



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:44 PM UTC

PDB ID : 6X5K / pdb_00006x5k
Title : Crystal structure of CODH/ACS with carbon monoxide bound to the A-cluster
Authors : Cohen, S.E.; Wittenborn, E.C.; Hendrickson, R.; Drennan, C.L.
Deposited on : 2020-05-26
Resolution : 2.47 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

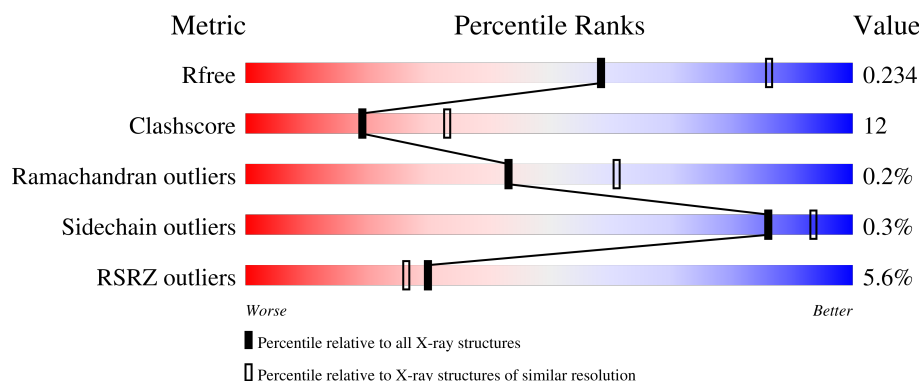
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.47 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	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (2.50-2.46)

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	674	% 76% 24%
1	B	674	% 77% 22%
1	C	674	2% 77% 22%
1	D	674	2% 76% 23% .
2	M	729	% 78% 22%

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Mol	Chain	Length	Quality of chain
2	N	729	% 78% 21%
2	O	729	26% 70% 29%
2	P	729	9% 74% 25%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	XCC	A	703	-	-	X	-
4	XCC	B	702	-	-	X	-
4	XCC	C	703	-	-	X	-
4	XCC	D	702	-	-	X	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 44026 atoms, of which 32 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.

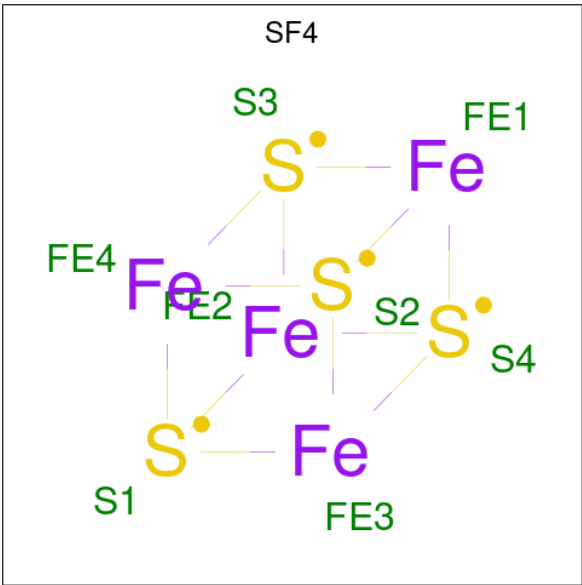
- Molecule 1 is a protein called Carbon monoxide dehydrogenase/acetyl-CoA synthase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	672	Total	C	N	O	S	0	3	0
			5109	3209	895	961	44			
1	B	672	Total	C	N	O	S	0	3	0
			5109	3209	895	961	44			
1	C	672	Total	C	N	O	S	0	3	0
			5109	3209	895	961	44			
1	D	672	Total	C	N	O	S	0	2	0
			5103	3206	894	960	43			

- Molecule 2 is a protein called Carbon monoxide dehydrogenase/acetyl-CoA synthase subunit alpha.

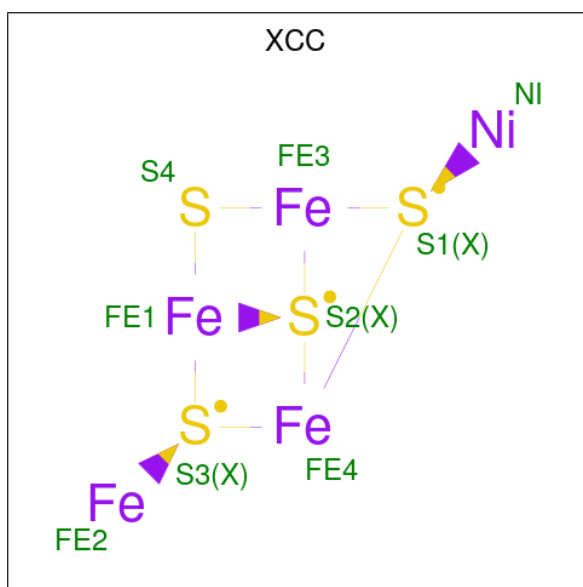
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	M	728	Total	C	N	O	S	0	0	0
			5740	3681	956	1068	35			
2	N	728	Total	C	N	O	S	0	0	0
			5740	3681	956	1068	35			
2	O	726	Total	C	N	O	S	0	0	0
			5723	3670	954	1065	34			
2	P	728	Total	C	N	O	S	0	0	0
			5740	3681	956	1068	35			

- Molecule 3 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄) (labeled as "Ligand of Interest" by depositor).



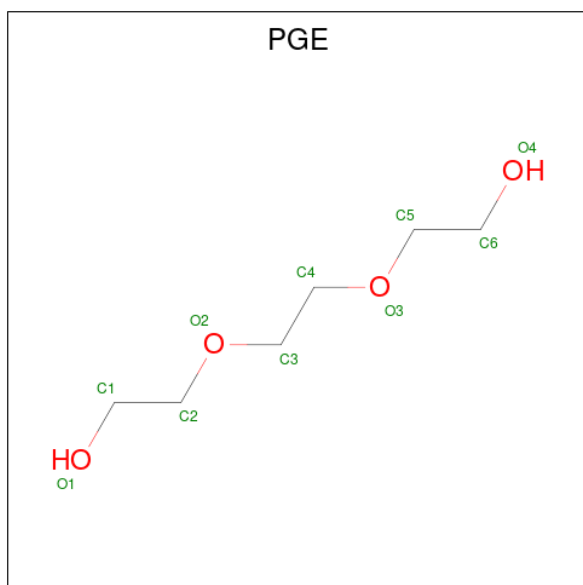
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Fe	S	0	0
			8	4	4		
3	A	1	Total	Fe	S	0	0
			8	4	4		
3	B	1	Total	Fe	S	0	0
			8	4	4		
3	C	1	Total	Fe	S	0	0
			8	4	4		
3	C	1	Total	Fe	S	0	0
			8	4	4		
3	D	1	Total	Fe	S	0	0
			8	4	4		
3	M	1	Total	Fe	S	0	0
			8	4	4		
3	N	1	Total	Fe	S	0	0
			8	4	4		
3	O	1	Total	Fe	S	0	0
			8	4	4		
3	P	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 4 is FE(4)-NI(1)-S(4) CLUSTER (CCD ID: XCC) (formula: Fe₄NiS₄).



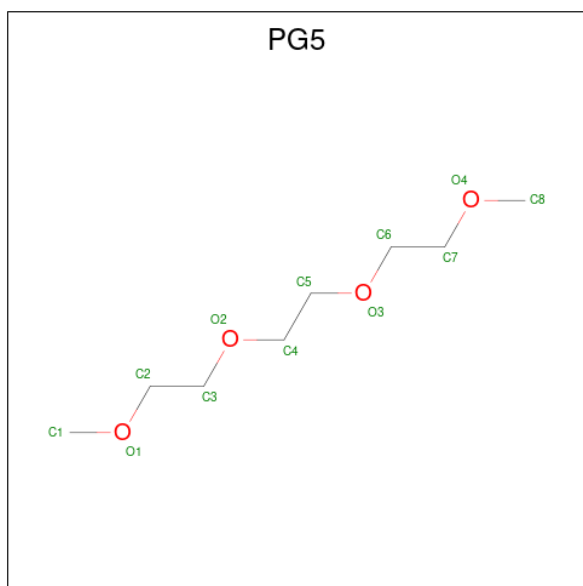
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	Fe	Ni	S	0	0
			9	4	1	4		
4	B	1	Total	Fe	Ni	S	0	0
			9	4	1	4		
4	C	1	Total	Fe	Ni	S	0	0
			9	4	1	4		
4	D	1	Total	Fe	Ni	S	0	0
			9	4	1	4		

- Molecule 5 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	0	0
			24	6	14	4		

- Molecule 6 is 1-METHOXY-2-[2-(2-METHOXY-ETHOXY)]-ETHANE (CCD ID: PG5) (formula: C₈H₁₈O₄).

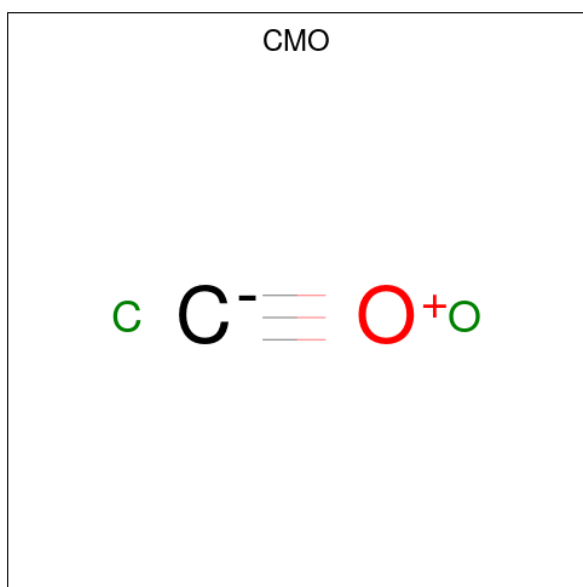


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	C	1	Total	C	H	O	0	0
			30	8	18	4		

- Molecule 7 is NICKEL (II) ION (CCD ID: NI) (formula: Ni) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	M	2	Total	Ni	0	0
			2	2		
7	N	2	Total	Ni	0	0
			2	2		
7	O	2	Total	Ni	0	0
			2	2		
7	P	2	Total	Ni	0	0
			2	2		

- Molecule 8 is CARBON MONOXIDE (CCD ID: CMO) (formula: CO) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	M	1	Total	C	O	0	0
			2	1	1		
8	N	1	Total	C	O	0	0
			2	1	1		
8	P	1	Total	C	O	0	0
			2	1	1		

- Molecule 9 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	N	1	Total	Mg	0	0
			1	1		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	84	Total	O	0	0
			84	84		
10	B	77	Total	O	0	0
			77	77		
10	C	44	Total	O	0	0
			44	44		
10	D	51	Total	O	0	0
			51	51		
10	M	76	Total	O	0	0
			76	76		

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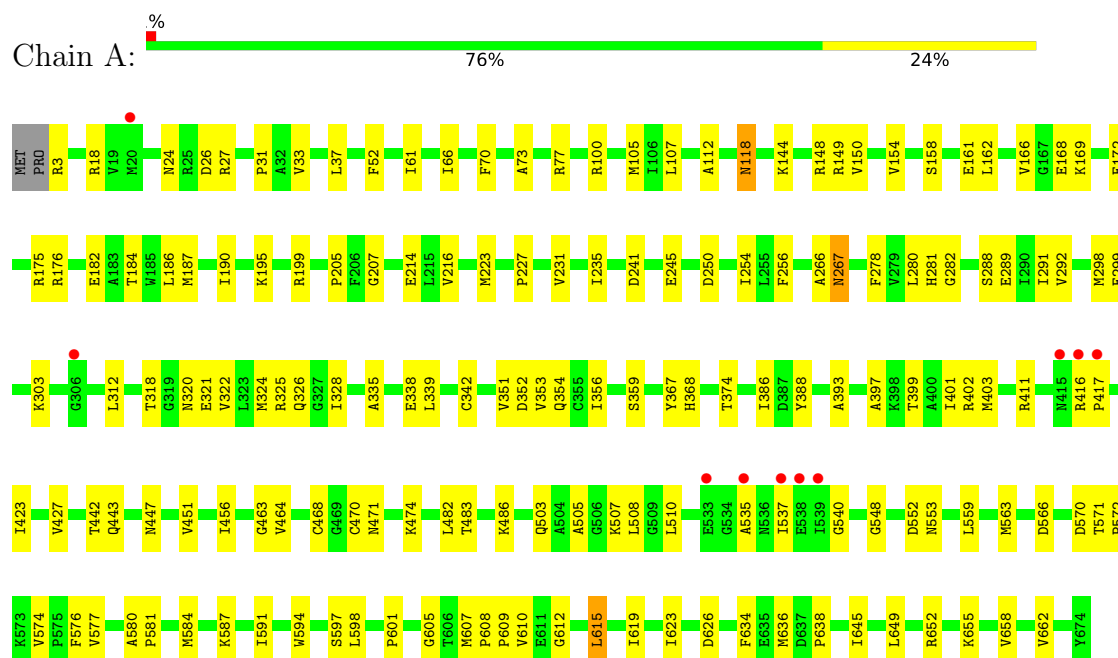
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	N	79	Total 79	O 79	0	0
10	O	12	Total 12	O 12	0	0
10	P	45	Total 45	O 45	0	0

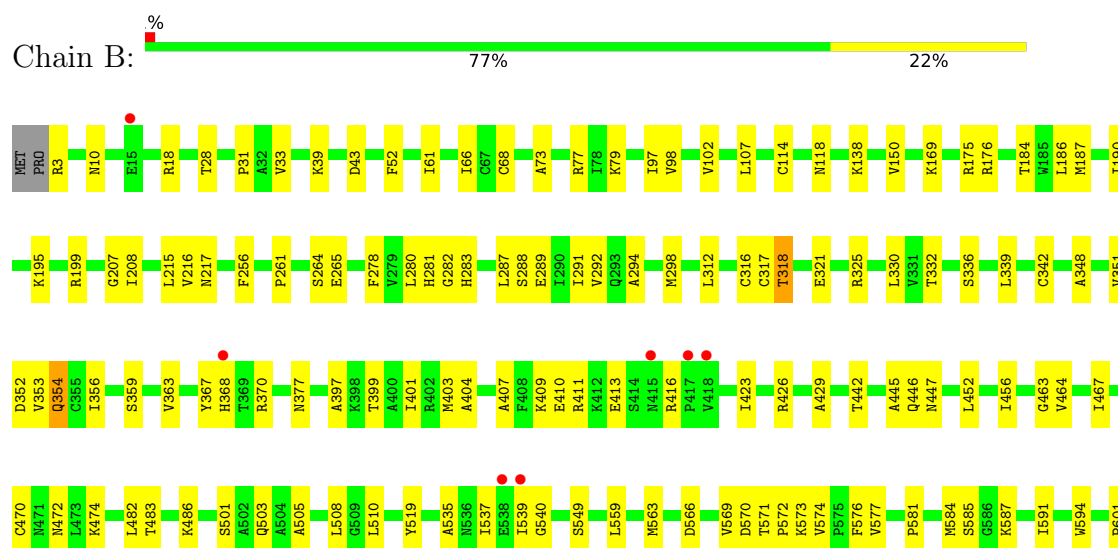
3 Residue-property plots [i](#)

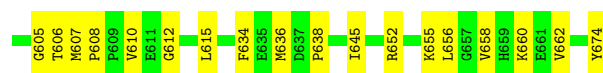
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Carbon monoxide dehydrogenase/acetyl-CoA synthase subunit beta

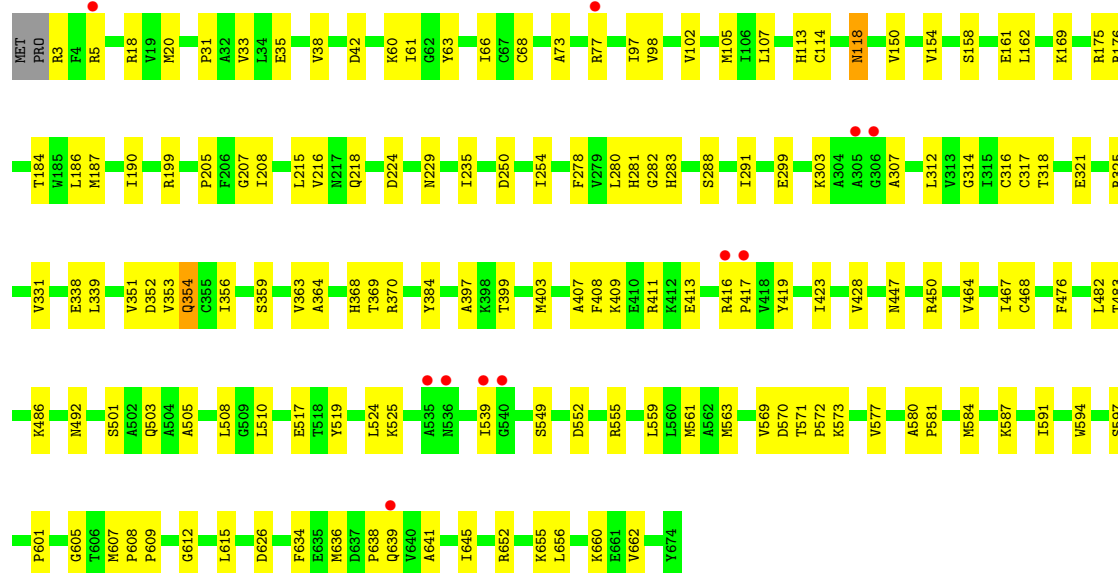
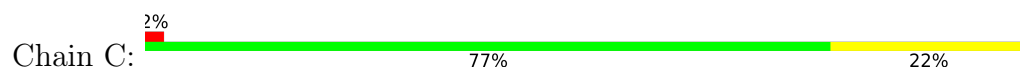


- Molecule 1: Carbon monoxide dehydrogenase/acetyl-CoA synthase subunit beta

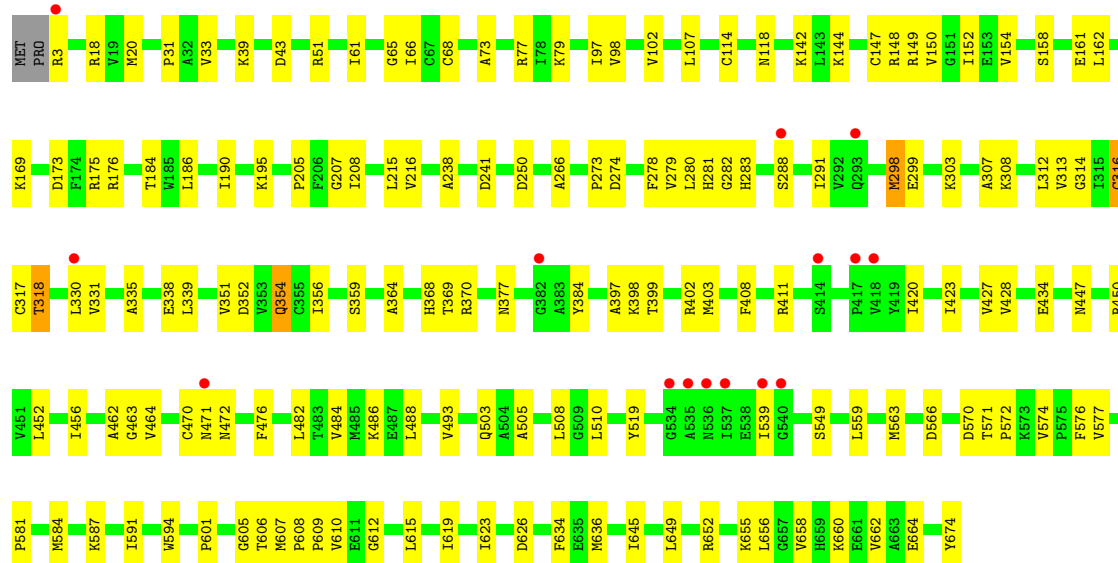
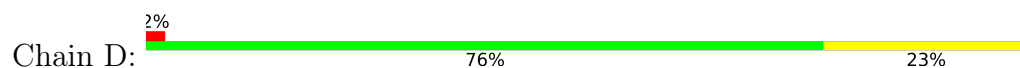




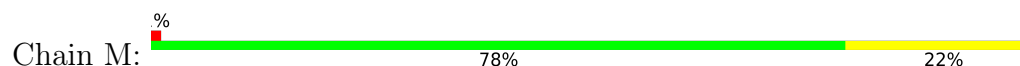
- Molecule 1: Carbon monoxide dehydrogenase/acetyl-CoA synthase subunit beta

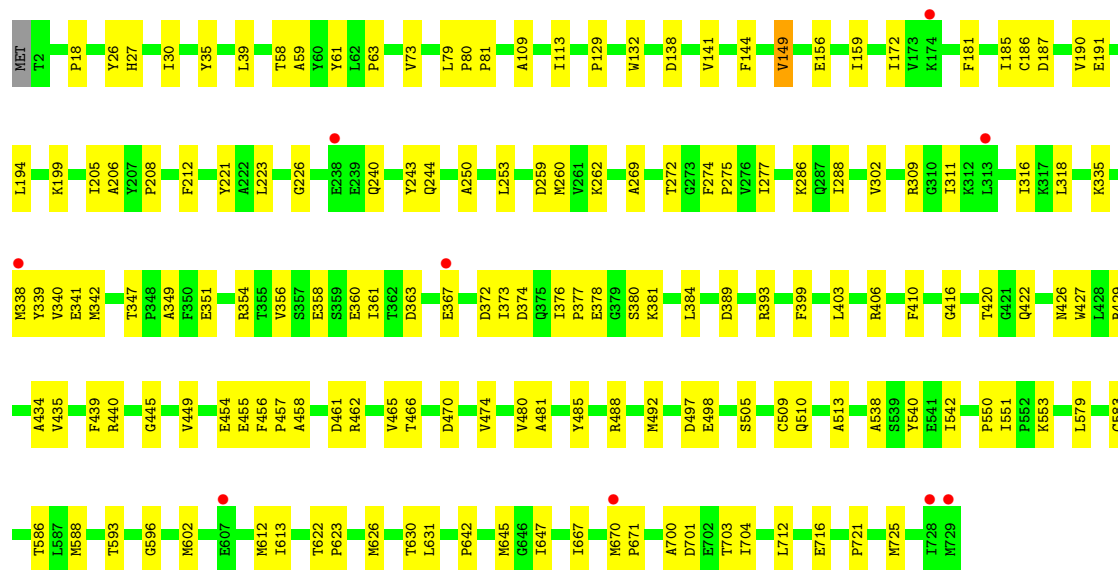


- Molecule 1: Carbon monoxide dehydrogenase/acetyl-CoA synthase subunit beta



- Molecule 2: Carbon monoxide dehydrogenase/acetyl-CoA synthase subunit alpha

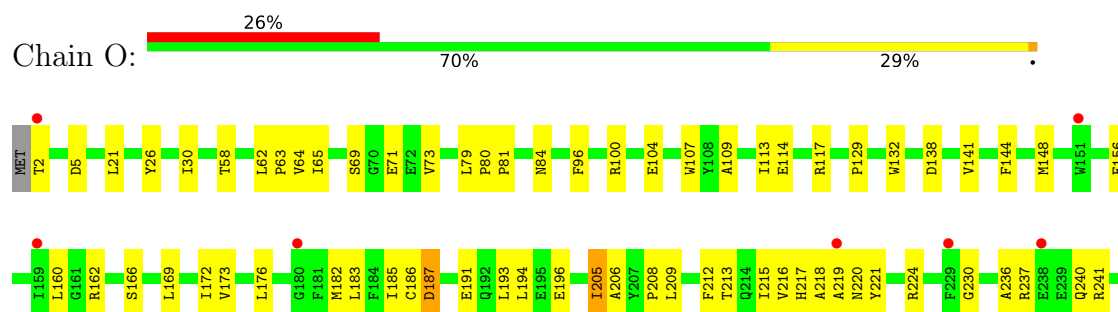


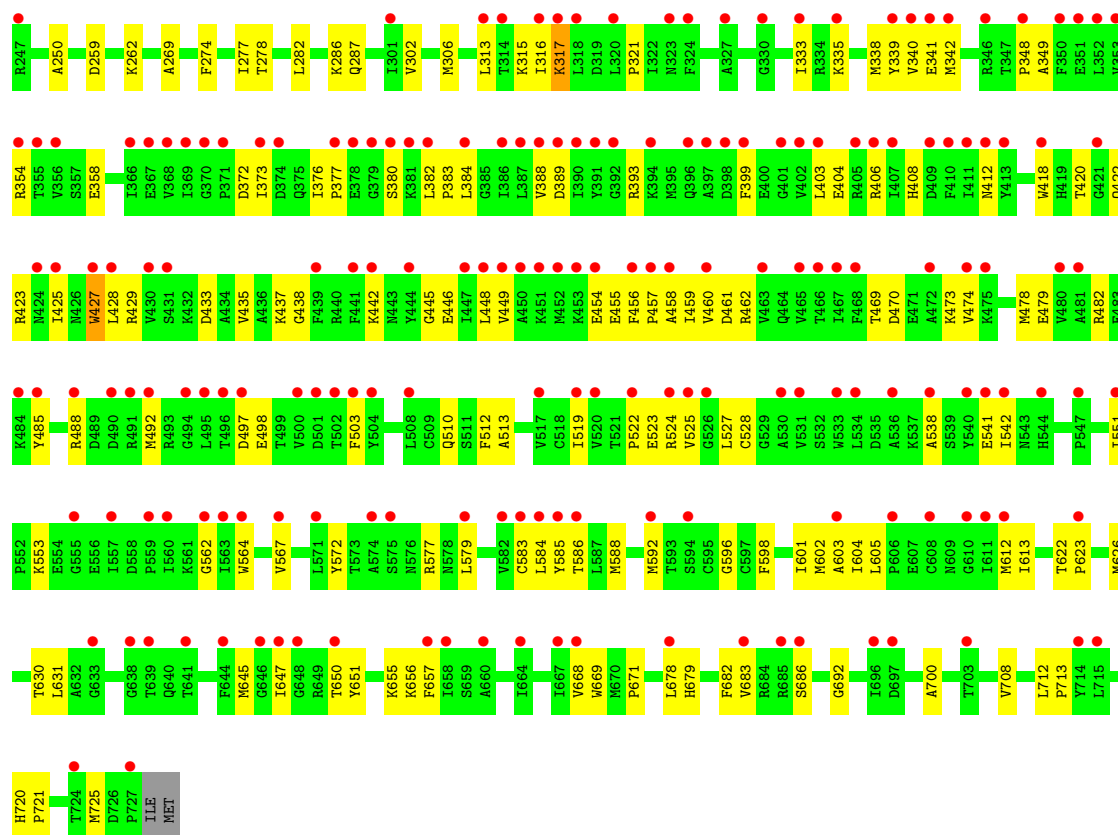


• Molecule 2: Carbon monoxide dehydrogenase/acetyl-CoA synthase subunit alpha

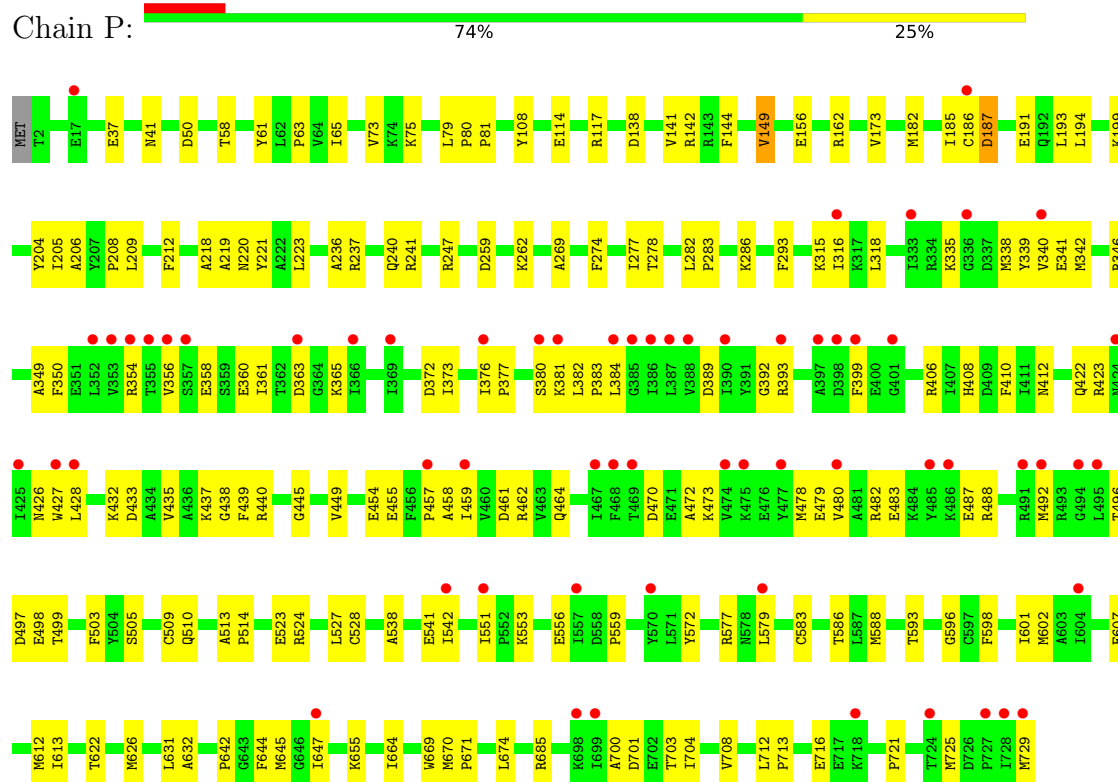
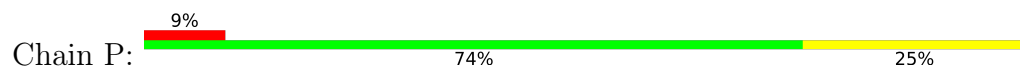


• Molecule 2: Carbon monoxide dehydrogenase/acetyl-CoA synthase subunit alpha





- Molecule 2: Carbon monoxide dehydrogenase/acetyl-CoA synthase subunit alpha



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	99.73Å 137.38Å 141.87Å 101.82° 109.08° 103.66°	Depositor
Resolution (Å)	49.63 – 2.47 49.63 – 2.47	Depositor EDS
% Data completeness (in resolution range)	86.9 (49.63-2.47) 79.3 (49.63-2.47)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.48Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, R_{free}	0.203 , 0.234 0.204 , 0.234	Depositor DCC
R_{free} test set	10245 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.154	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.6	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	0.93	EDS
Total number of atoms	44026	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% 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: XCC, CMO, PG5, NI, SF4, PGE, MG

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	0.09	0/5202	0.26	0/7048
1	B	0.10	0/5202	0.27	0/7048
1	C	0.10	0/5202	0.25	0/7048
1	D	0.09	0/5196	0.27	1/7040 (0.0%)
2	M	0.09	0/5874	0.27	0/7954
2	N	0.09	0/5874	0.26	0/7954
2	O	0.08	0/5857	0.24	0/7933
2	P	0.08	0/5874	0.24	0/7954
All	All	0.09	0/44281	0.26	1/59979 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	316	CYS	CB-CA-C	-6.85	107.64	115.79

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5109	0	5103	129	0
1	B	5109	0	5103	129	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	5109	0	5103	125	0
1	D	5103	0	5099	131	0
2	M	5740	0	5701	122	0
2	N	5740	0	5701	108	0
2	O	5723	0	5681	158	0
2	P	5740	0	5701	148	0
3	A	16	0	0	0	0
3	B	8	0	0	0	0
3	C	16	0	0	0	0
3	D	8	0	0	0	0
3	M	8	0	0	0	0
3	N	8	0	0	1	0
3	O	8	0	0	0	0
3	P	8	0	0	1	0
4	A	9	0	0	2	0
4	B	9	0	0	3	0
4	C	9	0	0	2	0
4	D	9	0	0	3	0
5	A	10	14	14	0	0
6	C	12	18	18	1	0
7	M	2	0	0	0	0
7	N	2	0	0	0	0
7	O	2	0	0	0	0
7	P	2	0	0	0	0
8	M	2	0	0	0	0
8	N	2	0	0	0	0
8	P	2	0	0	0	0
9	N	1	0	0	0	0
10	A	84	0	0	2	0
10	B	77	0	0	1	0
10	C	44	0	0	1	0
10	D	51	0	0	1	0
10	M	76	0	0	1	0
10	N	79	0	0	1	0
10	O	12	0	0	0	0
10	P	45	0	0	0	0
All	All	43994	32	43224	1015	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 12.

