
1:A:374:LEU:HD11	1:A:397:PHE:HB2	1.99	0.45
1:B:92:LEU:HG	1:B:140:GLU:HB2	1.98	0.45
1:A:68:PRO:HB2	1:A:138:THR:HG23	1.99	0.45
1:B:394:SER:O	1:B:398:ARG:HB3	2.14	0.45
1:A:72:GLY:HA3	1:A:135:TYR:CZ	2.52	0.45
1:B:343:LEU:HD13	1:B:431:VAL:HG13	1.99	0.45
1:B:60:LEU:HD11	1:B:70:LEU:HD13	1.97	0.45
1:A:161:TYR:HB3	1:A:231:ILE:HG23	1.99	0.44
1:B:111:THR:CG2	1:A:424:LYS:HB3	2.47	0.44
1:A:464:LEU:HD12	1:A:465:PRO:HD2	2.00	0.44
1:A:178:HIS:CG	1:A:179:PRO:HD2	2.53	0.44
1:A:238:ARG:HD3	1:A:256:TYR:CE1	2.53	0.44
1:B:164:SER:HB3	1:B:169:THR:N	2.33	0.44
1:B:464:LEU:HD12	1:B:465:PRO:HD2	1.99	0.43
1:A:59:HIS:HD2	1:A:188:ARG:O	2.01	0.43
1:A:342:TRP:CD1	1:A:382:LEU:HB3	2.53	0.43
1:B:142:SER:O	1:B:151:PRO:HD2	2.19	0.43
1:A:190:THR:HA	1:A:213:LYS:O	2.18	0.43
1:A:347:PHE:CZ	1:A:403:LEU:HD21	2.53	0.43
1:B:53:VAL:HG11	1:B:56:TYR:CZ	2.54	0.43
1:B:202:LEU:HD21	1:B:288:GLU:HB2	2.01	0.43
1:A:390:ILE:HG13	1:A:391:PHE:CD2	2.54	0.42
1:B:206:GLN:HA	1:B:214:THR:O	2.19	0.42
