



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 09:47 PM UTC

PDB ID : 6RL0 / pdb_00006rl0
Title : Recombinant Pseudomonas stutzeri nitrous oxide reductase, form I
Authors : Zhang, L.; Wuest, A.; Prasser, B.; Mueller, C.; Einsle, O.
Deposited on : 2019-04-30
Resolution : 1.78 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

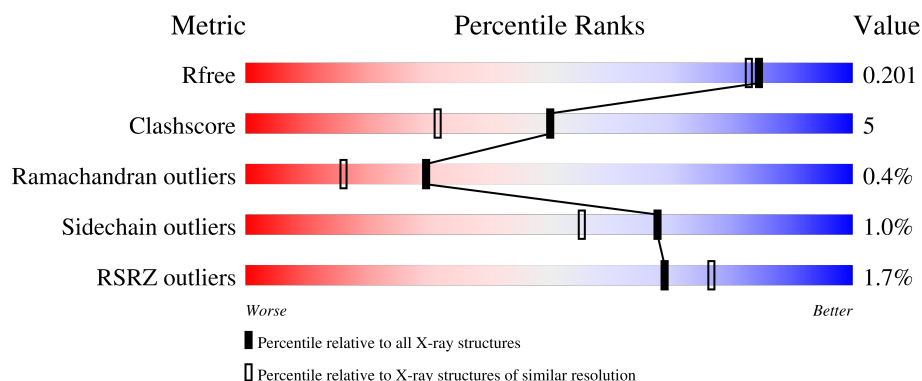
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

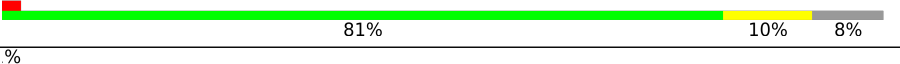
The reported resolution of this entry is 1.78 Å.

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	1365 (1.78-1.78)
Clashscore	190562	1395 (1.78-1.78)
Ramachandran outliers	187476	1382 (1.78-1.78)
Sidechain outliers	187428	1382 (1.78-1.78)
RSRZ outliers	180081	1365 (1.78-1.78)

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	638	 0% 80% 11% 8%
1	B	638	 2% 81% 10% 8%
1	C	638	 0% 80% 11% 9%
1	D	638	 2% 80% 10% 9%

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
3	CL	A	702	-	-	X	-
3	CL	C	702	-	-	X	-
7	FMT	D	706	-	-	X	-

2 Entry composition [i](#)

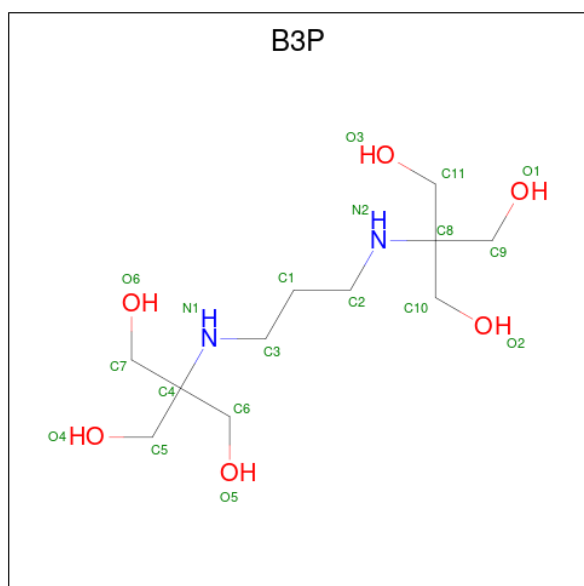
There are 10 unique types of molecules in this entry. The entry contains 19996 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Nitrous-oxide reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	584	Total	C	N	O	S	0	3	0
			4610	2914	794	870	32			
1	B	584	Total	C	N	O	S	0	4	0
			4615	2917	795	871	32			
1	C	582	Total	C	N	O	S	0	3	0
			4596	2907	791	866	32			
1	D	581	Total	C	N	O	S	0	2	0
			4581	2897	789	863	32			

- Molecule 2 is 2-[3-(2-HYDROXY-1,1-DIHYDROXYMETHYL-ETHYLAMINO)-PROPYLAMINO]-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: B3P) (formula: $C_{11}H_{26}N_2O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			19	11	2	6		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			19	11	2	6		
2	C	1	Total	C	N	O	0	0
			19	11	2	6		
2	D	1	Total	C	N	O	0	0
			19	11	2	6		

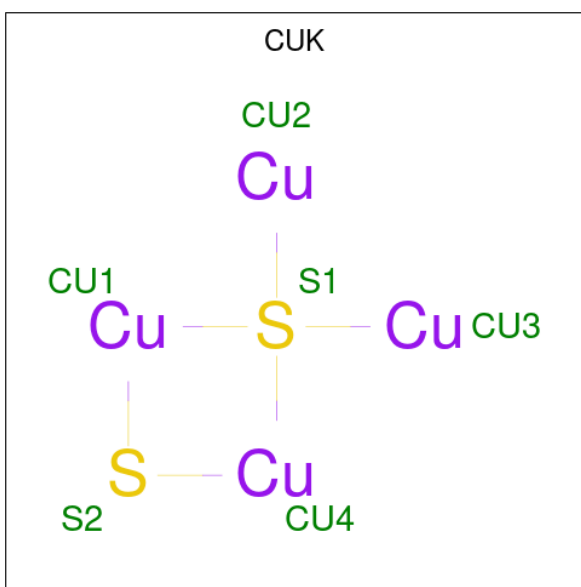
- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	B	1	Total	Cl	0	0
			1	1		
3	C	1	Total	Cl	0	0
			1	1		
3	D	1	Total	Cl	0	0
			1	1		

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

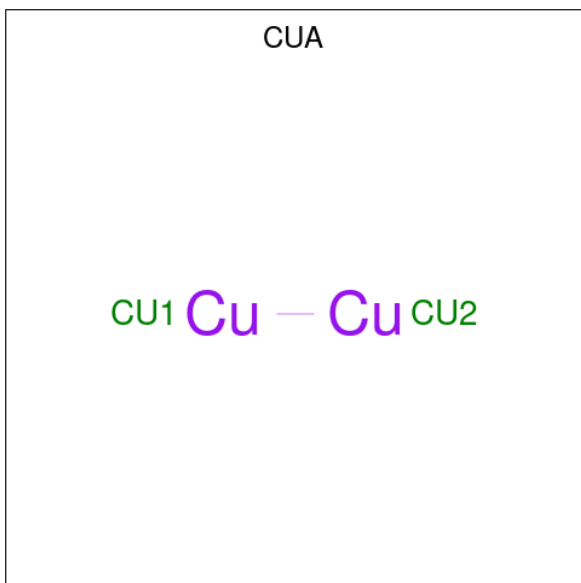
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ca	0	0
			1	1		
4	B	1	Total	Ca	0	0
			1	1		
4	C	1	Total	Ca	0	0
			1	1		
4	D	1	Total	Ca	0	0
			1	1		

- Molecule 5 is [4Cu:2S] cluster (CCD ID: CUK) (formula: Cu₄S₂) (labeled as "Ligand of Interest" by depositor).



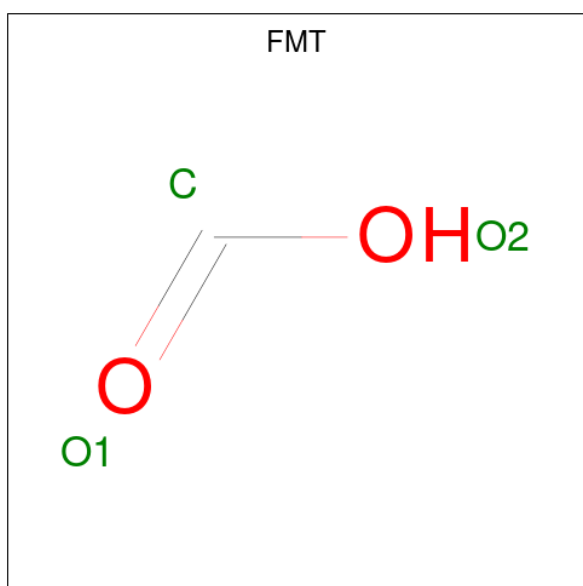
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	Cu	S	0	0
			6	4	2		
5	B	1	Total	Cu	S	0	0
			6	4	2		
5	C	1	Total	Cu	S	0	0
			6	4	2		
5	D	1	Total	Cu	S	0	0
			6	4	2		

- Molecule 6 is DINUCLEAR COPPER ION (CCD ID: CUA) (formula: Cu₂).



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

- Molecule 7 is FORMIC ACID (CCD ID: FMT) (formula: CH₂O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C O 3 1 2	0	0
7	B	1	Total C O 3 1 2	0	0
7	C	1	Total C O 3 1 2	0	0
7	C	1	Total C O 3 1 2	0	0
7	D	1	Total C O 3 1 2	0	0

- Molecule 8 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total K 1 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	B	1	Total K 1 1	0	0
8	C	1	Total K 1 1	0	0
8	D	1	Total K 1 1	0	0

- Molecule 9 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total Na 1 1	0	0
9	B	1	Total Na 1 1	0	0
9	C	1	Total Na 1 1	0	0
9	D	1	Total Na 1 1	0	0

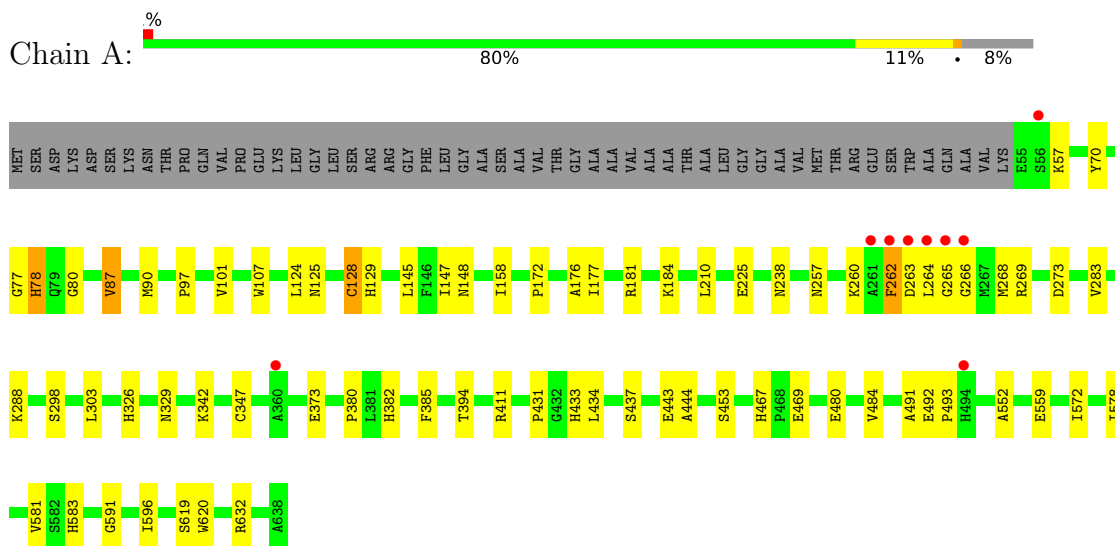
- Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	413	Total O 413 413	0	0
10	B	392	Total O 392 392	0	0
10	C	362	Total O 362 362	0	0
10	D	288	Total O 288 288	0	0

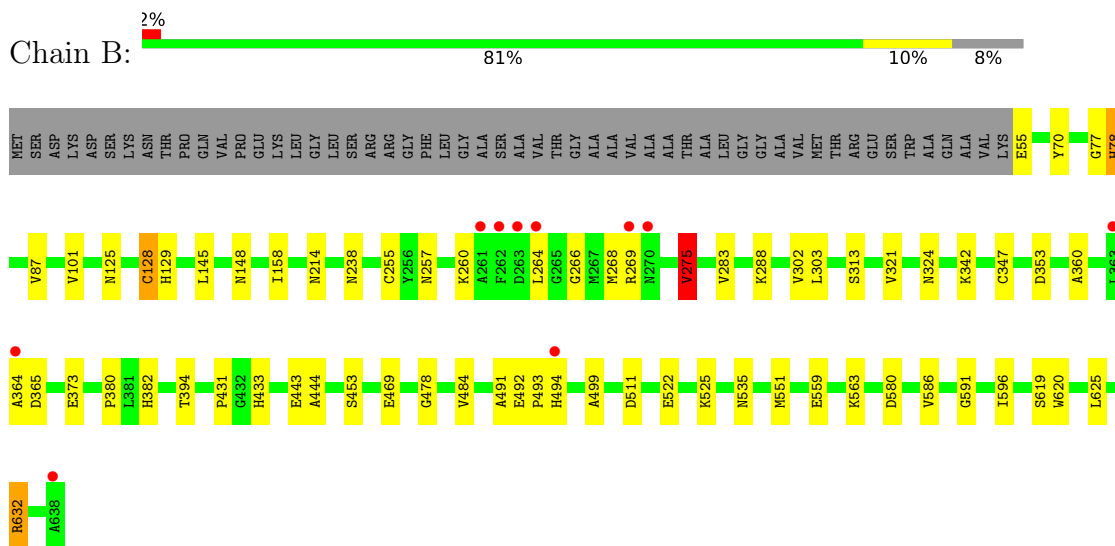
3 Residue-property plots

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.