All (1015) 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)
2:O:602:MET:HE2	2:O:647:ILE:HD13	1.48	0.94
2:P:602:MET:HE2	2:P:647:ILE:HD13	1.48	0.92
1:D:482:LEU:HD21	1:D:508:LEU:HD12	1.53	0.90
1:A:281:HIS:HB3	1:A:351:VAL:HG12	1.51	0.90
1:B:607:MET:HE3	1:B:608:PRO:HD2	1.52	0.90
1:A:150:VAL:HG11	1:A:169:LYS:HG2	1.56	0.87
2:P:497:ASP:O	2:P:553:LYS:NZ	2.07	0.86
2:M:602:MET:HE2	2:M:647:ILE:HD13	1.56	0.86
1:A:607:MET:HE3	1:A:608:PRO:HD2	1.56	0.85
1:B:535:ALA:HB3	1:B:537:ILE:HD11	1.58	0.85
1:A:442:THR:HG21	1:A:535:ALA:HB2	1.59	0.84
2:M:339:TYR:HD2	2:M:340:VAL:HG23	1.42	0.84
1:B:482:LEU:HD21	1:B:508:LEU:HD12	1.60	0.82
2:M:149:VAL:HG11	2:M:509:CYS:HA	1.62	0.81
1:B:150:VAL:HG11	1:B:169:LYS:HG2	1.61	0.80
2:N:339:TYR:HD2	2:N:340:VAL:HG23	1.46	0.80
2:P:655:LYS:HG2	2:P:685:ARG:HH12	1.46	0.80
1:D:61:ILE:HD13	1:D:77:ARG:HE	1.48	0.79
1:D:607:MET:HE3	1:D:608:PRO:HD2	1.64	0.78
2:O:479:GLU:HA	2:O:482:ARG:HD3	1.62	0.78
1:C:607:MET:HE3	1:C:608:PRO:HD2	1.63	0.78
1:A:368[A]:HIS:CE1	1:A:411:ARG:HD3	2.18	0.78
1:B:559:LEU:HG	1:B:563:MET:HE3	1.68	0.76
1:D:298:MET:HE1	1:D:402:ARG:HG2	1.66	0.76
1:D:150:VAL:HG11	1:D:169:LYS:HG2	1.66	0.76
1:C:482:LEU:HD21	1:C:508:LEU:HD12	1.67	0.76
2:P:149:VAL:HG11	2:P:509:CYS:HA	1.68	0.76
2:O:377:PRO:HG2	2:O:380:SER:HB3	1.68	0.75
1:B:662:VAL:HG21	2:N:194:LEU:HD13	1.69	0.75
1:A:535:ALA:HB3	1:A:537:ILE:HD11	1.68	0.74
1:C:66:ILE:HG23	1:C:612:GLY:HA3	1.68	0.74
1:C:291:ILE:HD13	1:C:397:ALA:HB1	1.69	0.74
2:P:338:MET:SD	2:P:341:GLU:HB2	2.27	0.74
2:M:335:LYS:HE3	2:M:429:ARG:HH22	1.51	0.74
1:C:368[B]:HIS:CE1	1:C:411:ARG:HD3	2.23	0.74
2:N:602:MET:HE2	2:N:647:ILE:HD13	1.68	0.74
2:O:478:MET:O	2:O:482:ARG:HG3	1.87	0.74
1:A:282:GLY:HA3	1:A:352:ASP:OD1	1.88	0.74
1:B:61:ILE:HD13	1:B:77:ARG:HE	1.53	0.73
1:D:31:PRO:HB2	1:D:423:ILE:HD13	1.70	0.73
1:D:368[A]:HIS:CE1	1:D:411:ARG:HD3	2.23	0.73
1:B:281:HIS:HB3	1:B:351:VAL:HG12	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:61:ILE:HD13	1:C:77:ARG:NE	2.04	0.73
2:M:339:TYR:CE1	2:M:378:GLU:HG2	2.24	0.73
1:D:281:HIS:HB3	1:D:351:VAL:HG12	1.71	0.73
1:C:150:VAL:HG11	1:C:169:LYS:HG2	1.71	0.72
2:N:602:MET:HE2	2:N:647:ILE:HG21	1.73	0.71
2:P:721:PRO:O	2:P:725:MET:HG3	1.89	0.71
1:A:442:THR:HG21	1:A:535:ALA:CB	2.20	0.71
1:D:291:ILE:HD13	1:D:397:ALA:HB1	1.71	0.71
1:A:537:ILE:HG22	1:A:540:GLY:H	1.55	0.70
2:M:721:PRO:O	2:M:725:MET:HG3	1.91	0.70
2:P:339:TYR:CD2	2:P:340:VAL:HG23	2.26	0.70
2:O:58:THR:HA	2:O:138:ASP:OD1	1.91	0.70
1:D:482:LEU:CD2	1:D:508:LEU:HD12	2.21	0.70
1:A:535:ALA:HB3	1:A:537:ILE:CD1	2.22	0.70
2:O:2:THR:OG1	2:O:5:ASP:OD1	2.07	0.70
2:N:149:VAL:HG11	2:N:509:CYS:HA	1.73	0.69
1:D:607:MET:HE1	1:D:634:PHE:CD1	2.27	0.69
1:B:281:HIS:O	1:B:351:VAL:HA	1.92	0.69
2:O:338:MET:SD	2:O:341:GLU:HB2	2.33	0.69
2:P:498:GLU:N	2:P:498:GLU:OE1	2.26	0.69
1:B:537:ILE:HG22	1:B:540:GLY:H	1.56	0.69
1:A:149:ARG:NH2	1:A:250:ASP:OD2	2.24	0.69
1:A:607:MET:HE1	1:A:634:PHE:CD1	2.28	0.69
1:B:33:VAL:HG13	1:B:339:LEU:HD12	1.74	0.69
2:O:269:ALA:HB1	2:O:274:PHE:HB2	1.73	0.69
1:B:368[B]:HIS:CE1	1:B:411:ARG:HD3	2.27	0.68
1:C:482:LEU:CD2	1:C:508:LEU:HD12	2.23	0.68
1:A:33:VAL:HG13	1:A:339:LEU:HD12	1.76	0.68
1:C:281:HIS:HB3	1:C:351:VAL:HG12	1.75	0.68
2:M:613:ILE:HB	2:M:670:MET:HG2	1.75	0.68
1:C:282:GLY:HA3	1:C:352:ASP:OD1	1.94	0.68
2:P:339:TYR:HD2	2:P:340:VAL:HG23	1.57	0.67
1:D:282:GLY:HA3	1:D:352:ASP:OD1	1.95	0.67
2:O:498:GLU:OE1	2:O:498:GLU:N	2.27	0.67
2:O:588:MET:HE1	2:O:604:ILE:HG21	1.76	0.67
1:A:281:HIS:O	1:A:351:VAL:HA	1.95	0.67
1:A:615:LEU:HD23	2:M:27:HIS:HB3	1.75	0.67
1:D:66:ILE:HG23	1:D:612:GLY:HA3	1.76	0.67
2:M:498:GLU:N	2:M:498:GLU:OE2	2.28	0.67
1:B:607:MET:HE1	1:B:634:PHE:CD1	2.28	0.67
2:P:488:ARG:O	2:P:492:MET:HG2	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:ILE:HG23	1:B:612:GLY:HA3	1.76	0.67
1:A:559:LEU:HG	1:A:563:MET:HE3	1.77	0.67
2:O:488:ARG:O	2:O:492:MET:HG2	1.96	0.66
1:A:280:LEU:HD13	1:A:288:SER:HB2	1.77	0.66
1:A:291:ILE:HD13	1:A:397:ALA:HB1	1.75	0.66
2:O:553:LYS:HE2	2:O:564:TRP:CZ2	2.31	0.66
1:B:442:THR:HG21	1:B:535:ALA:HB2	1.77	0.66
2:M:602:MET:HE2	2:M:647:ILE:HG21	1.78	0.66
1:B:482:LEU:CD2	1:B:508:LEU:HD12	2.26	0.66
2:N:377:PRO:HG2	2:N:380:SER:HB3	1.77	0.66
2:M:439:PHE:O	2:M:440:ARG:HD2	1.96	0.65
1:D:144:LYS:O	1:D:148:ARG:HG3	1.96	0.65
1:D:299:GLU:HG2	1:D:303:LYS:HE2	1.78	0.65
2:M:340:VAL:HG21	2:M:376:ILE:HD11	1.79	0.65
1:C:31:PRO:HB2	1:C:423:ILE:HD13	1.77	0.65
1:C:416:ARG:CD	1:C:417:PRO:HD2	2.27	0.65
2:P:613:ILE:O	2:P:671:PRO:HD3	1.95	0.65
2:P:363:ASP:OD1	2:P:462:ARG:HA	1.97	0.65
1:C:601:PRO:HD3	1:C:652:ARG:CZ	2.27	0.65
1:D:274:ASP:CB	1:D:308:LYS:HE2	2.26	0.65
1:D:470:CYS:HB3	4:D:702:XCC:S2	2.36	0.64
2:N:700:ALA:HA	2:N:704:ILE:HD12	1.78	0.64
1:B:176:ARG:O	1:B:207:GLY:HA3	1.97	0.64
1:D:559:LEU:HG	1:D:563:MET:HE3	1.80	0.64
1:D:591:ILE:HA	1:D:594:TRP:NE1	2.12	0.64
2:O:721:PRO:O	2:O:725:MET:HG3	1.98	0.64
1:B:288:SER:O	1:B:292:VAL:HG23	1.96	0.64
1:A:66:ILE:HG23	1:A:612:GLY:HA3	1.78	0.64
1:D:368[A]:HIS:CD2	1:D:411:ARG:HB2	2.33	0.64
2:P:346:ARG:HB3	2:P:381:LYS:HD2	1.80	0.64
2:P:579:LEU:HD11	2:P:593:THR:HG21	1.80	0.64
1:B:607:MET:CE	1:B:608:PRO:HD2	2.25	0.64
1:B:607:MET:HE3	1:B:608:PRO:CD	2.26	0.63
1:B:282:GLY:HA3	1:B:352:ASP:OD1	1.98	0.63
2:P:602:MET:HE2	2:P:647:ILE:HG21	1.79	0.63
1:C:591:ILE:HA	1:C:594:TRP:NE1	2.13	0.63
1:C:607:MET:CE	1:C:608:PRO:HD2	2.29	0.63
1:A:482:LEU:HD21	1:A:508:LEU:HD12	1.80	0.63
1:B:535:ALA:HB3	1:B:537:ILE:CD1	2.28	0.63
1:C:468:CYS:HB2	10:C:840:HOH:O	1.98	0.63
2:N:338:MET:SD	2:N:341:GLU:HB2	2.39	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:433:ASP:OD2	2:N:437:LYS:HE2	1.98	0.63
2:M:73:VAL:HG11	2:M:79:LEU:HD21	1.80	0.63
2:O:598:PHE:CE1	2:O:601:ILE:HD11	2.34	0.62
2:P:422:GLN:OE1	2:P:422:GLN:N	2.32	0.62
2:P:50:ASP:HA	2:P:75:LYS:HD2	1.79	0.62
1:A:591:ILE:HA	1:A:594:TRP:NE1	2.15	0.62
1:B:482:LEU:HD12	1:B:486:LYS:HE3	1.80	0.62
2:M:422:GLN:OE1	2:M:422:GLN:N	2.32	0.62
2:O:339:TYR:CD2	2:O:340:VAL:HG23	2.34	0.62
2:M:349:ALA:HA	2:M:384:LEU:O	2.00	0.62
2:O:339:TYR:CD2	2:O:435:VAL:HG21	2.35	0.62
2:P:510:GLN:HA	2:P:513:ALA:O	2.00	0.62
1:D:338:GLU:OE2	1:D:359:SER:OG	2.18	0.62
2:O:510:GLN:HA	2:O:513:ALA:O	1.99	0.62
1:A:482:LEU:HD12	1:A:486:LYS:HE3	1.80	0.61
2:M:505:SER:HB3	2:M:551:ILE:HD11	1.82	0.61
2:O:340:VAL:HG12	2:O:342:MET:SD	2.40	0.61
2:O:423:ARG:HB3	2:O:485:TYR:CZ	2.34	0.61
2:O:372:ASP:OD1	2:O:373:ILE:N	2.31	0.61
1:D:149:ARG:NH2	1:D:250:ASP:OD2	2.33	0.61
2:P:382:LEU:HD23	2:P:383:PRO:O	2.00	0.61
2:M:367:GLU:HG3	2:M:466:THR:HG23	1.82	0.61
2:O:349:ALA:HA	2:O:384:LEU:O	2.00	0.61
1:C:351:VAL:HB	1:C:356:ILE:HD13	1.82	0.61
2:N:358:GLU:HG3	2:N:462:ARG:NH1	2.15	0.61
1:D:68:CYS:HB2	1:D:97:ILE:HG23	1.83	0.61
2:N:63:PRO:HB2	2:N:220:ASN:HB2	1.83	0.61
2:P:58:THR:HA	2:P:138:ASP:OD1	1.99	0.61
1:A:3:ARG:HG2	10:M:939:HOH:O	2.01	0.61
1:B:107:LEU:HD21	1:B:215:LEU:HB3	1.82	0.61
2:O:572:TYR:CE2	2:O:577:ARG:HG2	2.36	0.61
1:A:33:VAL:HG13	1:A:339:LEU:CD1	2.31	0.61
1:D:281:HIS:O	1:D:351:VAL:HA	2.01	0.61
2:N:586:THR:HG22	2:N:588:MET:H	1.66	0.61
1:C:281:HIS:O	1:C:351:VAL:HA	2.00	0.61
2:O:382:LEU:HD23	2:O:383:PRO:O	2.01	0.61
2:P:701:ASP:OD1	2:P:703:THR:OG1	2.10	0.61
1:B:591:ILE:HA	1:B:594:TRP:NE1	2.15	0.60
1:C:154:VAL:HG13	1:C:162:LEU:HD11	1.83	0.60
2:O:348:PRO:HG2	2:O:383:PRO:HB3	1.83	0.60
2:N:602:MET:HE1	2:N:645:MET:HE3	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:338:GLU:OE2	1:C:359:SER:OG	2.20	0.60
1:C:601:PRO:HD3	1:C:652:ARG:NH2	2.17	0.60
1:D:176:ARG:O	1:D:207:GLY:HA3	2.01	0.60
1:D:607:MET:CE	1:D:608:PRO:HD2	2.30	0.60
2:O:186:CYS:O	2:O:187:ASP:HB2	2.01	0.60
1:D:447:ASN:ND2	1:D:450:ARG:HB2	2.17	0.60
1:A:570:ASP:OD2	1:A:571:THR:N	2.33	0.60
1:B:601:PRO:HD3	1:B:652:ARG:CZ	2.31	0.60
2:P:358:GLU:HG3	2:P:462:ARG:NH1	2.16	0.60
1:C:291:ILE:HD13	1:C:397:ALA:CB	2.31	0.59
1:D:601:PRO:HD3	1:D:652:ARG:CZ	2.32	0.59
2:N:354:ARG:NH2	2:N:480:VAL:HG11	2.16	0.59
2:O:358:GLU:HG3	2:O:462:ARG:NH1	2.17	0.59
1:A:235:ILE:HG23	1:A:597:SER:OG	2.03	0.59
2:N:58:THR:HA	2:N:138:ASP:OD1	2.01	0.59
2:N:187:ASP:OD2	2:N:262:LYS:HE3	2.02	0.59
2:N:349:ALA:HA	2:N:384:LEU:O	2.02	0.59
2:O:626:MET:HB3	2:O:630:THR:HB	1.84	0.59
2:P:372:ASP:OD1	2:P:373:ILE:N	2.32	0.59
1:B:655:LYS:HG3	2:N:191:GLU:HG2	1.85	0.59
1:C:584:MET:HG3	1:D:73:ALA:HB2	1.85	0.59
1:B:442:THR:HG21	1:B:535:ALA:CB	2.33	0.59
1:D:33:VAL:HG13	1:D:339:LEU:HD12	1.85	0.59
1:D:559:LEU:HG	1:D:563:MET:CE	2.32	0.59
2:M:339:TYR:CD2	2:M:340:VAL:HG23	2.31	0.59
2:O:492:MET:HE2	2:O:492:MET:HA	1.83	0.59
1:A:581:PRO:O	1:A:605:GLY:HA3	2.03	0.59
1:A:615:LEU:CD2	2:M:27:HIS:HB3	2.33	0.59
1:C:235:ILE:HG23	1:C:597:SER:OG	2.02	0.59
2:N:721:PRO:O	2:N:725:MET:HG3	2.02	0.59
2:P:356:VAL:HG21	2:P:361:ILE:HB	1.85	0.59
1:D:18:ARG:NE	1:D:636:MET:HE3	2.18	0.58
2:O:626:MET:HE3	2:O:631:LEU:HD21	1.86	0.58
2:P:354:ARG:HD2	2:P:389:ASP:OD2	2.03	0.58
2:P:479:GLU:OE1	2:P:482:ARG:HD2	2.03	0.58
1:B:31:PRO:HB2	1:B:423:ILE:HD13	1.86	0.58
1:A:154:VAL:HG13	1:A:162:LEU:HD11	1.85	0.58
1:B:447:ASN:ND2	1:B:566:ASP:OD2	2.24	0.58
2:M:338:MET:SD	2:M:341:GLU:HB2	2.43	0.58
2:P:269:ALA:HB1	2:P:274:PHE:HB2	1.85	0.58
1:B:18:ARG:NE	1:B:636:MET:HE3	2.19	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:VAL:HB	1:A:356:ILE:HD13	1.84	0.58
1:A:150:VAL:HG13	1:A:184:THR:HG21	1.85	0.58
2:P:199:LYS:HA	2:P:199:LYS:HE2	1.85	0.58
1:B:577:VAL:HG21	1:B:645:ILE:HG23	1.86	0.58
2:O:423:ARG:HD3	2:O:485:TYR:CD2	2.39	0.58
1:D:18:ARG:CD	1:D:636:MET:HE3	2.34	0.57
2:M:351:GLU:HB3	2:M:426:ASN:HD21	1.69	0.57
1:C:33:VAL:HG13	1:C:339:LEU:HD12	1.85	0.57
2:O:172:ILE:HG23	2:O:302:VAL:CG2	2.34	0.57
1:A:587:LYS:HE3	4:A:703:XCC:S4	2.44	0.57
1:C:61:ILE:HG21	1:C:77:ARG:HE	1.69	0.57
2:N:422:GLN:HG2	2:N:423:ARG:HG3	1.86	0.57
2:N:335:LYS:HE3	2:N:429:ARG:HH22	1.68	0.57
2:P:433:ASP:OD1	2:P:437:LYS:HE2	2.03	0.57
1:C:176:ARG:O	1:C:207:GLY:HA3	2.04	0.57
1:D:587:LYS:HE3	4:D:702:XCC:S4	2.45	0.57
2:N:393:ARG:HG2	2:N:461:ASP:OD2	2.05	0.57
1:A:416:ARG:HG3	1:A:417:PRO:HD2	1.86	0.57
1:C:607:MET:HE1	1:C:634:PHE:CD1	2.40	0.57
1:D:398:LYS:O	1:D:402:ARG:HG3	2.03	0.57
2:N:186:CYS:O	2:N:187:ASP:HB2	2.05	0.57
2:O:399:PHE:HZ	2:O:541:GLU:HG3	1.69	0.57
2:O:433:ASP:OD2	2:O:437:LYS:HE2	2.04	0.57
1:D:150:VAL:HG13	1:D:184:THR:HG21	1.86	0.57
1:D:317:CYS:HB3	1:D:549:SER:HB2	1.87	0.57
2:M:141:VAL:HG22	2:M:221:TYR:CE2	2.39	0.57
2:N:114:GLU:OE1	2:N:117:ARG:NH1	2.38	0.57
1:A:175:ARG:HB3	1:B:175:ARG:HB3	1.87	0.56
1:B:278:PHE:HB3	1:B:312:LEU:HD23	1.87	0.56
1:C:61:ILE:HD13	1:C:77:ARG:HE	1.69	0.56
1:C:303:LYS:HA	1:C:307:ALA:O	2.04	0.56
1:D:655:LYS:NZ	1:D:674:TYR:O	2.36	0.56
1:D:291:ILE:HD13	1:D:397:ALA:CB	2.34	0.56
1:C:559:LEU:HG	1:C:563:MET:CE	2.35	0.56
2:N:602:MET:CE	2:N:645:MET:HE3	2.36	0.56
1:A:144:LYS:O	1:A:148:ARG:HG3	2.06	0.56
1:B:351:VAL:HB	1:B:356:ILE:HD13	1.87	0.56
2:M:579:LEU:HD11	2:M:593:THR:HG21	1.88	0.56
2:P:583:CYS:HB2	2:P:586:THR:HG22	1.87	0.56
1:B:587:LYS:HE3	4:B:702:XCC:S4	2.46	0.56
1:A:559:LEU:HG	1:A:563:MET:CE	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:615:LEU:HD23	2:M:27:HIS:CB	2.36	0.56
1:B:470:CYS:HB3	4:B:702:XCC:S2	2.46	0.56
2:M:358:GLU:HG3	2:M:462:ARG:NH1	2.21	0.56
2:O:26:TYR:CZ	2:O:30:ILE:HD11	2.40	0.56
2:O:679:HIS:O	2:O:683:VAL:HG23	2.06	0.56
2:P:376:ILE:HG21	2:P:382:LEU:HD12	1.87	0.56
2:P:583:CYS:CB	2:P:586:THR:HG22	2.36	0.56
1:C:60:LYS:H	6:C:704:PG5:H83	1.71	0.56
2:O:712:LEU:HB3	2:O:713:PRO:HD3	1.86	0.56
2:M:354:ARG:HD2	2:M:389:ASP:OD2	2.05	0.56
2:N:478:MET:O	2:N:482:ARG:HG3	2.06	0.56
2:O:433:ASP:O	2:O:437:LYS:HG3	2.05	0.56
1:A:470:CYS:HB3	4:A:703:XCC:S2	2.46	0.55
2:N:498:GLU:OE2	2:N:498:GLU:N	2.38	0.55
1:B:370:ARG:HD3	1:B:407:ALA:HB2	1.87	0.55
1:C:68:CYS:HB2	1:C:97:ILE:HG23	1.88	0.55
1:A:456:ILE:HD13	1:A:463:GLY:HA2	1.88	0.55
1:C:581:PRO:O	1:C:605:GLY:HA3	2.06	0.55
2:O:422:GLN:N	2:O:422:GLN:OE1	2.39	0.55
1:A:61:ILE:HD13	1:A:77:ARG:HD2	1.87	0.55
1:A:150:VAL:CG1	1:A:169:LYS:HG2	2.32	0.55
1:A:292:VAL:HB	10:A:838:HOH:O	2.07	0.55
1:B:77:ARG:NH1	1:B:79:LYS:HE2	2.22	0.55
1:B:559:LEU:HG	1:B:563:MET:CE	2.36	0.55
1:C:587:LYS:HE3	4:C:703:XCC:S4	2.46	0.55
1:C:607:MET:HE3	1:C:608:PRO:CD	2.33	0.55
2:M:416:GLY:HA3	2:M:434:ALA:HB2	1.88	0.55
2:M:626:MET:HE3	2:M:631:LEU:HD21	1.88	0.55
2:N:69:SER:OG	2:N:71:GLU:OE1	2.23	0.55
2:P:670:MET:HE3	2:P:674:LEU:HD23	1.88	0.55
2:P:700:ALA:HA	2:P:704:ILE:HD12	1.87	0.55
2:M:583:CYS:CB	2:M:586:THR:HG22	2.37	0.55
1:C:447:ASN:ND2	1:C:450:ARG:HB2	2.22	0.55
1:D:335:ALA:HB2	1:D:471:ASN:HB3	1.89	0.55
2:P:602:MET:CE	2:P:645:MET:HE3	2.36	0.55
2:P:342:MET:HG3	2:P:382:LEU:O	2.07	0.55
1:C:559:LEU:HG	1:C:563:MET:HE3	1.88	0.54
2:O:69:SER:OG	2:O:71:GLU:OE1	2.24	0.54
1:C:150:VAL:HG13	1:C:184:THR:HG21	1.88	0.54
1:D:601:PRO:HD3	1:D:652:ARG:NH2	2.22	0.54
2:M:144:PHE:CE1	2:M:205:ILE:HG12	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:278:THR:HG21	2:O:282:LEU:HD11	1.89	0.54
1:C:158:SER:OG	1:C:161:GLU:HG3	2.07	0.54
2:O:317:LYS:NZ	2:O:317:LYS:HB3	2.22	0.54
1:D:158:SER:OG	1:D:161:GLU:HG3	2.07	0.54
1:B:399:THR:HG22	1:B:403:MET:HE3	1.89	0.54
2:N:492:MET:HE2	2:N:492:MET:HA	1.88	0.54
1:C:107:LEU:HD13	1:C:216:VAL:HG22	1.89	0.54
2:M:206:ALA:O	2:M:208:PRO:HD3	2.08	0.54
2:O:503:PHE:O	2:O:551:ILE:N	2.36	0.54
2:O:686:SER:HB3	2:O:692:GLY:O	2.08	0.54
1:A:321:GLU:OE1	1:A:325:ARG:NH1	2.34	0.54
2:P:114:GLU:OE1	2:P:117:ARG:NH1	2.41	0.54
1:D:577:VAL:HG13	1:D:649:LEU:CD2	2.38	0.54
2:O:114:GLU:OE1	2:O:117:ARG:NH1	2.41	0.54
2:O:182:MET:HA	2:O:205:ILE:HG22	1.90	0.54
1:A:176:ARG:O	1:A:207:GLY:HA3	2.08	0.54
2:N:712:LEU:O	2:N:716:GLU:HG3	2.08	0.54
2:O:156:GLU:HG3	2:O:182:MET:HB3	1.89	0.54
2:O:333:ILE:HD12	2:O:429:ARG:HB2	1.91	0.54
1:D:314:GLY:O	1:D:331:VAL:HG12	2.09	0.53
1:D:581:PRO:O	1:D:605:GLY:HA3	2.08	0.53
2:M:470:ASP:O	2:M:474:VAL:HG23	2.08	0.53
2:M:700:ALA:HA	2:M:704:ILE:HD12	1.89	0.53
2:P:199:LYS:HD3	2:P:204:TYR:OH	2.08	0.53
1:D:656:LEU:O	1:D:660:LYS:HG3	2.09	0.53
2:M:399:PHE:O	2:M:403:LEU:HD13	2.07	0.53
1:A:464:VAL:HG21	1:A:563:MET:HE1	1.89	0.53
1:A:607:MET:CE	1:A:608:PRO:HD2	2.33	0.53
1:D:486:LYS:HE2	1:D:519:TYR:CE2	2.43	0.53
1:D:570:ASP:OD2	1:D:571:THR:N	2.40	0.53
1:D:607:MET:HE3	1:D:608:PRO:CD	2.36	0.53
2:O:316:ILE:HD12	2:O:316:ILE:N	2.24	0.53
2:O:456:PHE:HB3	2:O:459:ILE:CG2	2.38	0.53
2:O:585:TYR:CE1	2:O:656:LYS:HB3	2.44	0.53
2:O:603:ALA:HB3	2:O:612:MET:HG3	1.90	0.53
1:A:607:MET:HE3	1:A:608:PRO:CD	2.35	0.53
2:O:141:VAL:HG22	2:O:221:TYR:CE2	2.43	0.53
2:P:602:MET:HE1	2:P:645:MET:HE3	1.90	0.53
1:C:607:MET:HE1	1:C:634:PHE:CG	2.44	0.53
2:M:372:ASP:OD1	2:M:373:ILE:N	2.39	0.53
2:P:186:CYS:O	2:P:187:ASP:HB2	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:187:ASP:OD2	2:P:262:LYS:HE3	2.09	0.53
2:P:439:PHE:O	2:P:440:ARG:HD2	2.08	0.53
1:A:291:ILE:HD13	1:A:397:ALA:CB	2.37	0.53
2:M:340:VAL:CG2	2:M:376:ILE:HD11	2.38	0.53
2:M:356:VAL:HG21	2:M:361:ILE:HB	1.91	0.53
2:N:439:PHE:O	2:N:440:ARG:HD2	2.09	0.53
1:A:601:PRO:HD3	1:A:652:ARG:CZ	2.39	0.53
2:M:187:ASP:OD2	2:M:262:LYS:HE3	2.09	0.53
2:O:404:GLU:OE1	2:O:423:ARG:HG2	2.08	0.53
1:A:505:ALA:HB1	1:A:510:LEU:HB2	1.89	0.53
2:N:354:ARG:HD2	2:N:389:ASP:OD2	2.08	0.53
2:N:402:VAL:HG13	2:N:535:ASP:OD1	2.09	0.53
2:O:21:LEU:HB2	2:O:286:LYS:HA	1.90	0.53
1:B:570:ASP:OD2	1:B:571:THR:N	2.41	0.53
2:P:588:MET:O	2:P:642:PRO:HB3	2.08	0.53
1:B:118:ASN:C	1:B:118:ASN:HD22	2.17	0.52
1:C:33:VAL:HG13	1:C:339:LEU:CD1	2.38	0.52
1:C:428:VAL:HG22	1:C:539:ILE:HD11	1.91	0.52
2:O:376:ILE:HG21	2:O:382:LEU:HD12	1.91	0.52
1:A:368[A]:HIS:CD2	1:A:411:ARG:HB2	2.45	0.52
1:C:370:ARG:HD2	1:C:384:TYR:CE2	2.44	0.52
1:B:482:LEU:HD13	1:B:510:LEU:HD11	1.92	0.52
1:C:18:ARG:NE	1:C:636:MET:HE3	2.24	0.52
1:A:577:VAL:HG13	1:A:649:LEU:CD2	2.40	0.52
2:O:524:ARG:NH1	2:O:584:LEU:O	2.41	0.52
2:P:141:VAL:HG22	2:P:221:TYR:CE2	2.44	0.52
2:P:626:MET:HE3	2:P:631:LEU:HD21	1.91	0.52
1:A:577:VAL:HG21	1:A:645:ILE:CG2	2.40	0.52
2:N:21:LEU:HB2	2:N:286:LYS:HA	1.92	0.52
2:P:438:GLY:HA2	2:P:440:ARG:NH1	2.24	0.52
2:P:612:MET:HE1	2:P:622:THR:HB	1.92	0.52
1:A:107:LEU:HD13	1:A:216:VAL:HG22	1.92	0.52
1:C:105:MET:HE1	1:C:609:PRO:HD2	1.90	0.52
1:D:107:LEU:HD13	1:D:216:VAL:HG22	1.91	0.52
2:M:58:THR:HA	2:M:138:ASP:OD1	2.10	0.52
2:N:206:ALA:O	2:N:208:PRO:HD3	2.09	0.52
2:N:422:GLN:N	2:N:422:GLN:OE1	2.43	0.52
2:P:350:PHE:HA	2:P:426:ASN:HD22	1.75	0.52
1:B:10:ASN:HB2	2:N:257:GLU:OE2	2.09	0.52
2:P:384:LEU:CD2	2:P:428:LEU:HD22	2.39	0.52
1:A:118:ASN:C	1:A:118:ASN:HD22	2.18	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:446:GLN:NE2	1:C:42:ASP:OD2	2.43	0.52
2:M:602:MET:HE1	2:M:645:MET:HE3	1.92	0.52
2:O:445:GLY:O	2:O:449:VAL:HG23	2.10	0.52
1:C:224:ASP:OD1	1:C:229:ASN:ND2	2.40	0.51
1:C:368[B]:HIS:CD2	1:C:407:ALA:HB1	2.45	0.51
1:D:447:ASN:ND2	1:D:566:ASP:OD2	2.25	0.51
1:C:607:MET:HE3	1:C:607:MET:HA	1.92	0.51
1:D:282:GLY:C	1:D:318:THR:HB	2.35	0.51
2:M:156:GLU:O	2:M:250:ALA:HA	2.11	0.51
1:B:150:VAL:CG1	1:B:169:LYS:HG2	2.34	0.51
2:O:622:THR:OG1	2:O:626:MET:O	2.25	0.51
1:D:77:ARG:NH1	1:D:79:LYS:HE2	2.24	0.51
2:O:187:ASP:OD2	2:O:262:LYS:HE3	2.11	0.51
1:A:468:CYS:HB2	10:A:857:HOH:O	2.10	0.51
1:C:577:VAL:HG21	1:C:645:ILE:HG23	1.91	0.51
1:D:77:ARG:HH11	1:D:79:LYS:HE2	1.75	0.51
2:N:510:GLN:HA	2:N:513:ALA:O	2.11	0.51
1:D:607:MET:HE3	1:D:607:MET:HA	1.93	0.51
1:C:186:LEU:CD2	1:C:205:PRO:HD2	2.41	0.51
1:C:587:LYS:O	1:C:591:ILE:HG13	2.10	0.51
1:D:571:THR:OG1	1:D:572:PRO:HD3	2.10	0.51
2:N:156:GLU:HG3	2:N:182:MET:HB3	1.93	0.51
2:N:406:ARG:HB3	2:N:410:PHE:CZ	2.46	0.51
2:P:373:ILE:HD13	2:P:435:VAL:HG22	1.93	0.51
1:D:370:ARG:HD2	1:D:384:TYR:CE2	2.46	0.51
1:D:456:ILE:HD13	1:D:463:GLY:HA2	1.93	0.51
2:M:377:PRO:HG2	2:M:380:SER:HB3	1.92	0.51
1:B:601:PRO:HD3	1:B:652:ARG:NH2	2.25	0.50
2:P:346:ARG:HB3	2:P:381:LYS:CD	2.41	0.50
2:P:423:ARG:O	2:P:426:ASN:ND2	2.44	0.50
1:D:577:VAL:HG21	1:D:645:ILE:HG23	1.92	0.50
2:M:186:CYS:O	2:M:187:ASP:HB2	2.10	0.50
2:P:655:LYS:HA	2:P:685:ARG:NH2	2.25	0.50
1:A:388:TYR:OH	1:A:393:ALA:HB1	2.12	0.50
1:B:61:ILE:HD13	1:B:77:ARG:NE	2.22	0.50
1:B:464:VAL:HG21	1:B:563:MET:HE1	1.94	0.50
1:C:368[B]:HIS:HD2	1:C:407:ALA:HB1	1.77	0.50
1:D:330:LEU:HD11	1:D:503:GLN:CG	2.42	0.50
2:P:338:MET:O	2:P:432:LYS:HE3	2.11	0.50
1:B:656:LEU:O	1:B:660:LYS:HG3	2.11	0.50
1:C:317:CYS:HB3	1:C:549:SER:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:291:ILE:HD13	1:B:397:ALA:HB1	1.93	0.50
1:B:316:CYS:N	1:B:503:GLN:HE22	2.09	0.50
2:M:602:MET:CE	2:M:645:MET:HE3	2.41	0.50
1:B:607:MET:HE3	1:B:607:MET:HA	1.93	0.50
1:C:5:ARG:HB2	2:O:162:ARG:NH2	2.26	0.50
2:M:613:ILE:O	2:M:671:PRO:HD3	2.11	0.50
2:N:384:LEU:CD2	2:N:428:LEU:HD22	2.41	0.50
2:O:156:GLU:HB2	2:O:250:ALA:HB2	1.94	0.50
2:O:470:ASP:HB3	2:O:473:LYS:CG	2.42	0.50
2:P:316:ILE:HG13	2:P:454:GLU:OE2	2.12	0.50
2:P:342:MET:HE2	2:P:384:LEU:HD22	1.94	0.50
2:P:439:PHE:C	2:P:440:ARG:HD2	2.37	0.50
2:M:172:ILE:HG23	2:M:302:VAL:CG2	2.42	0.50
2:N:505:SER:HB3	2:N:551:ILE:HD11	1.94	0.50
2:O:354:ARG:HD2	2:O:389:ASP:OD2	2.12	0.50
2:P:335:LYS:H	2:P:335:LYS:HD2	1.76	0.50
2:P:342:MET:HG2	2:P:384:LEU:HB2	1.92	0.50
2:P:393:ARG:HG2	2:P:461:ASP:OD2	2.12	0.50
1:A:548:GLY:HA3	1:A:552:ASP:OD2	2.12	0.50
2:M:354:ARG:NH2	2:M:480:VAL:HG11	2.27	0.50
2:M:583:CYS:HB2	2:M:586:THR:HG22	1.92	0.50
2:O:562:GLY:HA3	2:O:585:TYR:CD2	2.47	0.50
2:P:356:VAL:CG2	2:P:361:ILE:HB	2.40	0.50
2:P:478:MET:O	2:P:482:ARG:HG3	2.11	0.50
1:A:447:ASN:ND2	1:A:566:ASP:OD2	2.30	0.49
1:A:626:ASP:HB3	2:M:212:PHE:CG	2.47	0.49
1:B:33:VAL:HG13	1:B:339:LEU:CD1	2.40	0.49
1:B:77:ARG:HH11	1:B:79:LYS:HE2	1.76	0.49
1:D:384:TYR:CD1	2:O:84:ASN:HB3	2.46	0.49
2:P:206:ALA:O	2:P:208:PRO:HD3	2.11	0.49
2:P:602:MET:CE	2:P:647:ILE:HG21	2.42	0.49
1:C:316:CYS:N	1:C:503:GLN:HE22	2.11	0.49
1:D:505:ALA:HB1	1:D:510:LEU:HB2	1.94	0.49
2:P:173:VAL:HG21	2:P:193:LEU:HD21	1.93	0.49
2:M:701:ASP:OD1	2:M:703:THR:OG1	2.14	0.49
2:N:456:PHE:HA	2:N:542:ILE:HD13	1.95	0.49
1:C:505:ALA:HB1	1:C:510:LEU:HB2	1.93	0.49
1:C:569:VAL:HB	1:C:573:LYS:HD3	1.94	0.49
1:D:190:ILE:HG13	1:D:195:LYS:HG3	1.94	0.49
2:O:73:VAL:HG11	2:O:79:LEU:HD21	1.94	0.49
1:D:662:VAL:HG21	2:P:194:LEU:HD13	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:393:ARG:HG2	2:M:461:ASP:OD2	2.12	0.49
2:P:340:VAL:HG12	2:P:342:MET:SD	2.53	0.49
1:B:581:PRO:O	1:B:605:GLY:HA3	2.12	0.49
2:M:156:GLU:HB2	2:M:250:ALA:HB2	1.95	0.49
2:M:456:PHE:HA	2:M:542:ILE:HD13	1.94	0.49
2:O:335:LYS:NZ	2:O:429:ARG:HH22	2.10	0.49
2:O:583:CYS:HB2	2:O:586:THR:HG22	1.94	0.49
2:M:363:ASP:OD1	2:M:462:ARG:HA	2.13	0.49
2:N:602:MET:CE	2:N:647:ILE:HG21	2.42	0.49
2:O:524:ARG:NH2	2:O:645:MET:SD	2.85	0.49
1:A:443:GLN:HB2	1:A:451:VAL:HG21	1.93	0.49
1:C:280:LEU:HD13	1:C:288:SER:HB2	1.95	0.49
2:M:439:PHE:C	2:M:440:ARG:HD2	2.38	0.49
2:O:626:MET:HG2	2:O:630:THR:HG21	1.95	0.49
2:P:588:MET:HE2	2:P:729:MET:O	2.12	0.49
1:B:150:VAL:HG13	1:B:184:THR:HG21	1.94	0.49
1:B:321:GLU:OE1	1:B:325:ARG:NH1	2.41	0.49
1:C:63:TYR:OH	1:C:607:MET:HB3	2.13	0.49
1:C:150:VAL:HG12	1:C:169:LYS:HE2	1.94	0.49
1:C:482:LEU:HD12	1:C:486:LYS:HE3	1.94	0.49
2:M:109:ALA:O	2:M:113:ILE:HG13	2.13	0.49
2:O:523:GLU:OE1	2:O:655:LYS:HB2	2.13	0.49
1:D:118:ASN:C	1:D:118:ASN:HD22	2.21	0.48
2:M:588:MET:O	2:M:642:PRO:HB3	2.13	0.48
2:O:315:LYS:HA	2:O:454:GLU:OE2	2.13	0.48
2:P:259:ASP:OD1	2:P:262:LYS:HD2	2.12	0.48
2:M:492:MET:HA	2:M:492:MET:HE2	1.94	0.48
1:B:39:LYS:NZ	1:B:43:ASP:OD2	2.42	0.48
1:D:587:LYS:O	1:D:591:ILE:HG13	2.12	0.48
1:D:98:VAL:O	1:D:102:VAL:HG12	2.13	0.48
1:C:118:ASN:C	1:C:118:ASN:HD22	2.20	0.48
1:C:486:LYS:HE2	1:C:519:TYR:CE2	2.48	0.48
2:O:470:ASP:O	2:O:474:VAL:HG23	2.14	0.48
1:A:298:MET:HE2	1:A:401:ILE:HB	1.96	0.48
1:A:615:LEU:HD22	2:M:260:MET:SD	2.54	0.48
1:A:658:VAL:O	1:A:662:VAL:HG23	2.14	0.48
1:B:483:THR:HB	1:B:638:PRO:HB2	1.95	0.48
1:C:278:PHE:HB3	1:C:312:LEU:HD23	1.95	0.48
1:C:482:LEU:HD13	1:C:510:LEU:HD11	1.96	0.48
1:D:107:LEU:HD21	1:D:215:LEU:HB3	1.96	0.48
2:M:540:TYR:CD1	2:M:550:PRO:HD3	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:199:LYS:HD3	2:P:204:TYR:CZ	2.49	0.48
1:B:505:ALA:HB1	1:B:510:LEU:HB2	1.96	0.48
1:C:570:ASP:HB3	1:C:572:PRO:HD2	1.96	0.48
1:D:279:VAL:HG22	1:D:313:VAL:CG2	2.43	0.48
2:N:259:ASP:OD1	2:N:262:LYS:HD2	2.13	0.48
2:O:376:ILE:O	2:O:376:ILE:HD12	2.14	0.48
2:O:626:MET:HE3	2:O:631:LEU:CD2	2.44	0.48
2:P:360:GLU:O	2:P:360:GLU:HG2	2.13	0.48
1:A:482:LEU:CD2	1:A:508:LEU:HD12	2.44	0.48
1:B:577:VAL:HG21	1:B:645:ILE:CG2	2.43	0.48
2:N:603:ALA:HB3	2:N:612:MET:HG3	1.96	0.48
2:O:393:ARG:HG2	2:O:461:ASP:OD2	2.14	0.48
1:A:31:PRO:HB2	1:A:423:ILE:HD13	1.96	0.48
1:B:186:LEU:O	1:B:190:ILE:HG12	2.14	0.48
2:N:622:THR:HB	2:N:623:PRO:HD2	1.96	0.48
1:A:299:GLU:HG2	1:A:303:LYS:HE2	1.96	0.48
1:B:368[B]:HIS:CD2	1:B:411:ARG:HB2	2.48	0.48