1:B:235:LEU:HA	1:B:257:ALA:HA	2.01	0.42
1:B:371:TYR:HB2	1:B:464:LEU:HD13	2.00	0.42
1:B:86:ARG:HE	1:B:116:GLU:HB3	1.85	0.42
1:B:178:HIS:CG	1:B:179:PRO:HD2	2.55	0.42
1:A:61:THR:HB	1:A:71:SER:HB2	2.01	0.41
1:A:42:ILE:HG23	1:A:387:GLU:HG3	2.02	0.41
1:B:248:GLN:NE2	1:B:285:TYR:O	2.39	0.41
1:A:409:LYS:NZ	1:A:447:GLU:HB3	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:285:TYR:CD1	1:B:286:PRO:HD2	2.55	0.41
1:B:101:TRP:CH2	1:B:129:GLY:HA3	2.56	0.41
1:B:59:HIS:HD2	1:B:188:ARG:O	2.04	0.41
1:A:338:TRP:O	1:A:341:THR:HG23	2.21	0.41
1:B:198:ALA:HA	1:B:217:PHE:CZ	2.56	0.41
1:B:342:TRP:CD2	1:B:382:LEU:HD13	2.55	0.41
1:B:338:TRP:O	1:B:341:THR:HG23	2.21	0.40
1:A:285:TYR:CD1	1:A:286:PRO:HD2	2.56	0.40
1:A:350:TYR:OH	1:A:407:ARG:HD3	2.22	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	435/474 (92%)	416 (96%)	19 (4%)	0	100	100
1	B	432/474 (91%)	413 (96%)	19 (4%)	0	100	100
All	All	867/948 (92%)	829 (96%)	38 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	346/381 (91%)	342 (99%)	4 (1%)	71	80