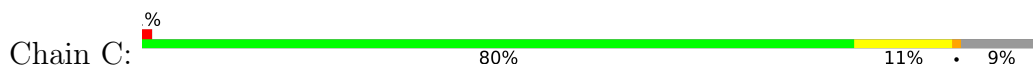
- Molecule 1: Nitrous-oxide reductase

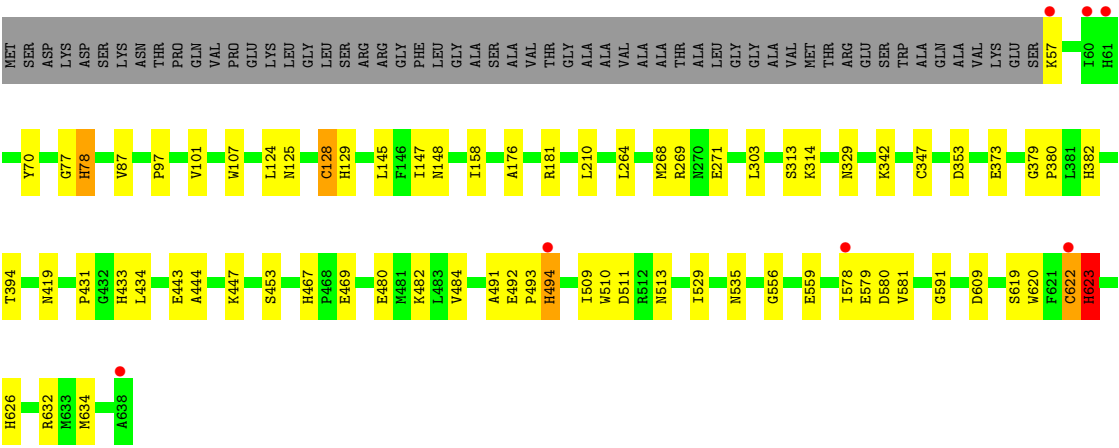


- Molecule 1: Nitrous-oxide reductase

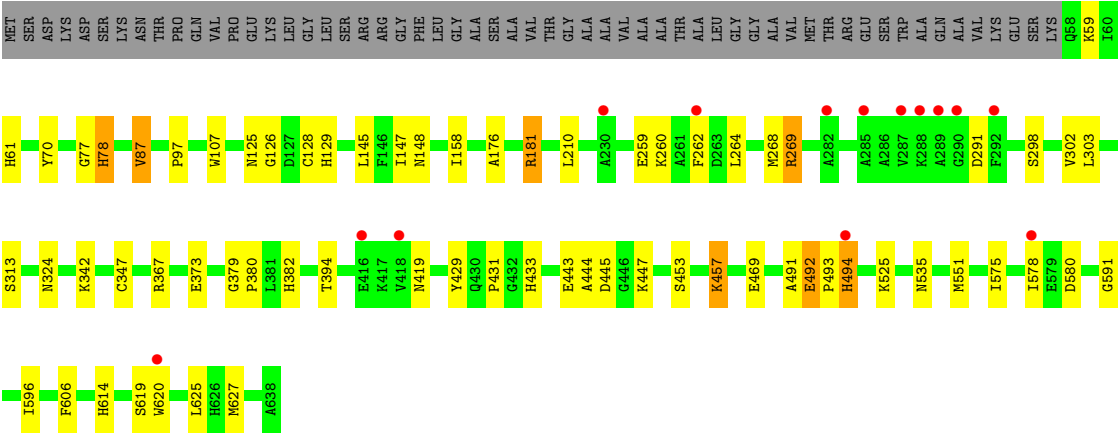
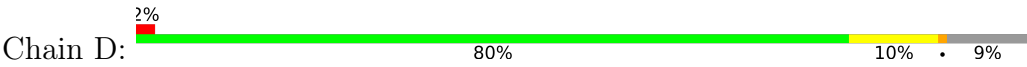


- Molecule 1: Nitrous-oxide reductase





● Molecule 1: Nitrous-oxide reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	108.92Å 73.37Å 136.34Å 90.00° 95.07° 90.00°	Depositor
Resolution (Å)	48.91 – 1.78 48.91 – 1.78	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.91-1.78) 100.0 (48.91-1.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 1.78Å)	Xtriage
Refinement program	REFMAC 5.8.0218	Depositor
R, R_{free}	0.157 , 0.192 0.167 , 0.201	Depositor DCC
R_{free} test set	10244 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	23.0	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 30.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	19996	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 39.59 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.1041e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: CL, K, NA, FMT, CUK, CUA, B3P, CA

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	1.32	12/4732 (0.3%)	1.10	14/6406 (0.2%)
1	B	1.29	11/4740 (0.2%)	1.09	13/6417 (0.2%)
1	C	1.32	5/4718 (0.1%)	1.12	12/6387 (0.2%)
1	D	1.28	12/4700 (0.3%)	1.06	7/6364 (0.1%)
All	All	1.30	40/18890 (0.2%)	1.09	46/25574 (0.2%)

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	324	ASN	C-O	9.69	1.28	1.23
1	A	128	CYS	C-O	7.69	1.32	1.23
1	A	87	VAL	CA-C	7.31	1.57	1.52
1	B	535	ASN	N-CA	-7.12	1.38	1.46
1	A	493	PRO	C-N	-7.07	1.23	1.33
1	B	619	SER	C-O	-6.62	1.15	1.24
1	A	87	VAL	N-CA	-6.59	1.40	1.46
1	D	492	GLU	N-CA	-6.49	1.36	1.46
1	C	128	CYS	C-O	-6.31	1.16	1.24
1	C	622	CYS	C-N	-6.14	1.25	1.33
1	C	493	PRO	C-N	-6.12	1.25	1.33
1	D	87	VAL	CA-C	6.06	1.56	1.52
1	D	494	HIS	N-CA	6.05	1.53	1.46
1	D	367	ARG	CZ-NH1	6.00	1.41	1.32
1	D	324	ASN	CA-C	5.93	1.57	1.53
1	D	298	SER	C-O	-5.72	1.17	1.23
1	A	572	ILE	C-O	-5.72	1.18	1.24
1	A	172	PRO	C-O	-5.62	1.17	1.23
1	D	181	ARG	CZ-NH2	5.60	1.40	1.33
1	C	529	ILE	C-O	-5.55	1.17	1.24
1	D	302	VAL	C-O	5.55	1.30	1.24
1	A	184	LYS	CA-C	5.46	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	494	HIS	N-CA	5.45	1.53	1.46
1	B	511	ASP	CB-CG	5.44	1.65	1.52
1	A	177	ILE	N-CA	-5.43	1.40	1.46
1	D	126	GLY	N-CA	-5.43	1.39	1.45
1	A	385	PHE	N-CA	5.41	1.52	1.46
1	A	437	SER	C-O	5.38	1.30	1.23
1	D	614	HIS	CA-C	-5.36	1.46	1.52
1	C	419	ASN	CG-OD1	5.33	1.33	1.23
1	B	302	VAL	C-O	5.27	1.30	1.24
1	B	586	VAL	N-CA	-5.19	1.40	1.46
1	D	535	ASN	C-O	5.18	1.30	1.23
1	B	499	ALA	CA-C	5.16	1.58	1.52
1	D	87	VAL	C-O	-5.13	1.19	1.23
1	A	298	SER	C-O	-5.08	1.18	1.23
1	B	128	CYS	CB-SG	-5.08	1.64	1.81
1	B	632	ARG	CD-NE	-5.07	1.39	1.46
1	B	493	PRO	C-N	-5.03	1.26	1.33
1	A	225	GLU	N-CA	-5.01	1.40	1.46

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	622	CYS	N-CA-C	-14.92	86.39	109.52
1	C	493	PRO	CA-C-N	-8.74	106.17	122.60
1	C	493	PRO	C-N-CA	-8.74	106.17	122.60
1	C	623	HIS	N-CA-C	8.68	129.28	110.80
1	D	493	PRO	CA-C-N	-8.37	103.72	121.80
1	D	493	PRO	C-N-CA	-8.37	103.72	121.80
1	B	493	PRO	CA-C-N	-8.26	103.96	121.80
1	B	493	PRO	C-N-CA	-8.26	103.96	121.80
1	A	493	PRO	O-C-N	-8.01	113.08	122.85
1	C	634	MET	CG-SD-CE	-7.89	83.55	100.90
1	B	493	PRO	O-C-N	-7.50	114.39	123.01
1	C	493	PRO	O-C-N	-7.49	114.40	123.01
1	B	364	ALA	N-CA-C	-7.31	96.64	108.41
1	A	493	PRO	CA-C-N	-7.25	106.19	121.64
1	A	493	PRO	C-N-CA	-7.25	106.19	121.64
1	B	275	VAL	N-CA-CB	7.03	119.43	110.99
1	D	491	ALA	CB-CA-C	-7.00	100.44	111.51
1	C	581	VAL	N-CA-C	6.81	113.49	106.21
1	B	128	CYS	CB-CA-C	-6.61	97.49	109.38
1	C	484	VAL	N-CA-C	6.51	119.47	112.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	491	ALA	N-CA-C	6.50	121.16	113.23
1	D	493	PRO	O-C-N	-6.47	115.57	123.01
1	C	494	HIS	N-CA-C	-6.36	104.22	112.68
1	D	269	ARG	CB-CA-C	-6.30	100.14	110.85
1	B	484	VAL	N-CA-C	5.97	120.41	113.42
1	B	563	LYS	CD-CE-NZ	5.96	130.97	111.90
1	B	353	ASP	CB-CA-C	-5.94	100.58	110.68
1	B	275	VAL	CB-CA-C	5.83	118.77	110.84
1	B	491	ALA	N-CA-C	5.78	119.63	112.58
1	A	262	PHE	CB-CA-C	5.77	120.57	110.64
1	C	609	ASP	CB-CG-OD2	-5.75	105.17	118.40
1	A	484	VAL	N-CA-C	5.71	119.13	113.53
1	B	275	VAL	CG1-CB-CG2	5.53	122.97	110.80
1	A	80	GLY	N-CA-C	5.52	122.09	114.92
1	D	491	ALA	N-CA-C	5.49	119.27	112.24
1	A	262	PHE	CA-C-N	5.45	130.83	122.93
1	A	262	PHE	C-N-CA	5.45	130.83	122.93
1	A	491	ALA	N-CA-C	5.36	119.10	112.24
1	A	128	CYS	CB-CA-C	-5.31	100.56	109.48
1	C	419	ASN	N-CA-CB	5.29	119.12	110.39
1	A	263	ASP	N-CA-CB	-5.24	102.25	110.69
1	A	273	ASP	N-CA-C	-5.17	104.54	108.78
1	B	551	MET	CG-SD-CE	-5.15	89.57	100.90
1	D	419	ASN	N-CA-CB	5.12	119.03	110.53
1	A	581	VAL	N-CA-C	5.11	111.67	106.21
1	A	57	LYS	CB-CA-C	-5.05	101.36	110.35

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	4610	0	4480	54	0
1	B	4615	0	4486	38	0
1	C	4596	0	4469	57	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	4581	0	4450	56	0
2	A	19	0	26	0	0
2	B	19	0	26	1	0
2	C	19	0	26	0	0
2	D	19	0	26	0	0
3	A	1	0	0	2	0
3	B	1	0	0	1	0
3	C	1	0	0	2	0
3	D	1	0	0	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	6	0	0	0	0
5	B	6	0	0	0	0
5	C	6	0	0	0	0
5	D	6	0	0	0	0
6	A	2	0	0	0	0
6	B	2	0	0	0	0
6	C	2	0	0	0	0
6	D	2	0	0	0	0
7	A	3	0	1	0	0
7	B	3	0	1	0	0
7	C	6	0	2	0	0
7	D	3	0	1	2	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
8	C	1	0	0	0	0
8	D	1	0	0	0	0
9	A	1	0	0	0	0
9	B	1	0	0	0	0
9	C	1	0	0	0	0
9	D	1	0	0	0	0
10	A	413	0	0	6	1
10	B	392	0	0	8	1
10	C	362	0	0	8	1
10	D	288	0	0	8	1
All	All	19996	0	17994	194	3

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 5.

