



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 11:46 AM UTC

PDB ID : 8OQP / pdb_00008oqp
Title : Structure of Mycobacterium tuberculosis beta-oxidation trifunctional enzyme
in complex with Fragment-M-76
Authors : Dalwani, S.; Wierenga, R.K.; Venkatesan, R.
Deposited on : 2023-04-12
Resolution : 2.18 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

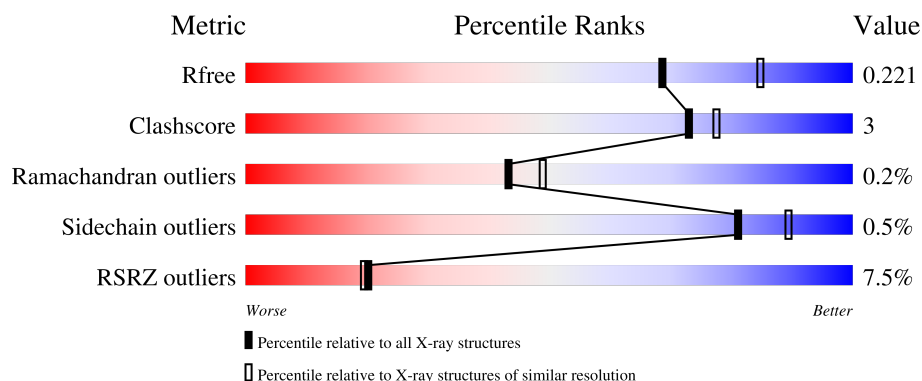
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.18 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	8975 (2.20-2.16)
Clashscore	190562	9786 (2.20-2.16)
Ramachandran outliers	187476	9664 (2.20-2.16)
Sidechain outliers	187428	9664 (2.20-2.16)
RSRZ outliers	180081	8979 (2.20-2.16)

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	736	7% 92% 6%
1	B	736	6% 92% 7%
2	C	403	8% 91% 8%
2	D	403	10% 91% 9%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 17630 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called 3-hydroxyacyl-CoA dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	724	Total	C	N	O	S	0	0	0
			5387	3410	925	1031	21			
1	B	726	Total	C	N	O	S	0	0	0
			5403	3419	929	1034	21			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	MET	-	initiating methionine	UNP O53872
A	-14	GLY	-	expression tag	UNP O53872
A	-13	SER	-	expression tag	UNP O53872
A	-12	SER	-	expression tag	UNP O53872
A	-11	HIS	-	expression tag	UNP O53872
A	-10	HIS	-	expression tag	UNP O53872
A	-9	HIS	-	expression tag	UNP O53872
A	-8	HIS	-	expression tag	UNP O53872
A	-7	HIS	-	expression tag	UNP O53872
A	-6	HIS	-	expression tag	UNP O53872
A	-5	SER	-	expression tag	UNP O53872
A	-4	GLN	-	expression tag	UNP O53872
A	-3	ASP	-	expression tag	UNP O53872
A	-2	PRO	-	expression tag	UNP O53872
A	-1	ASN	-	expression tag	UNP O53872
A	0	SER	-	expression tag	UNP O53872
B	-15	MET	-	initiating methionine	UNP O53872
B	-14	GLY	-	expression tag	UNP O53872
B	-13	SER	-	expression tag	UNP O53872
B	-12	SER	-	expression tag	UNP O53872
B	-11	HIS	-	expression tag	UNP O53872
B	-10	HIS	-	expression tag	UNP O53872
B	-9	HIS	-	expression tag	UNP O53872
B	-8	HIS	-	expression tag	UNP O53872
B	-7	HIS	-	expression tag	UNP O53872

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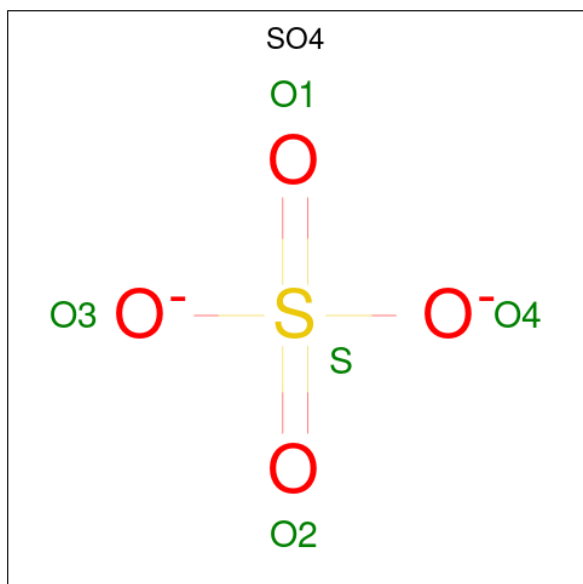
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	HIS	-	expression tag	UNP O53872
B	-5	SER	-	expression tag	UNP O53872
B	-4	GLN	-	expression tag	UNP O53872
B	-3	ASP	-	expression tag	UNP O53872
B	-2	PRO	-	expression tag	UNP O53872
B	-1	ASN	-	expression tag	UNP O53872
B	0	SER	-	expression tag	UNP O53872

- Molecule 2 is a protein called Putative acyltransferase Rv0859.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	402	Total	C	N	O	S	0	0	0
			2966	1853	525	573	15			
2	D	402	Total	C	N	O	S	0	0	0
			2966	1853	525	573	15			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

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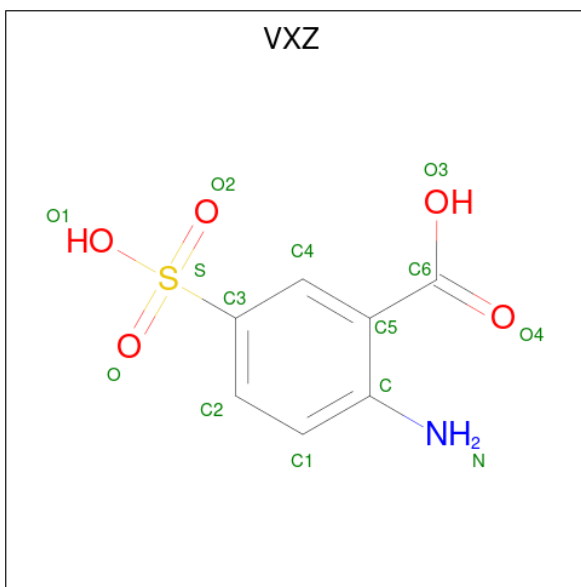
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is 2-azanyl-5-sulfo-benzoic acid (CCD ID: VXZ) (formula: C₇H₇NO₅S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			14	7	1	5	1		
5	A	1	Total	C	N	O	S	0	0
			14	7	1	5	1		
5	A	1	Total	C	N	O	S	0	0
			14	7	1	5	1		
5	A	1	Total	C	N	O	S	0	0
			14	7	1	5	1		
5	A	1	Total	C	N	O	S	0	0
			14	7	1	5	1		
5	B	1	Total	C	N	O	S	0	0
			14	7	1	5	1		
5	B	1	Total	C	N	O	S	0	0
			14	7	1	5	1		
5	B	1	Total	C	N	O	S	0	0
			14	7	1	5	1		
5	B	1	Total	C	N	O	S	0	0
			14	7	1	5	1		
5	C	1	Total	C	N	O	S	0	0
			14	7	1	5	1		
5	C	1	Total	C	N	O	S	0	0
			14	7	1	5	1		
5	D	1	Total	C	N	O	S	0	0
			14	7	1	5	1		
5	D	1	Total	C	N	O	S	0	0
			14	7	1	5	1		
5	D	1	Total	C	N	O	S	0	0
			14	7	1	5	1		

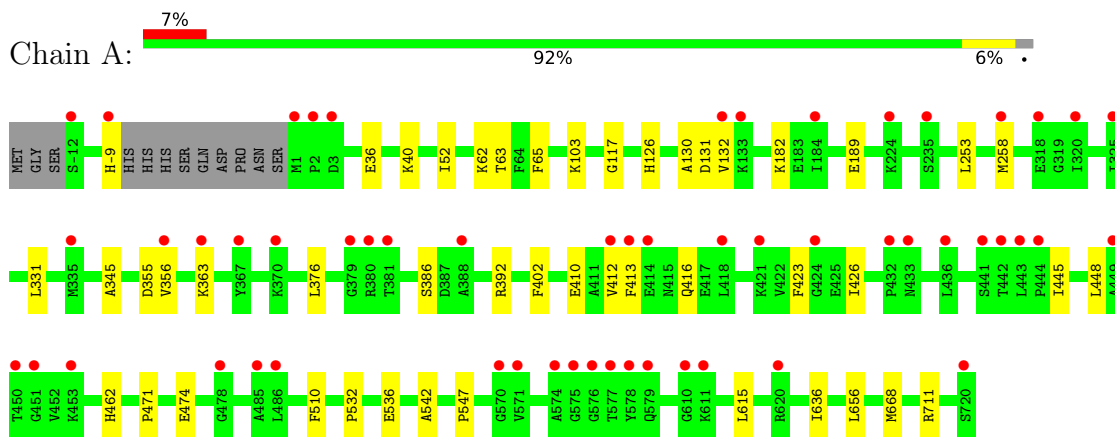
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	152	Total 152	O 152	0	0
6	B	173	Total 173	O 173	0	0
6	C	143	Total 143	O 143	0	0
6	D	107	Total 107	O 107	0	0

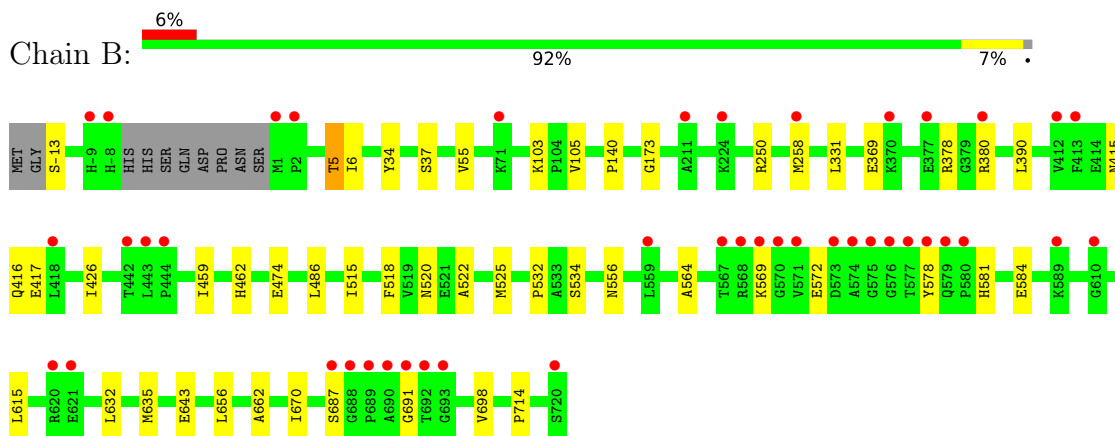
3 Residue-property plots

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

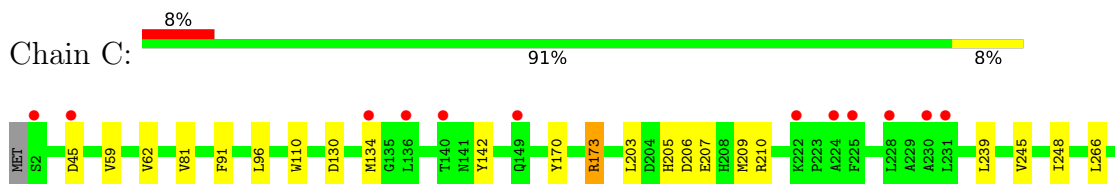
- Molecule 1: 3-hydroxyacyl-CoA dehydrogenase

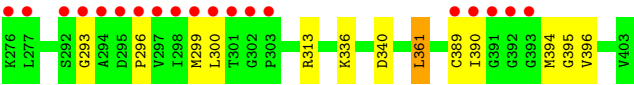


- Molecule 1: 3-hydroxyacyl-CoA dehydrogenase

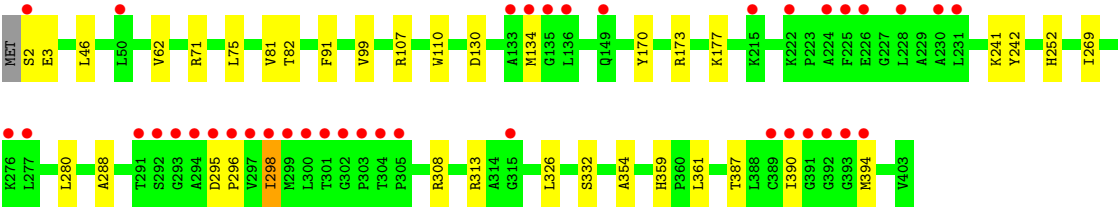
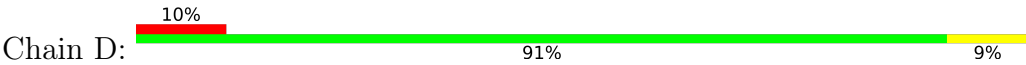


- Molecule 2: Putative acyltransferase Rv0859





● Molecule 2: Putative acyltransferase Rv0859



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	250.94Å 132.80Å 119.29Å 90.00° 110.46° 90.00°	Depositor
Resolution (Å)	69.73 – 2.18 69.73 – 2.18	Depositor EDS
% Data completeness (in resolution range)	76.3 (69.73-2.18) 76.4 (69.73-2.18)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 2.18Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.196 , 0.224 0.194 , 0.221	Depositor DCC
R_{free} test set	7114 reflections (3.74%)	wwPDB-VP
Wilson B-factor (Å ²)	46.5	Xtriage
Anisotropy	0.004	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17630	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.87% 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: GOL, SO4, VXZ

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	0.10	0/5487	0.29	0/7425
1	B	0.11	0/5504	0.28	0/7448
2	C	0.11	0/3011	0.32	0/4077
2	D	0.11	0/3011	0.32	0/4077
All	All	0.11	0/17013	0.30	0/23027

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [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	5387	0	5436	29	0
1	B	5403	0	5448	29	0
2	C	2966	0	2986	29	0
2	D	2966	0	2986	29	0
3	A	25	0	0	0	0
3	B	25	0	0	0	0
3	C	25	0	0	0	0
3	D	20	0	0	0	0
4	A	18	0	24	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	12	0	16	0	0
4	D	12	0	16	0	0
5	A	70	0	0	1	0
5	B	56	0	0	1	0
5	C	28	0	0	0	0
5	D	42	0	0	1	0
6	A	152	0	0	2	0
6	B	173	0	0	1	0
6	C	143	0	0	2	0
6	D	107	0	0	0	0
All	All	17630	0	16912	109	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 3.