2:O:206:ALA:O	2:O:208:PRO:HD3	2.14	0.48
2:O:236:ALA:O	2:O:240:GLN:HG2	2.14	0.48
1:D:33:VAL:HG13	1:D:339:LEU:CD1	2.44	0.47
1:D:274:ASP:HB2	1:D:308:LYS:HE2	1.95	0.47
2:O:321:PRO:HG3	2:O:438:GLY:O	2.14	0.47
2:O:377:PRO:HG2	2:O:380:SER:CB	2.42	0.47
2:O:527:LEU:HD12	2:O:528:CYS:N	2.29	0.47
2:P:156:GLU:HG3	2:P:182:MET:HB3	1.95	0.47
1:A:149:ARG:HH22	1:A:250:ASP:CG	2.21	0.47
1:A:150:VAL:HG12	1:A:169:LYS:HE2	1.95	0.47
1:A:186:LEU:O	1:A:190:ILE:HG12	2.14	0.47
1:B:409:LYS:O	1:B:413:GLU:HG3	2.14	0.47
1:C:113:HIS:NE2	1:C:587:LYS:HD2	2.29	0.47
1:C:517:GLU:OE2	1:C:525:LYS:HD2	2.14	0.47
2:P:572:TYR:CE2	2:P:577:ARG:HG2	2.48	0.47
2:M:59:ALA:HB2	2:M:630:THR:HA	1.96	0.47
1:A:278:PHE:HB3	1:A:312:LEU:HD23	1.96	0.47
2:N:579:LEU:HD11	2:N:593:THR:HG21	1.97	0.47
1:B:370:ARG:HH22	1:B:410:GLU:CD	2.22	0.47
1:C:73:ALA:CB	1:D:584:MET:HE3	2.45	0.47
1:C:571:THR:OG1	1:C:572:PRO:HD3	2.13	0.47
1:D:283:HIS:CE1	4:D:702:XCC:S3	3.07	0.47
1:D:330:LEU:HD23	1:D:330:LEU:C	2.39	0.47
2:N:129:PRO:O	2:N:132:TRP:HB2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:360:GLU:O	2:N:360:GLU:HG2	2.14	0.47
2:N:727:PRO:HB2	2:N:729:MET:HE2	1.95	0.47
2:O:470:ASP:HB3	2:O:473:LYS:HG2	1.95	0.47
2:P:505:SER:HB3	2:P:551:ILE:HD11	1.95	0.47
1:C:483:THR:CG2	1:C:638:PRO:HB2	2.44	0.47
2:N:439:PHE:C	2:N:440:ARG:HD2	2.39	0.47
2:O:538:ALA:O	2:O:542:ILE:HG13	2.14	0.47
2:P:315:LYS:HA	2:P:454:GLU:OE1	2.15	0.47
1:A:105:MET:HE3	1:A:609:PRO:HG3	1.95	0.47
1:B:107:LEU:HD13	1:B:216:VAL:HG22	1.96	0.47
1:C:584:MET:HE3	1:D:73:ALA:CB	2.45	0.47
1:D:150:VAL:CG1	1:D:169:LYS:HG2	2.40	0.47
2:M:199:LYS:HD2	2:M:199:LYS:HA	1.59	0.47
2:M:373:ILE:CD1	2:M:435:VAL:HG22	2.44	0.47
2:M:510:GLN:HA	2:M:513:ALA:O	2.14	0.47
2:N:340:VAL:HG12	2:N:342:MET:SD	2.55	0.47
2:N:372:ASP:OD1	2:N:373:ILE:N	2.42	0.47
2:P:354:ARG:NH2	2:P:480:VAL:HG11	2.29	0.47
1:C:250:ASP:O	1:C:254:ILE:HG13	2.15	0.47
1:D:607:MET:HE1	1:D:634:PHE:CG	2.50	0.47
1:D:655:LYS:HG2	2:P:191:GLU:HG2	1.97	0.47
2:M:457:PRO:O	2:M:458:ALA:HB3	2.15	0.47
2:O:144:PHE:O	2:O:148:MET:HG3	2.14	0.47
2:O:572:TYR:HA	2:O:579:LEU:O	2.14	0.47
2:O:585:TYR:CD1	2:O:656:LYS:HB3	2.50	0.47
2:P:73:VAL:HG11	2:P:79:LEU:HD21	1.97	0.47
1:A:182:GLU:OE2	1:A:199:ARG:HD3	2.14	0.47
1:B:452:LEU:HD21	1:B:464:VAL:HG21	1.96	0.47
1:C:150:VAL:CG1	1:C:169:LYS:HE2	2.44	0.47
1:C:468:CYS:O	1:C:580:ALA:HA	2.15	0.47
2:O:144:PHE:CE1	2:O:205:ILE:HG12	2.49	0.47
2:O:479:GLU:HA	2:O:482:ARG:CD	2.37	0.47
2:P:185:ILE:HD12	2:P:185:ILE:N	2.30	0.47
2:P:445:GLY:O	2:P:449:VAL:HG23	2.15	0.47
1:B:607:MET:HE1	1:B:634:PHE:CG	2.50	0.47
1:C:73:ALA:HB2	1:D:584:MET:HG3	1.97	0.47
2:O:96:PHE:O	2:O:100:ARG:HG3	2.15	0.47
2:O:373:ILE:CD1	2:O:435:VAL:HG22	2.45	0.47
2:P:712:LEU:O	2:P:716:GLU:HG3	2.15	0.47
1:B:190:ILE:HG13	1:B:195:LYS:HG3	1.96	0.46
1:D:330:LEU:HD11	1:D:503:GLN:CD	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:335:LYS:HD2	2:M:335:LYS:H	1.81	0.46
2:P:283:PRO:HD2	2:P:286:LYS:HD3	1.96	0.46
2:P:342:MET:CE	2:P:384:LEU:HD22	2.45	0.46
1:B:607:MET:HE2	1:B:610:VAL:CG2	2.45	0.46
1:D:154:VAL:HG13	1:D:162:LEU:HD11	1.98	0.46
2:N:500:VAL:HG21	2:N:522:PRO:HG2	1.96	0.46
2:N:701:ASP:OD1	2:N:703:THR:OG1	2.22	0.46
2:O:166:SER:OG	2:O:196:GLU:HG2	2.14	0.46
1:A:190:ILE:HG13	1:A:195:LYS:HG3	1.97	0.46
1:A:468:CYS:O	1:A:580:ALA:HA	2.14	0.46
1:C:314:GLY:O	1:C:331:VAL:HG12	2.15	0.46
2:O:26:TYR:CE2	2:O:30:ILE:HD11	2.51	0.46
2:O:376:ILE:HD12	2:O:376:ILE:C	2.40	0.46
2:P:377:PRO:HG2	2:P:380:SER:HB3	1.98	0.46
1:D:298:MET:HE1	1:D:402:ARG:CG	2.41	0.46
2:M:269:ALA:HB1	2:M:274:PHE:HB2	1.98	0.46
2:M:342:MET:HE2	2:M:384:LEU:HD22	1.96	0.46
2:N:236:ALA:O	2:N:240:GLN:HG2	2.15	0.46
2:O:176:LEU:HD21	2:O:302:VAL:HG13	1.98	0.46
2:O:185:ILE:N	2:O:185:ILE:HD12	2.31	0.46
1:B:3:ARG:HB2	10:N:954:HOH:O	2.14	0.46
1:B:18:ARG:CD	1:B:636:MET:HE3	2.46	0.46
1:C:98:VAL:O	1:C:102:VAL:HG12	2.16	0.46
1:C:416:ARG:CG	1:C:417:PRO:HD2	2.45	0.46
2:M:275:PRO:HD2	2:M:309:ARG:HD3	1.97	0.46
2:P:670:MET:CE	2:P:674:LEU:HG	2.46	0.46
1:A:399:THR:HG22	1:A:403:MET:HE3	1.97	0.46
1:B:68:CYS:HB2	1:B:97:ILE:HG23	1.96	0.46
1:C:662:VAL:HG21	2:O:194:LEU:HD13	1.97	0.46
2:M:185:ILE:HG22	2:M:190:VAL:HG23	1.97	0.46
1:D:147:CYS:HB3	1:D:152:ILE:HB	1.96	0.46
2:N:222:ALA:HB1	2:N:248:ILE:HG21	1.98	0.46
2:P:523:GLU:OE1	2:P:655:LYS:HB2	2.16	0.46
1:A:52:PHE:CZ	1:A:474:LYS:HA	2.51	0.46
1:A:607:MET:HE1	1:A:634:PHE:CG	2.51	0.46
1:D:273:PRO:HD3	1:D:420:ILE:HD12	1.97	0.46
1:D:559:LEU:O	1:D:563:MET:HG3	2.16	0.46
2:N:503:PHE:O	2:N:551:ILE:N	2.46	0.46
2:O:277:ILE:N	2:O:277:ILE:HD12	2.31	0.46
2:P:80:PRO:N	2:P:81:PRO:HD2	2.31	0.46
1:A:577:VAL:HG21	1:A:645:ILE:HG23	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:416:ARG:HD2	1:C:417:PRO:HD2	1.97	0.46
1:C:570:ASP:OD2	1:C:571:THR:N	2.48	0.46
1:D:452:LEU:HD21	1:D:464:VAL:HG21	1.98	0.46
2:M:149:VAL:CG1	2:M:509:CYS:HA	2.41	0.46
2:M:445:GLY:O	2:M:449:VAL:HG23	2.16	0.46
2:N:626:MET:HB3	2:N:630:THR:HB	1.98	0.46
2:P:607:GLU:HB2	2:P:708:VAL:HG11	1.97	0.46
1:B:636:MET:HE2	10:B:830:HOH:O	2.15	0.46
1:B:658:VAL:O	1:B:662:VAL:HG23	2.16	0.46
2:N:457:PRO:O	2:N:458:ALA:HB3	2.15	0.46
2:N:727:PRO:CG	2:N:729:MET:HE2	2.46	0.46
2:O:160:LEU:HD13	2:O:215:ILE:HD13	1.97	0.46
2:O:525:VAL:HG21	2:O:650:THR:HB	1.97	0.46
2:O:605:LEU:HD11	2:O:612:MET:HB3	1.98	0.46
2:O:613:ILE:O	2:O:671:PRO:HD3	2.16	0.46
1:A:353:VAL:O	1:A:354:GLN:HB2	2.16	0.45
1:A:570:ASP:HB3	1:A:572:PRO:HD2	1.98	0.45
1:B:317:CYS:HB3	1:B:549:SER:HB2	1.98	0.45
1:B:342:CYS:HA	1:B:367:TYR:CZ	2.51	0.45
1:C:656:LEU:O	1:C:660:LYS:HG3	2.16	0.45
1:D:51:ARG:HD3	10:D:810:HOH:O	2.15	0.45
1:D:266:ALA:HB2	1:D:427:VAL:HG21	1.98	0.45
1:D:472:ASN:ND2	1:D:606:THR:HG21	2.31	0.45
2:M:626:MET:HB3	2:M:630:THR:HB	1.98	0.45
1:A:150:VAL:HG12	1:A:150:VAL:O	2.15	0.45
1:A:607:MET:HE2	1:A:610:VAL:CG2	2.46	0.45
1:C:316:CYS:O	1:C:503:GLN:NE2	2.49	0.45
1:D:384:TYR:HD1	2:O:84:ASN:HB3	1.81	0.45
2:M:540:TYR:CE1	2:M:550:PRO:HD3	2.51	0.45
1:C:359:SER:O	1:C:363:VAL:HG23	2.15	0.45
2:M:316:ILE:HD12	2:M:316:ILE:C	2.41	0.45
2:M:712:LEU:O	2:M:716:GLU:HG3	2.15	0.45
2:N:269:ALA:HB1	2:N:274:PHE:HB2	1.98	0.45
2:O:333:ILE:HD12	2:O:418:TRP:HB3	1.98	0.45
2:P:335:LYS:HD2	2:P:335:LYS:N	2.31	0.45
2:P:392:GLY:HA3	2:P:459:ILE:O	2.15	0.45
2:P:479:GLU:HA	2:P:482:ARG:HD2	1.98	0.45
2:P:664:ILE:HD13	2:P:685:ARG:HG3	1.97	0.45
1:A:187:MET:HE2	1:A:199:ARG:NH1	2.32	0.45
1:A:250:ASP:O	1:A:254:ILE:HG13	2.17	0.45
1:B:537:ILE:HD12	1:B:537:ILE:N	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:602:MET:CE	2:M:647:ILE:HG21	2.46	0.45
2:N:59:ALA:HB2	2:N:630:THR:HA	1.99	0.45
2:O:669:TRP:HA	2:O:700:ALA:O	2.16	0.45
2:O:708:VAL:O	2:O:712:LEU:HB2	2.17	0.45
1:B:28:THR:HG21	1:B:33:VAL:HG12	1.99	0.45
2:M:356:VAL:CG2	2:M:361:ILE:HB	2.45	0.45
2:O:107:TRP:CZ3	2:O:219:ALA:HB1	2.52	0.45
1:A:70:PHE:CE1	1:A:100:ARG:HB3	2.52	0.45
1:A:266:ALA:HB2	1:A:427:VAL:HG21	1.98	0.45
1:C:175:ARG:HB3	1:D:175:ARG:HB3	1.97	0.45
1:D:577:VAL:HG21	1:D:645:ILE:CG2	2.46	0.45
2:P:524:ARG:NH2	2:P:645:MET:SD	2.87	0.45
1:A:18:ARG:NE	1:A:636:MET:HE3	2.32	0.45
1:B:280:LEU:HD13	1:B:288:SER:HB2	1.99	0.45
1:B:569:VAL:HB	1:B:573:LYS:HD3	1.99	0.45
1:C:307:ALA:HB2	1:C:408:PHE:CE2	2.52	0.45
2:P:373:ILE:CD1	2:P:435:VAL:HG22	2.46	0.45
1:A:318:THR:O	1:A:322:VAL:HG22	2.15	0.45
1:B:114[A]:CYS:SG	1:B:208:ILE:CG2	3.05	0.45
2:N:185:ILE:HD12	2:N:185:ILE:N	2.32	0.45
2:O:339:TYR:CE2	2:O:376:ILE:HD11	2.52	0.45
2:O:492:MET:HE2	2:O:492:MET:CA	2.47	0.45
2:P:376:ILE:HD12	2:P:376:ILE:C	2.42	0.45
2:P:399:PHE:HZ	2:P:541:GLU:HG3	1.82	0.45
2:P:496:THR:OG1	2:P:499:THR:HG23	2.17	0.45
1:B:291:ILE:HD13	1:B:397:ALA:CB	2.47	0.45
1:C:107:LEU:HD21	1:C:215:LEU:HB3	1.98	0.45
1:D:274:ASP:OD2	1:D:308:LYS:HE2	2.17	0.45
2:N:316:ILE:C	2:N:316:ILE:HD12	2.42	0.45
2:N:601:ILE:HG23	2:N:635:ILE:HD13	1.98	0.45
1:C:626:ASP:HB3	2:O:212:PHE:CG	2.52	0.45
2:O:173:VAL:HG13	2:O:183:LEU:CD1	2.47	0.45
1:B:607:MET:HE2	1:B:610:VAL:HG22	1.99	0.44
1:C:283:HIS:CE1	4:C:703:XCC:S3	3.10	0.44
1:C:615:LEU:C	1:C:615:LEU:HD23	2.42	0.44
1:D:298:MET:HE3	1:D:298:MET:HB3	1.57	0.44
2:M:342:MET:CE	2:M:384:LEU:HD22	2.47	0.44
2:M:159:ILE:HG12	2:M:253:LEU:HD12	1.98	0.44
2:N:80:PRO:N	2:N:81:PRO:HD2	2.32	0.44
2:N:354:ARG:HH21	2:N:480:VAL:HG11	1.81	0.44
1:A:607:MET:HE1	1:A:634:PHE:CE1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:35:TYR:CZ	2:M:39:LEU:HD11	2.52	0.44
2:M:406:ARG:HB3	2:M:410:PHE:CZ	2.52	0.44
2:N:445:GLY:HA2	2:N:465:VAL:HG11	1.99	0.44
2:O:213:THR:O	2:O:216:VAL:HG12	2.18	0.44
2:O:373:ILE:HD11	2:O:435:VAL:HG22	1.98	0.44
1:A:619:ILE:HA	1:A:623:ILE:HB	1.99	0.44
1:B:282:GLY:C	1:B:318:THR:HB	2.43	0.44
2:N:496:THR:HB	2:N:523:GLU:OE2	2.16	0.44
2:O:80:PRO:N	2:O:81:PRO:HD2	2.33	0.44
2:P:341:GLU:OE2	2:P:427:TRP:NE1	2.39	0.44
2:P:712:LEU:HB3	2:P:713:PRO:HD3	2.00	0.44
1:A:577:VAL:HG13	1:A:649:LEU:HD21	1.99	0.44
1:A:607:MET:HE2	1:A:610:VAL:HG22	1.99	0.44
1:B:445:ALA:HB3	1:C:38:VAL:HG13	1.99	0.44
1:B:559:LEU:O	1:B:563:MET:HG3	2.17	0.44
1:D:114[A]:CYS:SG	1:D:208:ILE:CG2	3.05	0.44
2:M:79:LEU:N	2:M:80:PRO:CD	2.81	0.44
2:O:79:LEU:N	2:O:80:PRO:CD	2.80	0.44
2:P:277:ILE:HD12	2:P:277:ILE:N	2.33	0.44
2:P:365:LYS:HB3	2:P:464:GLN:HG3	1.98	0.44
1:B:52:PHE:CZ	1:B:474:LYS:HA	2.52	0.44
1:B:397:ALA:O	1:B:401:ILE:HG13	2.18	0.44
1:C:351:VAL:HB	1:C:356:ILE:CD1	2.47	0.44
1:C:428:VAL:HG13	1:C:539:ILE:HD11	2.00	0.44
1:C:483:THR:HG23	1:C:639:GLN:OE1	2.17	0.44
1:D:280:LEU:HD13	1:D:288:SER:HB2	1.99	0.44
1:D:399:THR:HG22	1:D:403:MET:HE3	2.00	0.44
2:M:80:PRO:N	2:M:81:PRO:HD2	2.33	0.44
2:O:62:LEU:HD12	2:O:65:ILE:HD12	1.99	0.44
2:P:61:TYR:O	2:P:63:PRO:HD3	2.18	0.44
2:P:455:GLU:C	2:P:457:PRO:HD3	2.42	0.44
1:B:150:VAL:O	1:B:150:VAL:HG12	2.18	0.44
1:B:256:PHE:CG	1:B:289:GLU:HG3	2.52	0.44
1:B:486:LYS:HE2	1:B:519:TYR:CE2	2.52	0.44
1:D:619:ILE:HA	1:D:623:ILE:HB	2.00	0.44
1:D:626:ASP:HB3	2:P:212:PHE:CG	2.53	0.44
2:N:157:ALA:HB3	2:N:183:LEU:CD2	2.47	0.44
2:O:205:ILE:HG22	2:O:205:ILE:O	2.18	0.44
2:P:63:PRO:HB2	2:P:220:ASN:HB2	1.99	0.44
2:P:349:ALA:HA	2:P:384:LEU:O	2.18	0.44
2:P:556:GLU:OE2	2:P:559:PRO:HB3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:338:GLU:OE2	1:A:359:SER:OG	2.34	0.44
2:M:481:ALA:HB1	2:M:485:TYR:CZ	2.53	0.44
2:M:497:ASP:O	2:M:553:LYS:NZ	2.46	0.44
2:M:667:ILE:HG21	2:M:670:MET:SD	2.58	0.44
2:P:316:ILE:C	2:P:316:ILE:HD12	2.42	0.44
2:P:492:MET:HE2	2:P:492:MET:HA	1.98	0.44
1:B:294:ALA:O	1:B:298:MET:HG2	2.17	0.44
1:D:307:ALA:HB2	1:D:408:PHE:CE2	2.53	0.44
1:D:351:VAL:HB	1:D:356:ILE:HD13	1.99	0.44
2:N:340:VAL:CG2	2:N:376:ILE:HD11	2.48	0.44
2:N:488:ARG:O	2:N:492:MET:HG2	2.18	0.44
2:O:497:ASP:OD1	2:O:522:PRO:HD2	2.18	0.44
2:P:470:ASP:HB3	2:P:473:LYS:CG	2.48	0.44
1:C:467:ILE:HG22	1:C:501:SER:OG	2.17	0.43
1:B:467:ILE:HG22	1:B:501:SER:OG	2.17	0.43
2:N:439:PHE:CE1	2:N:443:ASN:HB2	2.54	0.43
1:B:283:HIS:CD2	1:B:317:CYS:HB2	2.53	0.43
1:D:354:GLN:HB3	1:D:587:LYS:NZ	2.33	0.43
2:M:26:TYR:CZ	2:M:30:ILE:HD11	2.53	0.43
2:M:185:ILE:N	2:M:185:ILE:HD12	2.34	0.43
2:M:339:TYR:CG	2:M:435:VAL:HG21	2.53	0.43
2:N:156:GLU:O	2:N:250:ALA:HA	2.18	0.43
2:N:340:VAL:HG21	2:N:376:ILE:HD11	2.00	0.43
2:O:237:ARG:O	2:O:241:ARG:HG3	2.19	0.43
2:P:144:PHE:CE1	2:P:205:ILE:HG12	2.53	0.43
2:P:670:MET:HE3	2:P:674:LEU:CD2	2.47	0.43
1:A:584:MET:HE3	1:B:73:ALA:CB	2.48	0.43
2:N:712:LEU:HB3	2:N:713:PRO:HD3	2.01	0.43
2:M:61:TYR:O	2:M:63:PRO:HD3	2.18	0.43
2:M:347:THR:CG2	2:M:381:LYS:HD2	2.48	0.43
1:A:335:ALA:HB2	1:A:471:ASN:HB3	2.00	0.43
1:D:364:ALA:HB1	1:D:369:THR:HB	2.00	0.43
2:M:335:LYS:HD2	2:M:335:LYS:N	2.33	0.43
2:P:483:GLU:O	2:P:487:GLU:HG3	2.19	0.43
2:O:388:VAL:HG21	2:O:448:LEU:HD13	2.00	0.43
2:P:79:LEU:N	2:P:80:PRO:CD	2.81	0.43
1:A:162:LEU:O	1:A:166:VAL:HG12	2.19	0.43
1:C:3:ARG:O	2:O:162:ARG:NH1	2.49	0.43
2:M:286:LYS:HE2	2:M:286:LYS:HB3	1.87	0.43
2:M:583:CYS:HB3	2:M:586:THR:HG22	2.00	0.43
2:N:335:LYS:H	2:N:335:LYS:HD2	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:247:ARG:NH2	2:P:514:PRO:HG3	2.34	0.43
2:P:277:ILE:HG13	2:P:293:PHE:CZ	2.54	0.43
2:P:316:ILE:HG13	2:P:454:GLU:CD	2.44	0.43
2:P:538:ALA:O	2:P:542:ILE:HG13	2.19	0.43
1:A:73:ALA:HB2	1:B:584:MET:HG3	2.01	0.43
1:A:483:THR:HB	1:A:638:PRO:HB2	2.00	0.43
1:A:503:GLN:O	1:A:507:LYS:HG3	2.19	0.43
1:C:416:ARG:HG3	1:C:417:PRO:HD2	2.01	0.43
2:N:172:ILE:HG23	2:N:302:VAL:CG2	2.49	0.43
2:N:538:ALA:O	2:N:542:ILE:HG13	2.19	0.43
2:N:613:ILE:O	2:N:671:PRO:HD3	2.19	0.43
2:O:420:THR:CG2	2:O:427:TRP:HB3	2.49	0.43
2:O:456:PHE:HB3	2:O:459:ILE:HG23	2.00	0.43
2:O:553:LYS:HG2	2:O:553:LYS:O	2.18	0.43
1:C:409:LYS:O	1:C:413:GLU:HG3	2.19	0.43
1:C:464:VAL:HG11	1:C:563:MET:HE1	2.01	0.43
1:D:150:VAL:O	1:D:150:VAL:HG12	2.18	0.43
2:M:318:LEU:HD12	2:M:318:LEU:N	2.33	0.43
2:N:275:PRO:HD2	2:N:309:ARG:HD3	2.00	0.43
2:O:63:PRO:HB2	2:O:220:ASN:HB2	2.01	0.43
2:O:80:PRO:HB2	2:O:81:PRO:HD3	2.00	0.43
2:O:408:HIS:O	2:O:412:ASN:ND2	2.51	0.43
2:P:37:GLU:OE1	2:P:41:ASN:ND2	2.42	0.43
1:B:368[A]:HIS:CD2	1:B:416:ARG:HD3	2.54	0.42
1:D:20:MET:HE2	1:D:476:PHE:CD1	2.53	0.42
2:M:181:PHE:CE2	2:M:311:ILE:HD13	2.54	0.42
2:M:488:ARG:O	2:M:492:MET:HG2	2.18	0.42
2:N:79:LEU:N	2:N:80:PRO:CD	2.82	0.42
2:O:64:VAL:HG23	2:O:220:ASN:HB3	2.01	0.42
2:O:209:LEU:HD22	2:O:217:HIS:HB2	2.01	0.42
2:O:230:GLY:HA2	2:O:512:PHE:O	2.18	0.42
2:O:425:ILE:O	2:O:425:ILE:HG22	2.18	0.42
1:B:348:ALA:CB	1:B:404:ALA:HB2	2.48	0.42
1:C:190:ILE:HG22	1:C:561:MET:HE3	2.01	0.42
1:D:462:ALA:HB3	1:D:493:VAL:HG22	2.01	0.42
1:D:570:ASP:HB3	1:D:572:PRO:HD2	2.00	0.42
2:M:420:THR:HG22	2:M:427:TRP:HB3	2.01	0.42
2:N:41:ASN:O	2:N:45:ARG:HG3	2.19	0.42
1:C:114[B]:CYS:SG	1:C:208:ILE:CG2	3.08	0.42
1:C:399:THR:HG22	1:C:403:MET:HE3	2.01	0.42
1:C:492:ASN:HA	1:C:524:LEU:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:455:GLU:C	2:M:457:PRO:HD3	2.44	0.42
2:N:382:LEU:HD12	2:N:383:PRO:HD2	2.00	0.42
2:N:588:MET:O	2:N:642:PRO:HB3	2.19	0.42
2:O:382:LEU:HD21	2:O:469:THR:HB	2.01	0.42
2:O:651:TYR:CZ	2:O:657:PHE:HA	2.54	0.42
2:P:528:CYS:HB3	3:P:801:SF4:S3	2.58	0.42
2:P:669:TRP:HA	2:P:700:ALA:O	2.19	0.42
1:A:553:ASN:ND2	1:A:591:ILE:HD13	2.34	0.42
2:O:519:ILE:HB	2:O:592:MET:HE2	2.01	0.42
1:A:374:THR:HA	1:A:386:ILE:O	2.20	0.42
1:B:287:LEU:O	1:B:291:ILE:HG12	2.20	0.42
1:D:186:LEU:CD2	1:D:205:PRO:HD2	2.49	0.42
2:M:259:ASP:OD1	2:M:262:LYS:HD2	2.19	0.42
2:M:269:ALA:HA	2:M:272:THR:HG22	2.02	0.42
2:M:612:MET:HE1	2:M:622:THR:CB	2.49	0.42
2:O:406:ARG:HG3	2:O:456:PHE:CZ	2.54	0.42
2:O:455:GLU:C	2:O:457:PRO:HD3	2.44	0.42
1:C:317:CYS:O	1:C:321:GLU:HG2	2.19	0.42
2:N:714:TYR:CZ	2:N:718:LYS:HE3	2.55	0.42
1:B:585:SER:OG	4:B:702:XCC:S4	2.74	0.42
1:C:150:VAL:HG13	1:C:184:THR:CG2	2.50	0.42
2:M:240:GLN:O	2:M:244:GLN:HG2	2.20	0.42
2:M:445:GLY:HA2	2:M:465:VAL:HG11	2.02	0.42
2:N:655:LYS:HG2	2:N:685:ARG:NH1	2.35	0.42
2:O:428:LEU:N	2:O:428:LEU:HD12	2.34	0.42
1:A:584:MET:HG3	1:B:73:ALA:HB2	2.01	0.42
1:A:662:VAL:HG21	2:M:194:LEU:HD13	2.00	0.42
1:B:98:VAL:O	1:B:102:VAL:HG12	2.20	0.42
1:B:359:SER:O	1:B:363:VAL:HG23	2.20	0.42
1:C:364:ALA:HB1	1:C:369:THR:HB	2.01	0.42
1:C:428:VAL:HG13	1:C:539:ILE:CD1	2.50	0.42
1:D:482:LEU:HD13	1:D:510:LEU:HD11	2.01	0.42
2:N:524:ARG:NH2	2:N:645:MET:SD	2.85	0.42
2:O:259:ASP:OD1	2:O:262:LYS:HD2	2.20	0.42
2:P:219:ALA:O	2:P:223:LEU:HG	2.20	0.42
2:P:236:ALA:O	2:P:240:GLN:HG2	2.20	0.42
1:A:591:ILE:HG23	1:A:594:TRP:CZ2	2.55	0.42
1:B:187:MET:HE2	1:B:199:ARG:NH1	2.35	0.42
1:B:256:PHE:CD2	1:B:289:GLU:HG3	2.55	0.42
1:C:321:GLU:OE1	1:C:325:ARG:NH1	2.40	0.42
2:M:226:GLY:HA2	2:M:243:TYR:CE1	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:602:MET:O	2:N:644:PHE:HA	2.20	0.42
1:A:112:ALA:HA	1:B:217:ASN:OD1	2.20	0.42
1:C:187:MET:HE2	1:C:199:ARG:NH1	2.35	0.42
1:C:655:LYS:HG3	2:O:191:GLU:HG2	2.02	0.42
1:D:142:LYS:NZ	1:D:434:GLU:OE2	2.47	0.42
2:M:141:VAL:HG22	2:M:221:TYR:CZ	2.54	0.42
2:N:383:PRO:O	2:N:469:THR:HA	2.20	0.42
2:N:406:ARG:HB3	2:N:410:PHE:CE2	2.55	0.42
2:N:470:ASP:O	2:N:474:VAL:HG23	2.19	0.42
2:O:129:PRO:O	2:O:132:TRP:HB2	2.20	0.42
2:O:668:VAL:HG12	2:O:720:HIS:CE1	2.55	0.42
2:P:470:ASP:HB3	2:P:473:LYS:HG2	2.01	0.42
1:A:326:GLN:HB2	1:A:328:ILE:HG12	2.01	0.41
1:C:299:GLU:HG2	1:C:303:LYS:HE2	2.02	0.41
1:C:591:ILE:HG23	1:C:594:TRP:CZ2	2.55	0.41
2:M:360:GLU:O	2:M:360:GLU:HG2	2.20	0.41
2:M:622:THR:HB	2:M:623:PRO:HD2	2.02	0.41
2:N:316:ILE:HG13	2:N:454:GLU:CD	2.44	0.41
1:A:223:MET:HE3	1:B:377:ASN:O	2.20	0.41
1:B:574:VAL:O	1:B:576:PHE:N	2.53	0.41
1:B:607:MET:HE1	1:B:634:PHE:CE1	2.54	0.41
1:C:552:ASP:CG	1:C:555:ARG:HE	2.26	0.41
2:M:18:PRO:HG3	2:M:288:ILE:HG12	2.02	0.41
2:O:585:TYR:CE1	2:O:656:LYS:HD2	2.55	0.41
2:O:622:THR:HB	2:O:623:PRO:HD2	2.01	0.41
1:A:374:THR:HB	1:A:388:TYR:HB2	2.02	0.41
1:C:353:VAL:O	1:C:354:GLN:HB2	2.20	0.41
2:M:223:LEU:HA	2:M:244:GLN:NE2	2.35	0.41
1:A:256:PHE:CD2	1:A:289:GLU:HG3	2.56	0.41
1:D:39:LYS:NZ	1:D:43:ASP:OD2	2.49	0.41
1:D:316:CYS:HB3	1:D:317:CYS:H	1.64	0.41
2:M:129:PRO:O	2:M:132:TRP:HB2	2.20	0.41
2:N:185:ILE:HG22	2:N:190:VAL:CG2	2.51	0.41
2:N:342:MET:HE2	2:N:384:LEU:HD22	2.02	0.41
2:N:403:LEU:HD21	2:N:460:VAL:CG2	2.49	0.41
2:P:503:PHE:O	2:P:551:ILE:N	2.46	0.41
1:A:266:ALA:O	1:A:267:ASN:CB	2.68	0.41
1:C:559:LEU:O	1:C:563:MET:HG3	2.20	0.41
1:C:641:ALA:O	1:C:645:ILE:HG13	2.21	0.41
1:D:484:VAL:O	1:D:488:LEU:HG	2.21	0.41
2:M:277:ILE:N	2:M:277:ILE:HD12	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:373:ILE:HG22	2:M:440:ARG:HG3	2.03	0.41
2:O:109:ALA:O	2:O:113:ILE:HG13	2.21	0.41
2:P:373:ILE:HG22	2:P:440:ARG:HG3	2.03	0.41
2:P:527:LEU:HD12	2:P:528:CYS:N	2.35	0.41
2:P:583:CYS:HB3	2:P:586:THR:HG22	2.02	0.41
2:P:598:PHE:CE1	2:P:601:ILE:HD11	2.56	0.41
1:A:158:SER:OG	1:A:161:GLU:HG3	2.20	0.41
1:A:227:PRO:O	1:A:231:VAL:HG23	2.21	0.41
2:N:132:TRP:HA	2:N:208:PRO:HB2	2.03	0.41
2:N:142:ARG:CZ	2:N:632:ALA:HB1	2.51	0.41
2:O:306:MET:HE1	2:O:313:LEU:HG	2.03	0.41
2:O:403:LEU:HD21	2:O:460:VAL:CG2	2.50	0.41
2:O:602:MET:HE2	2:O:647:ILE:HG21	2.03	0.41
2:P:142:ARG:CZ	2:P:632:ALA:HB1	2.50	0.41
2:P:602:MET:O	2:P:644:PHE:HA	2.19	0.41
1:B:572:PRO:HG2	1:B:674:TYR:CZ	2.55	0.41
1:D:607:MET:HE2	1:D:610:VAL:HG22	2.03	0.41
1:D:615:LEU:HD23	1:D:615:LEU:C	2.45	0.41
2:M:374:ASP:HB3	2:M:440:ARG:HH21	1.85	0.41
2:N:455:GLU:C	2:N:457:PRO:HD3	2.45	0.41
2:N:586:THR:HG23	2:N:661:ASP:OD1	2.20	0.41
2:O:63:PRO:HB3	2:O:224:ARG:HG3	2.03	0.41
2:O:80:PRO:HB2	2:O:81:PRO:CD	2.50	0.41
2:O:169:LEU:HG	2:O:193:LEU:HD21	2.01	0.41
2:O:282:LEU:CB	2:O:287:GLN:HG3	2.50	0.41
2:O:457:PRO:O	2:O:458:ALA:HB3	2.21	0.41
1:A:168:GLU:O	1:A:172:GLU:HG3	2.21	0.41
1:C:20:MET:HE2	1:C:476:PHE:CG	2.55	0.41
2:M:316:ILE:HG13	2:M:454:GLU:CD	2.46	0.41
2:O:104:GLU:OE1	2:O:237:ARG:HD2	2.20	0.41
2:O:442:LYS:O	2:O:446:GLU:HG3	2.21	0.41
1:A:24:ASN:O	1:A:27:ARG:HG2	2.20	0.41
1:A:537:ILE:HD12	1:A:537:ILE:N	2.36	0.41
1:B:138:LYS:HD2	1:B:256:PHE:HA	2.02	0.41
1:B:316:CYS:O	1:B:503:GLN:NE2	2.54	0.41
1:B:426:ARG:HH22	1:B:539:ILE:CG1	2.33	0.41
1:B:456:ILE:HD13	1:B:463:GLY:HA2	2.03	0.41
1:B:615:LEU:HD23	1:B:615:LEU:C	2.45	0.41
1:D:65:GLY:CA	1:D:609:PRO:HB2	2.51	0.41
1:D:238:ALA:O	1:D:241:ASP:HB3	2.21	0.41
1:D:574:VAL:O	1:D:576:PHE:N	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:583:CYS:HB2	2:M:586:THR:CG2	2.50	0.41
2:N:157:ALA:HB3	2:N:183:LEU:HD23	2.02	0.41
2:N:160:LEU:HD13	2:N:215:ILE:HD13	2.02	0.41
2:N:169:LEU:HG	2:N:193:LEU:HD21	2.03	0.41
2:N:456:PHE:HA	2:N:542:ILE:CD1	2.51	0.41
2:O:678:LEU:O	2:O:682:PHE:HB2	2.21	0.41
2:P:65:ILE:HG12	2:P:108:TYR:CD1	2.56	0.41
2:P:237:ARG:O	2:P:241:ARG:HG3	2.21	0.41
2:P:399:PHE:CZ	2:P:541:GLU:HG3	2.56	0.41
2:P:408:HIS:O	2:P:412:ASN:ND2	2.54	0.41
2:P:457:PRO:O	2:P:458:ALA:HB3	2.21	0.41
2:P:572:TYR:HA	2:P:579:LEU:O	2.21	0.41
1:A:26:ASP:HA	1:A:37:LEU:HD13	2.02	0.41
1:A:70:PHE:CZ	1:A:100:ARG:HB3	2.56	0.41
1:A:298:MET:HE1	1:A:402:ARG:HG2	2.03	0.41
1:A:320:ASN:O	1:A:324:MET:HG2	2.21	0.41
1:A:342:CYS:HA	1:A:367:TYR:CZ	2.55	0.41
1:A:574:VAL:O	1:A:576:PHE:N	2.51	0.41
1:B:280:LEU:HD21	1:B:291:ILE:HG21	2.03	0.41
1:B:353:VAL:O	1:B:354:GLN:HB2	2.19	0.41
1:B:570:ASP:HB3	1:B:572:PRO:HD2	2.03	0.41
1:D:660:LYS:O	1:D:664:GLU:HG3	2.21	0.41
2:M:80:PRO:HB2	2:M:81:PRO:HD3	2.03	0.41
2:O:497:ASP:OD2	2:O:585:TYR:OH	2.26	0.41
2:O:551:ILE:HG21	2:O:567:VAL:HG22	2.03	0.41
2:P:318:LEU:HD12	2:P:318:LEU:N	2.36	0.41
2:P:670:MET:HE3	2:P:674:LEU:HG	2.03	0.41
1:A:66:ILE:CG2	1:A:612:GLY:HA3	2.47	0.40
1:B:332:THR:OG1	1:B:336:SER:OG	2.20	0.40
1:B:351:VAL:HB	1:B:356:ILE:CD1	2.51	0.40
1:C:218:GLN:HG2	1:D:377:ASN:HB2	2.03	0.40
1:D:3:ARG:O	2:P:162:ARG:NH1	2.54	0.40
1:D:278:PHE:HB3	1:D:312:LEU:HD23	2.03	0.40
1:D:283:HIS:CD2	1:D:317:CYS:HB2	2.56	0.40
2:M:456:PHE:HD2	2:M:542:ILE:HD11	1.86	0.40
2:P:209:LEU:HD11	2:P:218:ALA:HB2	2.03	0.40
2:P:278:THR:HG21	2:P:282:LEU:HD11	2.04	0.40
2:P:393:ARG:HG2	2:P:393:ARG:H	1.65	0.40
1:A:186:LEU:CD2	1:A:205:PRO:HD2	2.50	0.40
1:A:214:GLU:HA	1:A:214:GLU:OE1	2.22	0.40
1:A:256:PHE:CG	1:A:289:GLU:HG3	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:261:PRO:HA	1:B:429:ALA:O	2.20	0.40
1:B:330:LEU:HD13	1:B:503:GLN:HG2	2.02	0.40
1:C:105:MET:HE1	1:C:609:PRO:CD	2.50	0.40
2:N:528:CYS:HB3	3:N:801:SF4:S3	2.61	0.40
2:P:670:MET:HE3	2:P:674:LEU:CG	2.51	0.40
1:A:241:ASP:O	1:A:245:GLU:HG3	2.20	0.40
1:B:28:THR:HG21	1:B:33:VAL:CG1	2.51	0.40
1:B:317:CYS:O	1:B:321:GLU:HG2	2.22	0.40
1:C:35:GLU:OE1	1:C:419:TYR:OH	2.30	0.40
1:D:279:VAL:HA	1:D:313:VAL:HG23	2.03	0.40
1:D:428:VAL:HG13	1:D:539:ILE:HD11	2.02	0.40
1:D:482:LEU:HD12	1:D:486:LYS:HE3	2.03	0.40
2:M:538:ALA:O	2:M:542:ILE:HG13	2.21	0.40
2:P:80:PRO:HB2	2:P:81:PRO:HD3	2.02	0.40
1:A:594:TRP:O	1:A:598:LEU:HG	2.20	0.40
1:A:655:LYS:HG3	2:M:191:GLU:HG2	2.03	0.40
1:B:472:ASN:ND2	1:B:606:THR:HG21	2.37	0.40
1:C:61:ILE:HD13	1:C:77:ARG:CZ	2.49	0.40
1:D:61:ILE:HD13	1:D:77:ARG:NE	2.26	0.40
2:N:149:VAL:CG1	2:N:509:CYS:HA	2.47	0.40
2:O:603:ALA:HB3	2:O:612:MET:CG	2.52	0.40
2:P:612:MET:HE1	2:P:622:THR:CB	2.51	0.40
1:B:264:SER:OG	1:B:265:GLU:N	2.54	0.40
1:D:65:GLY:HA3	1:D:609:PRO:HB2	2.04	0.40
1:D:173:ASP:O	1:D:208:ILE:HG13	2.21	0.40
1:D:316:CYS:O	1:D:503:GLN:NE2	2.55	0.40
1:D:658:VAL:O	1:D:662:VAL:HG23	2.22	0.40
2:O:209:LEU:HD11	2:O:218:ALA:HB2	2.02	0.40
2:P:406:ARG:HB3	2:P:410:PHE:CZ	2.56	0.40
2:P:470:ASP:OD2	2:P:472:ALA:HB3	2.21	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	673/674 (100%)	647 (96%)	25 (4%)	1 (0%)	48	66
1	B	673/674 (100%)	648 (96%)	24 (4%)	1 (0%)	48	66
1	C	673/674 (100%)	650 (97%)	22 (3%)	1 (0%)	48	66
1	D	672/674 (100%)	648 (96%)	23 (3%)	1 (0%)	48	66
2	M	726/729 (100%)	702 (97%)	23 (3%)	1 (0%)	48	66
2	N	726/729 (100%)	699 (96%)	25 (3%)	2 (0%)	36	53
2	O	724/729 (99%)	696 (96%)	24 (3%)	4 (1%)	21	35
2	P	726/729 (100%)	700 (96%)	24 (3%)	2 (0%)	36	53
All	All	5593/5612 (100%)	5390 (96%)	190 (3%)	13 (0%)	43	61