1	B	340/381 (89%)	338 (99%)	2 (1%)	86	91
All	All	686/762 (90%)	680 (99%)	6 (1%)	78	86

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	167	HIS
1	B	253	ARG
1	A	167	HIS
1	A	238	ARG
1	A	253	ARG
1	A	379	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	100	GLN
1	B	210	ASN
1	A	100	GLN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	$\# Z > 2$
3	LEU	A	502	2	5,8,8	0.21	0	6,10,10	0.46	0
3	LEU	B	502	-	5,8,8	0.22	0	6,10,10	0.40	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	LEU	A	502	2	-	0/4/8/8	-
3	LEU	B	502	-	-	0/4/8/8	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

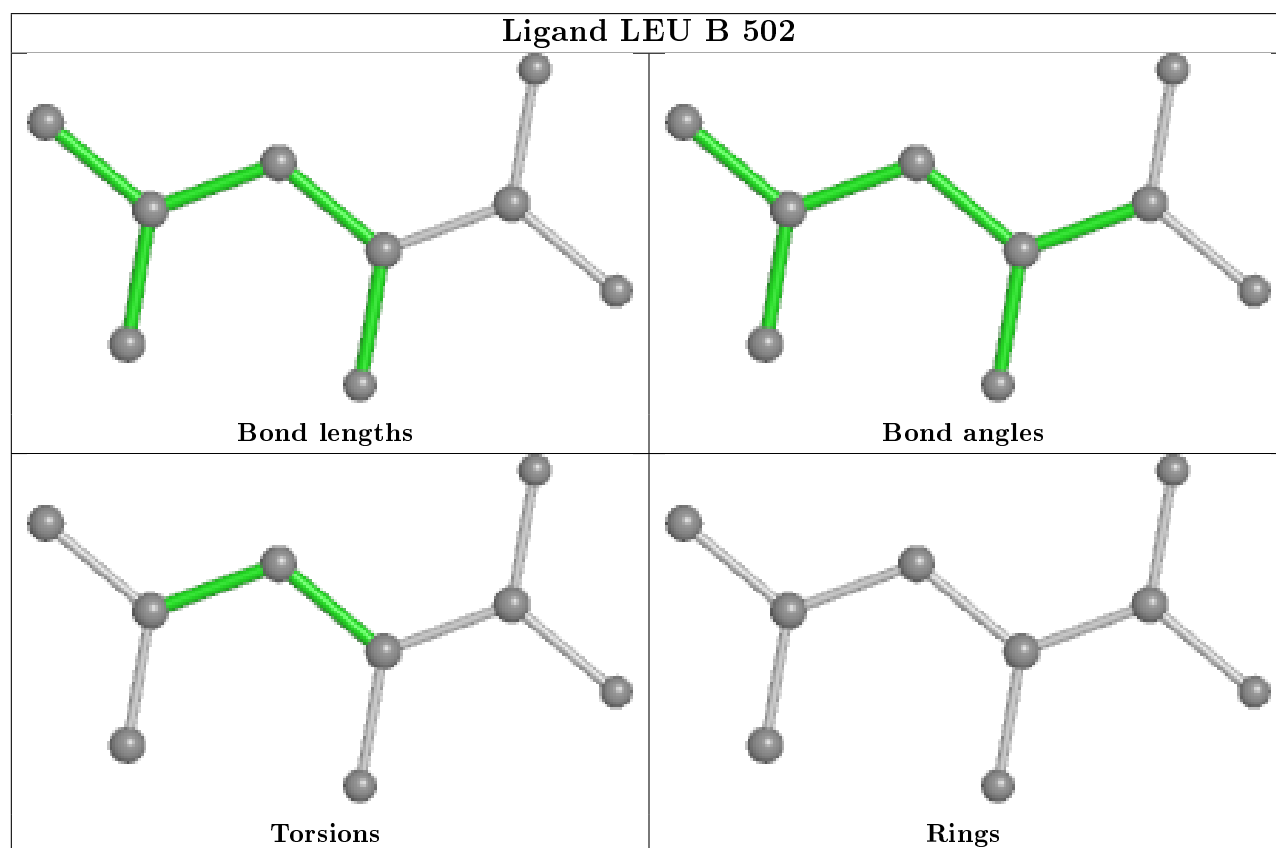
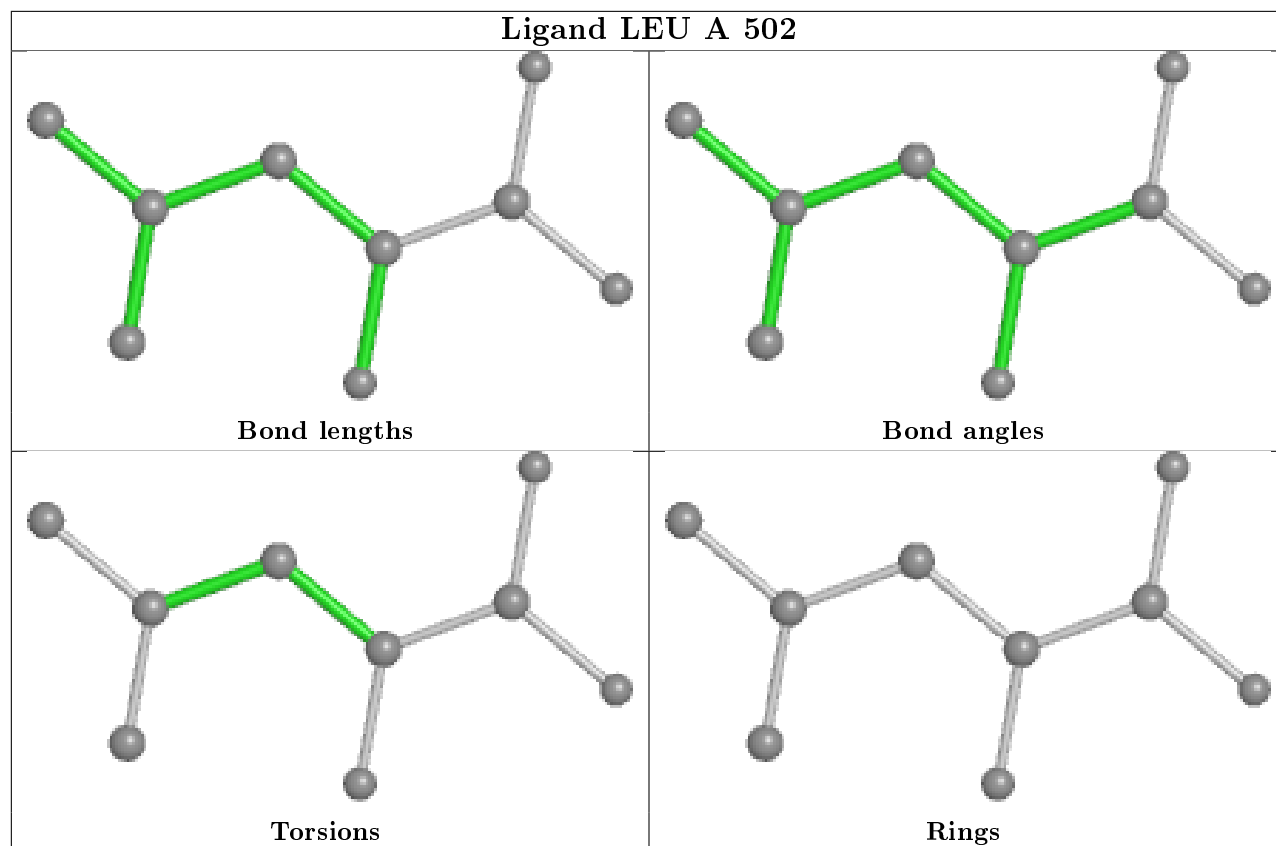
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	502	LEU	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	436/474 (91%)	0.84	27 (6%) 20 22	16, 25, 36, 54	0
1	B	433/474 (91%)	1.13	69 (15%) 1 1	20, 29, 41, 56	0
All	All	869/948 (91%)	0.99	96 (11%) 5 5	16, 26, 40, 56	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	145	PRO	7.7