All (109) 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:C:206:ASP:HB2	2:C:209:MET:HE2	1.73	0.70
1:B:378:ARG:HD2	1:B:380:ARG:HH12	1.60	0.67
2:C:62:VAL:HG12	2:D:62:VAL:HG12	1.77	0.66
1:A:412:VAL:HG23	1:A:413:PHE:H	1.60	0.66
1:A:103:LYS:NZ	6:A:901:HOH:O	2.30	0.64
2:C:336:LYS:NZ	2:C:340:ASP:OD2	2.33	0.62
2:C:96:LEU:HD23	2:C:396:VAL:HG13	1.82	0.61
1:A:711:ARG:HB3	4:A:804:GOL:H2	1.81	0.60
2:C:110:TRP:CD1	2:D:313:ARG:HD3	2.36	0.60
1:A:182:LYS:NZ	6:A:904:HOH:O	2.36	0.59
1:A:462:HIS:HB3	1:A:474:GLU:HB3	1.85	0.59
2:C:296:PRO:HD3	2:D:81:VAL:HG21	1.84	0.59
2:C:313:ARG:HD3	2:D:110:TRP:CD1	2.38	0.59
1:B:532:PRO:HB2	1:B:615:LEU:HD13	1.87	0.57
1:B:5:THR:HG22	1:B:37:SER:HB2	1.85	0.57
1:A:345:ALA:O	1:A:392:ARG:NH1	2.38	0.57
1:A:376:LEU:HD11	1:A:386:SER:HB2	1.86	0.57
1:B:250:ARG:NH1	2:C:142:TYR:O	2.38	0.56
1:B:369:GLU:HG2	1:B:390:LEU:HD13	1.87	0.56
1:B:632:LEU:HA	1:B:635:MET:HE3	1.88	0.55
2:D:390:ILE:HB	2:D:394:MET:HB2	1.89	0.55
2:C:81:VAL:HG11	2:D:296:PRO:HD3	1.87	0.55
2:D:252:HIS:HE1	2:D:332:SER:H	1.54	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:46:LEU:HD12	2:D:280:LEU:HD21	1.89	0.55
1:B:462:HIS:HB3	1:B:474:GLU:HB3	1.89	0.54
1:B:331:LEU:HD11	1:B:426:ILE:HD13	1.90	0.53
2:D:241:LYS:HE2	2:D:295:ASP:OD1	2.09	0.52
1:A:416:GLN:HG3	1:A:448:LEU:HD23	1.91	0.52
1:B:520:ASN:HB3	1:B:581:HIS:CE1	2.44	0.52
2:C:239:LEU:HD21	2:C:248:ILE:HG13	1.91	0.52
1:B:522:ALA:HA	1:B:525:MET:HE3	1.92	0.51
1:A:423:PHE:HA	1:A:426:ILE:HG22	1.92	0.51
2:C:299:MET:HB3	2:C:300:LEU:HD22	1.91	0.51
1:A:62:LYS:HG3	1:A:63:THR:HG23	1.93	0.51
2:D:62:VAL:HG11	2:D:130:ASP:HA	1.94	0.50
2:D:326:LEU:HD13	2:D:387:THR:HG23	1.93	0.50
2:D:99:VAL:HG13	2:D:269:ILE:HD11	1.93	0.49
1:A:532:PRO:HB2	1:A:615:LEU:HD13	1.94	0.49
2:C:91:PHE:HB2	2:C:390:ILE:HG23	1.94	0.49
2:D:242:TYR:OH	2:D:298:ILE:HB	2.13	0.48
1:B:534:SER:HB3	1:B:635:MET:HE1	1.94	0.48
1:B:698:VAL:HG13	1:B:714:PRO:HG3	1.94	0.48
1:A:445:ILE:HD13	1:A:448:LEU:HD12	1.96	0.48
1:B:5:THR:HG23	1:B:6:ILE:HG13	1.96	0.48
1:B:5:THR:HG22	1:B:37:SER:CB	2.43	0.48
2:C:293:GLY:H	2:D:82:THR:HG22	1.79	0.47
1:B:103:LYS:NZ	6:B:906:HOH:O	2.45	0.47
2:C:170:TYR:O	2:C:173:ARG:HG3	2.15	0.47
1:B:55:VAL:HB	1:B:105:VAL:HG22	1.97	0.47
2:C:62:VAL:HG11	2:C:130:ASP:HA	1.96	0.47
1:A:65:PHE:O	1:A:117:GLY:HA3	2.15	0.47
1:A:126:HIS:HE2	1:A:189:GLU:CD	2.20	0.47
2:C:390:ILE:HG21	2:C:394:MET:HE3	1.97	0.47
2:D:390:ILE:HG21	2:D:394:MET:HE3	1.97	0.47
1:A:510:PHE:CD1	1:A:656:LEU:HD11	2.50	0.47
1:A:253:LEU:HD13	1:A:258:MET:HB2	1.97	0.46
1:A:331:LEU:HB2	1:A:410:GLU:HA	1.98	0.46
2:C:45:ASP:N	2:C:45:ASP:OD1	2.45	0.46
1:A:711:ARG:HH11	4:A:804:GOL:H12	1.80	0.46
2:C:110:TRP:CZ2	2:D:288:ALA:HA	2.50	0.46
2:D:252:HIS:CE1	2:D:332:SER:H	2.32	0.46
1:B:415:ASN:OD1	1:B:417:GLU:HG2	2.16	0.46
1:B:569:LYS:HA	1:B:572:GLU:HG2	1.98	0.45
1:A:36:GLU:HG3	1:A:40:LYS:HE3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:ALA:O	1:A:132:VAL:N	2.47	0.45
1:B:459:ILE:HD11	1:B:486:LEU:HD12	1.99	0.45
1:B:632:LEU:HD12	1:B:635:MET:HE3	1.99	0.45
2:C:134:MET:SD	2:D:75:LEU:HD12	2.57	0.44
1:A:126:HIS:NE2	1:A:189:GLU:OE2	2.34	0.44
1:A:402:PHE:CB	1:A:426:ILE:HD11	2.47	0.44
1:B:564:ALA:HB1	1:B:584:GLU:OE2	2.17	0.44
2:D:173:ARG:NH2	2:D:177:LYS:HE3	2.32	0.44
1:A:363:LYS:HE2	1:A:363:LYS:HB3	1.82	0.43
2:C:205:HIS:HB2	6:C:734:HOH:O	2.18	0.43
2:C:59:VAL:HG21	2:C:361:LEU:HB3	2.01	0.43
1:A:536:GLU:CD	1:A:547:PRO:HB2	2.44	0.43
1:B:687:SER:HA	1:B:691:GLY:O	2.17	0.43
2:D:134:MET:HE3	2:D:134:MET:HB3	1.83	0.43
1:B:140:PRO:HA	1:B:173:GLY:HA3	2.01	0.43
1:A:471:PRO:HG2	1:A:668:MET:HB3	2.00	0.43
1:A:355:ASP:OD1	1:A:356:VAL:HG22	2.19	0.42
1:B:518:PHE:HB2	1:B:643:GLU:CD	2.44	0.42
2:C:203:LEU:HD11	2:C:207:GLU:HG3	2.01	0.42
2:D:390:ILE:HD12	2:D:394:MET:HB2	2.01	0.42
1:B:5:THR:HG21	1:B:34:TYR:HA	2.01	0.42
2:C:266:LEU:HD23	2:C:266:LEU:HA	1.88	0.42
2:C:390:ILE:HD12	2:C:394:MET:HB2	2.00	0.42
2:D:2:SER:HB3	2:D:107:ARG:O	2.20	0.42
2:D:170:TYR:O	2:D:173:ARG:HG2	2.20	0.42
1:B:515:ILE:HD12	1:B:670:ILE:HB	2.02	0.42
2:C:300:LEU:HD12	2:C:389:CYS:SG	2.59	0.42
1:A:376:LEU:HD11	1:A:386:SER:CB	2.48	0.42
5:A:807:VXZ:O4	5:A:807:VXZ:N	2.53	0.42
2:D:3:GLU:O	2:D:107:ARG:HG2	2.19	0.42
1:B:416:GLN:HE21	1:B:416:GLN:HB2	1.62	0.42
5:B:809:VXZ:O4	5:B:809:VXZ:N	2.53	0.41
2:C:210:ARG:NH1	6:C:609:HOH:O	2.52	0.41
1:A:542:ALA:HB2	1:A:636:ILE:HG23	2.03	0.41
5:D:507:VXZ:O4	5:D:507:VXZ:N	2.52	0.41
1:A:52:ILE:O	1:A:103:LYS:HD3	2.21	0.41
2:D:170:TYR:CD1	2:D:173:ARG:HD3	2.55	0.41
2:C:91:PHE:HB2	2:C:390:ILE:CG2	2.50	0.41
1:B:656:LEU:HD13	1:B:662:ALA:HB2	2.02	0.41
1:B:578:TYR:OH	1:B:584:GLU:OE2	2.35	0.41
2:D:91:PHE:HB2	2:D:390:ILE:CG2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:390:ILE:HD12	2:C:395:GLY:N	2.36	0.40
2:D:354:ALA:HB1	2:D:359:HIS:HB2	2.02	0.40
2:C:394:MET:HG2	2:D:71:ARG:CZ	2.51	0.40
2:D:91:PHE:HB2	2:D:390:ILE:HG23	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	720/736 (98%)	694 (96%)	25 (4%)	1 (0%)	48	55
1	B	722/736 (98%)	700 (97%)	21 (3%)	1 (0%)	48	55
2	C	400/403 (99%)	390 (98%)	9 (2%)	1 (0%)	36	39
2	D	400/403 (99%)	386 (96%)	13 (3%)	1 (0%)	36	39
All	All	2242/2278 (98%)	2170 (97%)	68 (3%)	4 (0%)	43	49

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	361	LEU
2	D	361	LEU
1	A	131	ASP
1	B	556	ASN

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	555/566 (98%)	554 (100%)	1 (0%)	87	94
1	B	557/566 (98%)	554 (100%)	3 (0%)	81	89
2	C	309/310 (100%)	307 (99%)	2 (1%)	78	87
2	D	309/310 (100%)	307 (99%)	2 (1%)	78	87
All	All	1730/1752 (99%)	1722 (100%)	8 (0%)	81	89

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-9	HIS
1	B	-13	SER
1	B	5	THR
1	B	258	MET
2	C	173	ARG
2	C	245	VAL
2	D	298	ILE
2	D	308	ARG

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

Mol	Chain	Res	Type
1	A	294	GLN
1	A	420	HIS
1	A	550	GLN
1	B	294	GLN
1	B	550	GLN
1	B	633	GLN
2	D	380	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

40 ligands are modelled in this entry.

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
5	VXZ	D	506	-	14,14,14	1.07	2 (14%)	21,21,21	0.88	2 (9%)
4	GOL	B	805	-	5,5,5	0.91	0	5,5,5	1.12	0
5	VXZ	B	809	-	14,14,14	1.22	3 (21%)	21,21,21	0.81	2 (9%)
5	VXZ	A	808	-	14,14,14	1.04	2 (14%)	21,21,21	0.90	2 (9%)
3	SO4	A	812	-	4,4,4	0.25	0	6,6,6	0.19	0
3	SO4	D	501	-	4,4,4	0.24	0	6,6,6	0.10	0
3	SO4	C	503	-	4,4,4	0.23	0	6,6,6	0.07	0
5	VXZ	B	806	-	14,14,14	1.07	2 (14%)	21,21,21	0.86	2 (9%)
3	SO4	C	507	-	4,4,4	0.23	0	6,6,6	0.15	0
4	GOL	A	806	-	5,5,5	0.87	0	5,5,5	1.12	0
3	SO4	A	813	-	4,4,4	0.24	0	6,6,6	0.11	0
3	SO4	C	501	-	4,4,4	0.23	0	6,6,6	0.09	0
3	SO4	B	802	-	4,4,4	0.23	0	6,6,6	0.08	0
3	SO4	D	503	-	4,4,4	0.23	0	6,6,6	0.09	0
4	GOL	D	505	-	5,5,5	0.89	0	5,5,5	1.05	0
5	VXZ	D	507	-	14,14,14	1.05	2 (14%)	21,21,21	0.86	2 (9%)
3	SO4	C	504	-	4,4,4	0.24	0	6,6,6	0.07	0
5	VXZ	A	807	-	14,14,14	1.06	2 (14%)	21,21,21	0.87	2 (9%)
5	VXZ	B	808	-	14,14,14	1.04	2 (14%)	21,21,21	0.90	2 (9%)
5	VXZ	A	811	-	14,14,14	1.23	3 (21%)	21,21,21	0.90	2 (9%)
5	VXZ	A	810	-	14,14,14	1.06	2 (14%)	21,21,21	0.88	2 (9%)
3	SO4	C	502	-	4,4,4	0.24	0	6,6,6	0.08	0
5	VXZ	C	506	-	14,14,14	1.07	2 (14%)	21,21,21	0.86	2 (9%)
4	GOL	A	805	-	5,5,5	0.88	0	5,5,5	1.11	0
3	SO4	B	810	-	4,4,4	0.25	0	6,6,6	0.16	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	B	801	-	4,4,4	0.22	0	6,6,6	0.11	0
4	GOL	D	504	-	5,5,5	0.94	0	5,5,5	1.14	0
3	SO4	A	801	-	4,4,4	0.25	0	6,6,6	0.10	0
3	SO4	A	803	-	4,4,4	0.24	0	6,6,6	0.08	0
3	SO4	B	803	-	4,4,4	0.24	0	6,6,6	0.07	0
4	GOL	A	804	-	5,5,5	0.91	0	5,5,5	1.08	0
5	VXZ	B	807	-	14,14,14	1.05	2 (14%)	21,21,21	0.89	2 (9%)
5	VXZ	A	809	-	14,14,14	1.06	2 (14%)	21,21,21	0.88	2 (9%)
3	SO4	D	502	-	4,4,4	0.23	0	6,6,6	0.08	0
5	VXZ	C	505	-	14,14,14	1.07	2 (14%)	21,21,21	0.87	2 (9%)
3	SO4	D	509	-	4,4,4	0.26	0	6,6,6	0.05	0
4	GOL	B	804	-	5,5,5	0.93	0	5,5,5	1.01	0
3	SO4	A	802	-	4,4,4	0.24	0	6,6,6	0.08	0
5	VXZ	D	508	-	14,14,14	1.05	2 (14%)	21,21,21	0.87	2 (9%)
3	SO4	B	811	-	4,4,4	0.23	0	6,6,6	0.07	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	VXZ	D	506	-	-	2/10/10/10	0/1/1/1
4	GOL	B	805	-	-	0/4/4/4	-
5	VXZ	B	809	-	-	0/10/10/10	0/1/1/1
5	VXZ	A	808	-	-	0/10/10/10	0/1/1/1
5	VXZ	B	806	-	-	0/10/10/10	0/1/1/1
4	GOL	A	806	-	-	2/4/4/4	-
4	GOL	D	505	-	-	4/4/4/4	-
5	VXZ	D	507	-	-	0/10/10/10	0/1/1/1
5	VXZ	A	807	-	-	0/10/10/10	0/1/1/1
5	VXZ	B	808	-	-	0/10/10/10	0/1/1/1
5	VXZ	A	811	-	-	0/10/10/10	0/1/1/1
5	VXZ	A	810	-	-	0/10/10/10	0/1/1/1
5	VXZ	C	506	-	-	0/10/10/10	0/1/1/1
4	GOL	A	805	-	-	1/4/4/4	-
4	GOL	D	504	-	-	0/4/4/4	-
5	VXZ	B	807	-	-	0/10/10/10	0/1/1/1
4	GOL	A	804	-	-	4/4/4/4	-
5	VXZ	A	809	-	-	0/10/10/10	0/1/1/1
5	VXZ	C	505	-	-	1/10/10/10	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	804	-	-	2/4/4/4	-
5	VXZ	D	508	-	-	0/10/10/10	0/1/1/1