All (194) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:271[B]:GLU:OE2	10:C:801:HOH:O	1.70	1.08
1:B:255:CYS:HB2	10:B:987:HOH:O	1.52	1.07
1:C:314:LYS:HD2	10:C:1131:HOH:O	1.58	1.02
1:B:288:LYS:HE2	10:B:1118:HOH:O	1.65	0.95
1:A:578:ILE:HG12	10:A:965:HOH:O	1.71	0.90
1:B:288:LYS:CE	10:B:1118:HOH:O	2.19	0.88
1:C:57:LYS:HG2	10:C:867:HOH:O	1.77	0.84
1:B:238[A]:ASN:OD1	1:B:257:ASN:HB3	1.80	0.82
1:A:238[B]:ASN:OD1	1:A:257:ASN:HB3	1.83	0.79
1:B:382[B]:HIS:CE1	1:B:433:HIS:HD1	2.00	0.78
1:A:469[A]:GLU:HG3	1:A:492:GLU:HA	1.66	0.78
1:B:382[A]:HIS:CE1	1:B:433:HIS:CE1	2.75	0.75
1:A:480:GLU:CD	10:A:816:HOH:O	2.29	0.75
1:C:382[B]:HIS:CE1	1:C:433:HIS:CE1	2.74	0.74
1:D:382[A]:HIS:CE1	1:D:433:HIS:HD1	2.05	0.73
1:D:606:PHE:HA	7:D:706:FMT:H	1.70	0.73
1:B:469[B]:GLU:HG3	1:B:492:GLU:HA	1.70	0.73
1:A:264:LEU:HD23	1:A:268:MET:HE1	1.71	0.73
1:A:382[A]:HIS:CE1	1:A:433:HIS:CE1	2.77	0.73
1:B:360:ALA:O	10:B:801:HOH:O	2.06	0.72
1:C:129:HIS:H	1:C:148:ASN:HD21	1.37	0.72
1:C:509:ILE:O	1:C:509:ILE:HD12	1.91	0.71
1:A:619:SER:O	1:D:78:HIS:CE1	2.44	0.71
1:C:469[A]:GLU:HG3	1:C:492:GLU:HA	1.71	0.71
1:B:283:VAL:HG13	1:B:303:LEU:HD13	1.73	0.70
1:C:579:GLU:O	10:C:802:HOH:O	2.09	0.70
1:A:382[B]:HIS:CE1	3:A:702:CL:CL	2.81	0.70
1:D:129:HIS:H	1:D:148:ASN:HD21	1.41	0.69
1:C:509:ILE:HD12	1:C:509:ILE:C	2.17	0.69
1:A:283:VAL:HG13	1:A:303:LEU:HD13	1.74	0.69
1:A:129:HIS:H	1:A:148:ASN:HD21	1.40	0.68
1:D:469[A]:GLU:HG3	1:D:492:GLU:HA	1.75	0.68
1:B:129:HIS:H	1:B:148:ASN:HD21	1.44	0.66
1:C:453:SER:O	1:C:469[B]:GLU:HG3	1.96	0.65
1:C:622:CYS:O	1:C:626:HIS:HB2	1.97	0.64
1:D:128:CYS:HB3	1:D:148:ASN:O	1.97	0.64
1:A:266:GLY:HA2	1:A:269:ARG:CZ	2.27	0.64
1:A:264:LEU:O	1:A:268:MET:HE2	1.97	0.64
1:B:288:LYS:HE3	10:B:1118:HOH:O	1.94	0.61
1:D:578:ILE:CD1	10:D:961:HOH:O	2.46	0.61
1:C:382[A]:HIS:CE1	1:C:433:HIS:HD1	2.18	0.61
1:D:578:ILE:HG13	10:D:961:HOH:O	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:264:LEU:HG	1:C:268:MET:HE2	1.82	0.61
1:A:382[B]:HIS:HE1	3:A:702:CL:CL	2.21	0.60
1:B:264:LEU:HG	1:B:268:MET:CE	2.32	0.60
1:A:260:LYS:O	1:A:262:PHE:CE1	2.55	0.60
1:C:129:HIS:CE1	1:C:494:HIS:CE1	2.90	0.59
1:A:210:LEU:HB2	1:A:262:PHE:CE1	2.37	0.59
1:D:129:HIS:CE1	1:D:494:HIS:CE1	2.90	0.59
1:A:559:GLU:OE1	1:A:632:ARG:NH2	2.36	0.58
1:C:382[A]:HIS:CE1	3:C:702:CL:CL	2.93	0.58
1:A:78:HIS:CE1	1:D:619:SER:O	2.57	0.58
1:A:264:LEU:HD23	1:A:268:MET:CE	2.33	0.58
1:D:210:LEU:HB2	1:D:262:PHE:CE2	2.37	0.58
1:A:382[B]:HIS:CE1	1:A:433:HIS:HD1	2.22	0.58
1:C:511:ASP:HB3	1:C:513:ASN:OD1	2.03	0.58
1:B:522:GLU:OE1	1:B:525:LYS:NZ	2.33	0.57
1:D:429:TYR:CE2	1:D:457:LYS:HE2	2.40	0.57
1:A:552:ALA:HB2	1:A:583:HIS:HE1	1.68	0.56
1:B:78:HIS:CE1	1:C:619:SER:O	2.59	0.56
1:D:494:HIS:HE1	10:D:824:HOH:O	1.88	0.56
1:C:128:CYS:HB2	1:C:147:ILE:HG12	1.88	0.56
1:C:329:ASN:HD21	1:C:382[A]:HIS:CE1	2.24	0.56
1:C:382[A]:HIS:HD2	1:C:434:LEU:H	1.53	0.55
1:A:411:ARG:NH2	10:A:806:HOH:O	2.37	0.55
1:C:382[A]:HIS:CE1	1:C:433:HIS:ND1	2.76	0.54
1:C:509:ILE:C	1:C:509:ILE:CD1	2.81	0.54
1:C:509:ILE:HD13	1:C:510:TRP:O	2.07	0.54
1:C:453:SER:O	1:C:469[B]:GLU:CG	2.55	0.54
1:D:210:LEU:HB3	1:D:262:PHE:CZ	2.44	0.53
1:C:580:ASP:HB3	10:C:802:HOH:O	2.07	0.53
1:A:329:ASN:HD21	1:A:382[B]:HIS:CE1	2.27	0.53
1:D:382[B]:HIS:NE2	1:D:433:HIS:CE1	2.77	0.53
1:B:264:LEU:CD1	1:B:268:MET:HE2	2.39	0.52
1:B:382[B]:HIS:CD2	3:B:702:CL:CL	2.99	0.52
1:A:77:GLY:HA2	1:A:128:CYS:O	2.10	0.52
1:A:181:ARG:HH22	1:A:382[B]:HIS:HE1	1.57	0.52
1:C:128:CYS:HB3	1:C:148:ASN:O	2.10	0.52
1:A:382[B]:HIS:HD2	1:A:434:LEU:H	1.58	0.51
1:D:494:HIS:CE1	10:D:824:HOH:O	2.64	0.51
1:D:578:ILE:HD12	10:D:961:HOH:O	2.07	0.50
1:B:266:GLY:HA2	1:B:269:ARG:CZ	2.41	0.50
1:A:145:LEU:HD22	1:A:158:ILE:HD12	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:259:GLU:HG2	1:D:269:ARG:HB2	1.93	0.49
1:A:380:PRO:HA	1:A:394:THR:O	2.13	0.49
1:D:551:MET:HE2	1:D:575:ILE:HD11	1.94	0.49
1:A:210:LEU:HB3	1:A:262:PHE:CZ	2.48	0.49
1:D:210:LEU:CB	1:D:262:PHE:CE2	2.95	0.49
1:B:275:VAL:HG13	1:B:321:VAL:HG21	1.95	0.49
1:A:431:PRO:HA	1:A:453:SER:HA	1.95	0.49
1:D:129:HIS:H	1:D:148:ASN:ND2	2.10	0.49
1:A:467:HIS:HB2	1:D:580:ASP:OD1	2.12	0.48
1:D:145:LEU:HD22	1:D:158:ILE:HD12	1.93	0.48
1:A:382[B]:HIS:CE1	1:A:433:HIS:ND1	2.81	0.48
1:D:578:ILE:CG1	10:D:961:HOH:O	2.58	0.48
1:B:55:GLU:HA	1:B:55:GLU:OE1	2.13	0.48
1:D:380:PRO:HA	1:D:394:THR:O	2.13	0.48
1:D:382[B]:HIS:CE1	1:D:433:HIS:CE1	3.01	0.48
1:D:77:GLY:HA2	1:D:128:CYS:O	2.13	0.48
1:D:453:SER:O	1:D:469[B]:GLU:HG3	2.13	0.48
1:A:443:GLU:O	1:A:444:ALA:C	2.57	0.48
1:C:70:TYR:OH	1:C:447:LYS:HD3	2.13	0.48
1:C:480:GLU:OE1	1:C:482:LYS:NZ	2.47	0.48
1:B:431:PRO:HA	1:B:453:SER:HA	1.96	0.47
1:A:268:MET:HE3	1:A:268:MET:HB2	1.73	0.47
1:A:288:LYS:HD2	10:A:801:HOH:O	2.15	0.47
1:B:443:GLU:O	1:B:444:ALA:C	2.57	0.47
1:C:431:PRO:HA	1:C:453:SER:HA	1.97	0.47
1:D:347:CYS:O	1:D:373:GLU:HA	2.15	0.47
1:D:431:PRO:HA	1:D:453:SER:HA	1.96	0.47
1:B:380:PRO:HA	1:B:394:THR:O	2.15	0.47
1:B:625:LEU:HD13	1:C:264:LEU:HD11	1.96	0.47
1:A:480:GLU:HG3	10:A:816:HOH:O	2.15	0.47
1:D:264:LEU:HG	1:D:268:MET:HE2	1.97	0.46
2:B:701:B3P:O5	2:B:701:B3P:H32	2.15	0.46
1:D:525:LYS:HE2	10:D:1069:HOH:O	2.15	0.46
1:A:326:HIS:CE1	1:D:627:MET:HE3	2.50	0.46
1:C:382[A]:HIS:HE1	3:C:702:CL:CL	2.35	0.46
1:D:443:GLU:O	1:D:444:ALA:C	2.58	0.46
1:C:129:HIS:H	1:C:148:ASN:ND2	2.09	0.46
1:D:382[B]:HIS:CE1	1:D:433:HIS:CG	3.03	0.46
1:B:77:GLY:HA2	1:B:128:CYS:O	2.15	0.46
1:C:77:GLY:HA2	1:C:128:CYS:O	2.16	0.46
1:B:70:TYR:HB2	1:B:87:VAL:HB	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:443:GLU:O	1:C:444:ALA:C	2.59	0.45
1:D:59:LYS:HG3	1:D:61:HIS:H	1.81	0.45
1:A:129:HIS:N	1:A:148:ASN:HD21	2.11	0.45
1:C:380:PRO:HA	1:C:394:THR:O	2.15	0.45
1:D:382[B]:HIS:CE1	1:D:433:HIS:CD2	3.04	0.45
1:B:260:LYS:HE3	10:B:1125:HOH:O	2.15	0.45
1:C:578:ILE:H	1:C:578:ILE:HD12	1.82	0.45
1:A:347:CYS:O	1:A:373:GLU:HA	2.16	0.45
1:A:382[B]:HIS:CD2	1:A:433:HIS:HD1	2.34	0.45
1:B:129:HIS:H	1:B:148:ASN:ND2	2.13	0.45
1:D:260:LYS:O	1:D:262:PHE:CE2	2.70	0.45
1:C:271[B]:GLU:CD	10:C:801:HOH:O	2.43	0.45
1:B:145:LEU:HD22	1:B:158:ILE:HD12	1.99	0.45
1:A:148:ASN:HD22	1:A:148:ASN:H	1.65	0.45
1:B:580:ASP:OD1	1:C:467:HIS:HB2	2.17	0.44
1:C:148:ASN:H	1:C:148:ASN:HD22	1.66	0.44
1:D:382[A]:HIS:CD2	3:D:702:CL:CL	3.07	0.44
1:B:125:ASN:OD1	1:C:591:GLY:HA2	2.17	0.44
1:A:210:LEU:HD13	1:A:262:PHE:CD2	2.52	0.44
1:B:55:GLU:HG3	1:B:478:GLY:HA3	1.98	0.44
1:C:101:VAL:HG21	1:C:124:LEU:HD22	2.00	0.44
1:C:353:ASP:HB2	10:C:893:HOH:O	2.18	0.44
1:D:382[B]:HIS:ND1	1:D:433:HIS:HA	2.33	0.44
1:A:181:ARG:HH22	1:A:382[B]:HIS:CE1	2.36	0.44
1:A:264:LEU:HD21	1:D:625:LEU:HD13	1.98	0.44
1:B:214:ASN:ND2	10:B:813:HOH:O	2.51	0.44
1:C:347:CYS:O	1:C:373:GLU:HA	2.17	0.44
1:D:148:ASN:HD22	1:D:148:ASN:H	1.66	0.43
1:C:97:PRO:HB2	1:C:107:TRP:CD1	2.53	0.43
1:A:125:ASN:OD1	1:D:591:GLY:HA2	2.18	0.43
1:D:70:TYR:HB2	1:D:87:VAL:HB	2.00	0.43
1:D:379:GLY:N	1:D:380:PRO:HD3	2.33	0.43
1:D:627:MET:HE1	10:D:1044:HOH:O	2.18	0.43
1:A:210:LEU:HD13	1:A:262:PHE:CE2	2.54	0.43
1:B:129:HIS:N	1:B:148:ASN:HD21	2.13	0.43
1:A:265:GLY:HA2	1:A:268:MET:CE	2.49	0.43
1:B:347:CYS:O	1:B:373:GLU:HA	2.19	0.43
1:D:606:PHE:HA	7:D:706:FMT:C	2.45	0.43
1:D:59:LYS:HE3	1:D:61:HIS:O	2.19	0.43
1:A:210:LEU:CB	1:A:262:PHE:CZ	3.02	0.42
1:C:181:ARG:HH22	1:C:382[A]:HIS:HE1	1.66	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:622:CYS:O	1:C:623:HIS:O	2.37	0.42
1:D:128:CYS:HB2	1:D:147:ILE:HG12	2.01	0.42
1:C:382[A]:HIS:CD2	1:C:433:HIS:HD1	2.36	0.42
1:A:70:TYR:HB2	1:A:87:VAL:HB	2.00	0.42
1:C:559:GLU:HA	1:C:632:ARG:O	2.19	0.42
1:C:379:GLY:N	1:C:380:PRO:HD3	2.35	0.42
1:C:78:HIS:CE1	1:C:492:GLU:HB3	2.54	0.42
1:B:632:ARG:HD3	10:B:988:HOH:O	2.19	0.41
1:A:97:PRO:HB2	1:A:107:TRP:CD1	2.55	0.41
1:A:264:LEU:CD2	1:A:268:MET:HE1	2.45	0.41
1:A:101:VAL:HG21	1:A:124:LEU:HD22	2.03	0.41
1:A:559:GLU:HA	1:A:632:ARG:O	2.21	0.41
1:C:269:ARG:HD2	10:C:815:HOH:O	2.20	0.41
1:A:591:GLY:HA2	1:D:125:ASN:OD1	2.20	0.41
1:B:148:ASN:H	1:B:148:ASN:HD22	1.68	0.41
1:C:303:LEU:HD22	1:C:303:LEU:N	2.35	0.41
1:C:535:ASN:OD1	1:C:556:GLY:HA3	2.20	0.41
1:C:382[A]:HIS:CE1	1:C:433:HIS:CE1	3.09	0.41
1:D:97:PRO:HB2	1:D:107:TRP:CD1	2.56	0.41
1:D:264:LEU:HG	1:D:268:MET:CE	2.51	0.41
1:A:90:MET:HG3	10:A:1067:HOH:O	2.21	0.41
1:B:559:GLU:HA	1:B:632:ARG:O	2.21	0.41
1:C:145:LEU:HD22	1:C:158:ILE:HD12	2.03	0.41
1:B:591:GLY:HA2	1:C:125:ASN:OD1	2.21	0.40
1:D:181:ARG:HH22	1:D:382[A]:HIS:HE2	1.70	0.40
1:D:303:LEU:HD22	1:D:303:LEU:N	2.37	0.40
1:D:453:SER:O	1:D:469[B]:GLU:CG	2.68	0.40
1:C:70:TYR:HB2	1:C:87:VAL:HB	2.04	0.40
1:D:445:ASP:O	1:D:447:LYS:HD2	2.21	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:C:1102:HOH:O	10:C:1136:HOH:O[1_565]	1.29	0.91
10:A:803:HOH:O	10:D:1060:HOH:O[1_565]	1.39	0.81
10:B:801:HOH:O	10:B:1080:HOH:O[2_555]	2.14	0.06