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	N	187	ASP
2	O	187	ASP
1	A	267	ASN
2	M	596	GLY
2	N	596	GLY
2	O	596	GLY
2	P	187	ASP
2	P	596	GLY
2	O	427	TRP
1	B	354	GLN
1	C	354	GLN
1	D	354	GLN
2	O	205	ILE

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	544/543 (100%)	542 (100%)	2 (0%)	84	92
1	B	544/543 (100%)	543 (100%)	1 (0%)	87	94
1	C	544/543 (100%)	542 (100%)	2 (0%)	84	92
1	D	543/543 (100%)	541 (100%)	2 (0%)	84	92
2	M	610/611 (100%)	609 (100%)	1 (0%)	87	94
2	N	610/611 (100%)	608 (100%)	2 (0%)	86	93
2	O	608/611 (100%)	607 (100%)	1 (0%)	87	94
2	P	610/611 (100%)	609 (100%)	1 (0%)	87	94
All	All	4613/4616 (100%)	4601 (100%)	12 (0%)	86	93

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	118	ASN
1	A	615	LEU
1	B	318	THR
1	C	118	ASN
1	C	318	THR
1	D	298	MET
1	D	318	THR
2	M	149	VAL
2	N	149	VAL
2	N	405	ARG
2	O	317	LYS
2	P	149	VAL

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	275	GLN
1	A	454	GLN
1	A	622	GLN
1	B	293	GLN
1	B	454	GLN
1	B	479	ASN
1	B	503	GLN
1	B	622	GLN
1	C	293	GLN
1	C	389	GLN

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Mol	Chain	Res	Type
1	C	479	ASN
1	C	503	GLN
1	C	622	GLN
1	D	122	HIS
1	D	454	GLN
1	D	479	ASN
1	D	565	ASN
1	D	622	GLN
2	M	426	ASN
2	M	549	GLN
2	N	581	GLN
2	O	515	ASN
2	O	576	ASN
2	O	618	HIS
2	P	426	ASN
2	P	581	GLN
2	P	590	ASN
2	P	618	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 28 ligands modelled in this entry, 9 are monoatomic - leaving 19 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	SF4	A	702	1	0,12,12	-	-	-		
8	CMO	P	804	7	0,1,1	-	-	-		
6	PG5	C	704	-	11,11,11	0.19	0	10,10,10	0.28	0
3	SF4	O	900	2	0,12,12	-	-	-		
3	SF4	D	701	1	0,12,12	-	-	-		
3	SF4	M	801	2	0,12,12	-	-	-		
5	PGE	A	704	-	9,9,9	0.36	0	8,8,8	0.20	0
8	CMO	M	804	-	0,1,1	-	-	-		
3	SF4	P	801	2	0,12,12	-	-	-		
3	SF4	A	701	1	0,12,12	-	-	-		
8	CMO	N	804	7	0,1,1	-	-	-		
4	XCC	D	702	1	0,11,11	-	-	-		
3	SF4	C	702	1	0,12,12	-	-	-		
3	SF4	N	801	2	0,12,12	-	-	-		
3	SF4	B	701	1	0,12,12	-	-	-		
3	SF4	C	701	1	0,12,12	-	-	-		
4	XCC	A	703	1,10	0,11,11	-	-	-		
4	XCC	C	703	1,10	0,11,11	-	-	-		
4	XCC	B	702	1,10	0,11,11	-	-	-		

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	SF4	A	702	1	-	-	0/6/5/5
6	PG5	C	704	-	-	4/9/9/9	-
3	SF4	N	801	2	-	-	0/6/5/5
3	SF4	O	900	2	-	-	0/6/5/5
3	SF4	A	701	1	-	-	0/6/5/5
3	SF4	D	701	1	-	-	0/6/5/5
3	SF4	M	801	2	-	-	0/6/5/5
3	SF4	C	701	1	-	-	0/6/5/5
4	XCC	D	702	1	-	-	0/3/3/3
5	PGE	A	704	-	-	2/7/7/7	-
4	XCC	A	703	1,10	-	-	0/3/3/3
4	XCC	C	703	1,10	-	-	0/3/3/3
3	SF4	C	702	1	-	-	0/6/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SF4	B	701	1	-	-	0/6/5/5
4	XCC	B	702	1,10	-	-	0/3/3/3
3	SF4	P	801	2	-	-	0/6/5/5

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

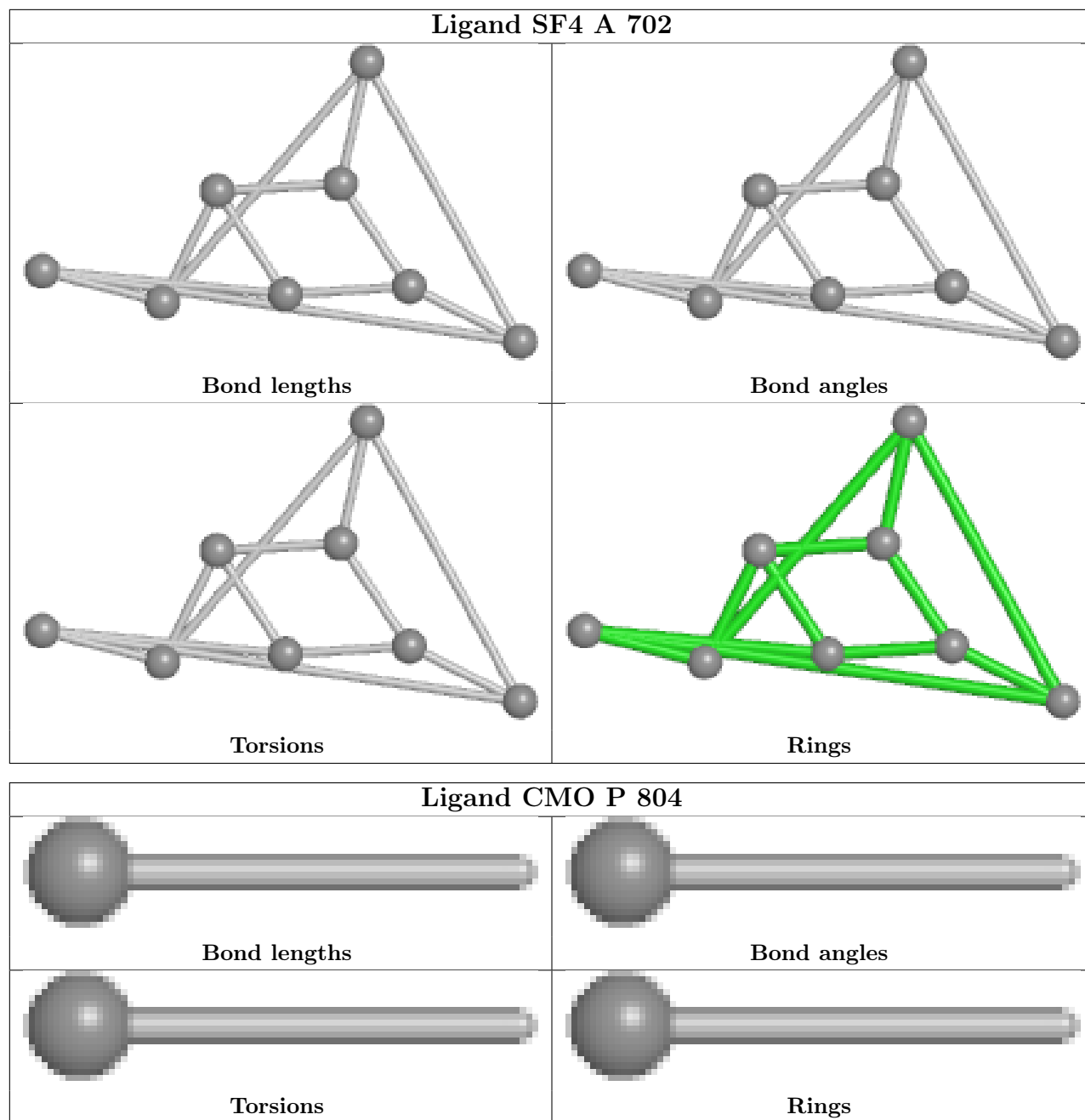
Mol	Chain	Res	Type	Atoms
5	A	704	PGE	O3-C5-C6-O4
5	A	704	PGE	O1-C1-C2-O2
6	C	704	PG5	C3-C2-O1-C1
6	C	704	PG5	O1-C2-C3-O2
6	C	704	PG5	O2-C4-C5-O3
6	C	704	PG5	O3-C6-C7-O4

There are no ring outliers.