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	809	VXZ	O4-C6	2.94	1.31	1.22
5	B	808	VXZ	O4-C6	2.89	1.31	1.22
5	C	506	VXZ	O4-C6	2.89	1.31	1.22
5	D	506	VXZ	O4-C6	2.88	1.31	1.22
5	A	809	VXZ	O4-C6	2.88	1.31	1.22
5	C	505	VXZ	O4-C6	2.88	1.31	1.22
5	B	806	VXZ	O4-C6	2.87	1.31	1.22
5	D	507	VXZ	O4-C6	2.85	1.31	1.22
5	A	811	VXZ	O4-C6	2.85	1.31	1.22
5	A	810	VXZ	O4-C6	2.85	1.31	1.22
5	B	807	VXZ	O4-C6	2.84	1.31	1.22
5	A	807	VXZ	O4-C6	2.83	1.31	1.22
5	A	808	VXZ	O4-C6	2.83	1.31	1.22
5	D	508	VXZ	O4-C6	2.83	1.31	1.22
5	C	505	VXZ	O3-C6	-2.59	1.22	1.30
5	B	806	VXZ	O3-C6	-2.59	1.22	1.30
5	A	809	VXZ	O3-C6	-2.57	1.22	1.30
5	D	506	VXZ	O3-C6	-2.57	1.22	1.30
5	C	506	VXZ	O3-C6	-2.57	1.22	1.30
5	A	811	VXZ	O2-S	2.56	1.56	1.43
5	B	809	VXZ	O3-C6	-2.56	1.22	1.30
5	A	807	VXZ	O3-C6	-2.56	1.22	1.30
5	B	807	VXZ	O3-C6	-2.56	1.22	1.30
5	D	507	VXZ	O3-C6	-2.55	1.22	1.30
5	A	810	VXZ	O3-C6	-2.55	1.22	1.30
5	B	808	VXZ	O3-C6	-2.55	1.22	1.30
5	D	508	VXZ	O3-C6	-2.55	1.22	1.30
5	A	808	VXZ	O3-C6	-2.54	1.23	1.30
5	A	811	VXZ	O3-C6	-2.51	1.23	1.30
5	B	809	VXZ	O2-S	2.31	1.54	1.43

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	808	VXZ	O4-C6-C5	-3.14	114.47	121.97
5	A	808	VXZ	O4-C6-C5	-3.13	114.50	121.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	811	VXZ	O4-C6-C5	-3.12	114.52	121.97
5	B	807	VXZ	O4-C6-C5	-3.09	114.58	121.97
5	A	807	VXZ	O4-C6-C5	-3.06	114.66	121.97
5	A	810	VXZ	O4-C6-C5	-3.06	114.68	121.97
5	D	506	VXZ	O4-C6-C5	-3.05	114.70	121.97
5	D	508	VXZ	O4-C6-C5	-3.04	114.72	121.97
5	A	809	VXZ	O4-C6-C5	-3.03	114.75	121.97
5	C	505	VXZ	O4-C6-C5	-3.02	114.75	121.97
5	B	806	VXZ	O4-C6-C5	-3.00	114.81	121.97
5	D	507	VXZ	O4-C6-C5	-2.99	114.83	121.97
5	C	506	VXZ	O4-C6-C5	-2.98	114.86	121.97
5	B	809	VXZ	O4-C6-C5	-2.78	115.33	121.97
5	A	811	VXZ	O3-C6-C5	2.51	122.41	115.28
5	B	808	VXZ	O3-C6-C5	2.45	122.24	115.28
5	A	808	VXZ	O3-C6-C5	2.43	122.19	115.28
5	B	807	VXZ	O3-C6-C5	2.41	122.13	115.28
5	A	809	VXZ	O3-C6-C5	2.38	122.06	115.28
5	D	506	VXZ	O3-C6-C5	2.38	122.05	115.28
5	A	810	VXZ	O3-C6-C5	2.38	122.03	115.28
5	D	508	VXZ	O3-C6-C5	2.36	121.99	115.28
5	C	505	VXZ	O3-C6-C5	2.35	121.97	115.28
5	A	807	VXZ	O3-C6-C5	2.35	121.95	115.28
5	D	507	VXZ	O3-C6-C5	2.33	121.92	115.28
5	B	806	VXZ	O3-C6-C5	2.31	121.85	115.28
5	C	506	VXZ	O3-C6-C5	2.31	121.85	115.28
5	B	809	VXZ	O3-C6-C5	2.26	121.70	115.28

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	804	GOL	O1-C1-C2-C3
4	A	804	GOL	C1-C2-C3-O3
4	A	806	GOL	O1-C1-C2-C3
4	D	505	GOL	O1-C1-C2-C3
4	D	505	GOL	C1-C2-C3-O3
4	A	804	GOL	O2-C2-C3-O3
4	B	804	GOL	O1-C1-C2-C3
4	A	806	GOL	O1-C1-C2-O2
4	D	505	GOL	O2-C2-C3-O3
4	B	804	GOL	O1-C1-C2-O2
4	D	505	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
4	A	804	GOL	O1-C1-C2-O2
4	A	805	GOL	O2-C2-C3-O3
5	D	506	VXZ	C-C5-C6-O4
5	D	506	VXZ	C-C5-C6-O3
5	C	505	VXZ	C-C5-C6-O4

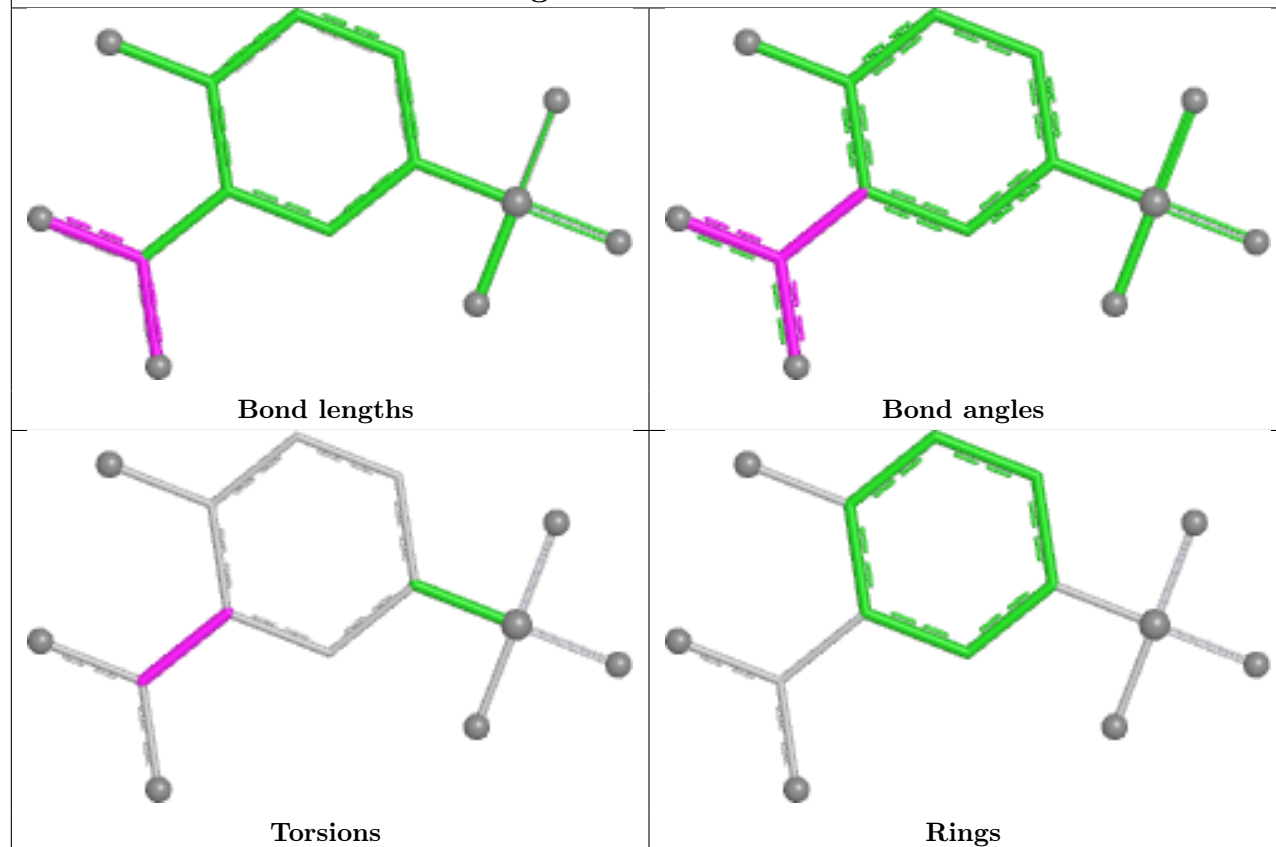
There are no ring outliers.

4 monomers are involved in 5 short contacts:

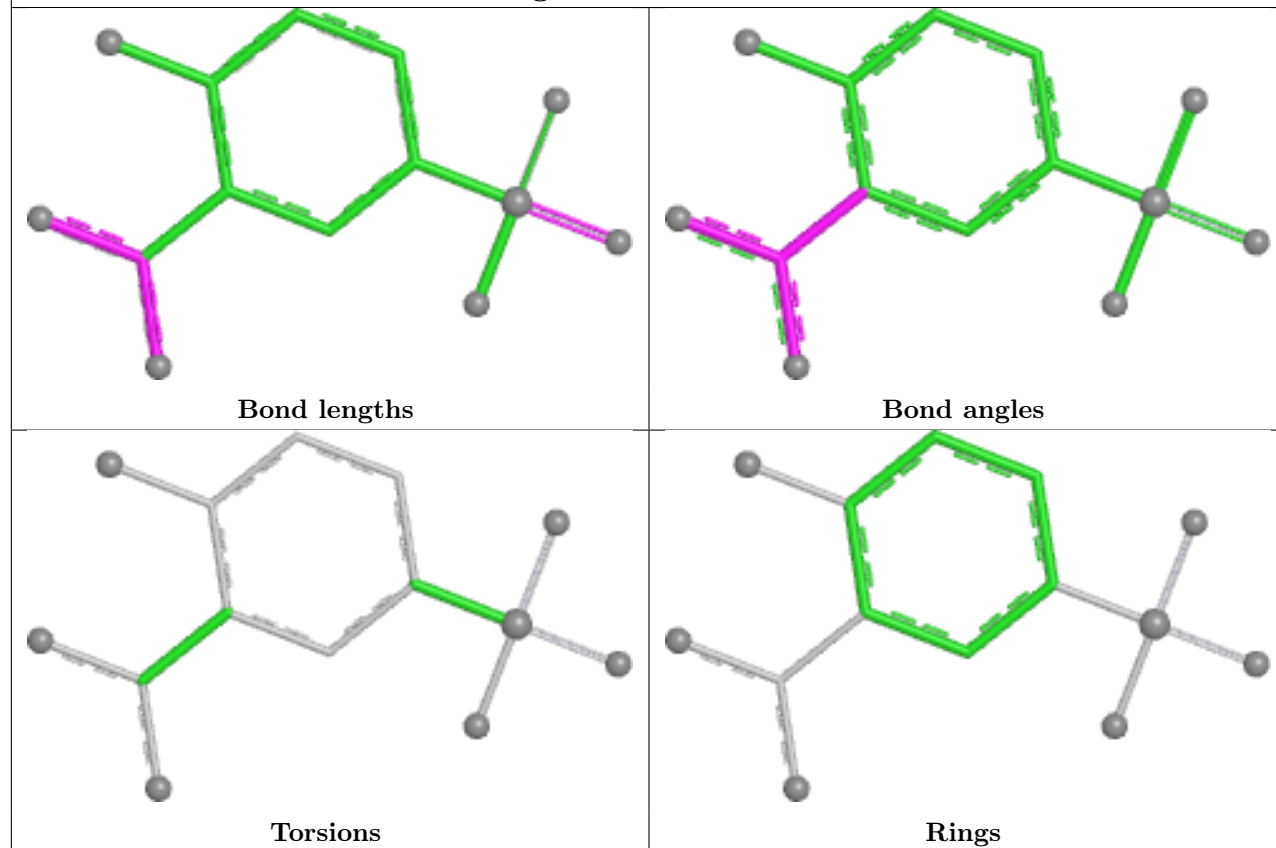
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	809	VXZ	1	0
5	D	507	VXZ	1	0
5	A	807	VXZ	1	0
4	A	804	GOL	2	0