1	B	460	THR	5.2
1	B	461	LEU	5.2
1	B	462	PRO	5.2
1	B	372	ALA	5.0
1	A	145	PRO	5.0
1	B	379	SER	4.8
1	B	381	PRO	4.7
1	B	112	ALA	4.7
1	B	463	PRO	4.2
1	A	68	PRO	4.1
1	B	36	GLN	3.9
1	B	207	THR	3.9
1	B	368	ALA	3.8
1	B	369	ASP	3.8
1	B	393	ALA	3.7
1	A	247	THR	3.7
1	B	459	ARG	3.6
1	A	144	ASP	3.6
1	A	460	THR	3.6
1	B	208	GLU	3.6
1	B	210	ASN	3.4
1	B	464	LEU	3.3
1	A	470	ALA	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	130	ARG	3.3
1	B	204	THR	3.2
1	B	247	THR	3.2
1	B	397	PHE	3.2
1	A	67	GLU	3.2
1	B	388	GLU	3.2
1	B	142	SER	2.9
1	B	263	GLY	2.9
1	B	68	PRO	2.8
1	B	220	ARG	2.8
1	B	37	SER	2.7
1	A	374	LEU	2.7
1	A	238	ARG	2.7
1	A	38	VAL	2.7
1	A	139	PRO	2.6
1	A	378	PRO	2.6
1	B	260	LEU	2.6
1	B	262	ALA	2.6
1	B	246	GLY	2.6
1	B	458	GLY	2.5
1	A	193	ALA	2.5
1	A	356	ALA	2.5
1	A	157	ALA	2.5
1	A	267	PRO	2.5
1	A	373	ARG	2.5
1	B	55	HIS	2.5
1	A	388	GLU	2.5
1	B	218	THR	2.4
1	B	378	PRO	2.4
1	B	194	SER	2.4
1	B	251	HIS	2.4
1	B	261	PRO	2.4