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	585/638 (92%)	560 (96%)	23 (4%)	2 (0%)	36	22
1	B	586/638 (92%)	561 (96%)	23 (4%)	2 (0%)	36	22
1	C	583/638 (91%)	560 (96%)	20 (3%)	3 (0%)	24	11
1	D	581/638 (91%)	558 (96%)	21 (4%)	2 (0%)	36	22
All	All	2335/2552 (92%)	2239 (96%)	87 (4%)	9 (0%)	30	16

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	623	HIS
1	A	342	LYS
1	B	342	LYS
1	B	365	ASP
1	C	342	LYS
1	D	342	LYS
1	C	176	ALA
1	D	176	ALA
1	A	176	ALA

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	502/537 (94%)	498 (99%)	4 (1%)	73	63
1	B	503/537 (94%)	497 (99%)	6 (1%)	63	49

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	500/537 (93%)	496 (99%)	4 (1%)	73	63
1	D	498/537 (93%)	492 (99%)	6 (1%)	63	49
All	All	2003/2148 (93%)	1983 (99%)	20 (1%)	68	55

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	HIS
1	A	147	ILE
1	A	596	ILE
1	A	620	TRP
1	B	78	HIS
1	B	101	VAL
1	B	275	VAL
1	B	313	SER
1	B	596	ILE
1	B	620	TRP
1	C	78	HIS
1	C	210	LEU
1	C	313	SER
1	C	620	TRP
1	D	78	HIS
1	D	291	ASP
1	D	313	SER
1	D	457	LYS
1	D	596	ILE
1	D	620	TRP

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

Mol	Chain	Res	Type
1	A	115	HIS
1	A	129	HIS
1	A	148	ASN
1	A	214	ASN
1	A	232	GLN
1	A	390	ASN
1	A	583	HIS
1	B	148	ASN
1	B	232	GLN

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Mol	Chain	Res	Type
1	B	390	ASN
1	B	599	GLN
1	C	61	HIS
1	C	115	HIS
1	C	129	HIS
1	C	148	ASN
1	C	211	GLN
1	C	390	ASN
1	C	494	HIS
1	D	129	HIS
1	D	148	ASN
1	D	214	ASN
1	D	232	GLN
1	D	241	ASN
1	D	281	HIS
1	D	390	ASN
1	D	467	HIS
1	D	494	HIS
1	D	599	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 33 ligands modelled in this entry, 16 are monoatomic - leaving 17 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	$\# Z > 2$
7	FMT	C	707	-	2,2,2	1.07	0	1,1,1	1.19	0
6	CUA	A	705	1	0,1,1	-	-	-		
6	CUA	D	705	1	0,1,1	-	-	-		
5	CUK	D	704	1	0,6,6	-	-	-		
6	CUA	C	705	1	0,1,1	-	-	-		
2	B3P	A	701	-	18,18,18	1.19	2 (11%)	23,23,23	1.84	8 (34%)
5	CUK	C	704	1	0,6,6	-	-	-		
2	B3P	D	701	-	18,18,18	0.79	0	23,23,23	1.68	4 (17%)
7	FMT	C	706	-	2,2,2	0.31	0	1,1,1	0.72	0
5	CUK	B	704	1	0,6,6	-	-	-		
7	FMT	A	706	-	2,2,2	0.27	0	1,1,1	0.85	0
7	FMT	B	706	-	2,2,2	0.06	0	1,1,1	0.26	0
7	FMT	D	706	-	2,2,2	0.14	0	1,1,1	0.66	0
6	CUA	B	705	1	0,1,1	-	-	-		
2	B3P	C	701	-	18,18,18	0.89	0	23,23,23	1.74	3 (13%)
2	B3P	B	701	-	18,18,18	0.92	0	23,23,23	1.59	4 (17%)
5	CUK	A	704	1	0,6,6	-	-	-		

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
5	CUK	D	704	1	-	-	0/1/1/1
2	B3P	A	701	-	-	2/28/28/28	-
5	CUK	C	704	1	-	-	0/1/1/1
2	B3P	D	701	-	-	3/28/28/28	-
5	CUK	B	704	1	-	-	0/1/1/1
2	B3P	C	701	-	-	2/28/28/28	-
2	B3P	B	701	-	-	1/28/28/28	-
5	CUK	A	704	1	-	-	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	B3P	C5-C4	3.31	1.57	1.53
2	A	701	B3P	C6-C4	2.02	1.55	1.53

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	701	B3P	C2-N2-C8	4.93	123.39	116.17
2	C	701	B3P	C2-N2-C8	4.74	123.11	116.17
2	B	701	B3P	C3-N1-C4	4.51	122.77	116.17
2	A	701	B3P	C6-C4-C5	3.56	117.83	110.02
2	A	701	B3P	C3-N1-C4	3.44	121.20	116.17
2	D	701	B3P	C3-N1-C4	3.38	121.11	116.17
2	B	701	B3P	C1-C3-N1	3.29	118.94	110.98
2	C	701	B3P	C3-N1-C4	2.98	120.52	116.17
2	D	701	B3P	C1-C2-N2	2.94	118.09	110.98
2	A	701	B3P	C6-C4-N1	-2.79	100.68	109.02
2	B	701	B3P	O4-C5-C4	-2.68	106.23	111.68
2	A	701	B3P	C1-C2-N2	2.67	117.44	110.98
2	A	701	B3P	C1-C3-N1	2.42	116.83	110.98
2	D	701	B3P	O3-C11-C8	-2.21	107.19	111.68
2	A	701	B3P	C5-C4-N1	2.14	115.40	109.02
2	A	701	B3P	O2-C10-C8	-2.13	107.34	111.68
2	B	701	B3P	C11-C8-C10	-2.09	105.43	110.02
2	C	701	B3P	C6-C4-N1	2.03	115.08	109.02
2	A	701	B3P	C7-C4-C6	-2.03	105.58	110.02

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	701	B3P	O2-C10-C8-C9
2	C	701	B3P	N2-C8-C9-O1
2	A	701	B3P	C7-C4-C6-O5
2	B	701	B3P	C7-C4-C6-O5
2	A	701	B3P	N1-C4-C6-O5
2	D	701	B3P	O2-C10-C8-N2
2	C	701	B3P	C10-C8-C9-O1
2	D	701	B3P	C9-C8-N2-C2