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

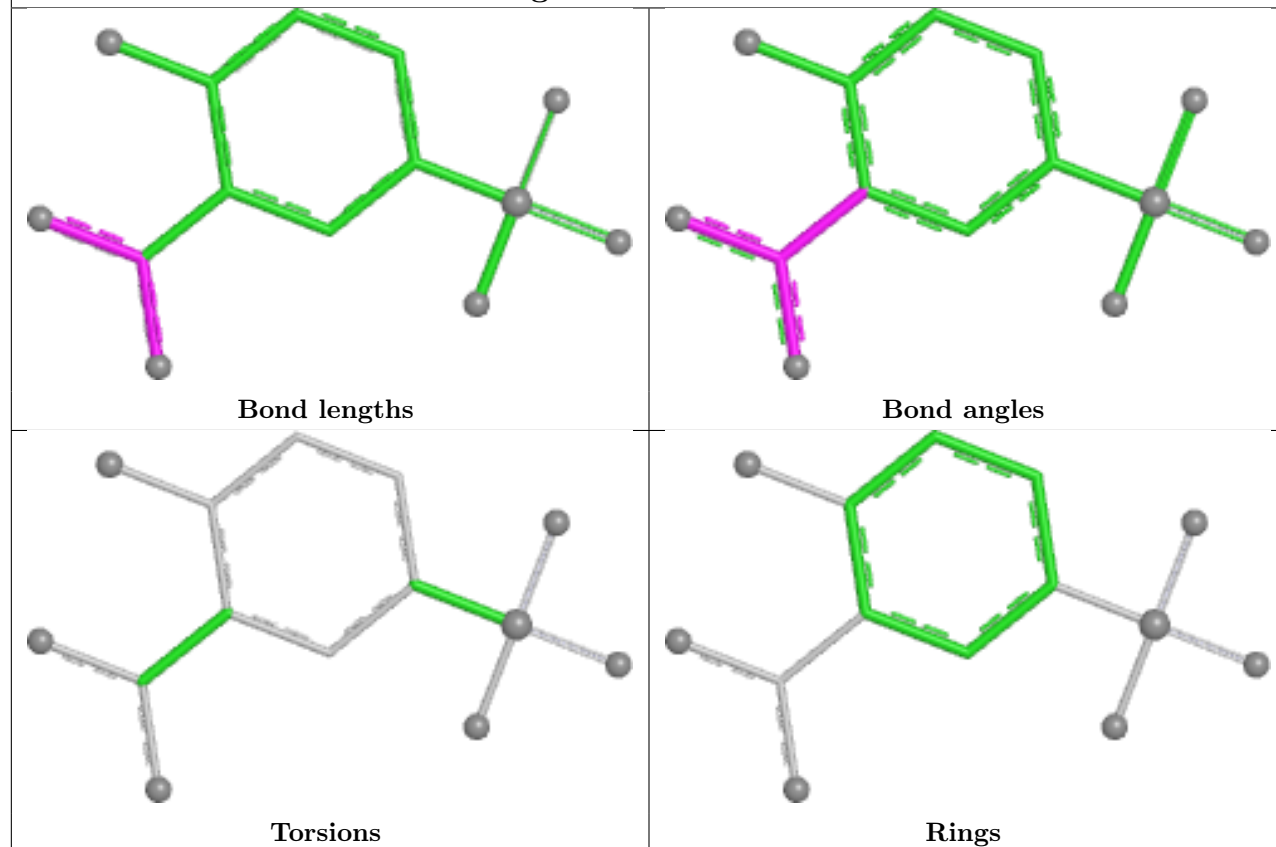
Ligand VXZ D 506



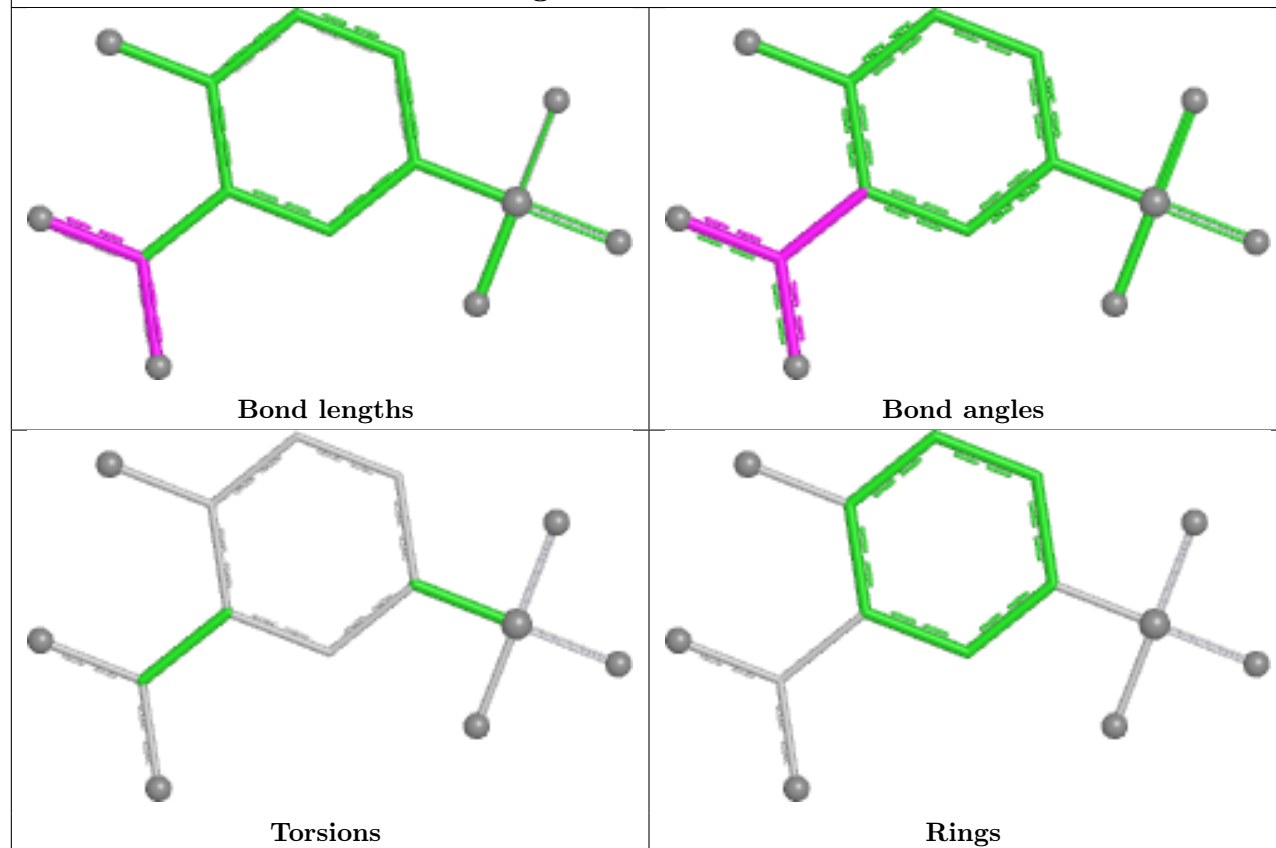
Ligand VXZ B 809



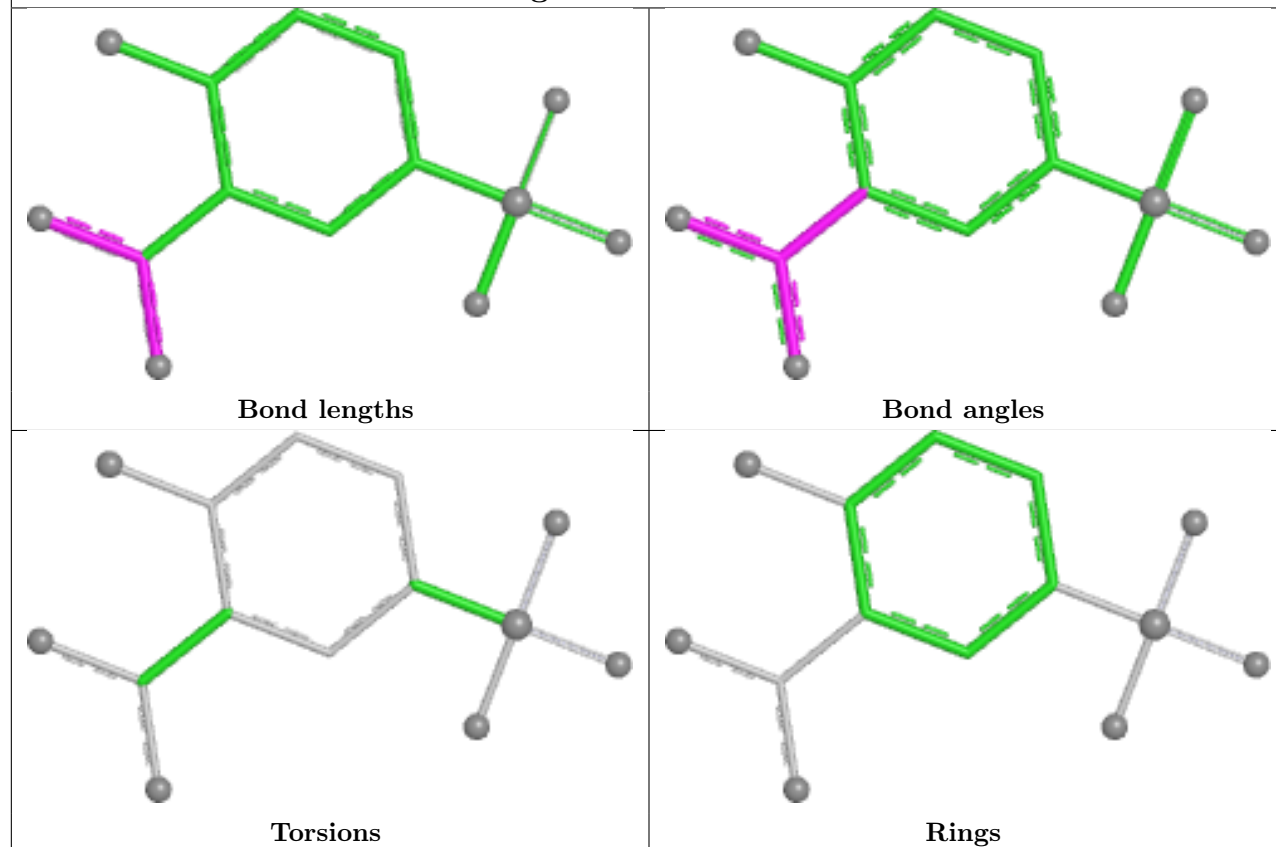
Ligand VXZ A 808



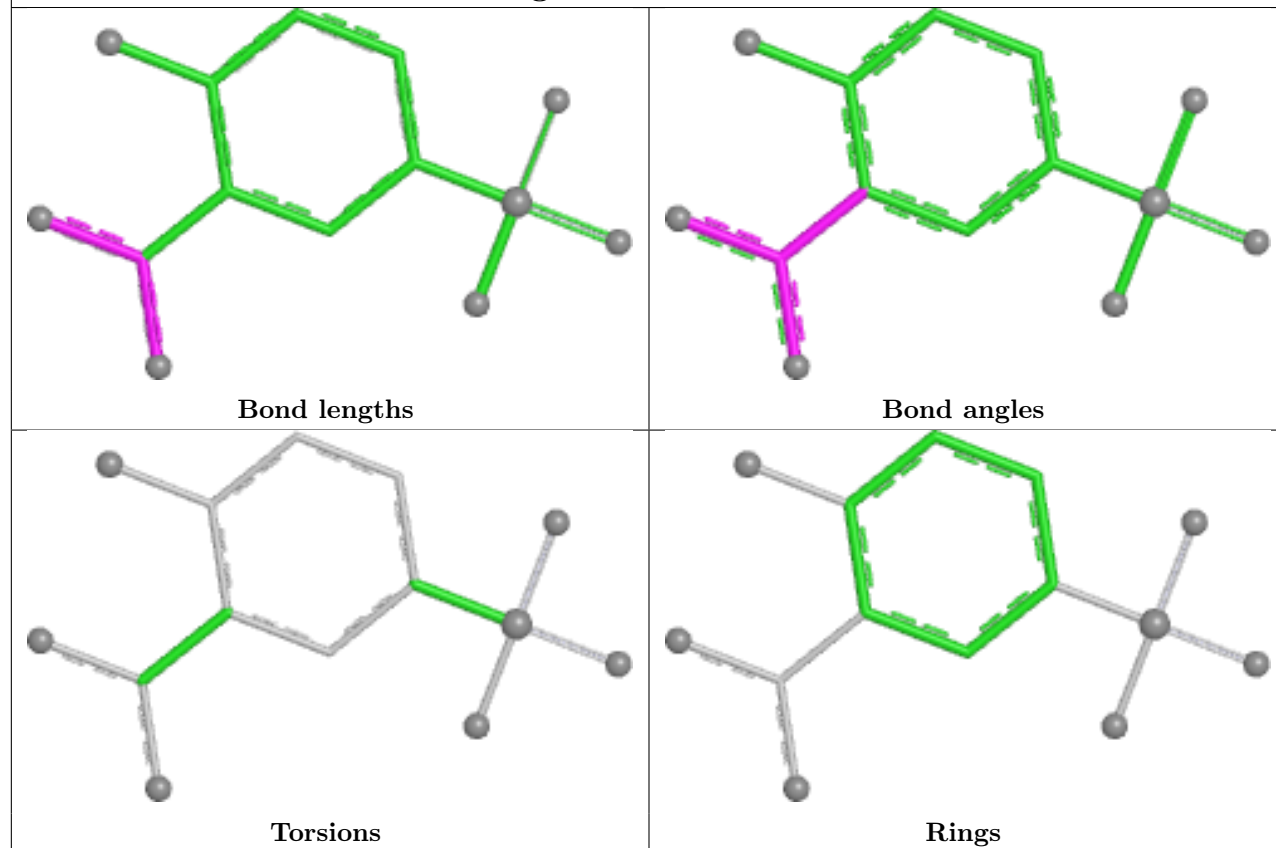
Ligand VXZ B 806



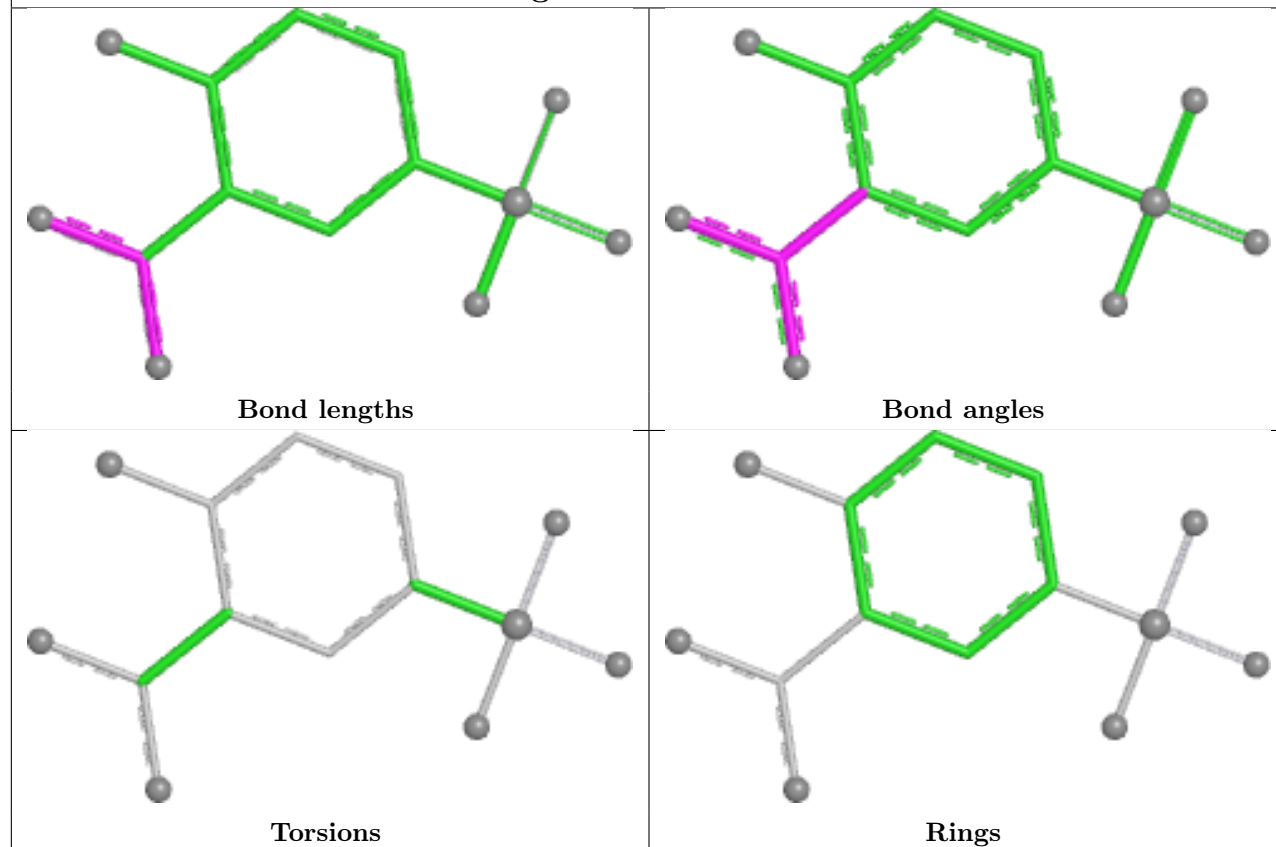
Ligand VXZ D 507



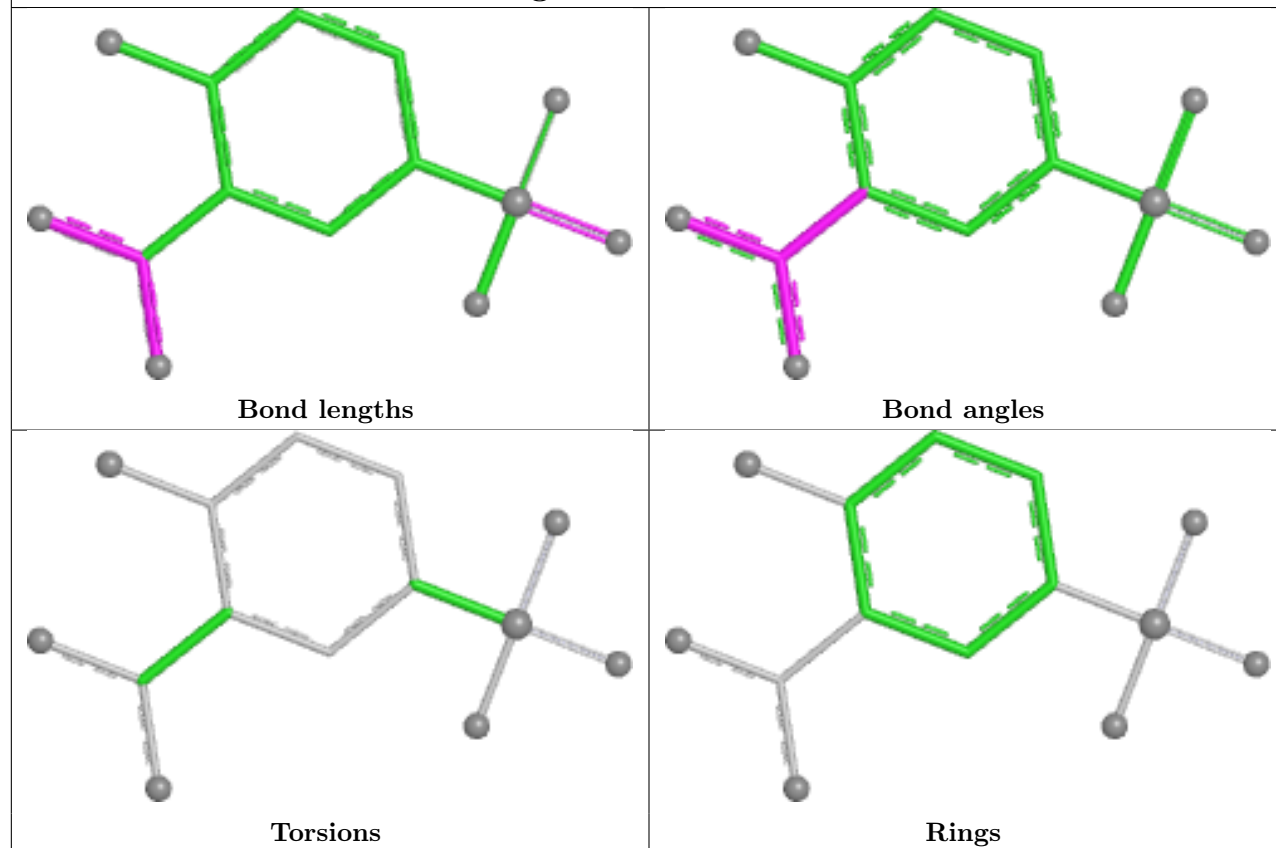
Ligand VXZ A 807