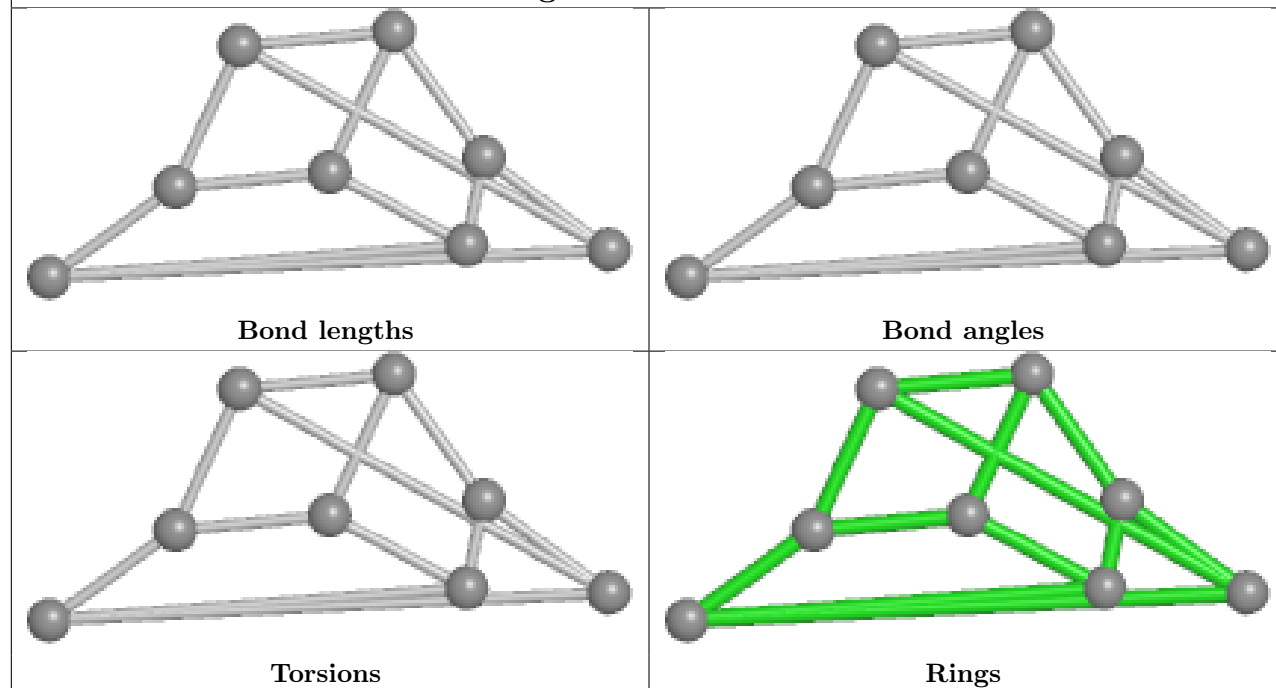
7 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	704	PG5	1	0
3	P	801	SF4	1	0
4	D	702	XCC	3	0
3	N	801	SF4	1	0
4	A	703	XCC	2	0
4	C	703	XCC	2	0
4	B	702	XCC	3	0

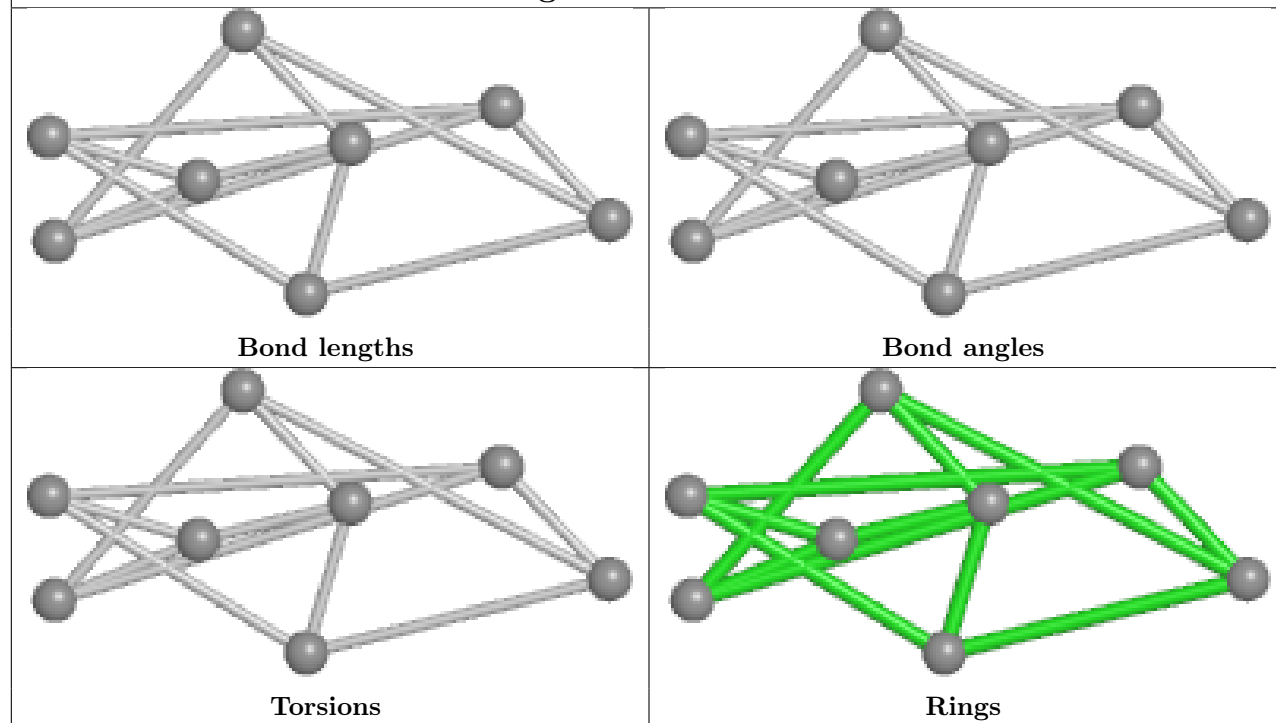
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

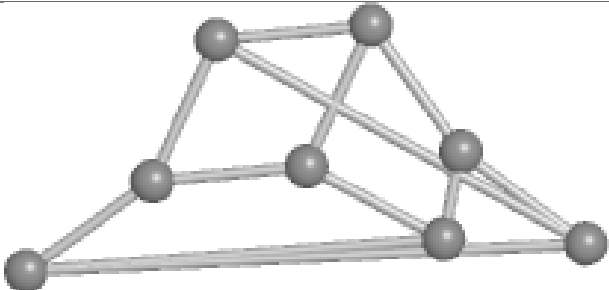
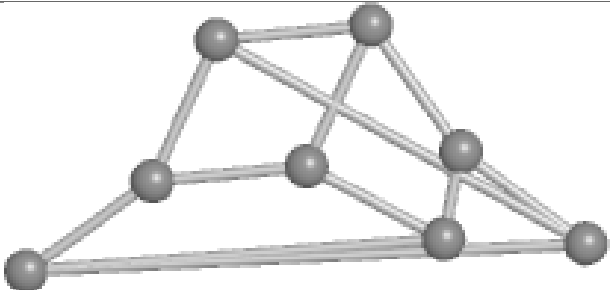
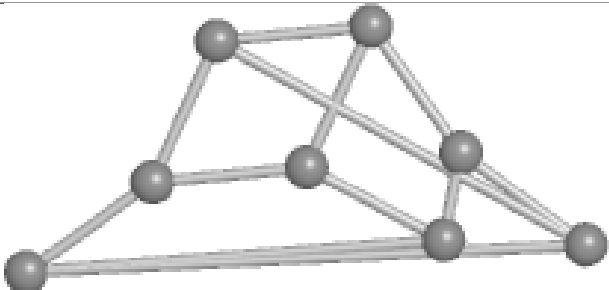
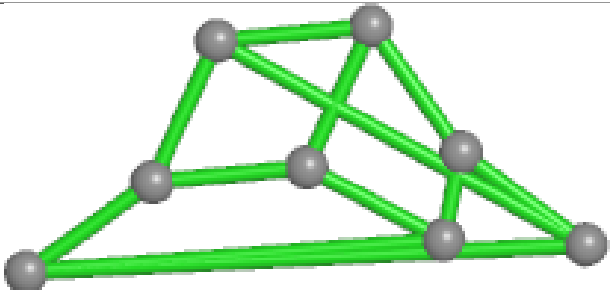


Ligand SF4 O 900

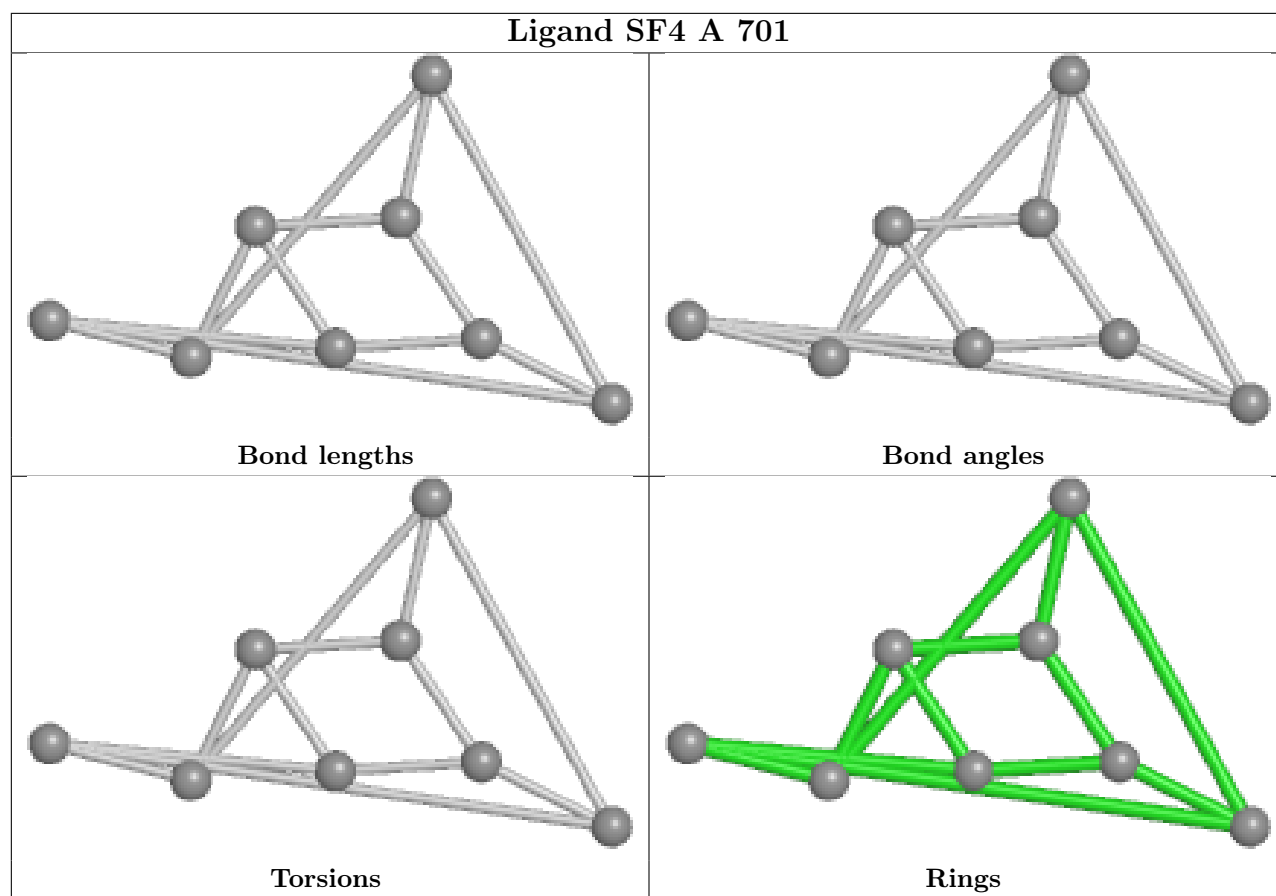
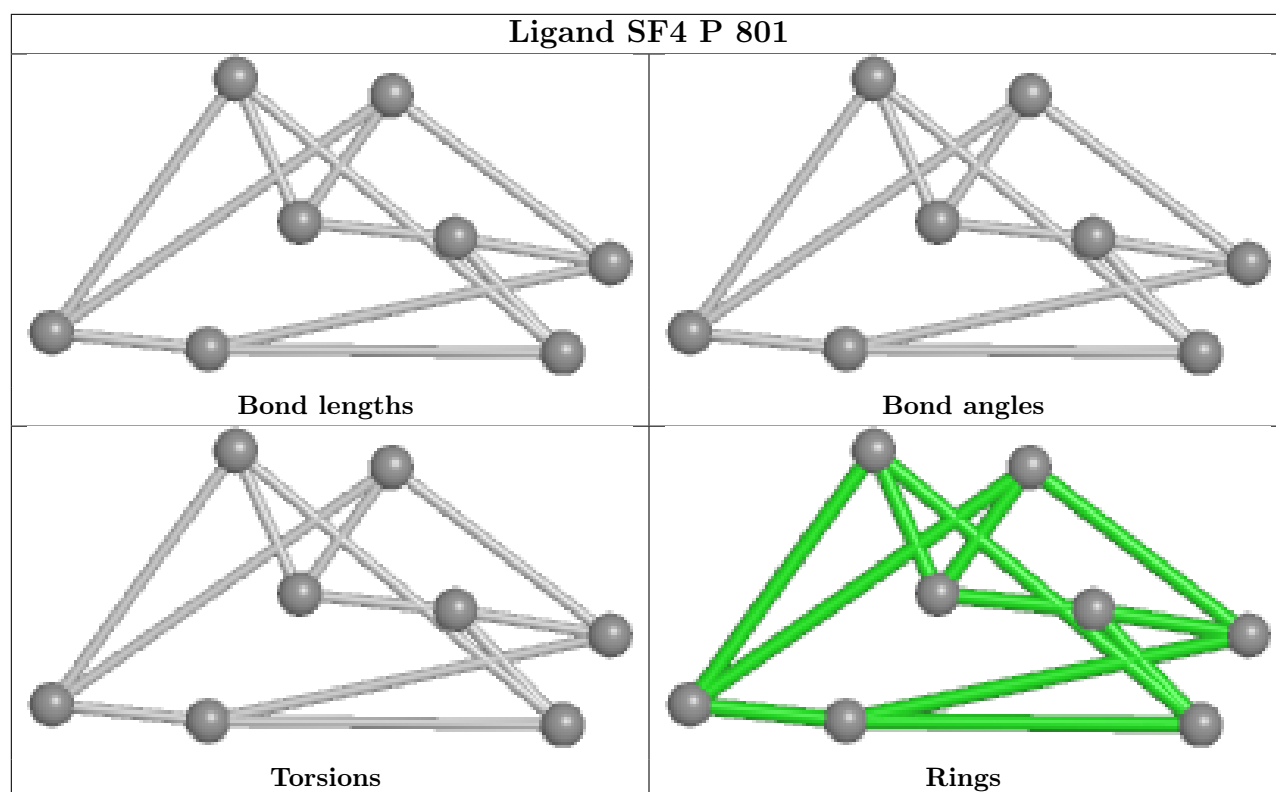


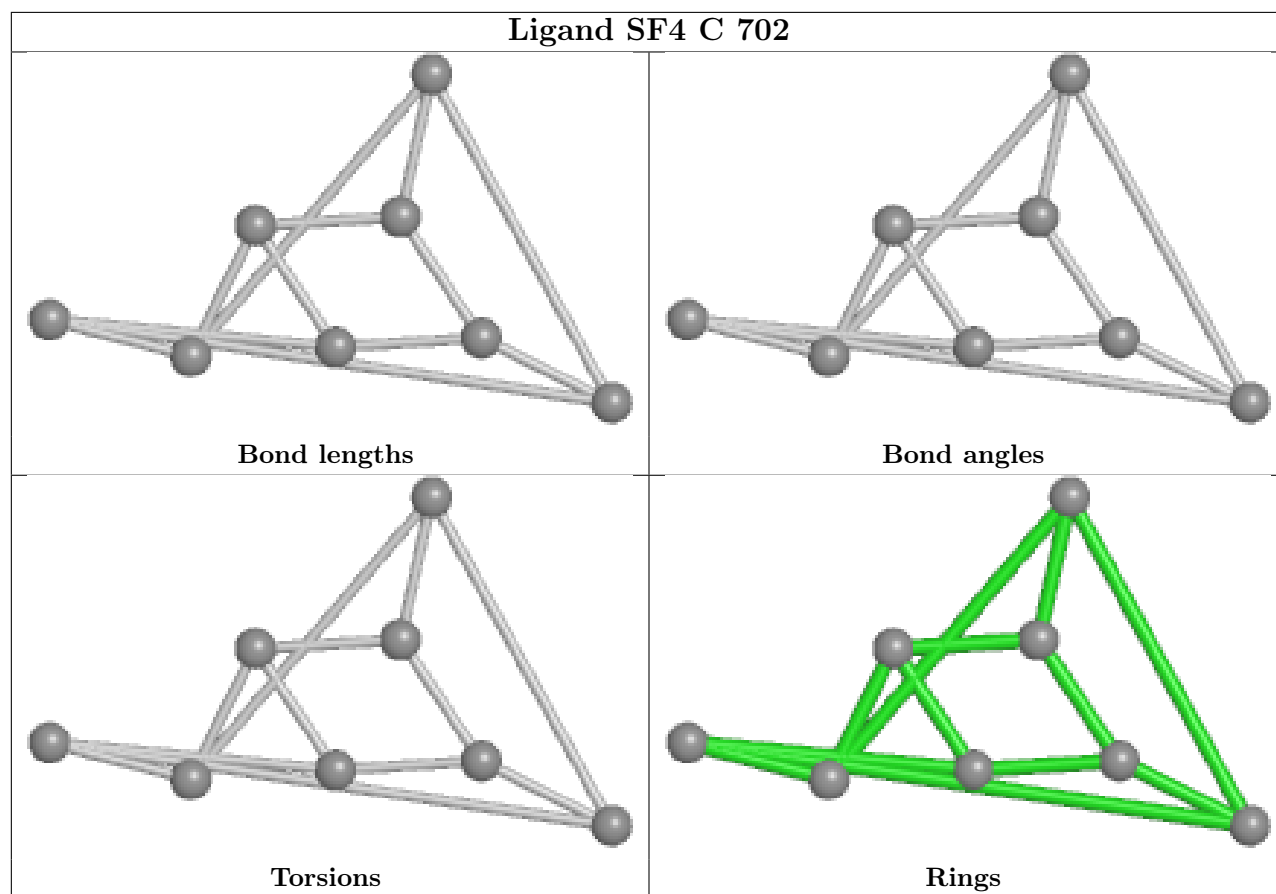
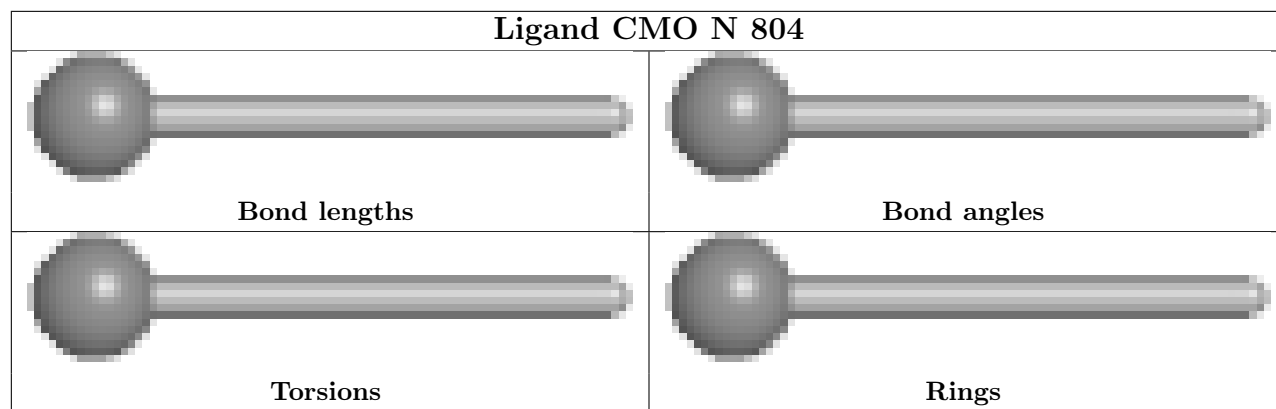
Ligand SF4 D 701

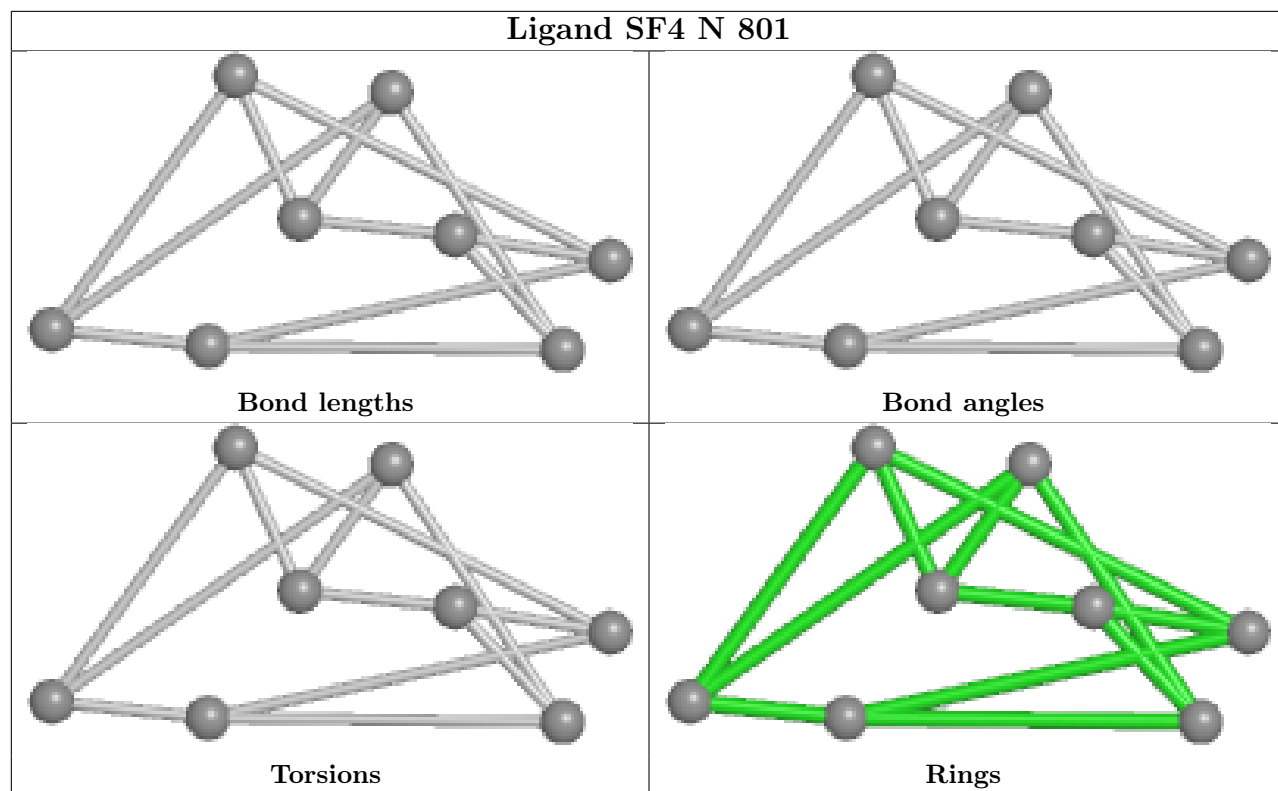


Ligand SF4 M 801	
	Bond lengths
	Bond angles
	Torsions
	Rings

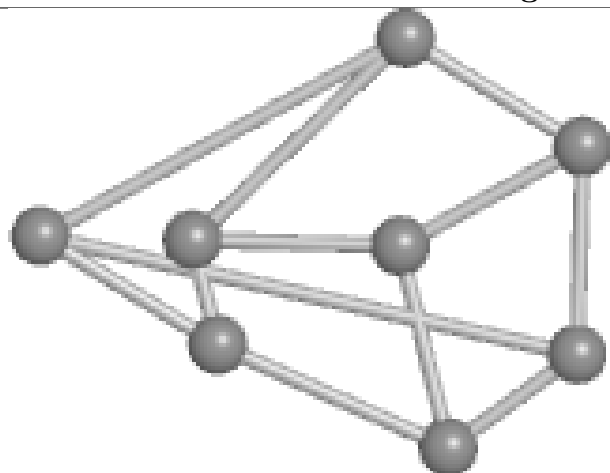
Ligand CMO M 804	
	Bond lengths
	Bond angles
	Torsions
	Rings