1	A	236	GLU	2.4
1	B	239	THR	2.4
1	B	146	GLY	2.4
1	B	451	THR	2.4
1	B	214	THR	2.3
1	B	81	ARG	2.3
1	A	334	THR	2.3
1	B	212	LEU	2.3
1	B	380	ARG	2.3
1	B	467	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	111	THR	2.3
1	B	61	THR	2.2
1	A	143	GLY	2.2
1	B	54	GLN	2.2
1	B	390	ILE	2.2
1	B	59	HIS	2.2
1	A	311	PRO	2.2
1	B	175	CYS	2.2
1	B	457	GLU	2.2
1	B	365	ALA	2.2
1	B	87	ILE	2.2
1	B	148	PRO	2.2
1	B	466	GLU	2.1
1	A	39	GLY	2.1
1	B	383	ARG	2.1
1	B	203	PHE	2.1
1	B	139	PRO	2.1
1	B	202	LEU	2.1
1	B	281	TRP	2.1
1	B	406	LEU	2.1
1	A	382	LEU	2.1
1	A	209	ARG	2.1
1	B	213	LYS	2.1
1	B	104	GLN	2.0
1	B	354	LEU	2.0
1	A	265	THR	2.0
1	A	390	ILE	2.0
1	B	63	PRO	2.0
1	B	38	VAL	2.0
1	B	156	GLU	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 carbohydrates in this entry.

6.4 Ligands [i](#)

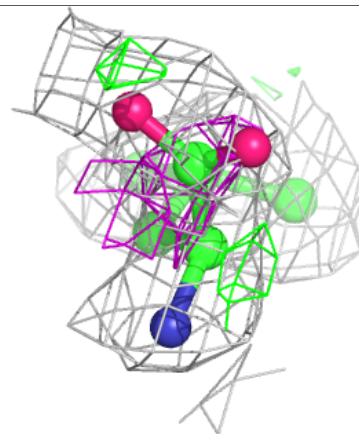
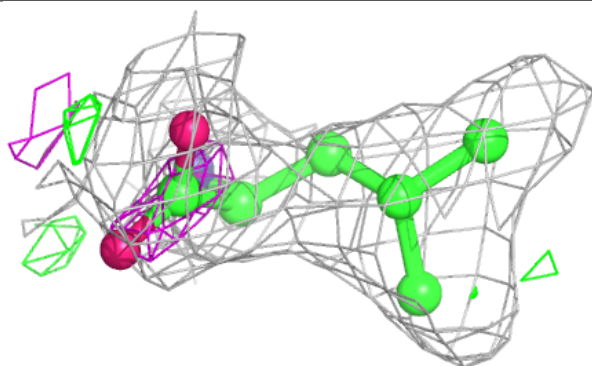
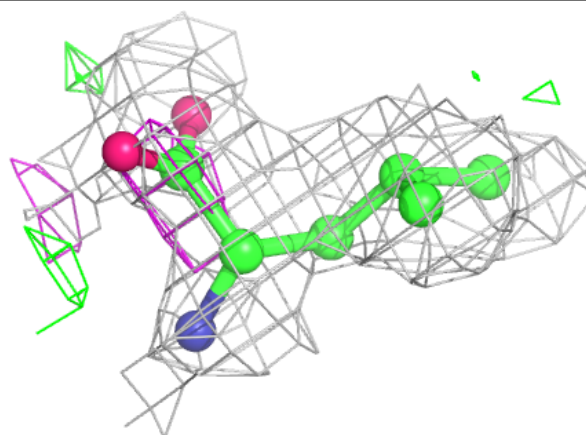
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	LEU	B	502	9/9	0.72	0.27	18,22,25,29	0