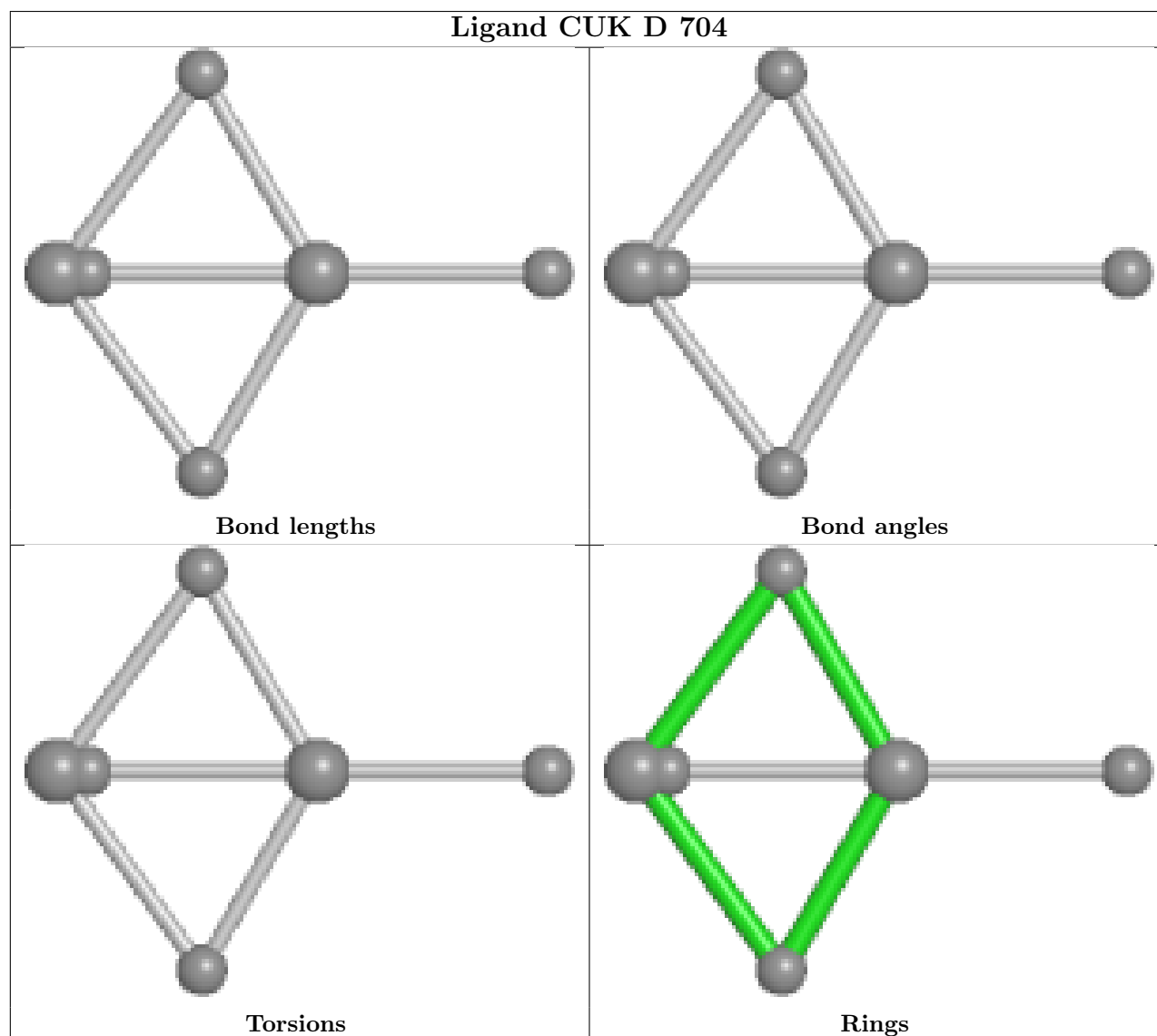
There are no ring outliers.

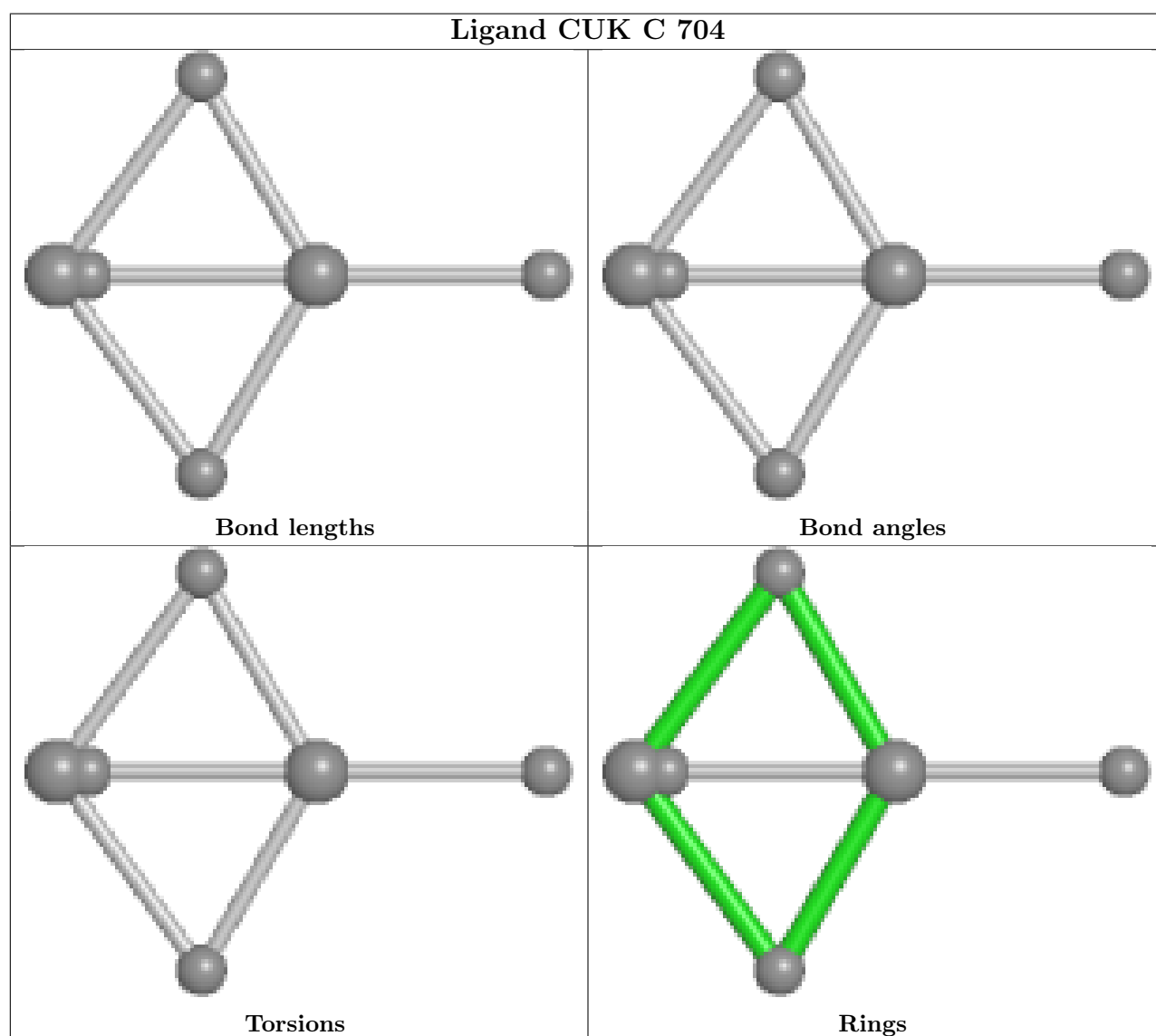
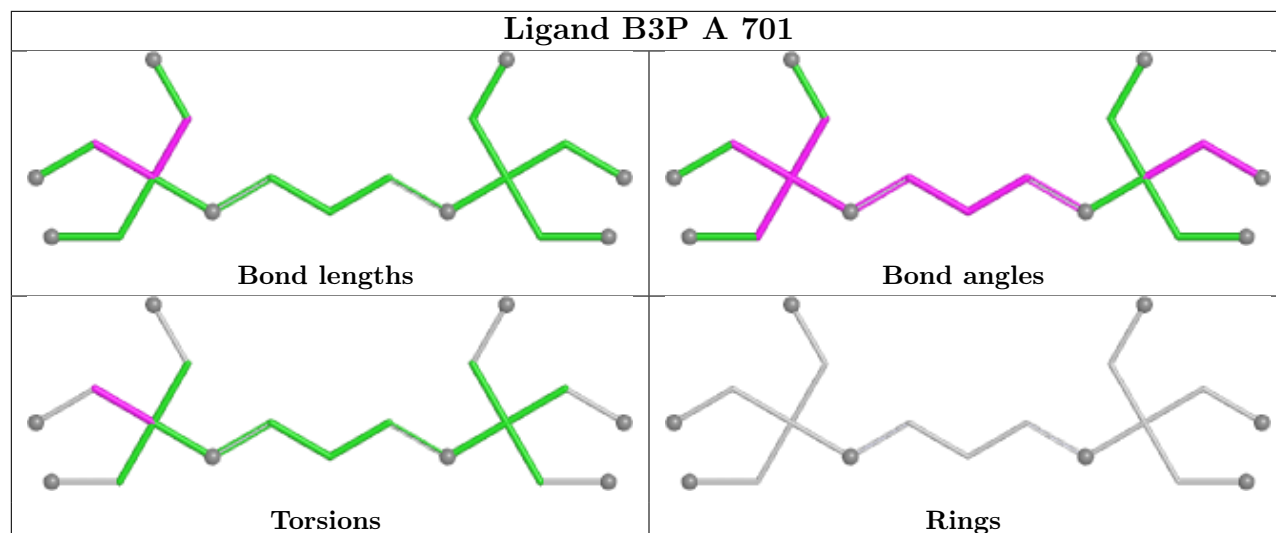
2 monomers are involved in 3 short contacts:

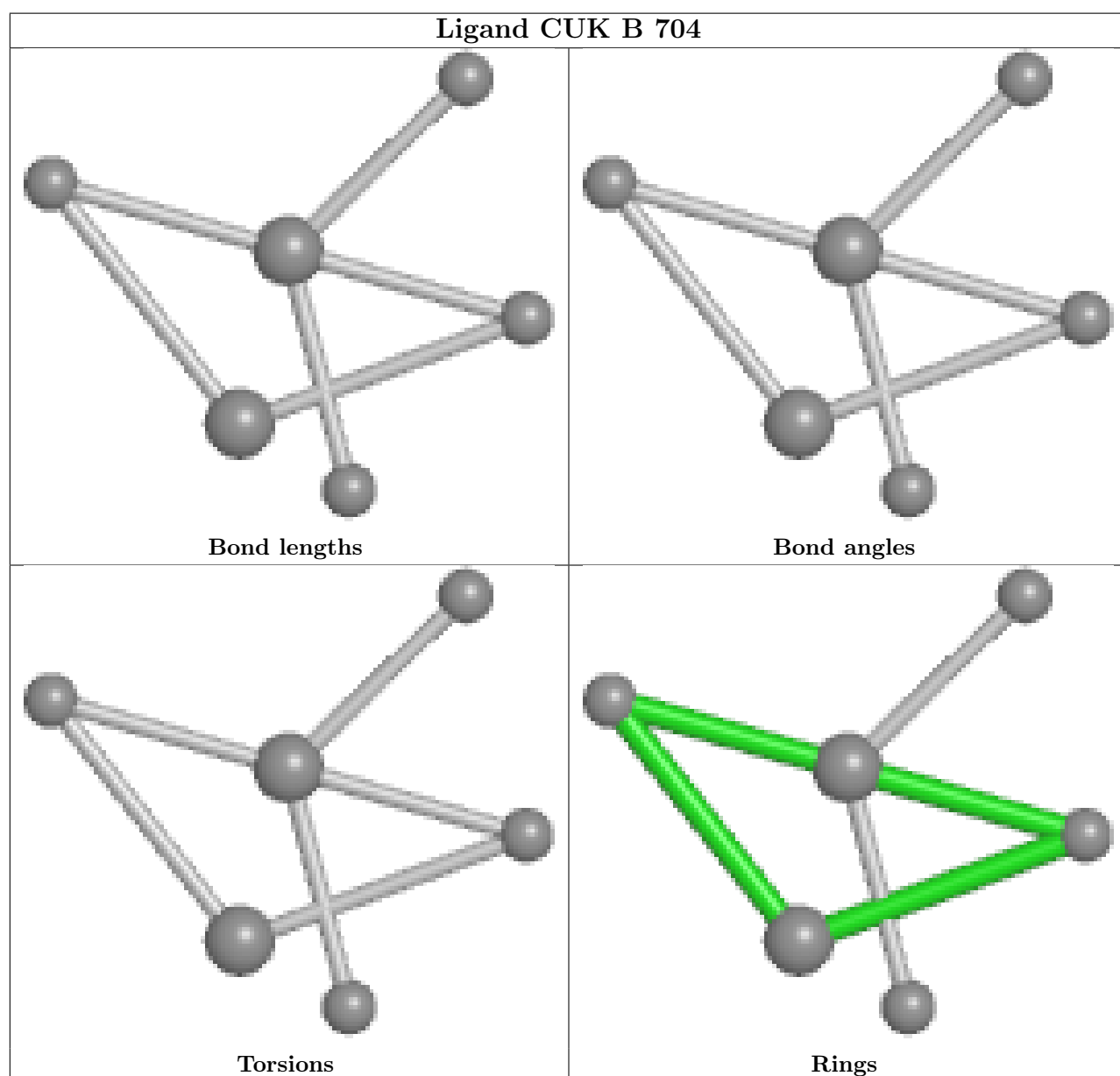
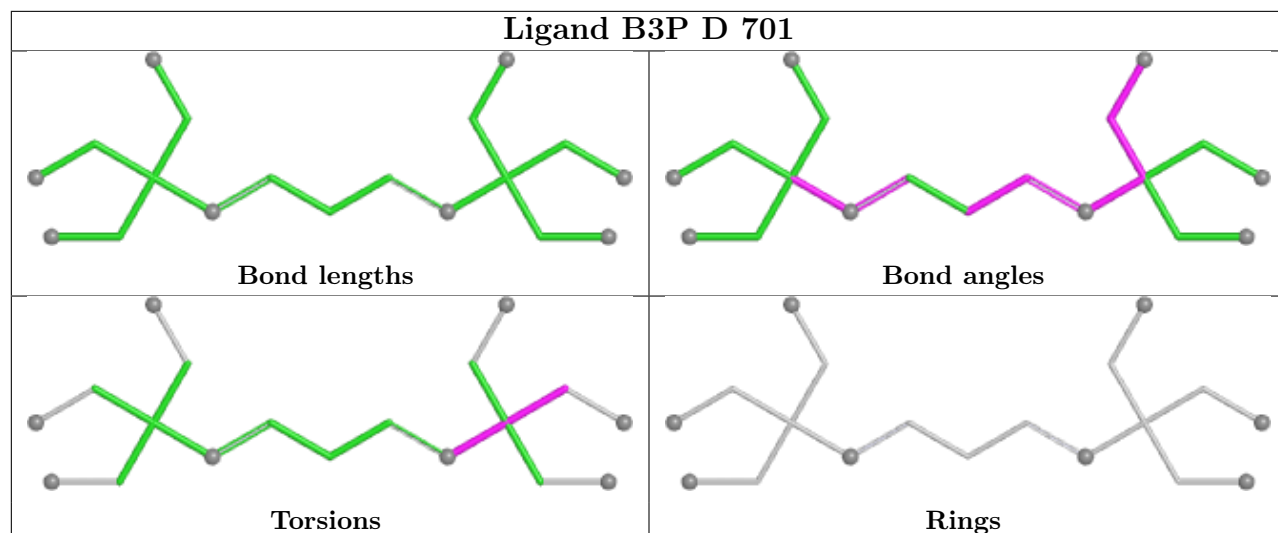
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	D	706	FMT	2	0
2	B	701	B3P	1	0

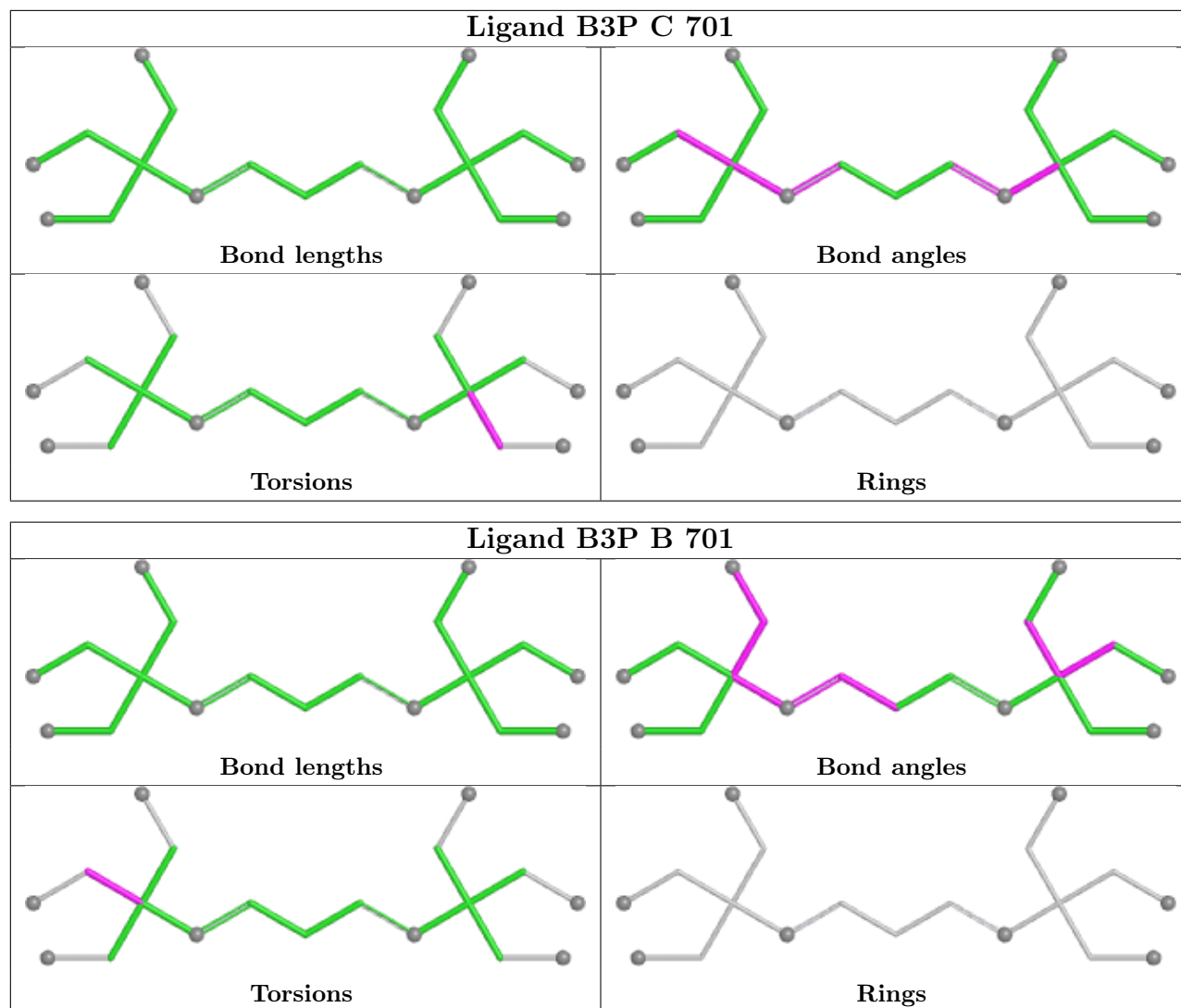
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

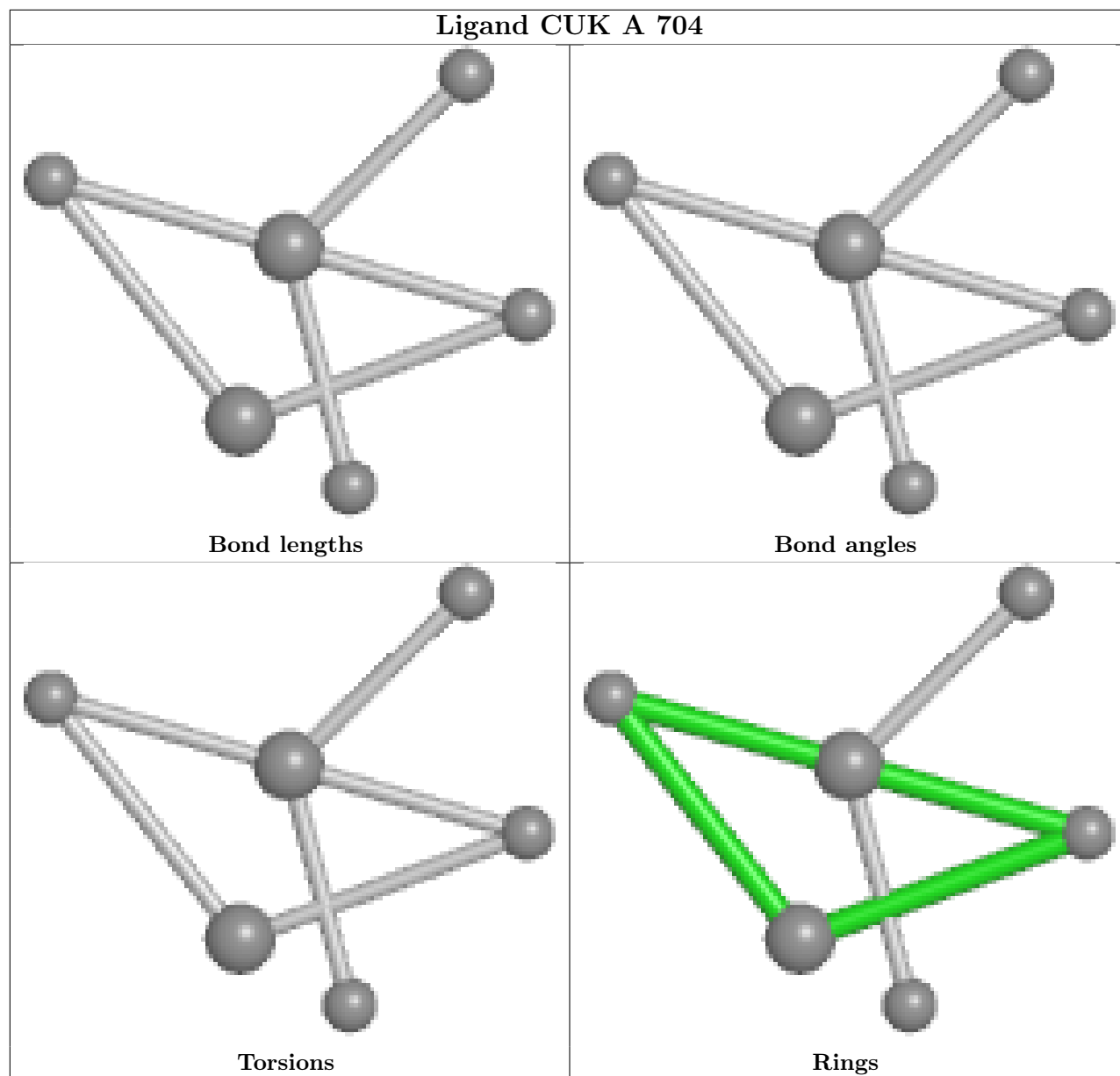
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 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 [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	584/638 (91%)	-0.17	9 (1%) 72 78	12, 24, 46, 75	3 (0%)
1	B	584/638 (91%)	-0.17	10 (1%) 69 77	14, 25, 47, 83	4 (0%)
1	C	582/638 (91%)	-0.14	7 (1%) 76 82	11, 25, 46, 67	3 (0%)
1	D	581/638 (91%)	0.01	14 (2%) 59 68	13, 28, 54, 81	2 (0%)
All	All	2331/2552 (91%)	-0.12	40 (1%) 69 77	11, 26, 49, 83	12 (0%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	262	PHE	5.9
1	B	262	PHE	4.8
1	D	289	ALA	4.3
1	A	265	GLY	3.8
1	A	261	ALA	3.7
1	A	263	ASP	3.4
1	A	264	LEU	3.4
1	D	262	PHE	3.4
1	D	288	LYS	3.2
1	A	494	HIS	3.0
1	B	261	ALA	2.8
1	B	364	ALA	2.8
1	C	60	ILE	2.8
1	C	578	ILE	2.8
1	D	292	PHE	2.8
1	B	494	HIS	2.8
1	D	287	VAL	2.8
1	D	282	ALA	2.7
1	D	285	ALA	2.6
1	B	263	ASP	2.6
1	C	57	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	418	VAL	2.5
1	B	264	LEU	2.4
1	A	56	SER	2.3
1	B	270	ASN	2.3
1	C	622	CYS	2.3
1	D	494	HIS	2.2
1	C	61	HIS	2.2
1	C	638	ALA	2.2
1	D	578	ILE	2.2
1	D	290	GLY	2.2
1	D	230	ALA	2.1
1	B	269	ARG	2.1
1	C	494	HIS	2.1
1	D	416	GLU	2.1
1	B	363	LEU	2.1
1	D	620	TRP	2.1
1	A	266	GLY	2.0
1	A	360	ALA	2.0
1	B	638	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	FMT	C	707	3/3	0.80	0.12	35,35,39,41	0
5	CUK	A	704	6/6	0.87	0.17	26,39,46,50	6
5	CUK	B	704	6/6	0.88	0.16	28,43,46,48	6
7	FMT	A	706	3/3	0.93	0.12	26,26,27,39	0