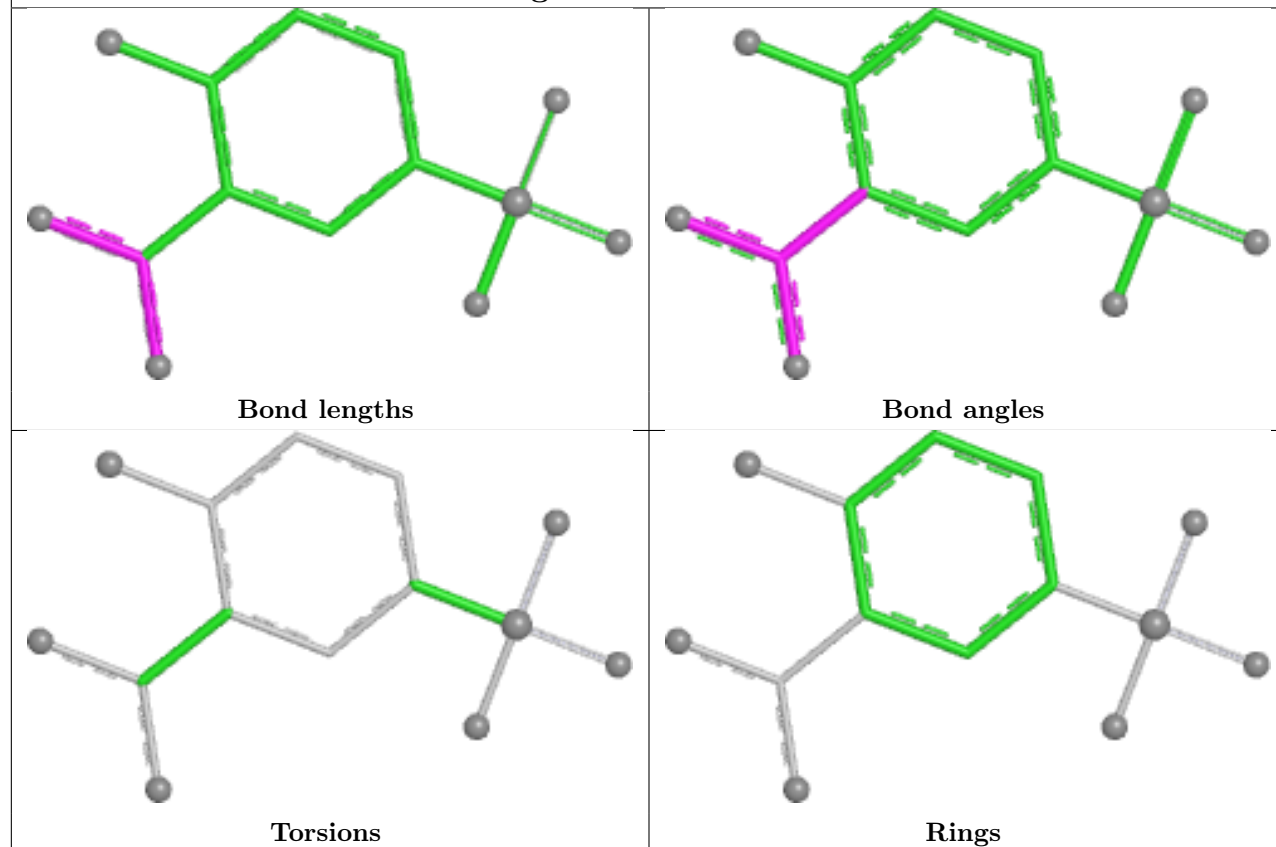
Ligand VXZ B 808



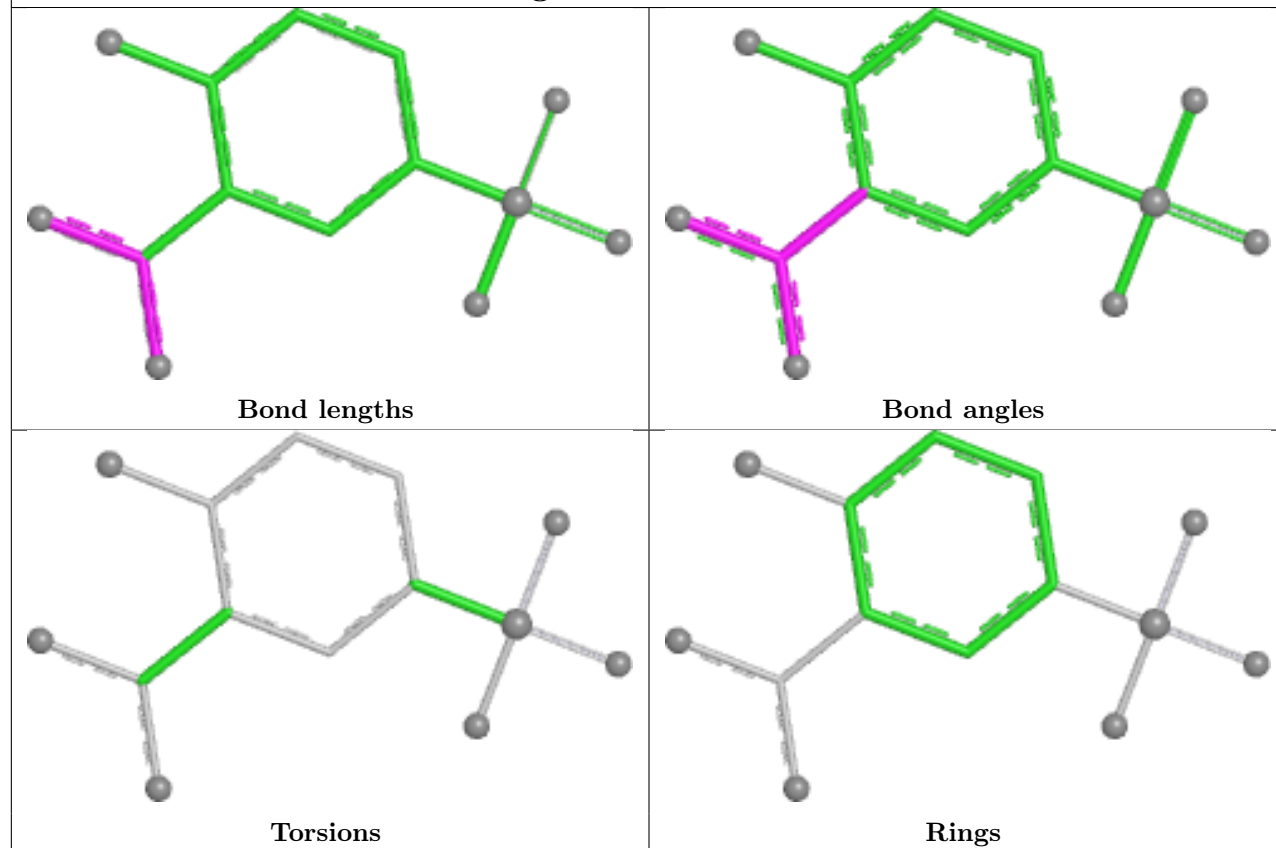
Ligand VXZ A 811



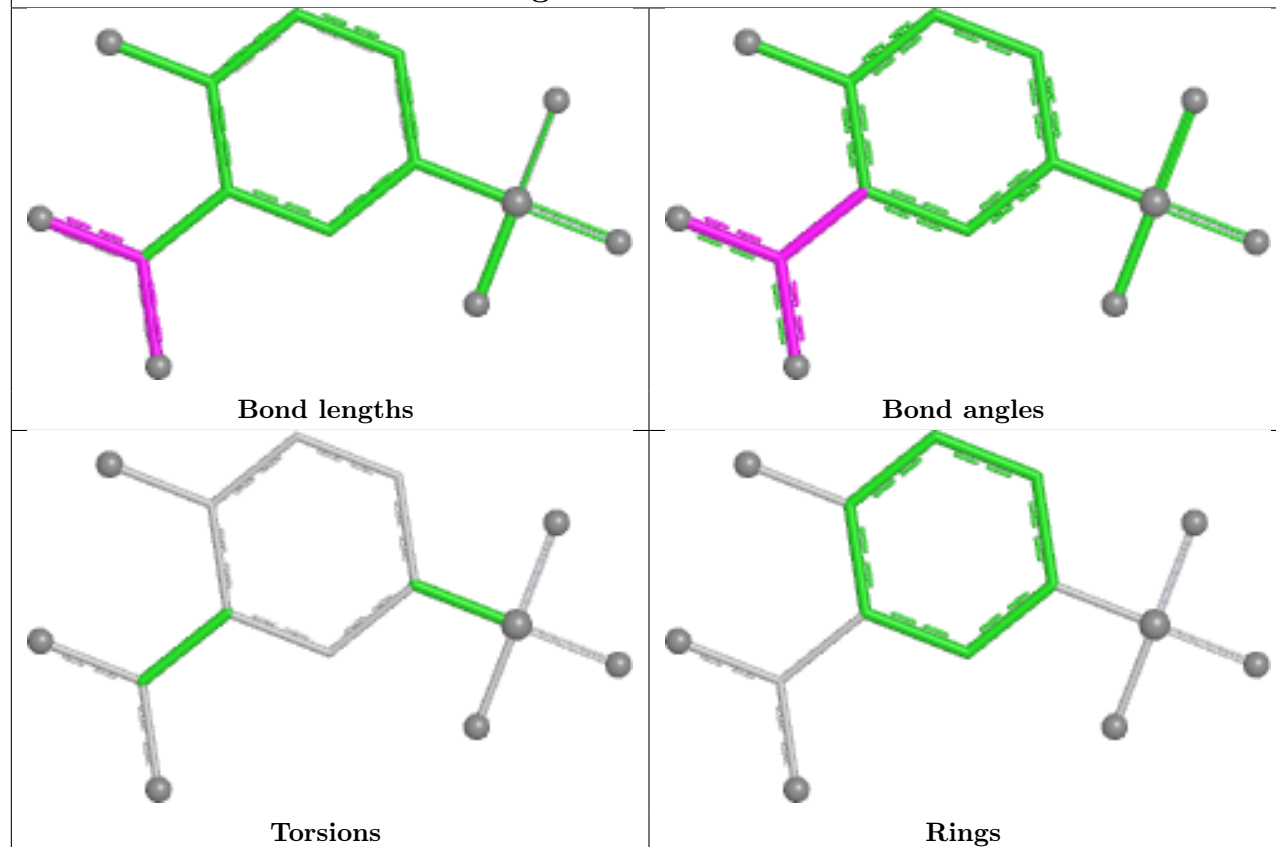
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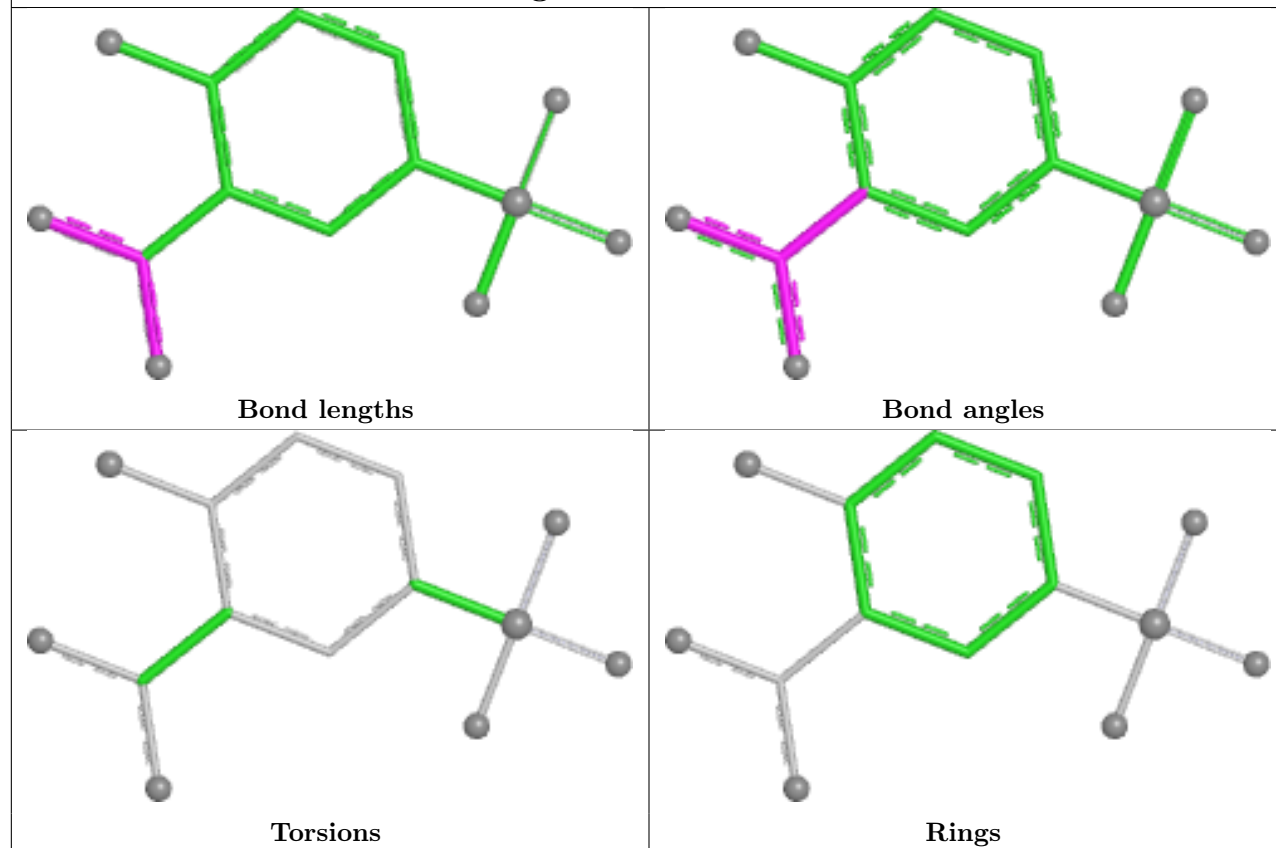
Ligand VXZ C 506



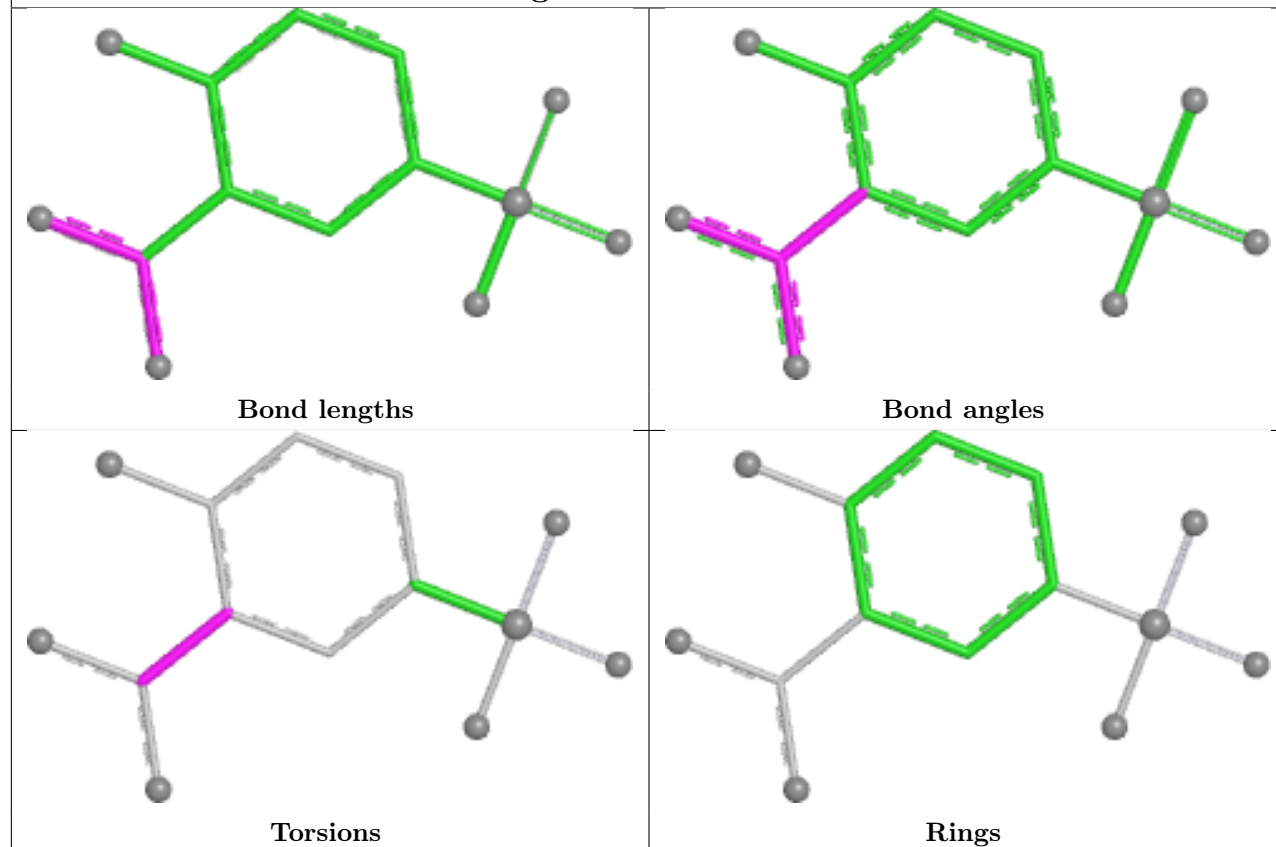
Ligand VXZ B 807



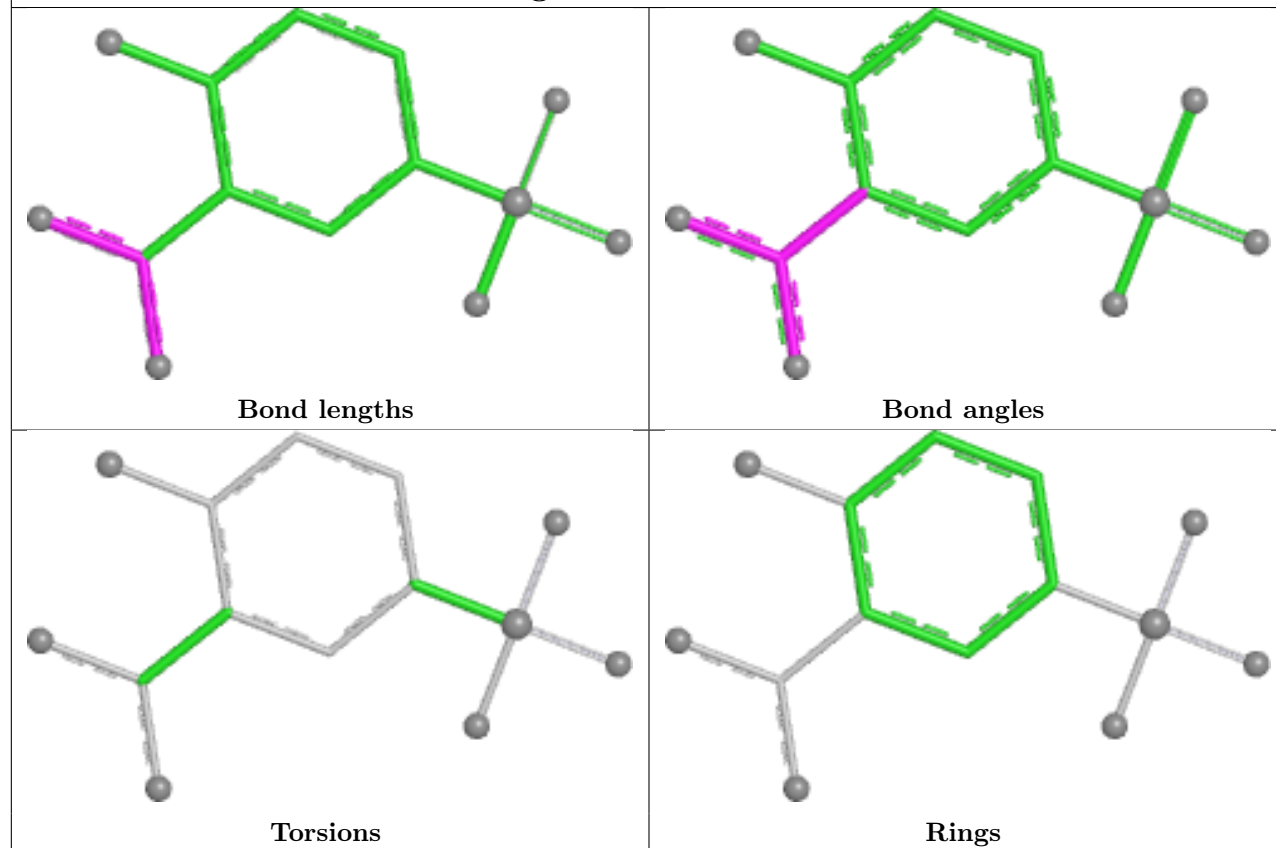
Ligand VXZ A 809



Ligand VXZ C 505



Ligand VXZ D 508



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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	724/736 (98%)	0.39	55 (7%) 20 19	32, 51, 87, 137	0
1	B	726/736 (98%)	0.26	43 (5%) 28 27	32, 48, 79, 135	0
2	C	402/403 (99%)	0.28	31 (7%) 19 18	33, 45, 77, 113	0
2	D	402/403 (99%)	0.31	39 (9%) 13 12	30, 46, 80, 140	0
All	All	2254/2278 (98%)	0.32	168 (7%) 20 19	30, 48, 82, 140	0