3	LEU	A	502	9/9	0.82	0.23	21,24,26,29	0
4	NA	A	503	1/1	0.91	0.25	10,10,10,10	0
4	NA	B	503	1/1	0.92	0.21	18,18,18,18	0
2	ZN	B	501	1/1	0.99	0.05	22,22,22,22	0
2	ZN	A	501	1/1	0.99	0.07	18,18,18,18	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

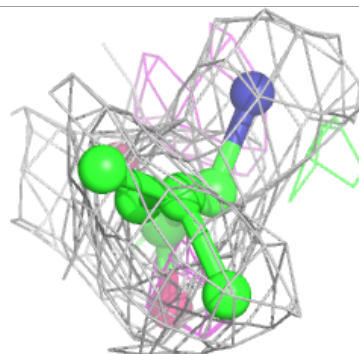
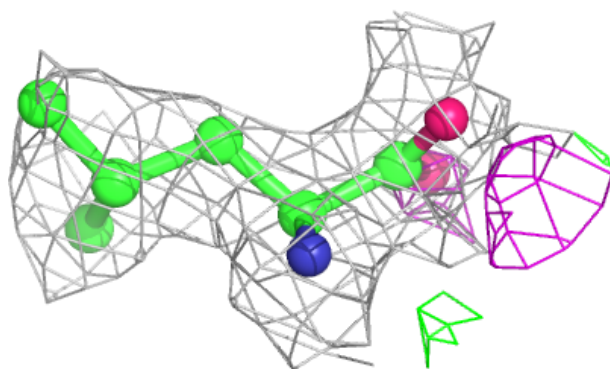
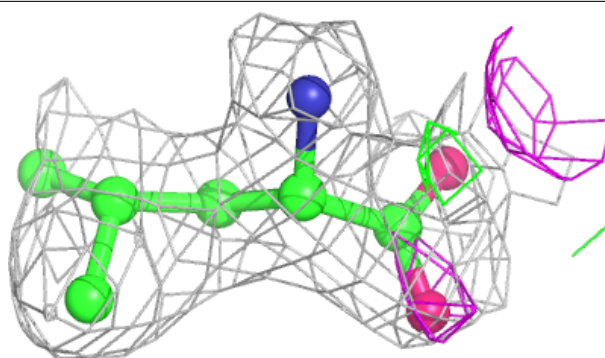
Electron density around LEU B 502:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



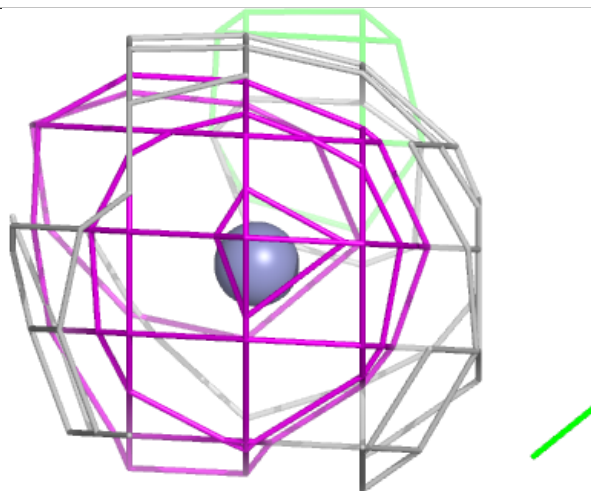
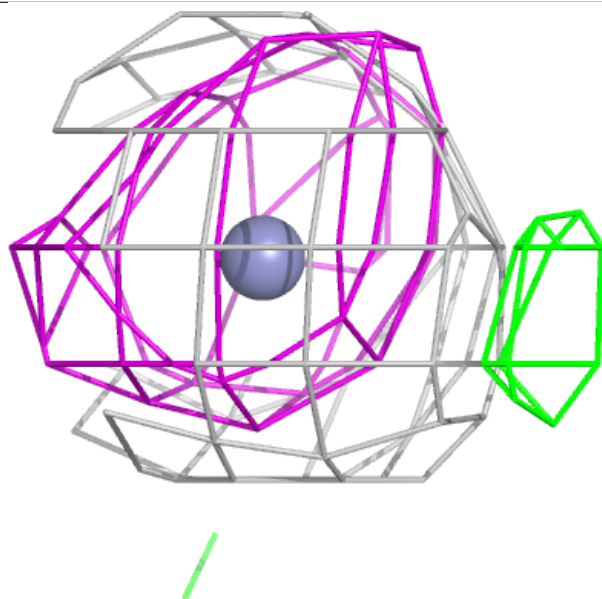
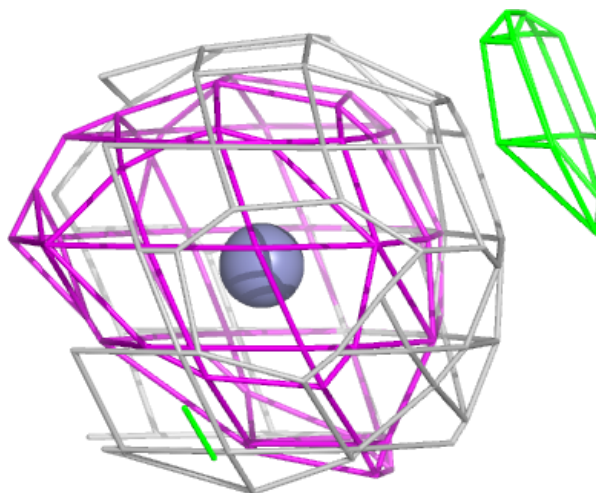
Electron density around LEU A 502:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



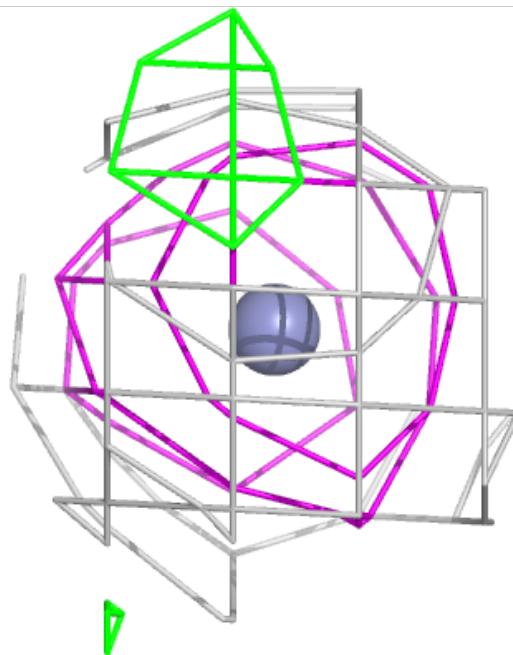
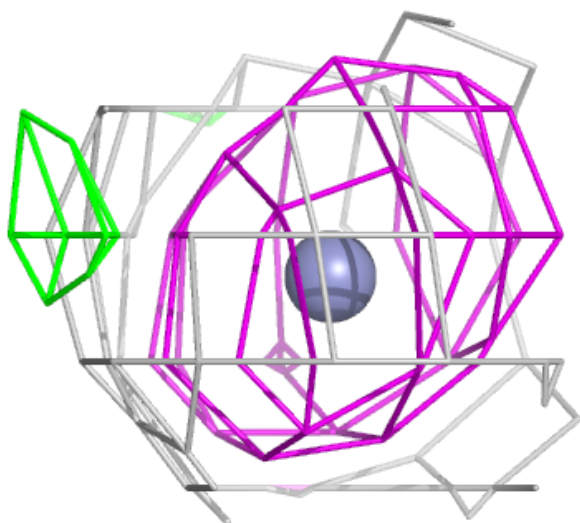
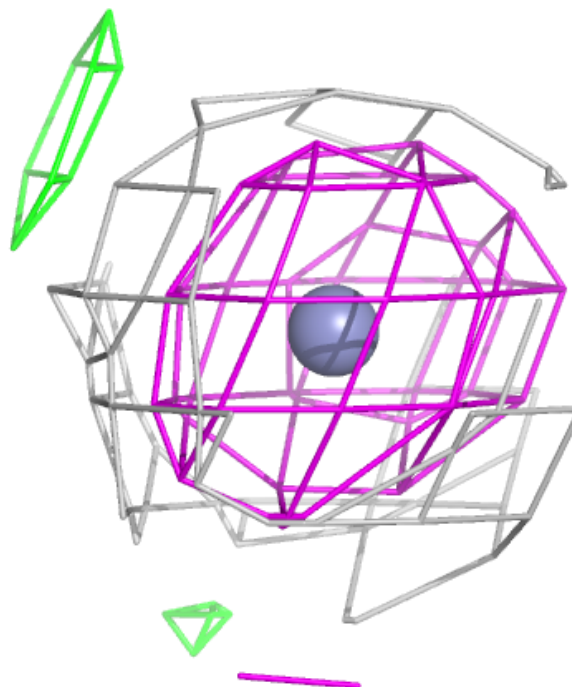
Electron density around ZN B 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ZN A 501:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.