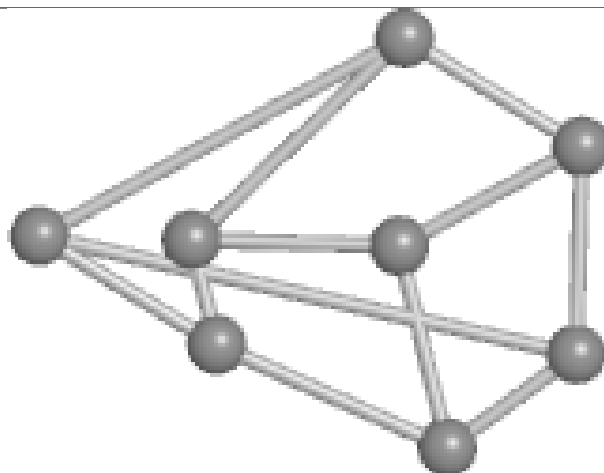




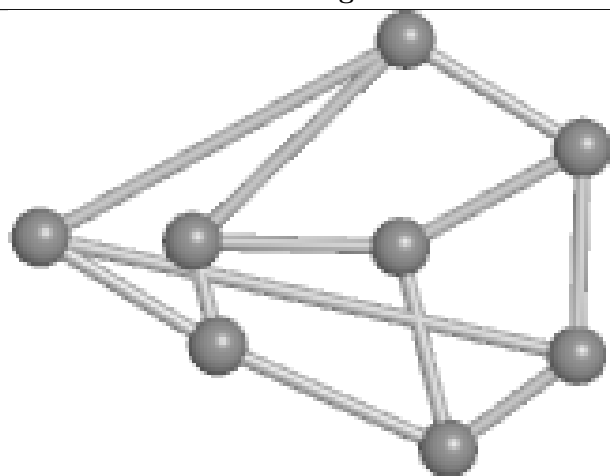
Ligand SF4 B 701



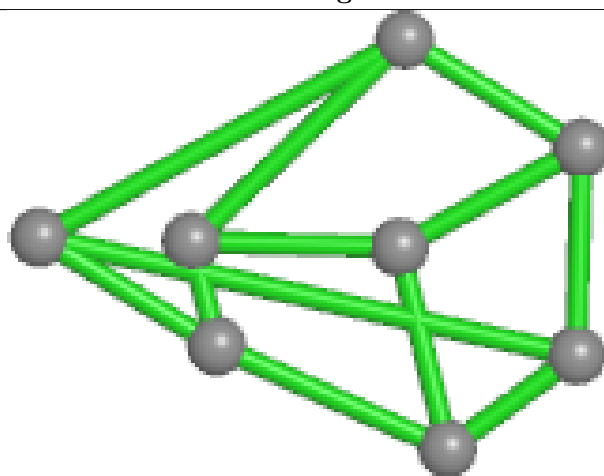
Bond lengths



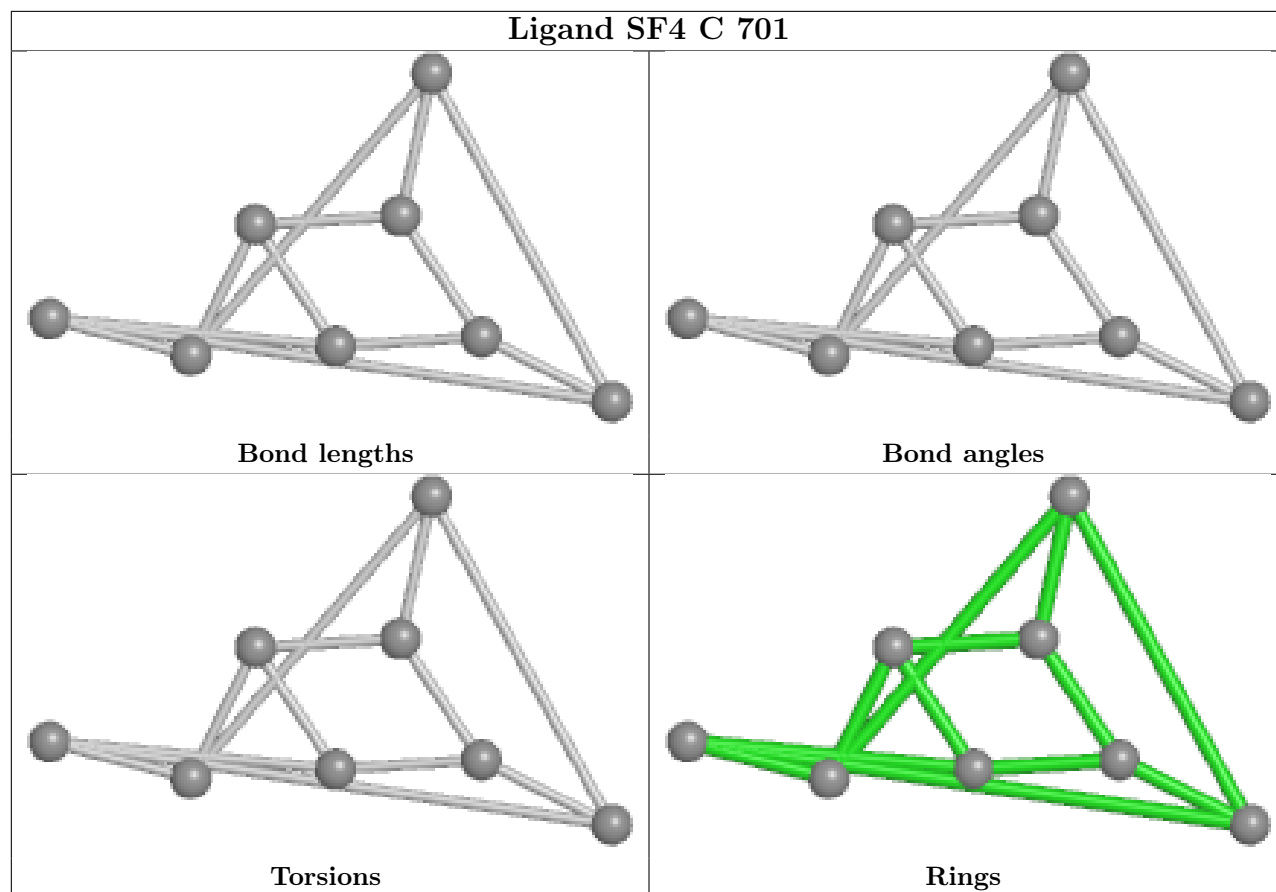
Bond angles



Torsions



Rings



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	672/674 (99%)	-0.09	10 (1%) 72 69	16, 35, 55, 127	3 (0%)
1	B	672/674 (99%)	-0.12	7 (1%) 79 77	16, 35, 54, 121	3 (0%)
1	C	672/674 (99%)	0.11	11 (1%) 70 68	19, 40, 61, 109	3 (0%)
1	D	672/674 (99%)	0.19	15 (2%) 62 59	20, 44, 67, 116	2 (0%)
2	M	728/729 (99%)	0.12	9 (1%) 76 74	23, 43, 71, 124	0
2	N	728/729 (99%)	0.13	9 (1%) 76 74	20, 41, 72, 107	0
2	O	726/729 (99%)	1.36	193 (26%) 1 1	37, 87, 163, 238	0
2	P	728/729 (99%)	0.66	62 (8%) 16 14	28, 57, 117, 194	0
All	All	5598/5612 (99%)	0.31	316 (5%) 30 26	16, 43, 112, 238	11 (0%)

All (316) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	O	316	ILE	6.6
2	N	729	MET	5.1
2	O	525	VAL	5.1
2	O	366	ILE	5.1
2	M	729	MET	5.0
1	C	540	GLY	5.0
2	O	399	PHE	4.9
2	O	449	VAL	4.8
2	O	465	VAL	4.8
1	D	535	ALA	4.7
2	O	425	ILE	4.5
2	O	318	LEU	4.5
1	C	77	ARG	4.4
2	O	481	ALA	4.3
1	D	540	GLY	4.3
2	P	353	VAL	4.3

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Mol	Chain	Res	Type	RSRZ
2	P	542	ILE	4.2
2	O	472	ALA	4.2
2	O	402	VAL	4.2
2	O	450	ALA	4.2
2	O	463	VAL	4.1
2	P	385	GLY	4.1
2	P	480	VAL	4.1
1	D	539	ILE	4.0
2	O	648	GLY	4.0
2	P	469	THR	4.0
2	O	411	ILE	3.9
2	O	444	TYR	3.9
2	P	492	MET	3.9
2	O	447	ILE	3.9
2	O	557	ILE	3.9
2	O	460	VAL	3.9
2	O	407	ILE	3.8
1	B	538	GLU	3.8
2	P	474	VAL	3.7
1	C	539	ILE	3.7
2	O	355	THR	3.7
2	O	583	CYS	3.7
2	O	441	PHE	3.7
2	O	480	VAL	3.7
2	P	475	LYS	3.7
2	P	494	GLY	3.6
2	O	608	CYS	3.6
1	D	534	GLY	3.6
2	O	563	ILE	3.6
2	O	667	ILE	3.6
2	O	504	TYR	3.6
2	O	403	LEU	3.6
2	O	381	LYS	3.5
2	O	485	TYR	3.5
1	C	535	ALA	3.5
2	O	384	LEU	3.5
2	O	367	GLU	3.5
2	O	612	MET	3.5
2	O	603	ALA	3.5
2	O	582	VAL	3.5
2	O	448	LEU	3.4
2	O	430	VAL	3.4

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Mol	Chain	Res	Type	RSRZ
2	O	428	LEU	3.4
2	O	696	ILE	3.4
2	O	567	VAL	3.4
2	O	658	ILE	3.3
2	O	475	LYS	3.3
2	O	571	LEU	3.3
2	O	368	VAL	3.3
2	P	459	ILE	3.2
2	O	534	LEU	3.2
2	O	519	ILE	3.2
2	O	606	PRO	3.2
2	P	399	PHE	3.2
2	O	544	HIS	3.2
2	O	427	TRP	3.1
1	D	536	ASN	3.1
2	P	384	LEU	3.1
2	M	728	ILE	3.1
2	O	390	ILE	3.1
2	O	467	ILE	3.1
2	O	584	LEU	3.1
2	O	370	GLY	3.1
2	P	491	ARG	3.1
2	P	728	ILE	3.1
2	O	594	SER	3.0
2	O	517	VAL	3.0
2	O	551	ILE	3.0
2	N	373	ILE	3.0
2	O	542	ILE	3.0
2	O	610	GLY	3.0
2	O	520	VAL	3.0
2	P	356	VAL	3.0
2	O	409	ASP	3.0
2	O	452	MET	2.9
2	P	336	GLY	2.9
2	O	353	VAL	2.9
2	O	683	VAL	2.9
2	O	559	PRO	2.9
2	O	533	TRP	2.9
2	O	439	PHE	2.9
2	N	2	THR	2.9
2	O	497	ASP	2.9
2	O	647	ILE	2.9

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Mol	Chain	Res	Type	RSRZ
2	O	371	PRO	2.9
2	O	397	ALA	2.9
1	D	330	LEU	2.9
2	O	352	LEU	2.9
2	O	373	ILE	2.8
2	P	457	PRO	2.8
1	A	535	ALA	2.8
2	O	657	PHE	2.8
2	O	394	LYS	2.8
2	O	466	THR	2.8
2	P	397	ALA	2.8
1	C	5	ARG	2.8
2	O	494	GLY	2.8
2	O	638	GLY	2.8
2	P	381	LYS	2.8
2	O	496	THR	2.8
2	O	715	LEU	2.8
2	O	238	GLU	2.8
2	P	604	ILE	2.7
2	N	357	SER	2.7
2	O	686	SER	2.7
2	O	492	MET	2.7
1	C	416	ARG	2.7
2	O	354	ARG	2.7
2	O	389	ASP	2.7
1	A	537	ILE	2.7
1	D	417	PRO	2.7
1	C	536	ASN	2.7
2	O	378	GLU	2.7
2	O	664	ILE	2.7
2	O	247	ARG	2.7
2	O	474	VAL	2.7
2	O	685	ARG	2.7
2	P	340	VAL	2.7
2	P	579	LEU	2.7
1	B	417	PRO	2.6
2	P	333	ILE	2.6
2	P	495	LEU	2.6
2	P	316	ILE	2.6
1	A	416	ARG	2.6
2	O	313	LEU	2.6
2	O	340	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
2	O	356	VAL	2.6
1	C	639	GLN	2.6
2	P	390	ILE	2.6
2	M	313	LEU	2.6
2	O	382	LEU	2.6
1	B	368[A]	HIS	2.6
2	O	377	PRO	2.6
2	P	354	ARG	2.6
2	P	727	PRO	2.6
2	O	333	ILE	2.6
2	O	410	PHE	2.6
1	B	415	ASN	2.6
2	O	639	THR	2.6
2	O	641	THR	2.6
1	C	306	GLY	2.5
2	O	342	MET	2.5
2	P	363	ASP	2.5
2	N	316	ILE	2.5
2	O	560	ILE	2.5
2	O	611	ILE	2.5
2	O	454	GLU	2.5
1	D	3	ARG	2.5
2	O	633	GLY	2.5
2	P	398	ASP	2.5
2	O	392	GLY	2.5
2	O	398	ASP	2.5
2	O	324	PHE	2.5
2	O	495	LEU	2.5
2	O	388	VAL	2.5
2	O	526	GLY	2.5
2	O	387	LEU	2.5
2	P	468	PHE	2.5
2	O	531	VAL	2.5
2	P	424	ASN	2.5
2	O	536	ALA	2.5
2	O	502	THR	2.5
2	O	592	MET	2.5
2	O	180	GLY	2.4
2	P	401	GLY	2.4
2	O	386	ILE	2.4
2	P	557	ILE	2.4
2	P	698	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
2	P	428	LEU	2.4
2	O	468	PHE	2.4
2	O	668	VAL	2.4
1	D	288	SER	2.4
2	O	317	LYS	2.4
1	A	539	ILE	2.4
2	N	728	ILE	2.4
2	O	413	TYR	2.4
2	P	387	LEU	2.4
2	O	456	PHE	2.4
1	A	417	PRO	2.4
2	N	389	ASP	2.4
2	O	346	ARG	2.4
2	O	151	TRP	2.4
2	P	427	TRP	2.4
2	M	174	LYS	2.4
2	M	238	GLU	2.4
2	M	607	GLU	2.4
2	O	522	PRO	2.4
1	D	414	SER	2.4
2	O	405	ARG	2.4
2	O	501	ASP	2.4
2	O	451	LYS	2.4
2	O	458	ALA	2.4
2	O	418	TRP	2.4
2	O	586	THR	2.3
2	O	697	ASP	2.3
2	O	442	LYS	2.3
1	A	538	GLU	2.3
2	O	351	GLU	2.3
2	O	457	PRO	2.3
2	O	724	THR	2.3
2	O	320	LEU	2.3
2	P	647	ILE	2.3
2	P	486	LYS	2.3
2	O	339	TYR	2.3
2	O	391	TYR	2.3
2	P	477	TYR	2.3
2	N	367	GLU	2.3
1	C	305	ALA	2.3
2	O	219	ALA	2.3
2	O	348	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
2	O	547	PRO	2.3
2	O	660	ALA	2.3
1	A	306	GLY	2.3
2	O	579	LEU	2.3
2	P	485	TYR	2.3
2	O	727	PRO	2.3
2	P	393	ARG	2.3
2	O	314	THR	2.3
2	O	555	GLY	2.3
2	O	431	SER	2.3
2	N	704	ILE	2.3
2	O	678	LEU	2.3
2	P	551	ILE	2.3
2	M	670	MET	2.3
2	O	229	PHE	2.3
2	O	341	GLU	2.3
2	O	406	ARG	2.3
2	O	488	ARG	2.3
2	O	327	ALA	2.2
2	P	355	THR	2.2
2	P	724	THR	2.2
2	O	159	ILE	2.2
2	O	380	SER	2.2
2	O	453	LYS	2.2
2	P	352	LEU	2.2
2	P	380	SER	2.2
2	P	718	LYS	2.2
2	O	491	ARG	2.2
2	O	524	ARG	2.2
2	O	421	GLY	2.2
2	P	729	MET	2.2
2	O	538	ALA	2.2
1	A	20	MET	2.2
2	O	541	GLU	2.2
2	P	17	GLU	2.2
2	O	503	PHE	2.2
2	O	335	LYS	2.2
2	M	338	MET	2.2
2	O	2	THR	2.2
2	O	650	THR	2.2
2	O	562	GLY	2.2
2	M	367	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
2	O	350	PHE	2.2
1	D	537	ILE	2.1
2	O	714	TYR	2.1
2	P	376	ILE	2.1
2	O	484	LYS	2.1
2	O	623	PRO	2.1
2	O	574	ALA	2.1
2	O	330	GLY	2.1
2	O	646	GLY	2.1
2	O	301	ILE	2.1
1	D	418	VAL	2.1
2	O	412	ASN	2.1
2	O	401	GLY	2.1
1	B	15	GLU	2.1
1	B	539	ILE	2.1
2	P	186	CYS	2.1
1	C	417	PRO	2.1
2	O	500	VAL	2.1
2	P	388	VAL	2.1
2	O	424	ASN	2.1
2	O	564	TRP	2.1
1	A	533	GLU	2.1
2	O	703	THR	2.1
2	P	366	ILE	2.1
2	P	369	ILE	2.1
2	P	699	ILE	2.1
2	O	374	ASP	2.1
2	O	540	TYR	2.1
2	P	570	TYR	2.1
2	O	575	SER	2.1
2	P	357	SER	2.1
1	B	418	VAL	2.0
2	O	644	PHE	2.0
1	A	415	ASN	2.0
1	D	471	ASN	2.0
1	D	382	GLY	2.0
2	O	369	ILE	2.0
2	P	386	ILE	2.0
2	P	425	ILE	2.0
2	O	585	TYR	2.0
1	D	293	GLN	2.0
2	O	396	GLN	2.0

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Mol	Chain	Res	Type	RSRZ
2	O	508	LEU	2.0
2	O	530	ALA	2.0
2	O	323	ASN	2.0
2	O	379	GLY	2.0
2	P	467	ILE	2.0
2	O	490	ASP	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.

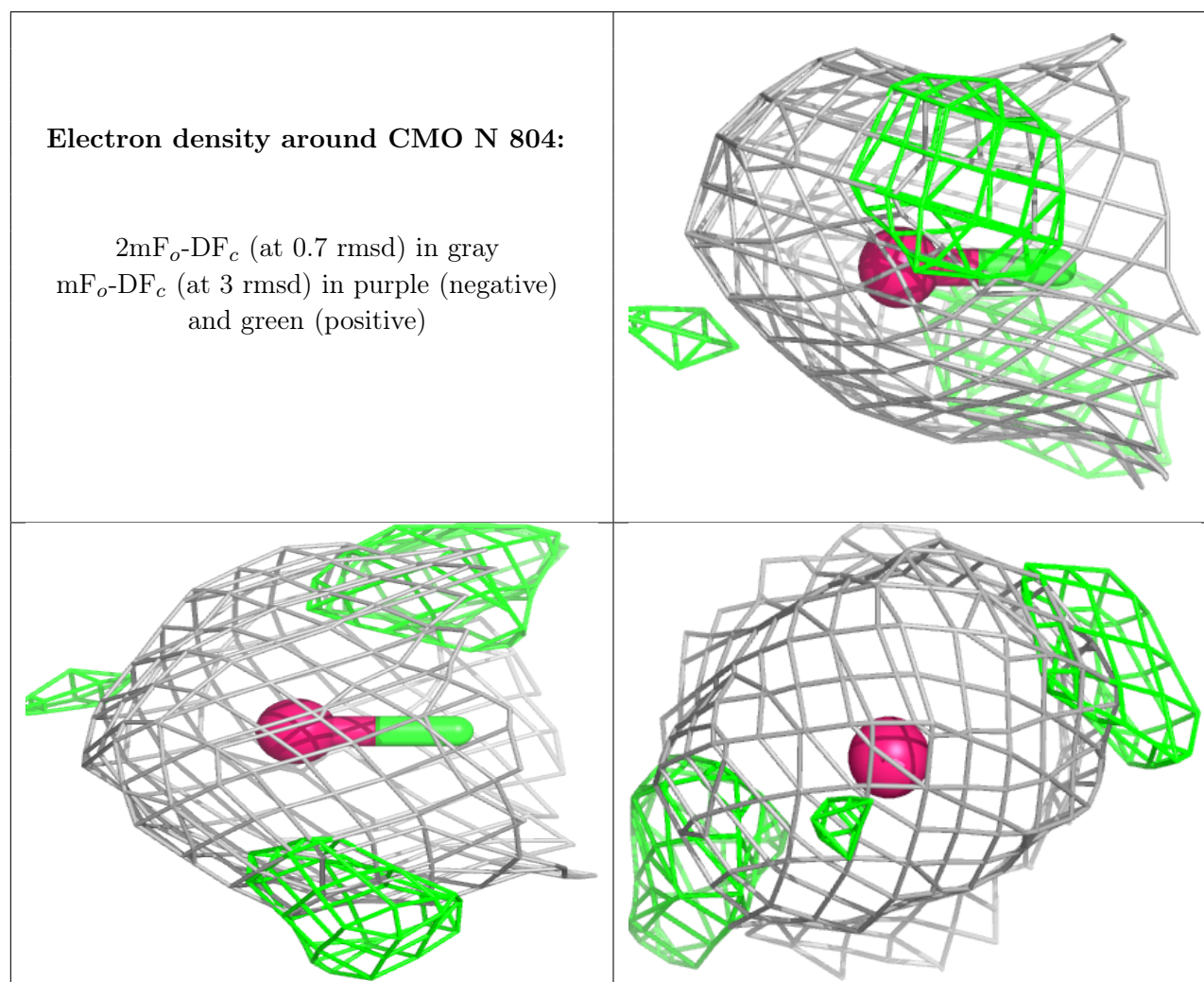
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	PG5	C	704	12/12	0.84	0.17	44,58,71,73	0
5	PGE	A	704	10/10	0.88	0.16	19,42,52,54	0
8	CMO	N	804	2/2	0.89	0.13	40,40,40,42	0
8	CMO	M	804	2/2	0.90	0.18	43,43,43,44	0
8	CMO	P	804	2/2	0.94	0.12	61,61,61,63	0
3	SF4	O	900	8/8	0.96	0.05	58,74,80,88	0
4	XCC	D	702	9/9	0.97	0.05	50,58,74,84	0
4	XCC	B	702	9/9	0.97	0.05	41,53,70,82	0
4	XCC	C	703	9/9	0.97	0.05	45,58,73,104	0
9	MG	N	805	1/1	0.97	0.04	43,43,43,43	0
7	NI	M	802	1/1	0.98	0.04	48,48,48,48	0
7	NI	O	901	1/1	0.98	0.07	75,75,75,75	0
4	XCC	A	703	9/9	0.98	0.04	43,52,63,69	0
3	SF4	C	702	8/8	0.99	0.03	35,38,41,43	0
3	SF4	D	701	8/8	0.99	0.04	36,38,44,45	0
3	SF4	M	801	8/8	0.99	0.03	28,30,34,36	0
3	SF4	N	801	8/8	0.99	0.04	27,32,33,35	0

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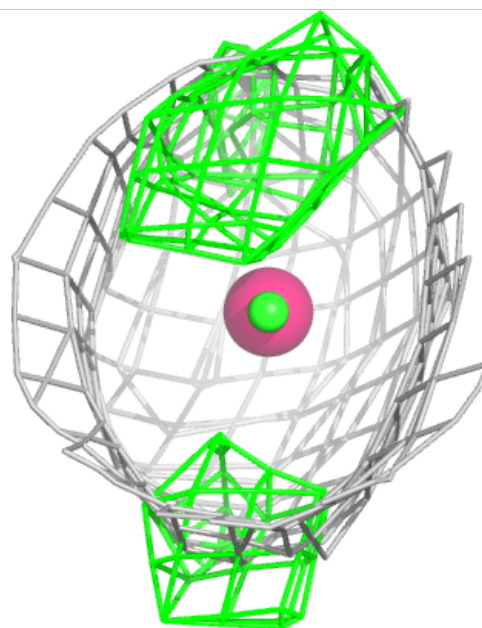
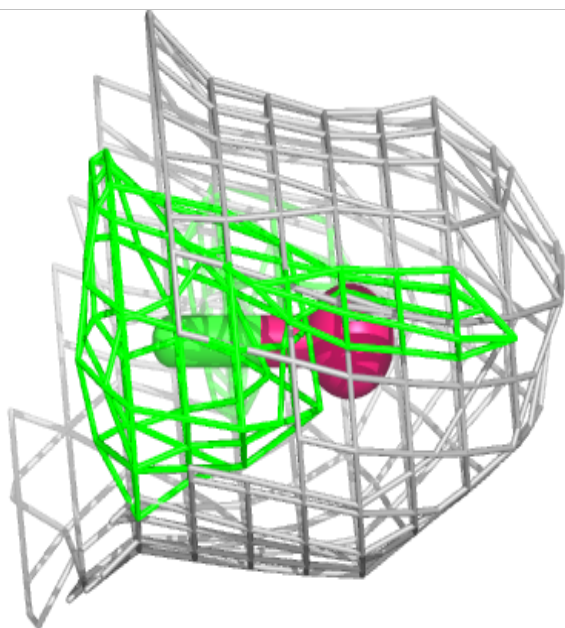
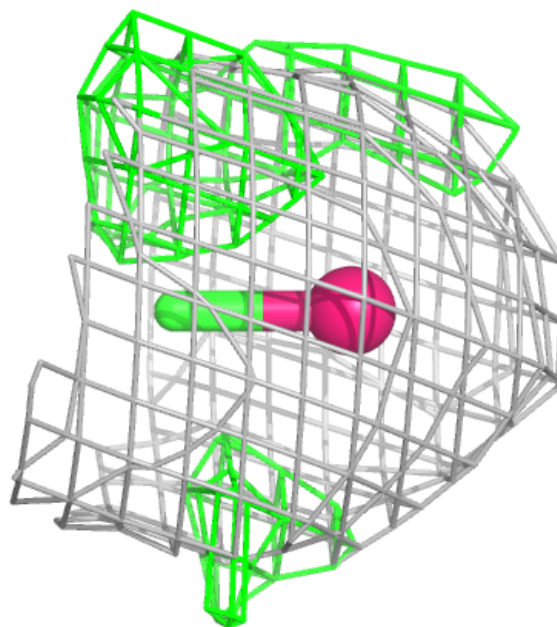
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	NI	M	803	1/1	0.99	0.03	31,31,31,31	0
7	NI	N	802	1/1	0.99	0.03	43,43,43,43	0
7	NI	N	803	1/1	0.99	0.04	31,31,31,31	0
3	SF4	A	701	8/8	0.99	0.02	21,26,27,29	0
7	NI	O	902	1/1	0.99	0.04	90,90,90,90	0
7	NI	P	802	1/1	0.99	0.03	61,61,61,61	0
7	NI	P	803	1/1	0.99	0.03	44,44,44,44	0
3	SF4	P	801	8/8	0.99	0.03	45,49,53,55	0
3	SF4	A	702	8/8	0.99	0.04	24,26,28,31	0
3	SF4	B	701	8/8	0.99	0.03	24,27,32,34	0
3	SF4	C	701	8/8	0.99	0.04	35,40,42,44	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.



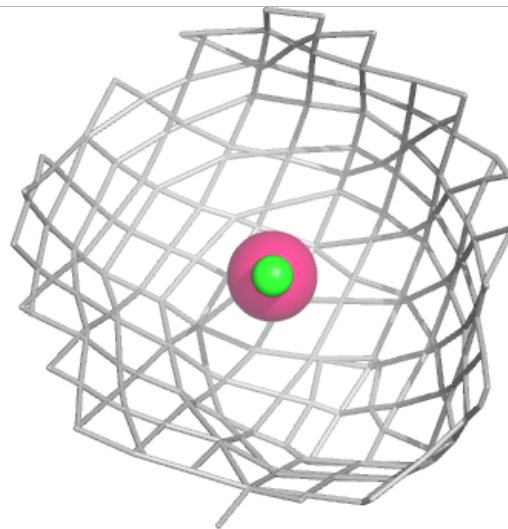
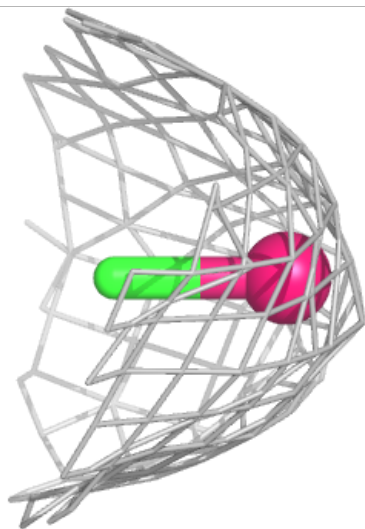
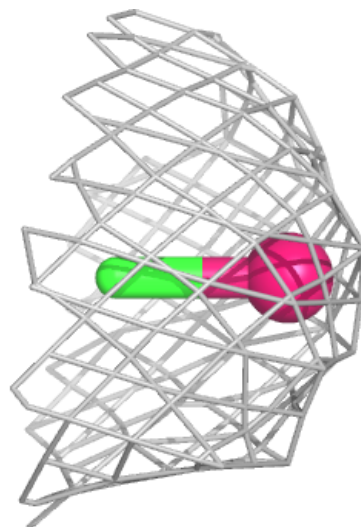
Electron density around CMO M 804:

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



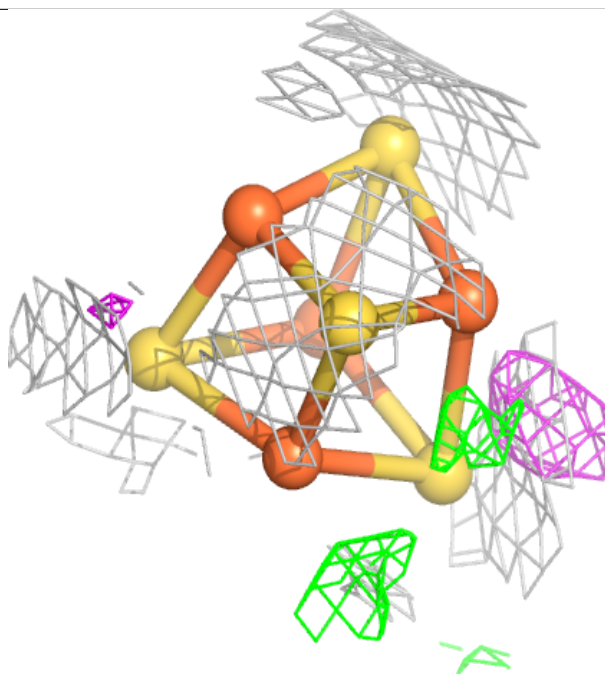
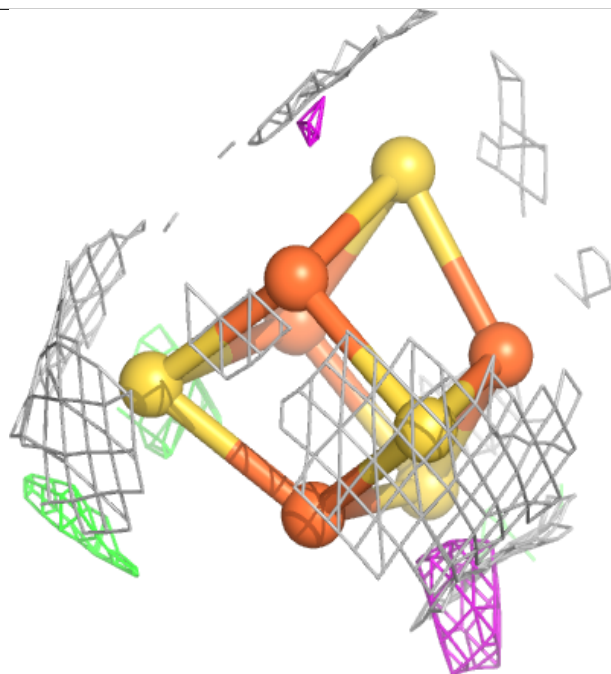
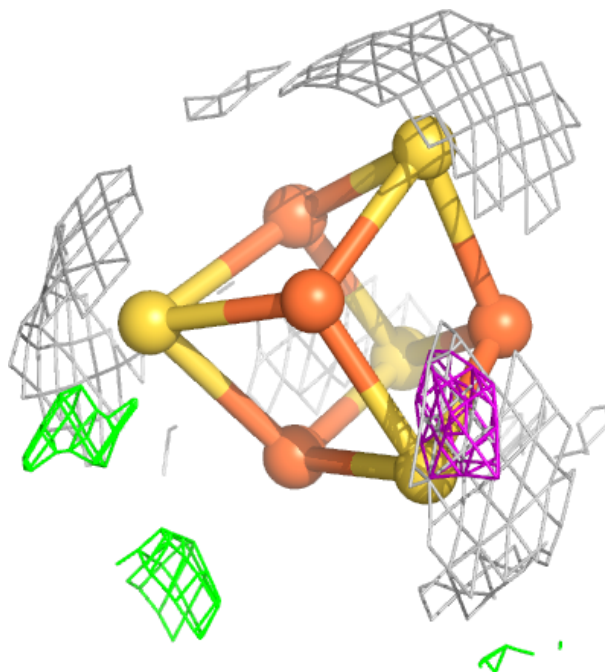
Electron density around CMO P 804:

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



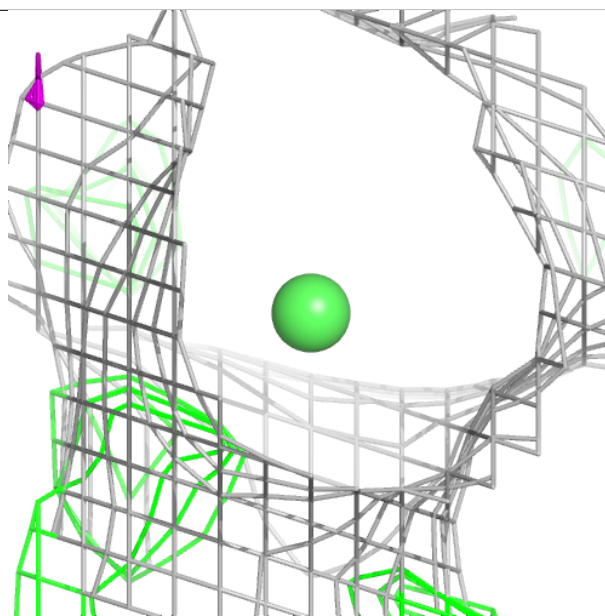
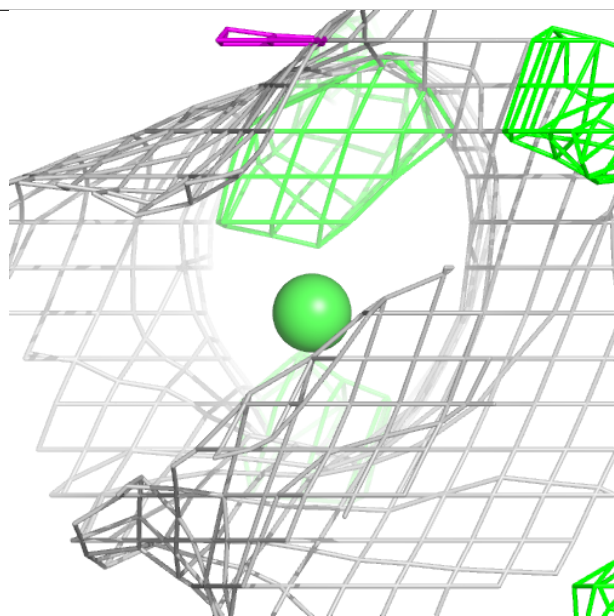
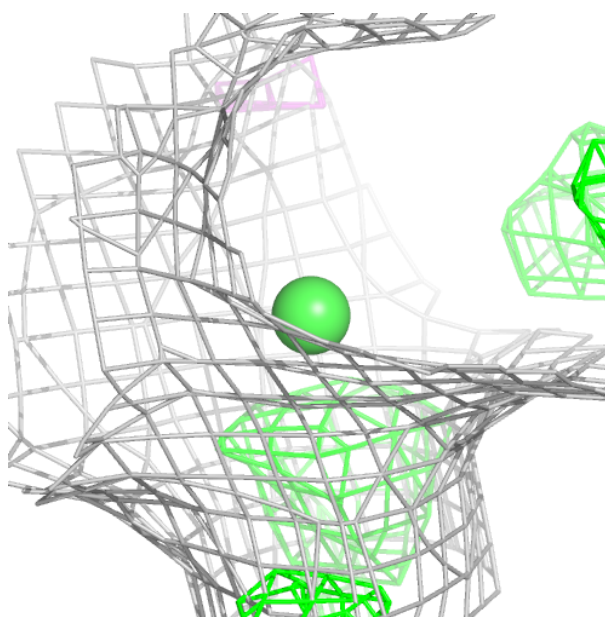
Electron density around SF4 O 900:

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and green (positive)



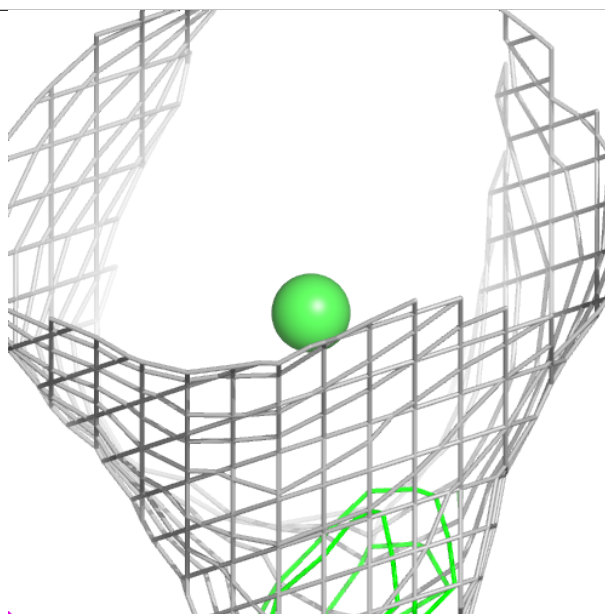
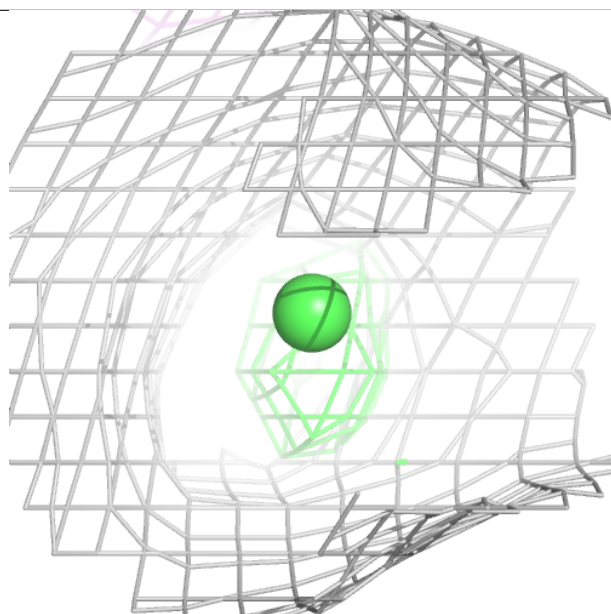
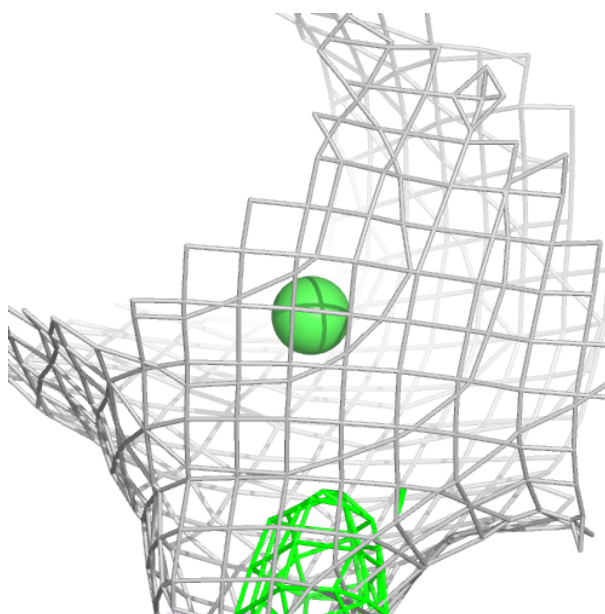
Electron density around NI M 802:

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



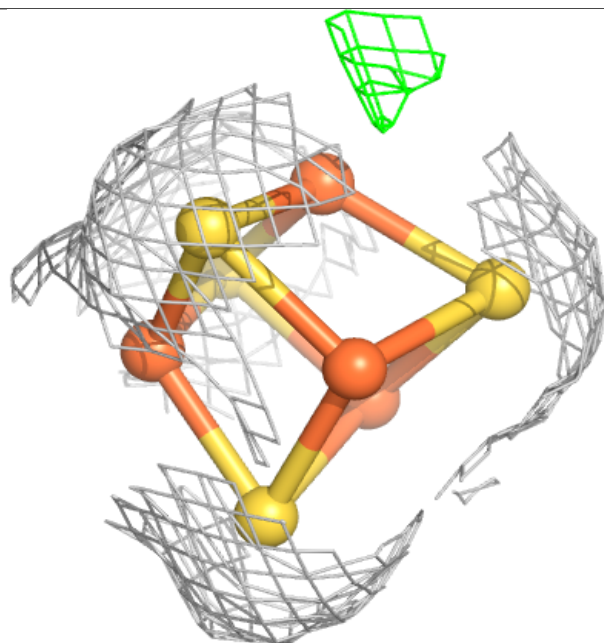
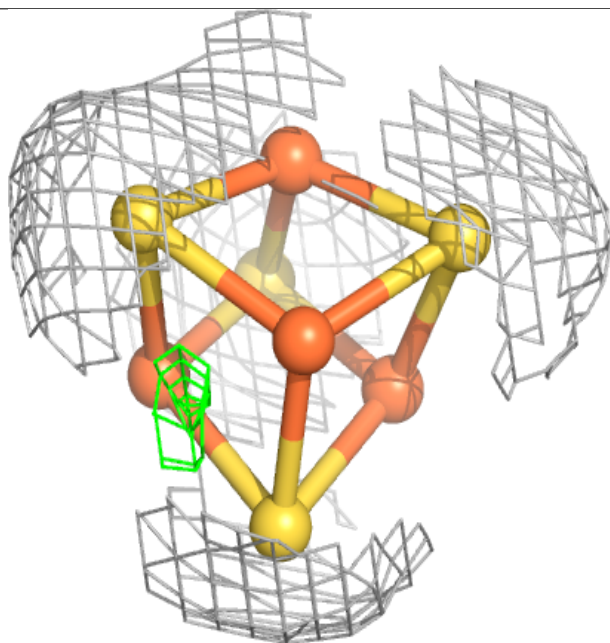
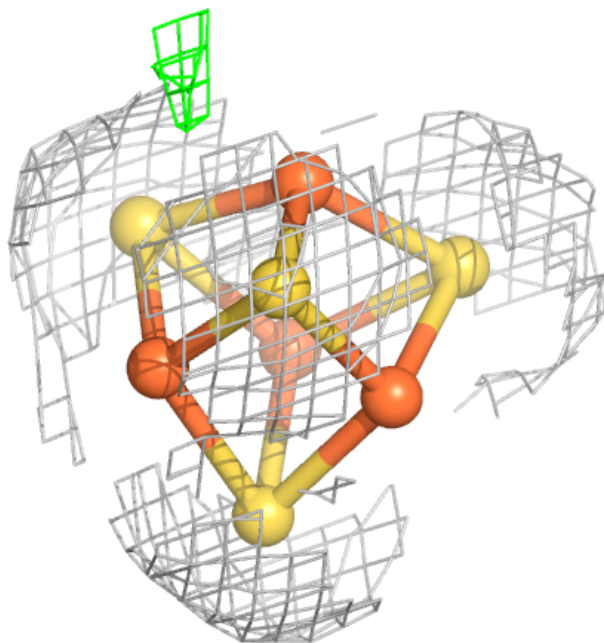
Electron density around NI O 901:

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and green (positive)



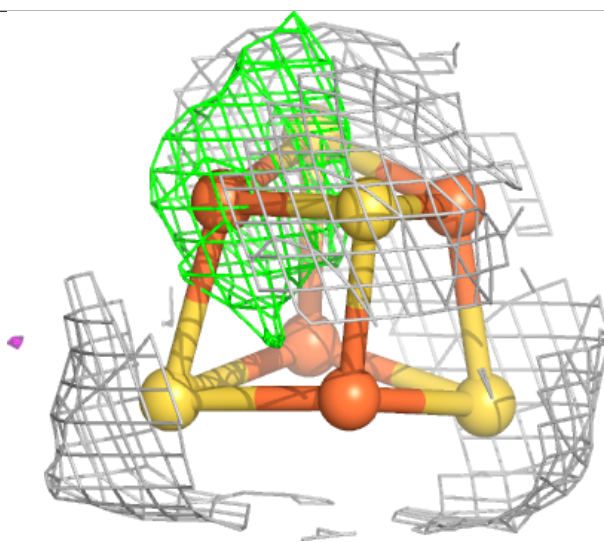
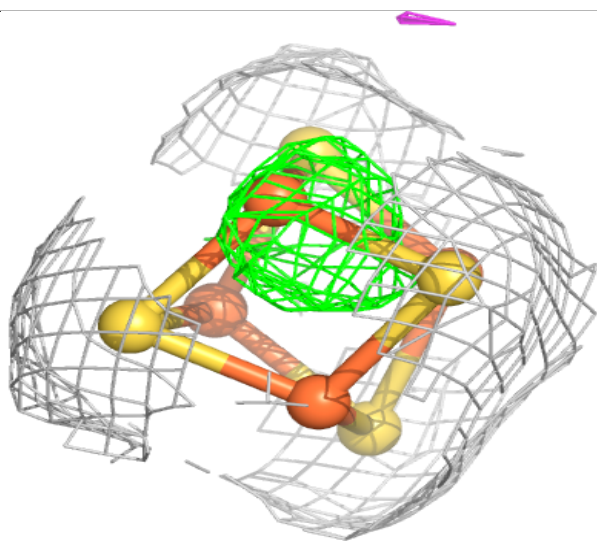
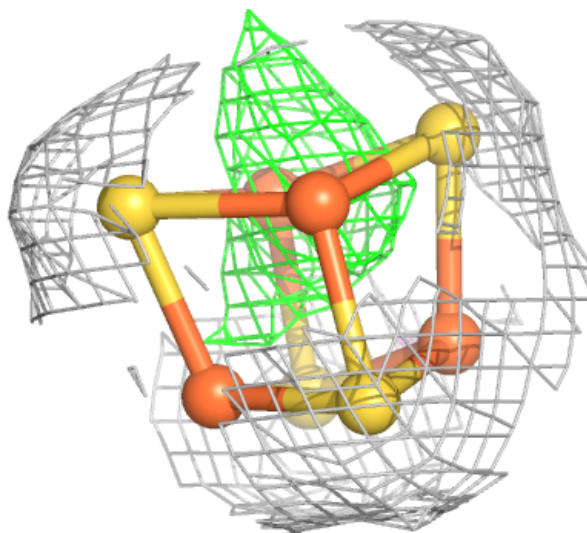
Electron density around SF4 C 702:

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



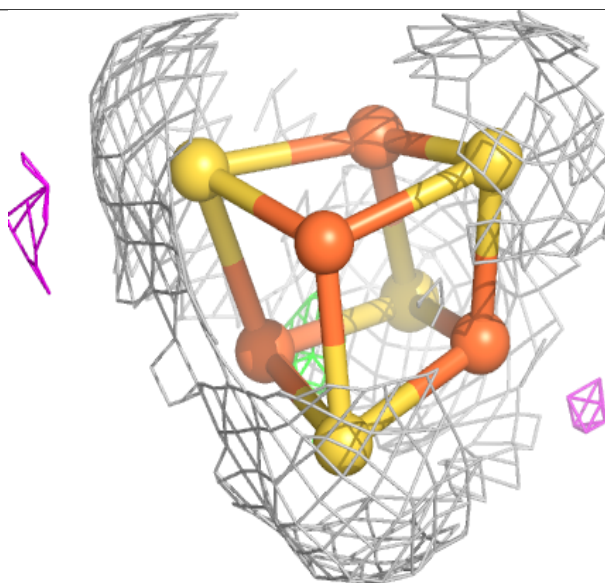
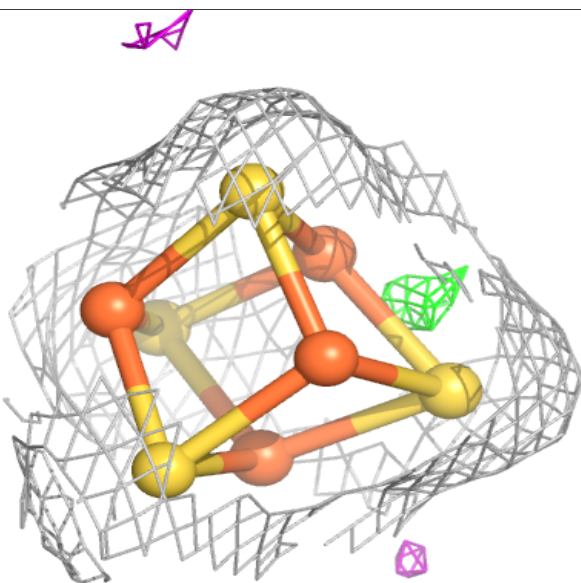
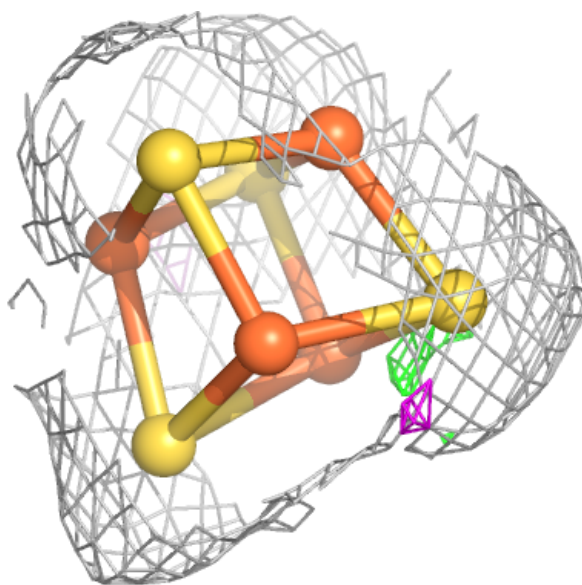
Electron density around SF4 D 701:

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and green (positive)



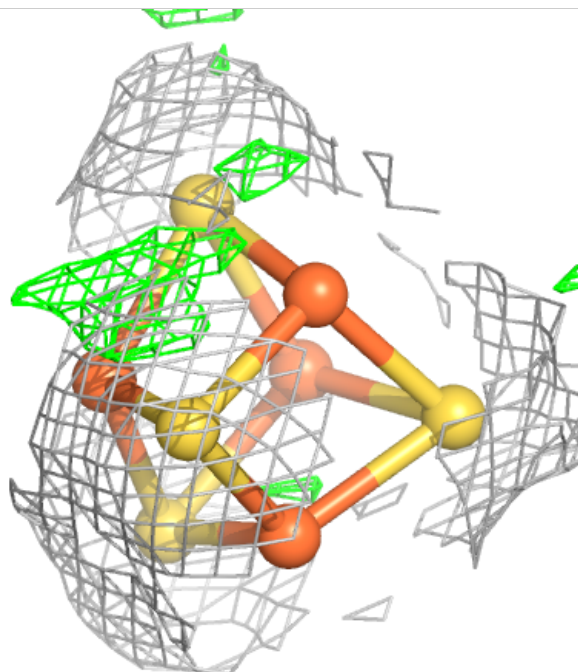
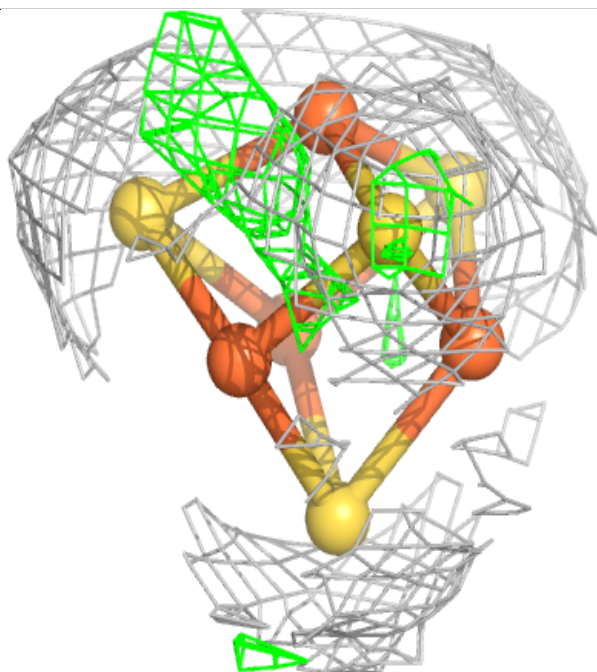
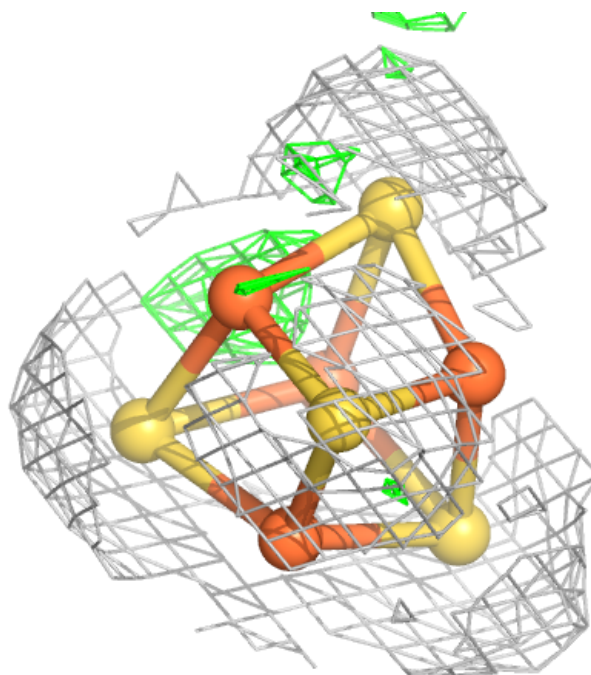
Electron density around SF4 M 801:

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



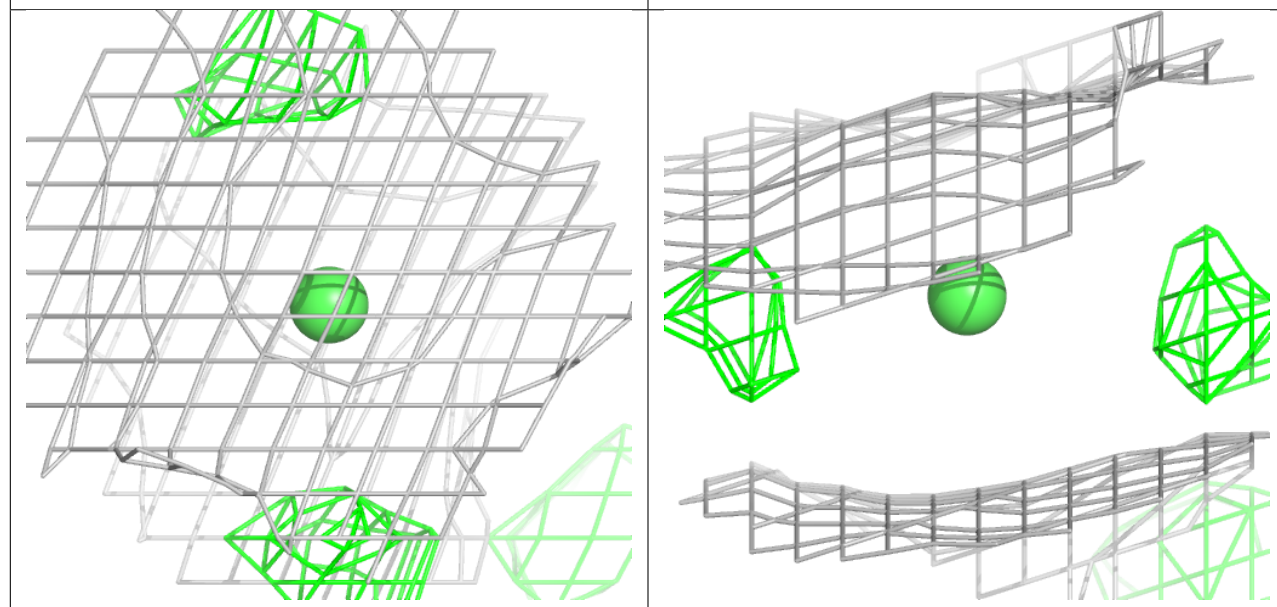
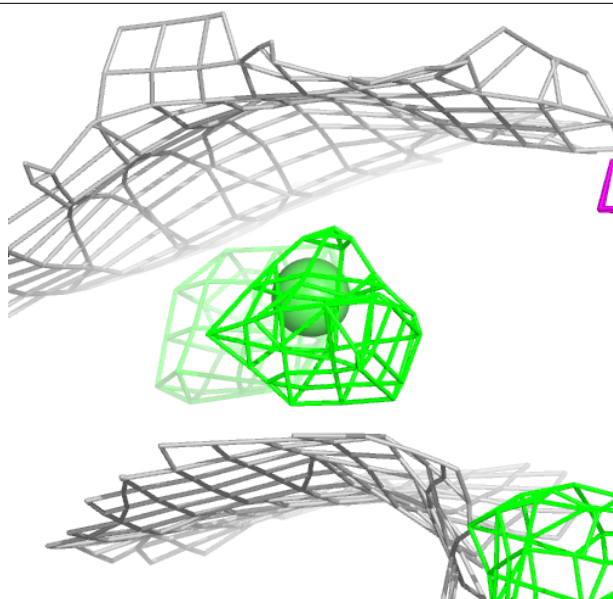
Electron density around SF4 N 801:

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



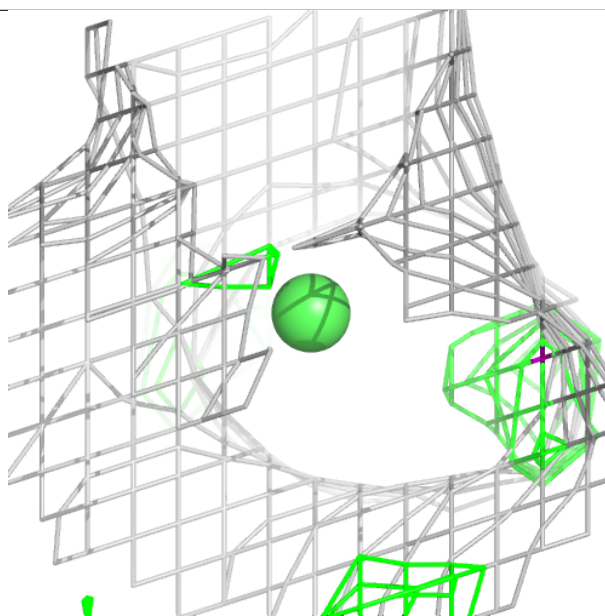
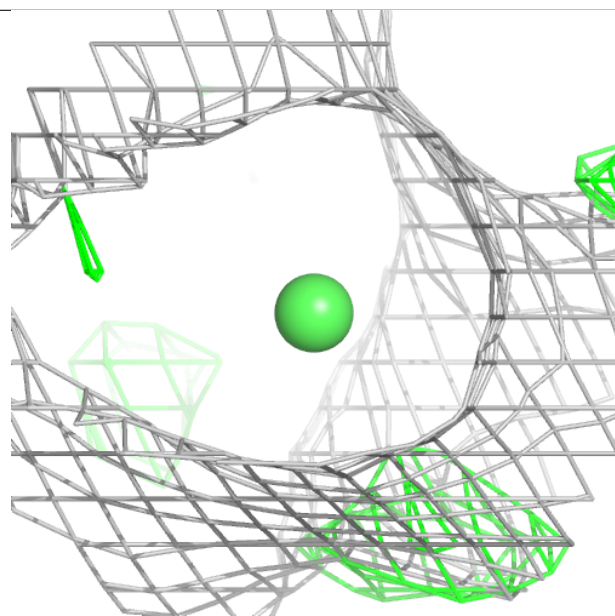
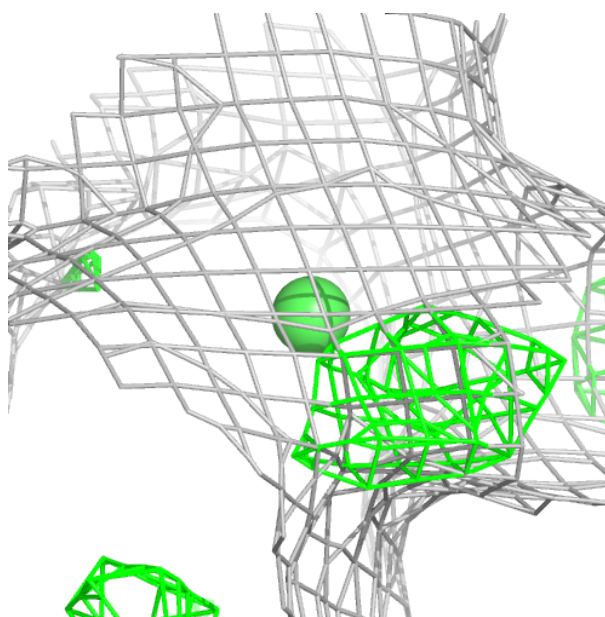
Electron density around NI M 803:

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



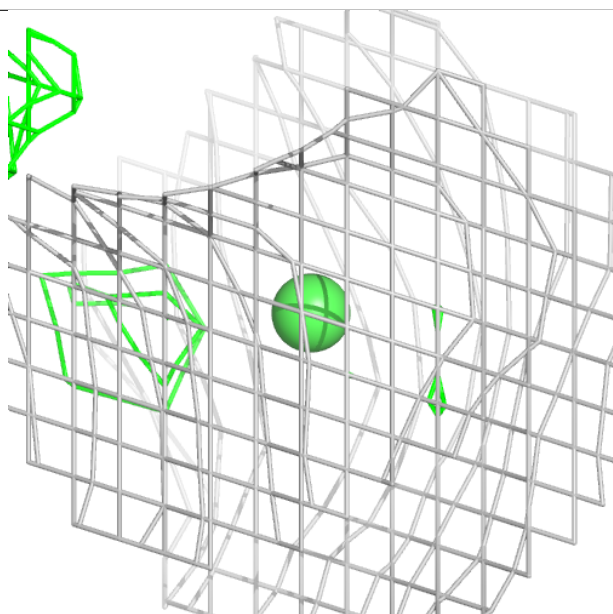
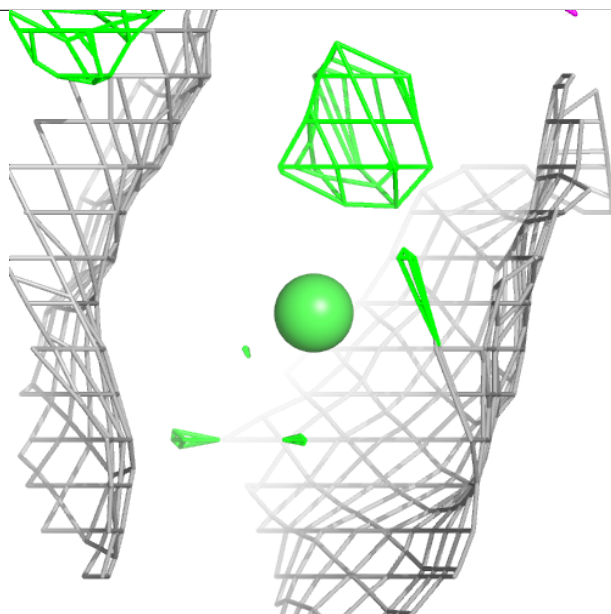
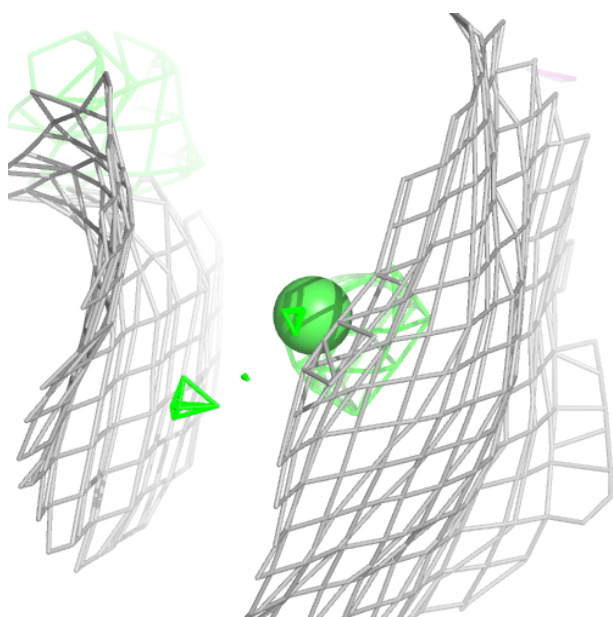
Electron density around NI N 802:

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and green (positive)



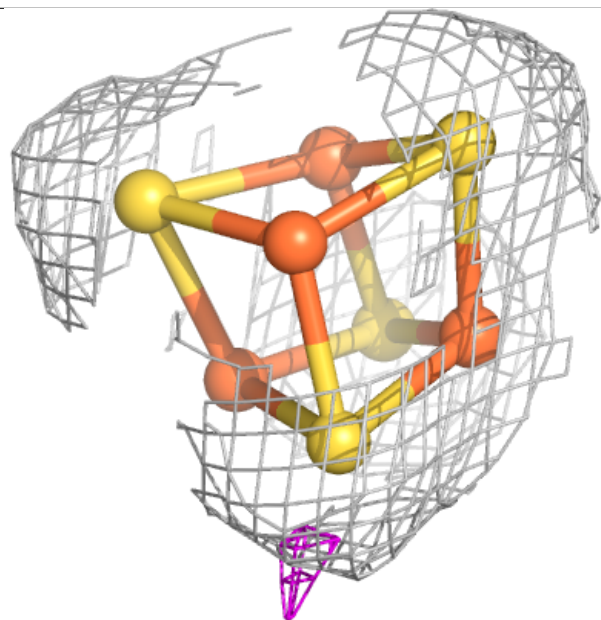
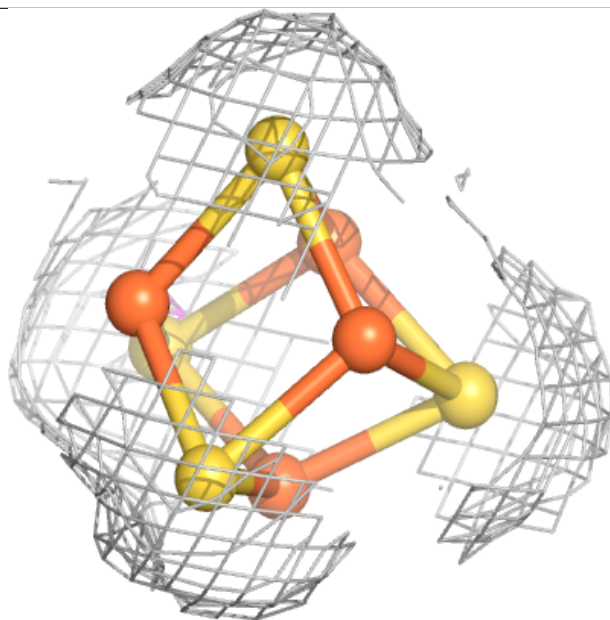
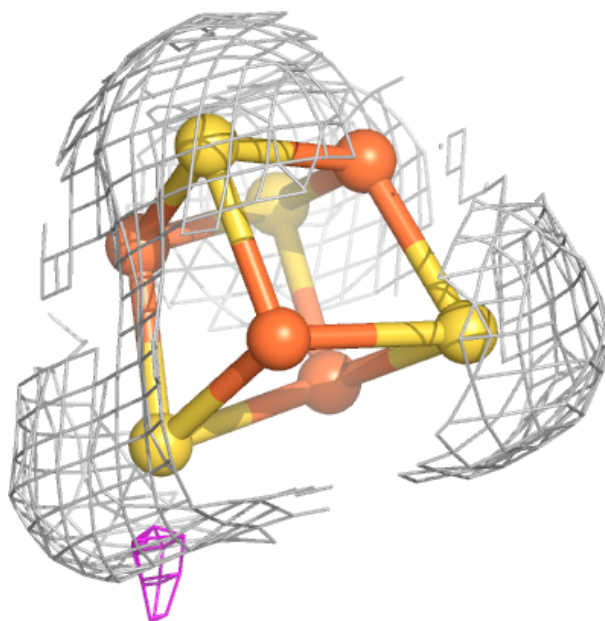
Electron density around NI N 803:

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and green (positive)



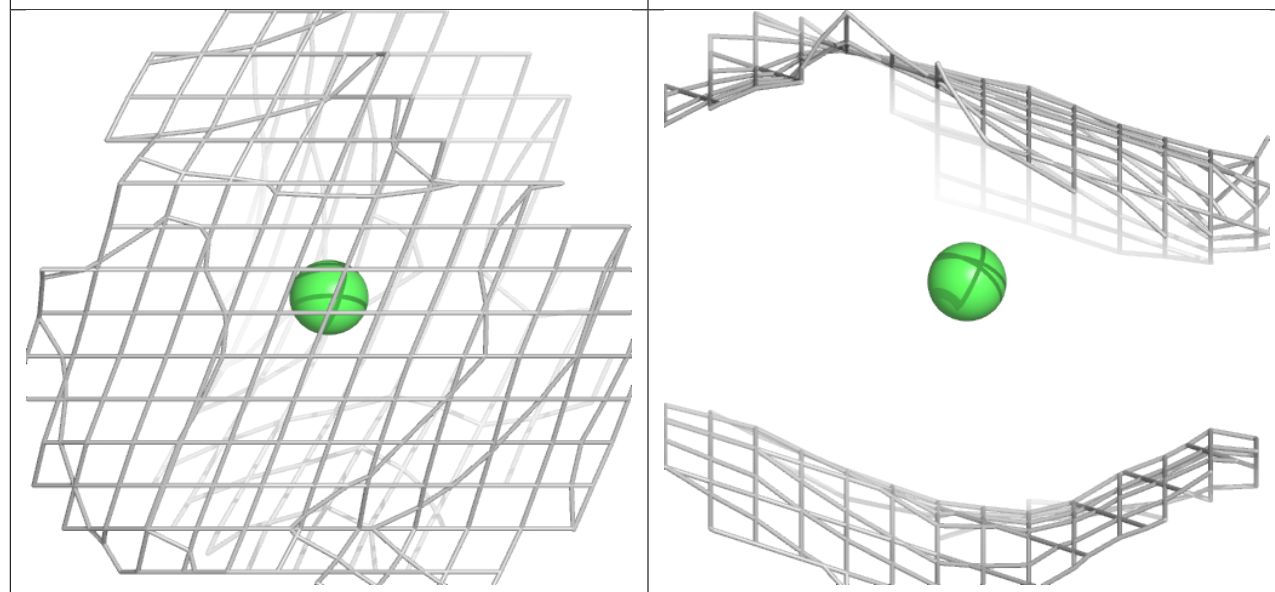
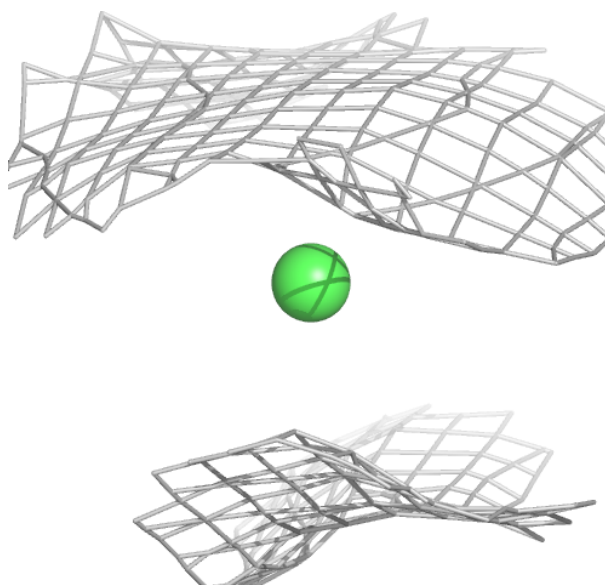
Electron density around SF4 A 701:

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



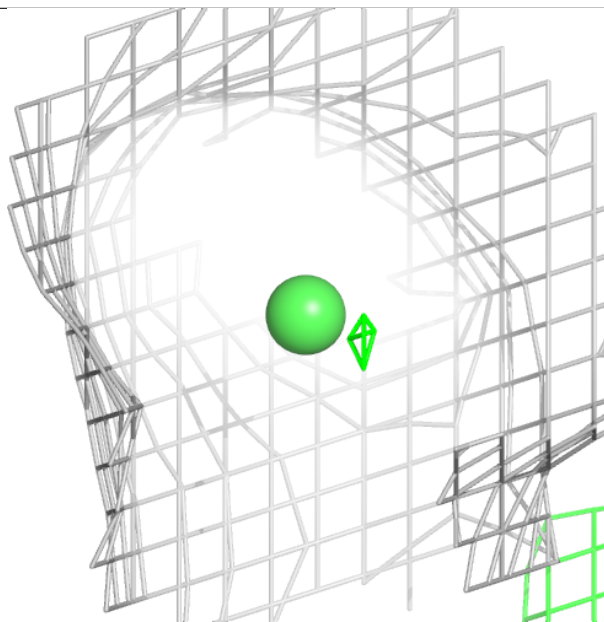
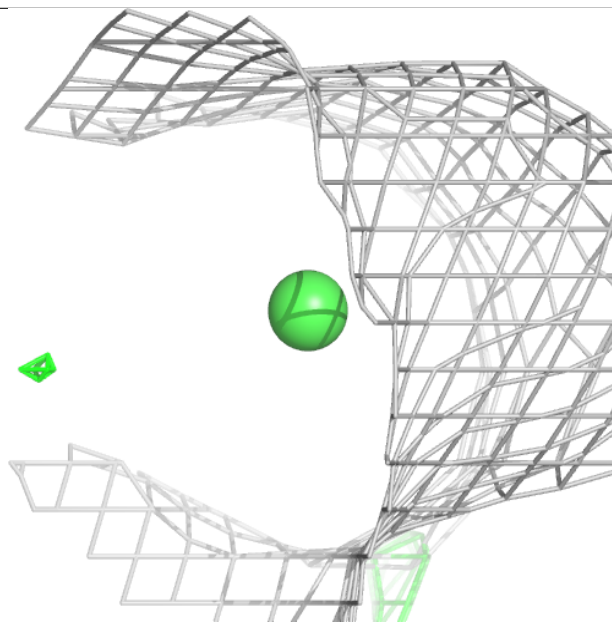
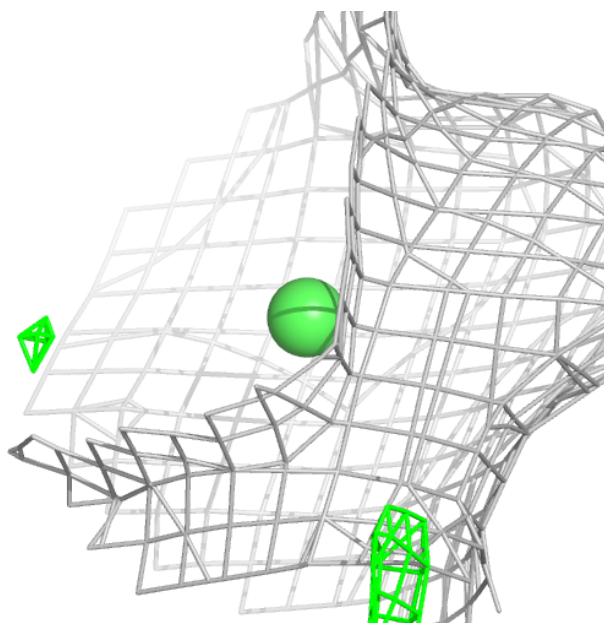
Electron density around NI O 902:

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



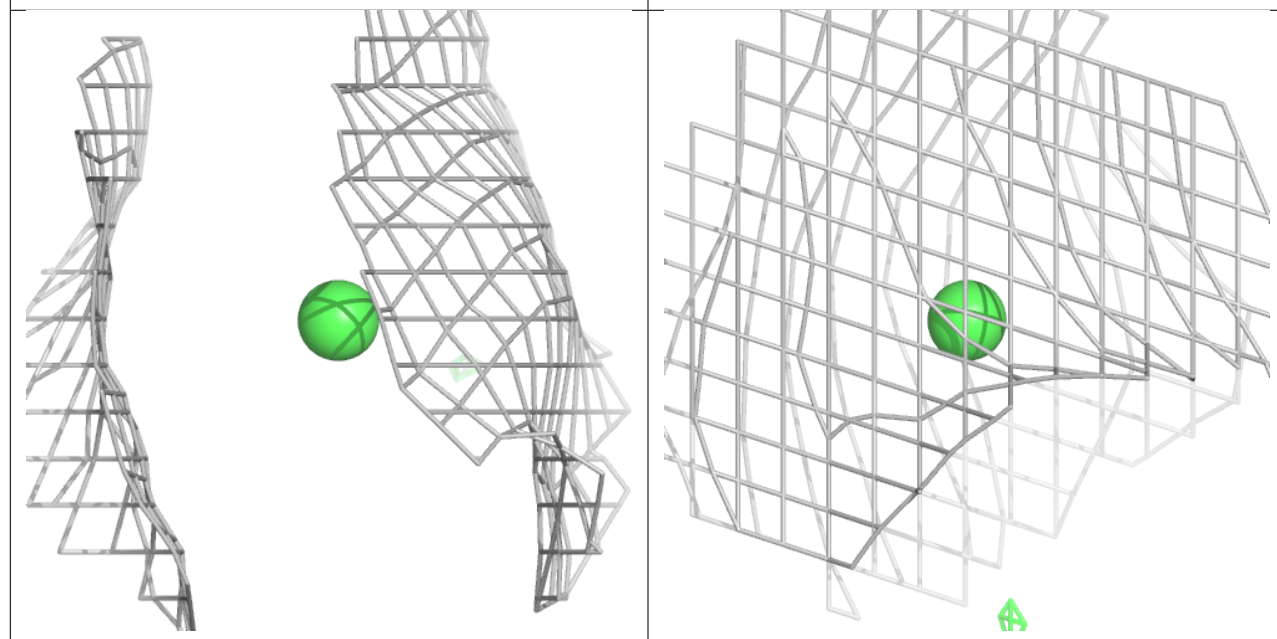
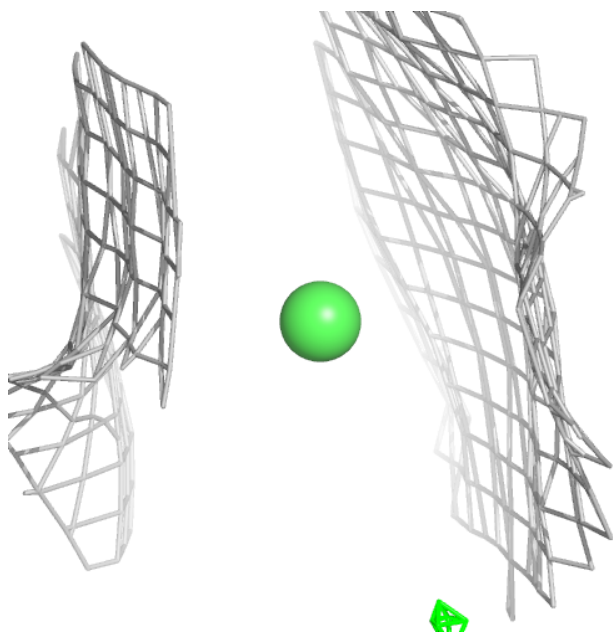
Electron density around NI P 802:

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and green (positive)



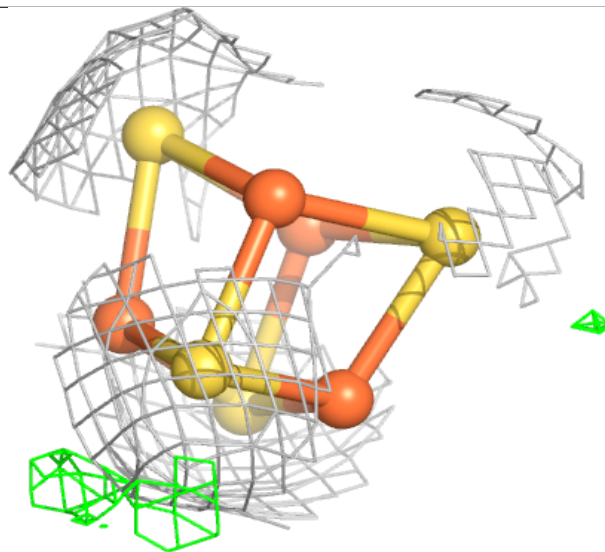
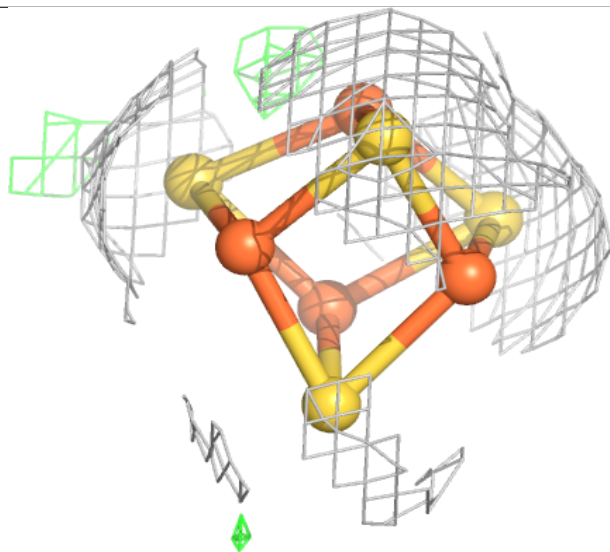
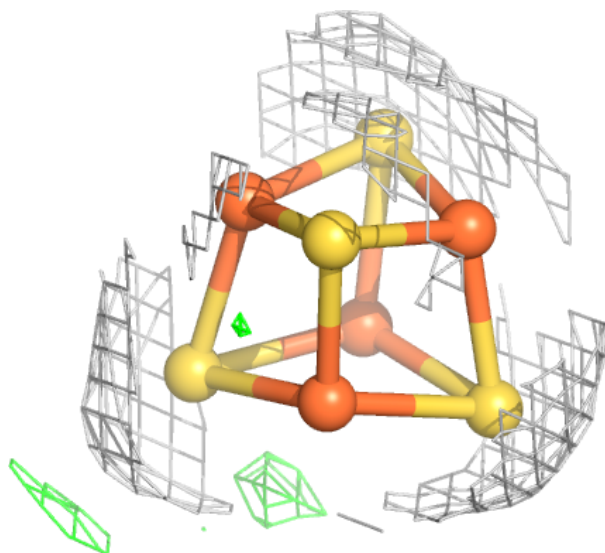
Electron density around NI P 803:

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and green (positive)



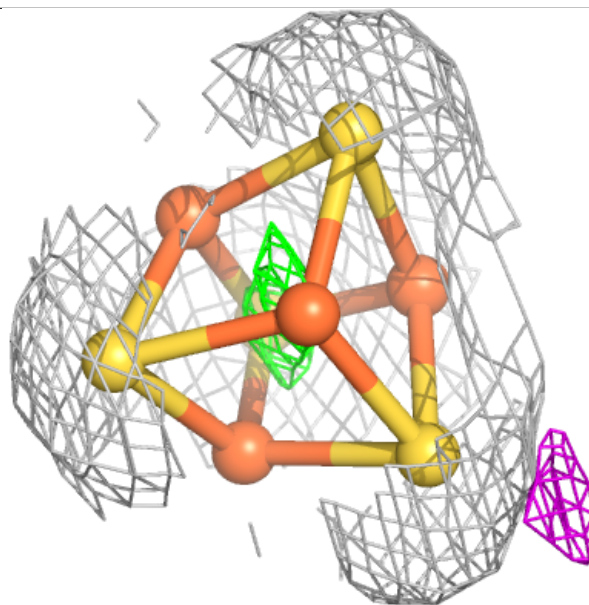
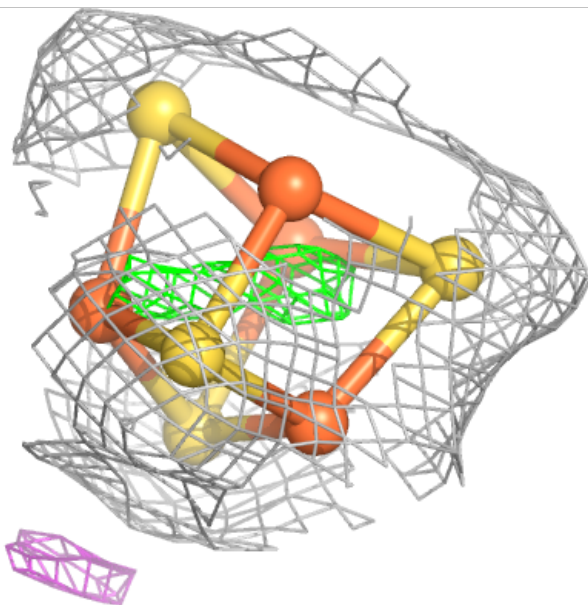
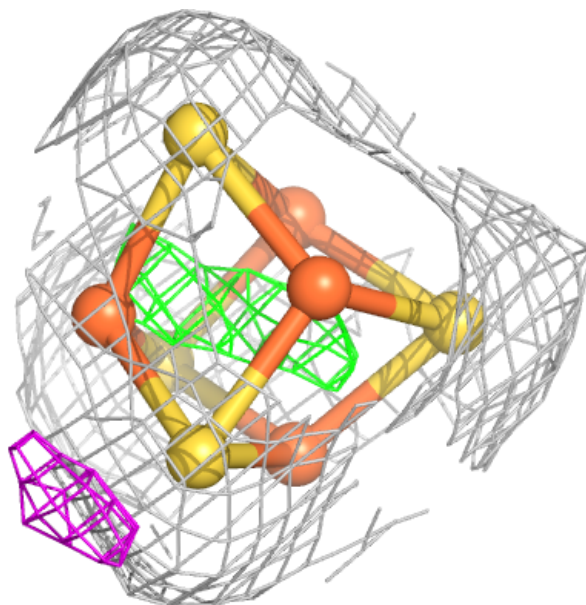
Electron density around SF4 P 801:

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and green (positive)



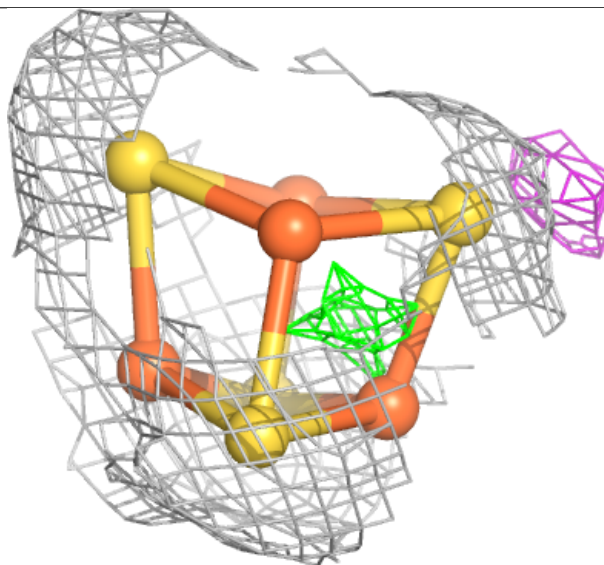
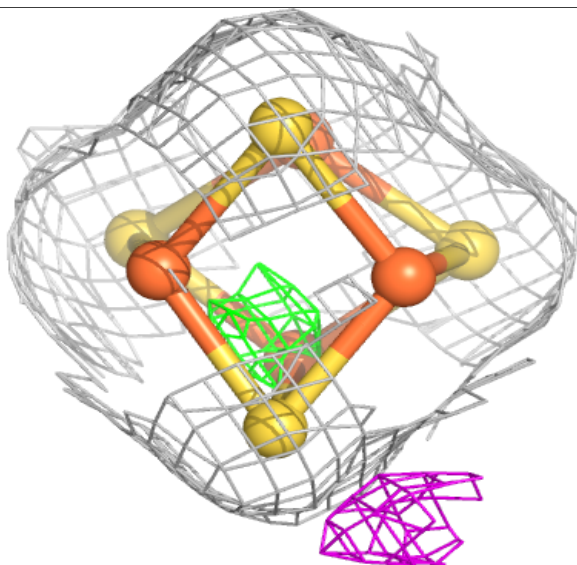
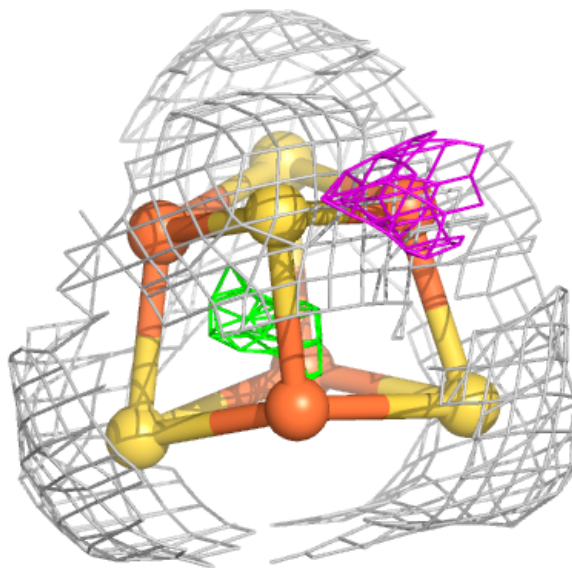
Electron density around SF4 A 702:

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



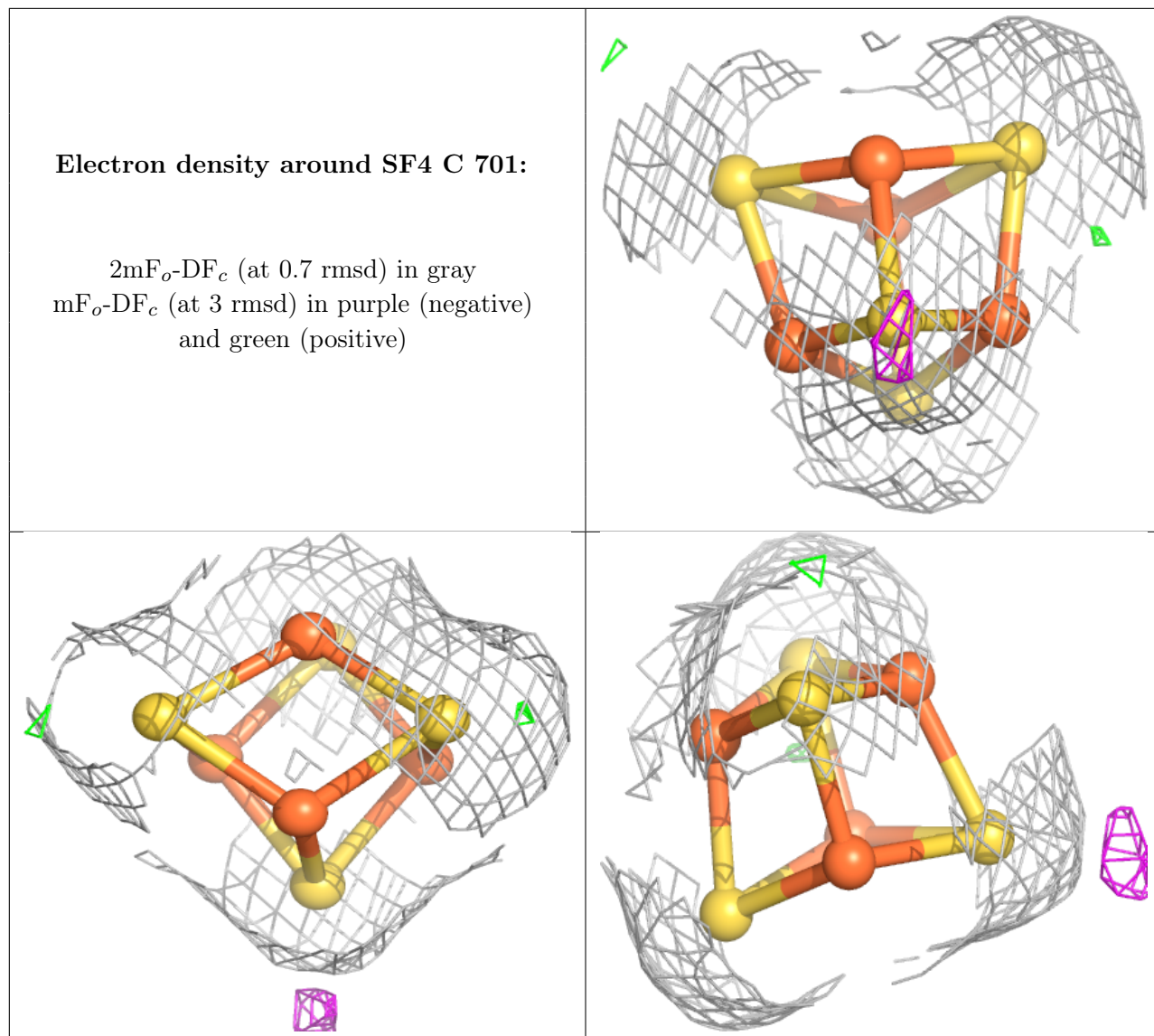
Electron density around SF4 B 701:

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



Electron density around SF4 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.