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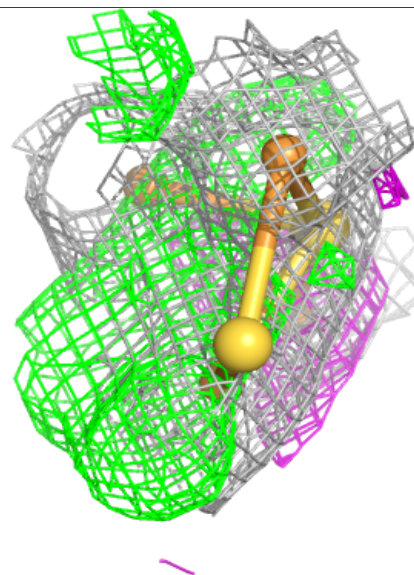
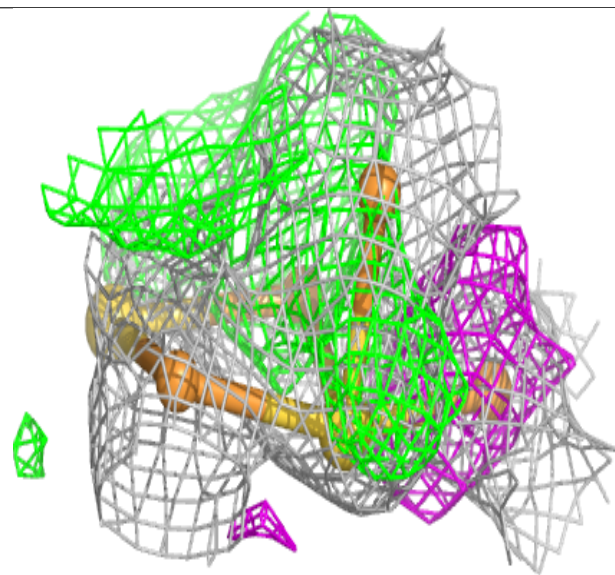
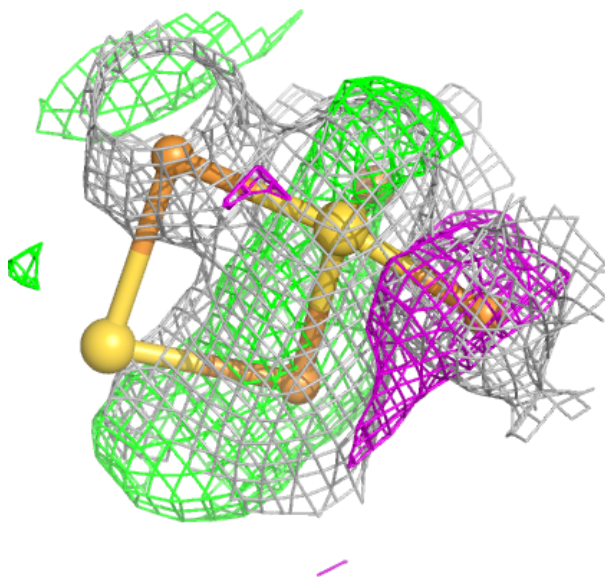
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	B3P	D	701	19/19	0.93	0.07	22,24,29,30	0
2	B3P	C	701	19/19	0.94	0.07	22,23,30,42	0
5	CUK	C	704	6/6	0.94	0.09	23,29,31,38	6
5	CUK	D	704	6/6	0.94	0.09	25,32,35,36	6
2	B3P	A	701	19/19	0.94	0.07	21,24,31,45	0
7	FMT	B	706	3/3	0.94	0.15	22,22,28,37	0
2	B3P	B	701	19/19	0.94	0.07	21,23,29,32	0
7	FMT	D	706	3/3	0.94	0.13	26,26,28,42	0
9	NA	D	708	1/1	0.94	0.07	34,34,34,34	1
8	K	A	707	1/1	0.95	0.08	34,34,34,34	1
9	NA	C	709	1/1	0.95	0.08	30,30,30,30	0
7	FMT	C	706	3/3	0.95	0.09	30,30,31,36	0
8	K	B	707	1/1	0.96	0.07	35,35,35,35	1
9	NA	B	708	1/1	0.96	0.05	26,26,26,26	0
8	K	C	708	1/1	0.97	0.07	30,30,30,30	1
3	CL	B	702	1/1	0.98	0.04	28,28,28,28	1
6	CUA	A	705	2/2	0.98	0.03	25,25,25,29	0
8	K	D	707	1/1	0.98	0.07	30,30,30,30	1
4	CA	D	703	1/1	0.99	0.04	26,26,26,26	0
6	CUA	B	705	2/2	0.99	0.03	24,24,24,29	0
6	CUA	C	705	2/2	0.99	0.03	26,26,26,33	0
6	CUA	D	705	2/2	0.99	0.03	28,28,28,34	0
3	CL	A	702	1/1	0.99	0.02	22,22,22,22	1
9	NA	A	708	1/1	0.99	0.02	27,27,27,27	0
3	CL	D	702	1/1	0.99	0.03	26,26,26,26	1
4	CA	A	703	1/1	0.99	0.03	27,27,27,27	1
4	CA	B	703	1/1	0.99	0.05	27,27,27,27	1
3	CL	C	702	1/1	1.00	0.02	26,26,26,26	1
4	CA	C	703	1/1	1.00	0.04	22,22,22,22	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.

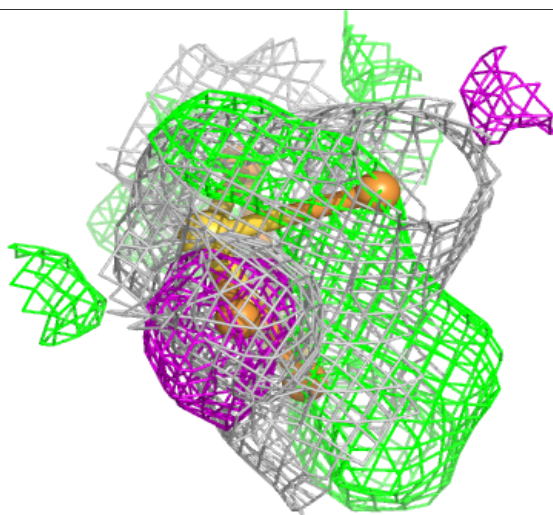
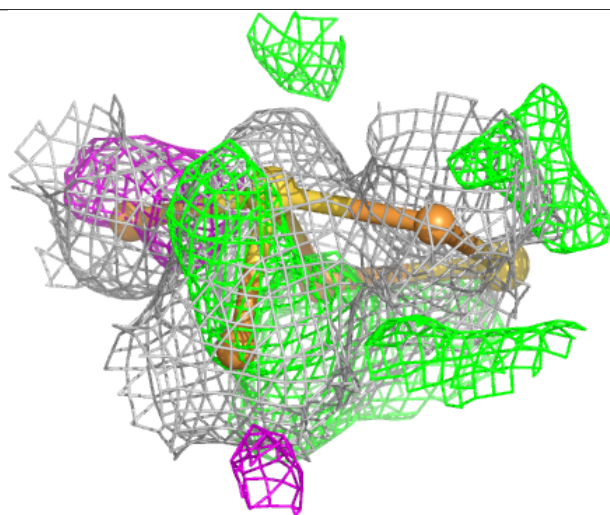
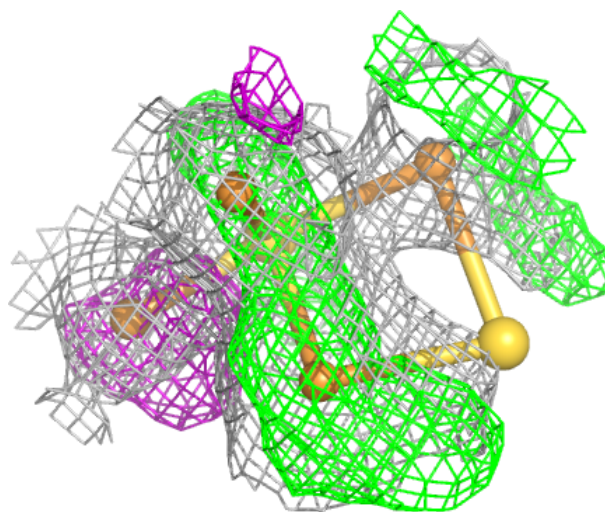
Electron density around CUK A 704:

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



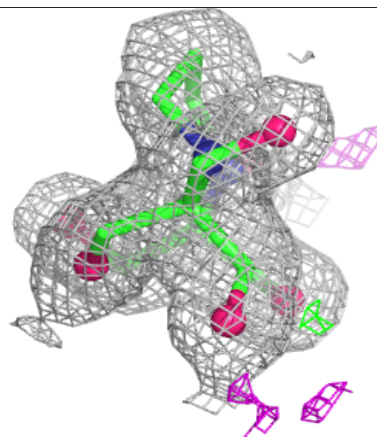
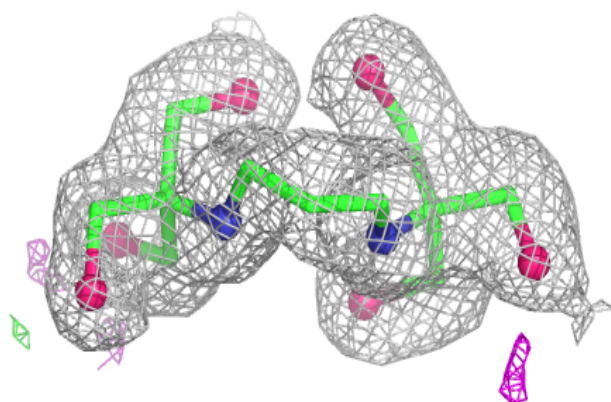
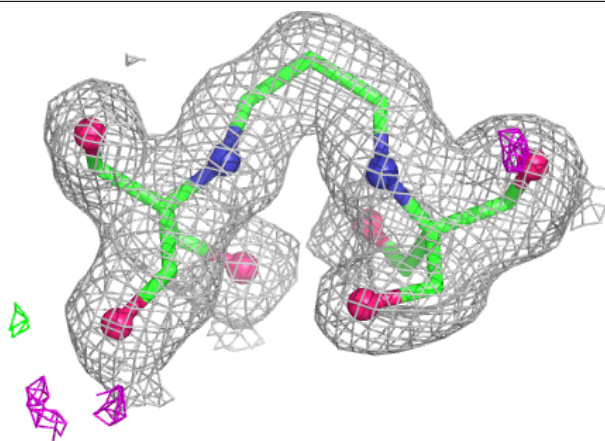
Electron density around CUK B 704:

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



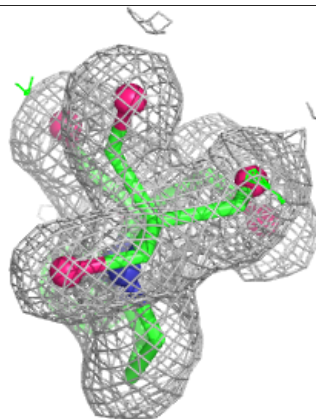
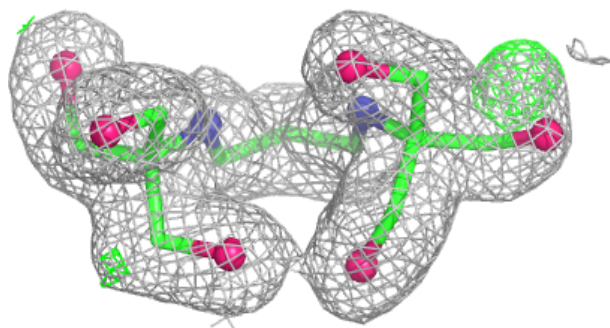
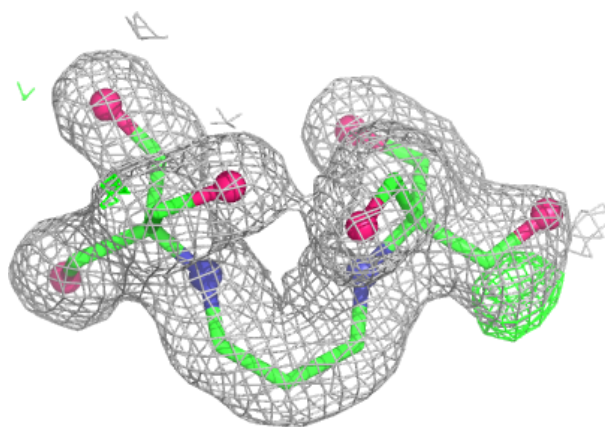
Electron density around B3P D 701:

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



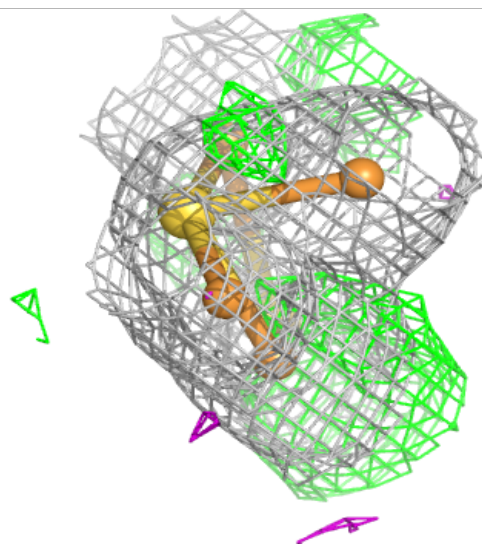
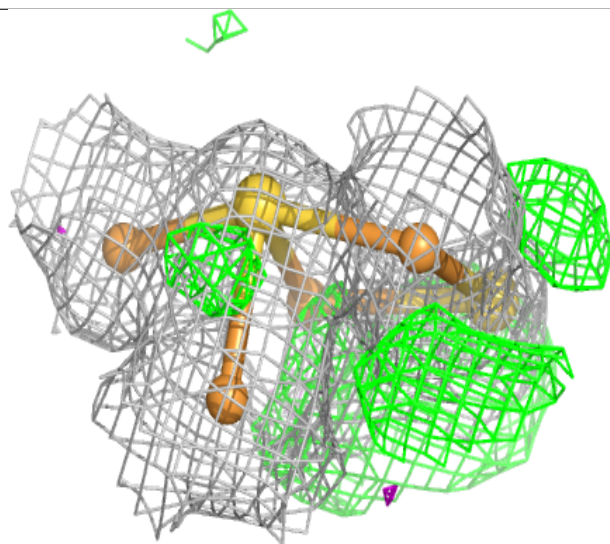
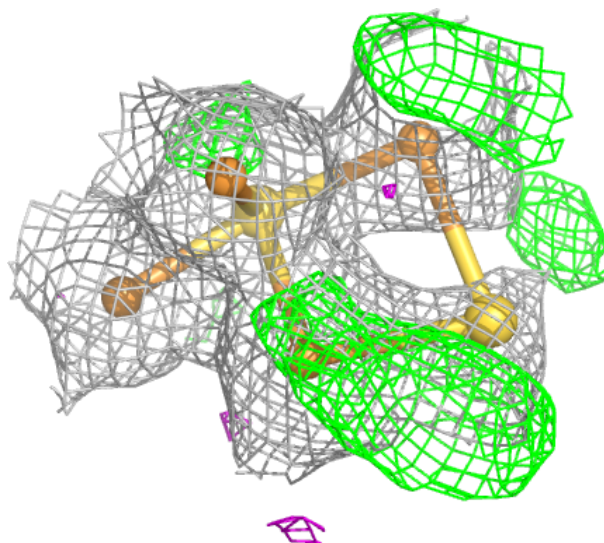
Electron density around B3P C 701:

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



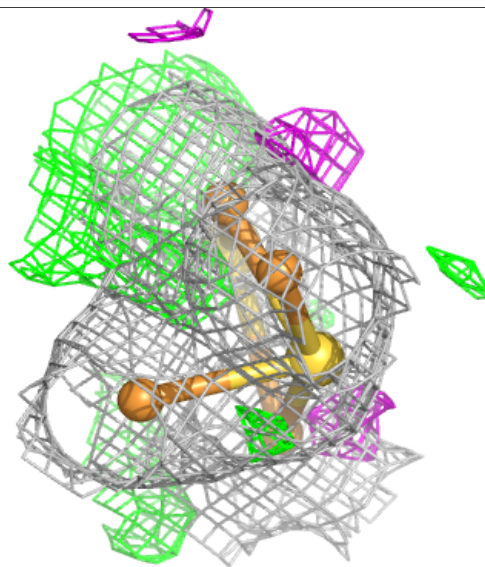
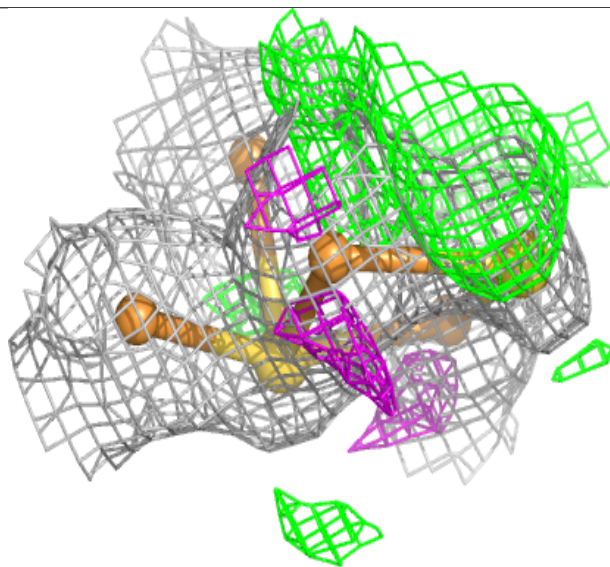
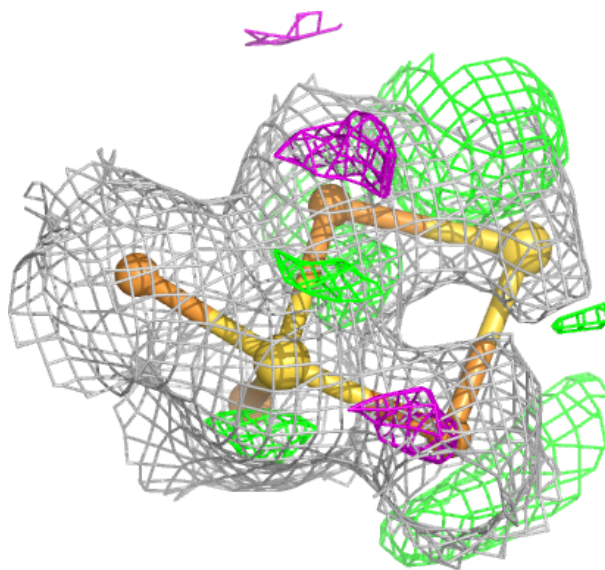
Electron density around CUK C 704:

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and green (positive)



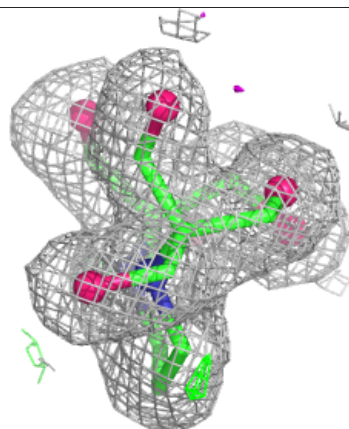
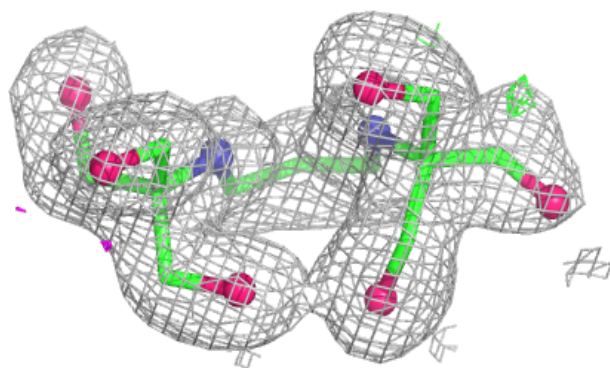
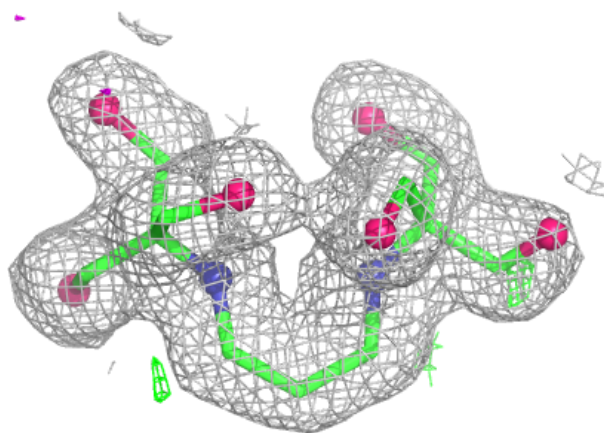
Electron density around CUK D 704:

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

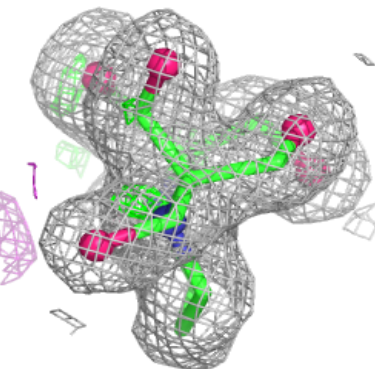
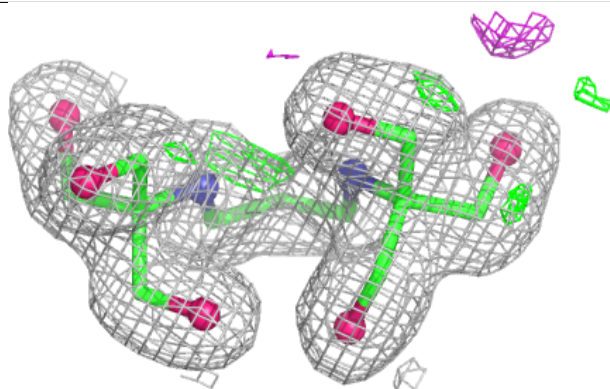
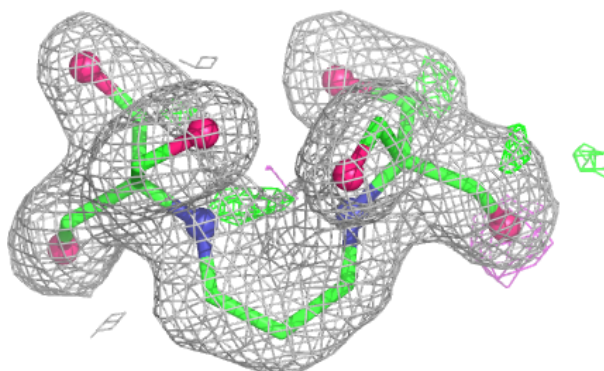


Electron density around B3P A 701:

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

**Electron density around B3P B 701:**

$2mF_o-DF_c$ (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.