All (168) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	294	ALA	8.4
2	D	390	ILE	6.8
1	A	574	ALA	6.7
2	C	294	ALA	6.1
2	D	135	GLY	6.0
2	D	302	GLY	6.0
2	D	136	LEU	5.8
2	C	293	GLY	5.7
1	A	413	PHE	5.6
2	D	296	PRO	5.5
2	C	390	ILE	5.5
2	D	293	GLY	5.4
2	C	225	PHE	5.4
2	D	297	VAL	5.3
1	B	720	SER	5.3
2	C	302	GLY	5.2
2	D	391	GLY	5.1
1	B	689	PRO	5.1
1	A	2	PRO	5.1
2	C	300	LEU	5.0
1	A	571	VAL	5.0

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Mol	Chain	Res	Type	RSRZ
1	A	720	SER	5.0
2	D	228	LEU	4.9
2	C	292	SER	4.9
1	B	413	PHE	4.9
1	B	571	VAL	4.8
1	B	577	THR	4.8
2	D	301	THR	4.8
2	C	301	THR	4.6
1	B	690	ALA	4.6
1	B	576	GLY	4.6
2	D	392	GLY	4.6
1	A	575	GLY	4.5
2	C	136	LEU	4.5
2	C	389	CYS	4.5
1	A	576	GLY	4.3
2	D	298	ILE	4.2
2	D	389	CYS	4.2
1	B	578	TYR	4.2
1	B	692	THR	4.2
2	D	300	LEU	4.1
2	D	225	PHE	4.1
1	B	-8	HIS	3.9
1	B	688	GLY	3.9
2	D	295	ASP	3.8
2	C	296	PRO	3.8
2	D	393	GLY	3.8
1	A	258	MET	3.8
1	B	1	MET	3.8
2	C	134	MET	3.8
2	D	2	SER	3.7
2	D	231	LEU	3.6
2	D	133	ALA	3.6
1	B	258	MET	3.5
1	B	687	SER	3.5
2	D	299	MET	3.5
1	A	414	GLU	3.5
2	D	303	PRO	3.5
1	B	-9	HIS	3.5
2	C	303	PRO	3.4
1	B	442	THR	3.4
1	A	451	GLY	3.4
1	B	568	ARG	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	367	TYR	3.4
1	A	-9	HIS	3.3
1	A	412	VAL	3.3
1	B	71	LYS	3.3
1	A	577	THR	3.3
1	B	570	GLY	3.3
1	A	1	MET	3.3
1	B	412	VAL	3.3
1	A	224	LYS	3.3
1	A	442	THR	3.2
2	C	276	LYS	3.2
1	B	575	GLY	3.2
1	A	444	PRO	3.2
1	B	569	LYS	3.2
2	D	134	MET	3.2
1	A	579	GLN	3.2
1	B	418	LEU	3.2
1	A	418	LEU	3.1
2	C	391	GLY	3.1
2	C	295	ASP	3.1
1	B	444	PRO	3.0
1	B	574	ALA	3.0
1	A	379	GLY	3.0
2	C	228	LEU	3.0
2	C	231	LEU	3.0
1	A	381	THR	3.0
1	A	432	PRO	2.9
2	C	392	GLY	2.9
2	C	393	GLY	2.9
2	D	222	LYS	2.9
1	B	579	GLN	2.9
1	B	580	PRO	2.9
2	C	297	VAL	2.8
2	C	298	ILE	2.8
2	D	291	THR	2.8
1	B	224	LYS	2.8
1	A	610	GLY	2.8
1	A	356	VAL	2.8
1	B	443	LEU	2.7
1	B	573	ASP	2.7
1	A	363	LYS	2.7
2	D	277	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	133	LYS	2.7
2	D	276	LYS	2.7
1	A	478	GLY	2.6
1	A	443	LEU	2.6
2	C	222	LYS	2.6
1	A	370	LYS	2.6
1	A	453	LYS	2.5
1	A	-12	SER	2.5
1	B	2	PRO	2.5
2	C	224	ALA	2.5
1	B	691	GLY	2.5
1	A	421	LYS	2.5
2	C	149	GLN	2.5
2	D	304	THR	2.5
2	C	2	SER	2.4
2	C	230	ALA	2.4
2	C	299	MET	2.4
2	D	50	LEU	2.4
1	A	424	GLY	2.4
1	A	325	ILE	2.4
1	A	578	TYR	2.4
1	B	567	THR	2.4
1	A	436	LEU	2.4
1	B	559	LEU	2.4
2	D	224	ALA	2.3
1	A	184	ILE	2.3
1	A	132	VAL	2.3
1	A	335	MET	2.3
1	A	611	LYS	2.3
1	A	318	GLU	2.3
2	C	277	LEU	2.3
1	B	370	LYS	2.3
2	C	140	THR	2.3
2	D	305	PRO	2.3
1	A	486	LEU	2.3
1	B	377	GLU	2.2
1	A	320	ILE	2.2
1	A	450	THR	2.2
2	C	45	ASP	2.2
1	A	485	ALA	2.2
2	D	394	MET	2.2
2	D	230	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
2	D	149	GLN	2.2
1	B	620	ARG	2.2
1	A	3	ASP	2.2
1	B	589	LYS	2.2
1	B	211	ALA	2.1
1	A	620	ARG	2.1
1	A	235	SER	2.1
1	A	441	SER	2.1
1	A	433	ASN	2.1
1	A	380	ARG	2.1
1	A	449	ALA	2.1
2	D	292	SER	2.1
2	D	226	GLU	2.1
2	D	215	LYS	2.1
1	A	388	ALA	2.1
1	B	610	GLY	2.1
1	B	693	GLY	2.1
1	B	621	GLU	2.0
1	B	380	ARG	2.0
1	A	570	GLY	2.0
2	D	315	GLY	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	A	802	5/5	0.63	0.14	62,83,97,99	5
3	SO4	C	501	5/5	0.68	0.13	102,108,121,130	0

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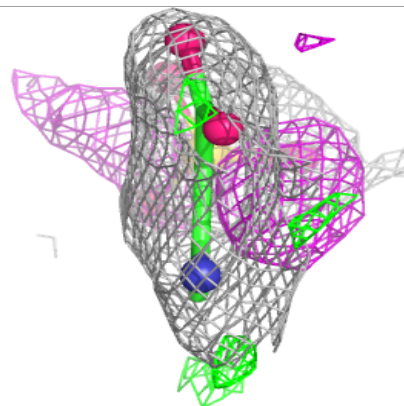
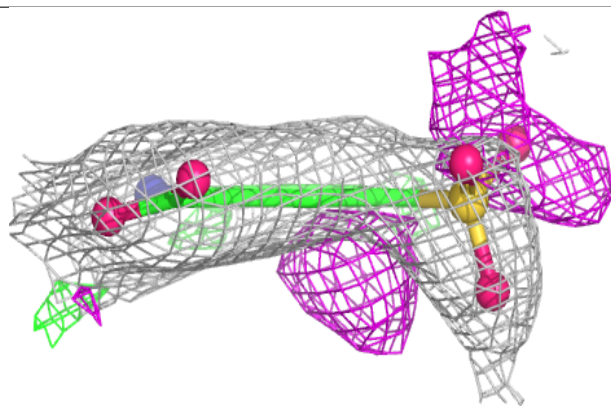
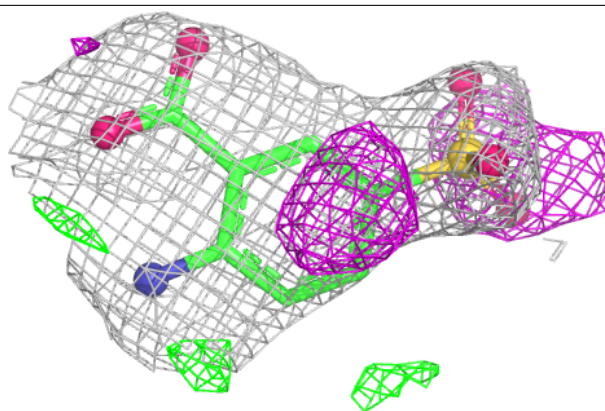
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	B	803	5/5	0.77	0.14	55,59,80,81	5
4	GOL	A	806	6/6	0.77	0.23	71,72,82,87	0
5	VXZ	A	810	14/14	0.77	0.19	66,82,125,138	0
4	GOL	D	504	6/6	0.79	0.18	64,67,70,71	0
3	SO4	A	813	5/5	0.79	0.12	61,70,104,137	0
5	VXZ	D	508	14/14	0.79	0.14	62,73,95,95	0
3	SO4	B	811	5/5	0.80	0.11	65,85,100,128	0
3	SO4	D	502	5/5	0.80	0.10	73,83,110,113	0
3	SO4	C	507	5/5	0.81	0.11	75,81,107,118	0
3	SO4	B	810	5/5	0.81	0.14	78,78,94,95	0
3	SO4	C	503	5/5	0.82	0.10	87,91,97,124	0
3	SO4	D	509	5/5	0.82	0.15	96,100,105,109	5
3	SO4	A	812	5/5	0.83	0.14	72,80,97,100	0
3	SO4	B	802	5/5	0.83	0.10	73,80,83,85	0
3	SO4	C	504	5/5	0.84	0.08	81,86,98,103	0
3	SO4	A	803	5/5	0.84	0.09	73,91,94,96	0
3	SO4	B	801	5/5	0.85	0.12	58,67,77,78	0
5	VXZ	B	808	14/14	0.85	0.12	66,80,98,99	0
5	VXZ	C	506	14/14	0.85	0.12	57,68,107,108	0
4	GOL	D	505	6/6	0.85	0.16	55,61,63,68	0
4	GOL	B	804	6/6	0.86	0.19	76,84,90,96	0
4	GOL	A	804	6/6	0.86	0.17	50,59,67,75	0
5	VXZ	A	811	14/14	0.86	0.11	64,72,85,92	0
5	VXZ	D	507	14/14	0.88	0.09	60,76,107,112	0
5	VXZ	B	809	14/14	0.89	0.11	53,68,82,85	0
5	VXZ	A	809	14/14	0.89	0.12	45,62,74,76	0
5	VXZ	C	505	14/14	0.90	0.11	65,79,87,90	0
3	SO4	A	801	5/5	0.90	0.09	56,66,76,85	0
4	GOL	A	805	6/6	0.91	0.11	42,44,54,54	0
5	VXZ	D	506	14/14	0.91	0.11	63,78,83,94	0
4	GOL	B	805	6/6	0.92	0.13	37,45,58,61	0
5	VXZ	B	807	14/14	0.93	0.12	57,71,82,83	0
3	SO4	D	501	5/5	0.94	0.07	71,76,84,91	0
3	SO4	D	503	5/5	0.95	0.06	64,64,71,75	0
5	VXZ	A	807	14/14	0.96	0.10	55,69,82,84	0
5	VXZ	B	806	14/14	0.96	0.08	53,62,69,71	0
3	SO4	C	502	5/5	0.97	0.05	63,65,73,76	0
5	VXZ	A	808	14/14	0.97	0.07	47,54,62,65	0

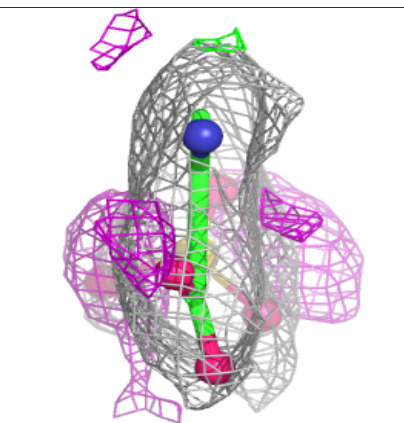
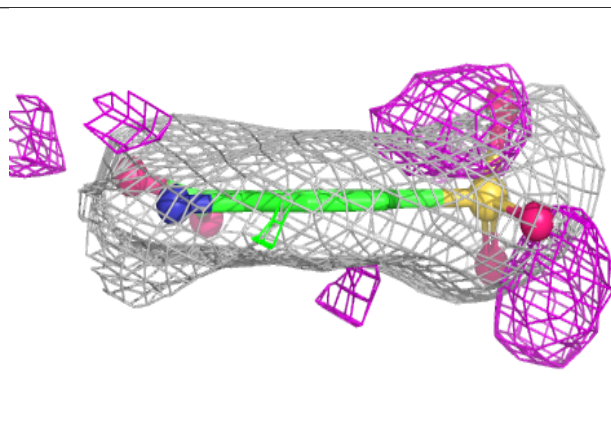
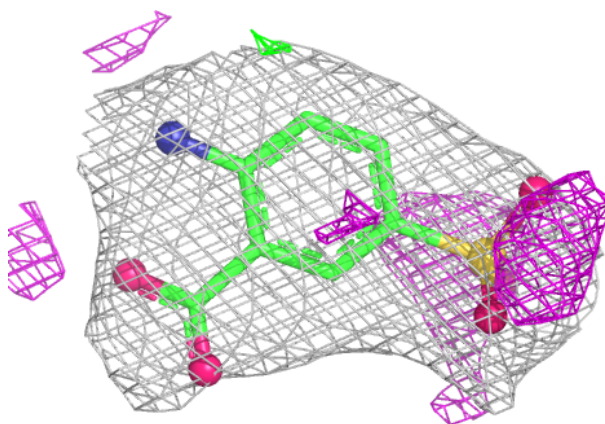
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around VXZ A 810:

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

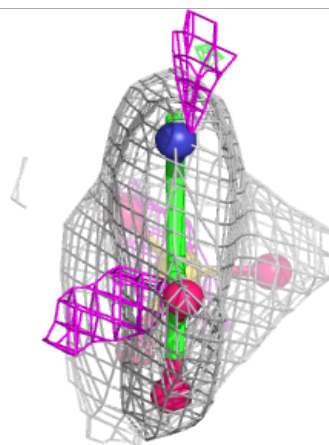
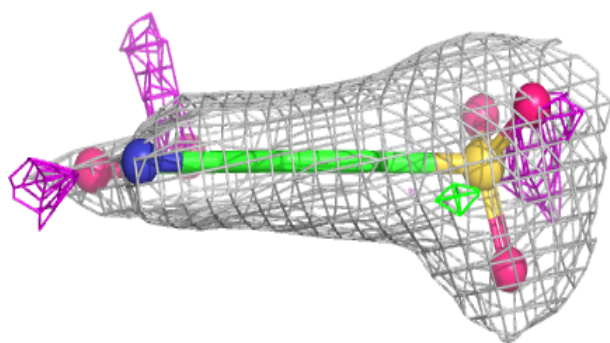
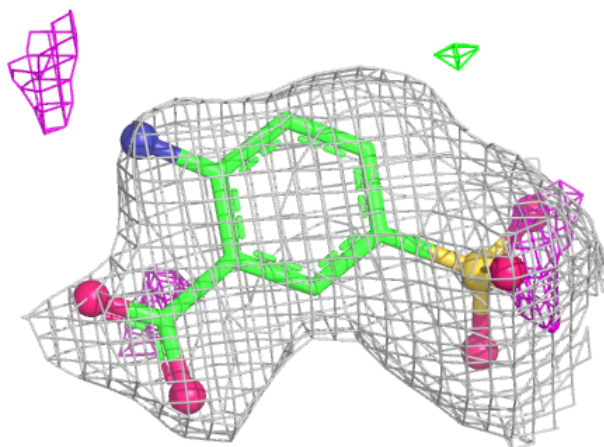
**Electron density around VXZ D 508:**

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



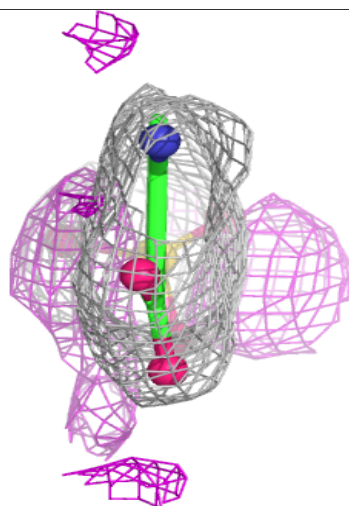
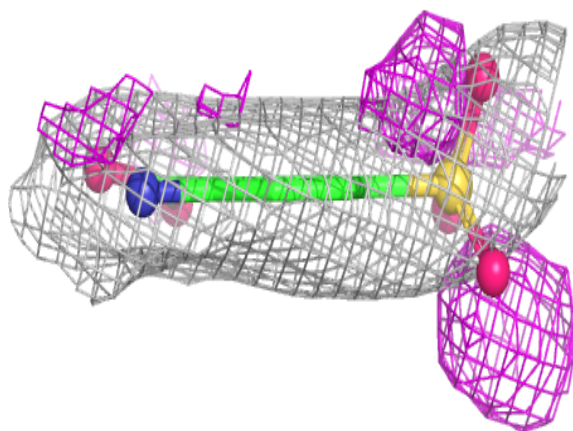
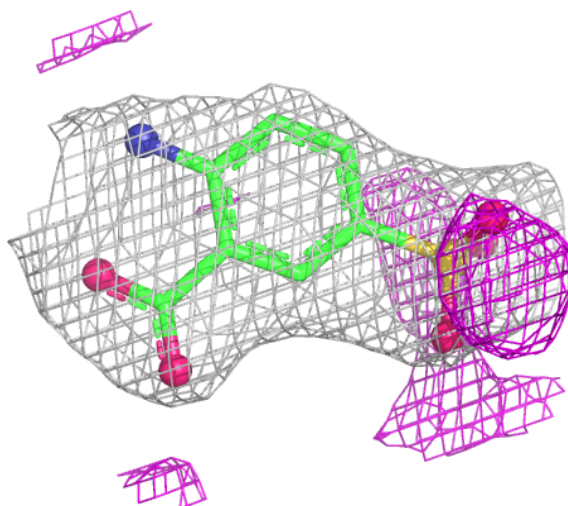
Electron density around VXZ B 808:

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



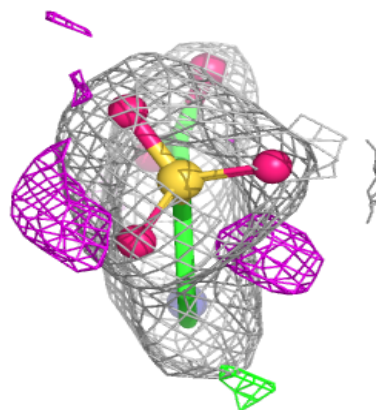
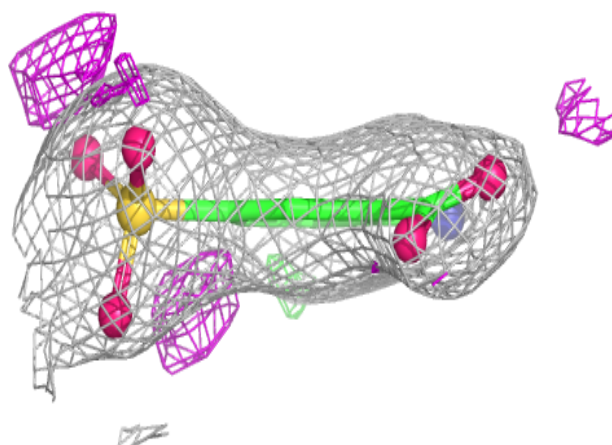
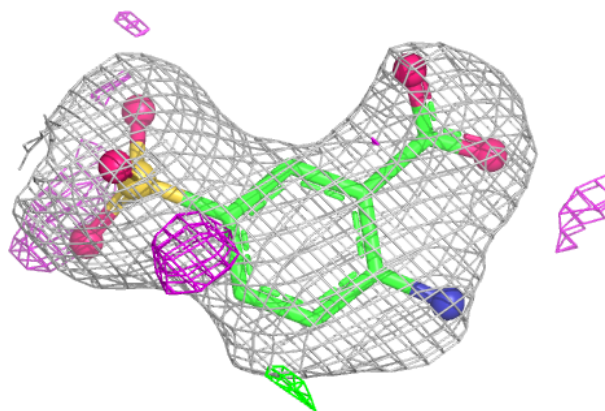
Electron density around VXZ C 506:

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



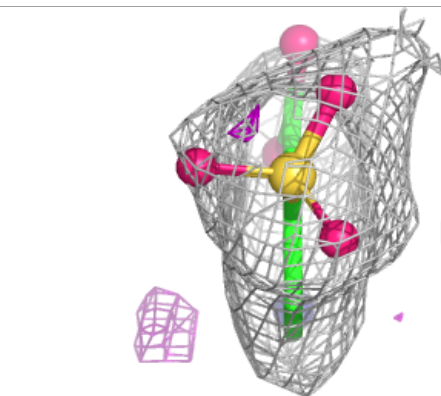
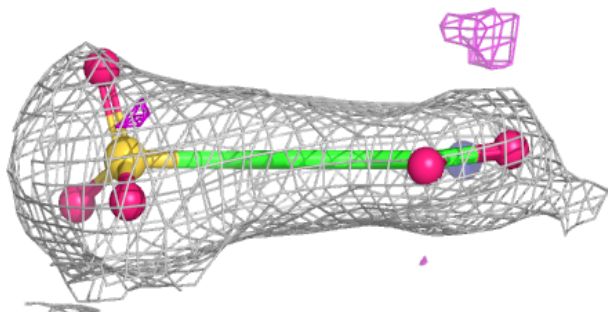
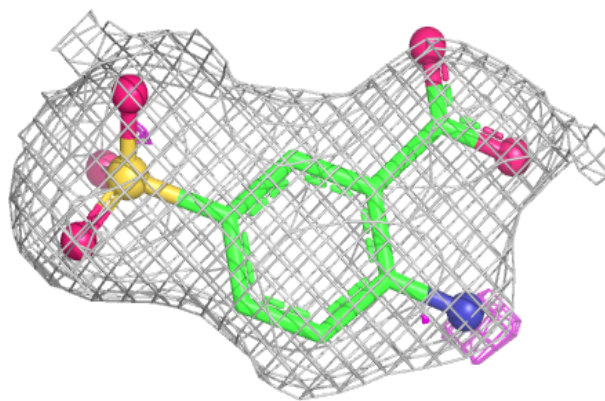
Electron density around VXZ A 811:

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

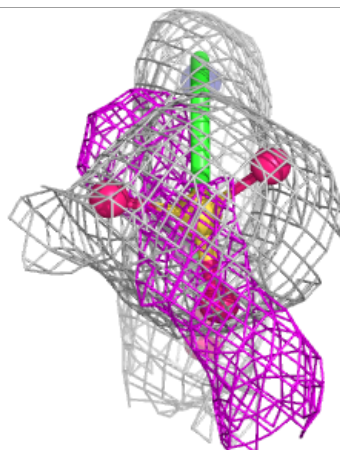
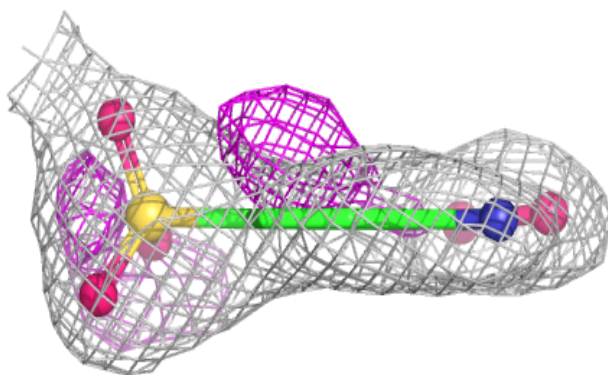
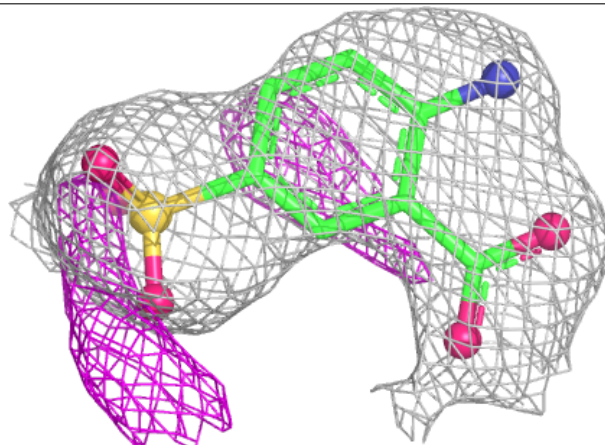


Electron density around VXZ D 507:

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

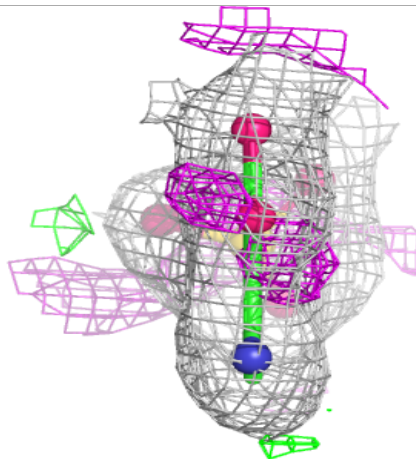
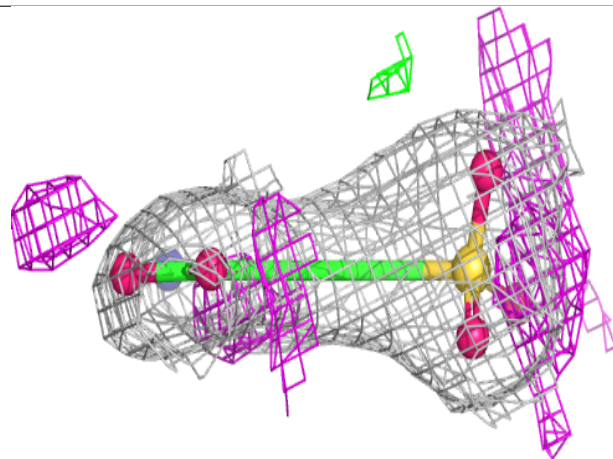
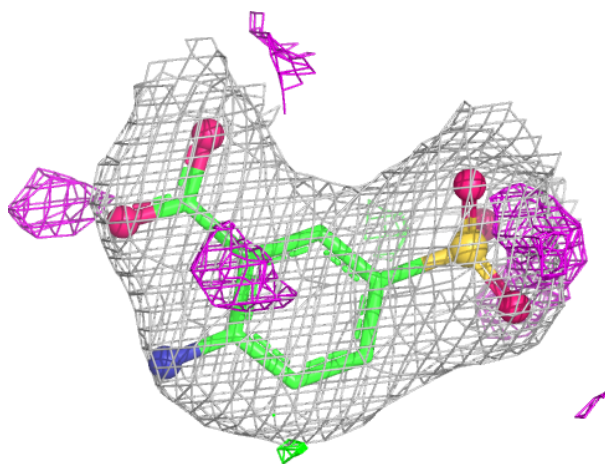
**Electron density around VXZ B 809:**

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and green (positive)



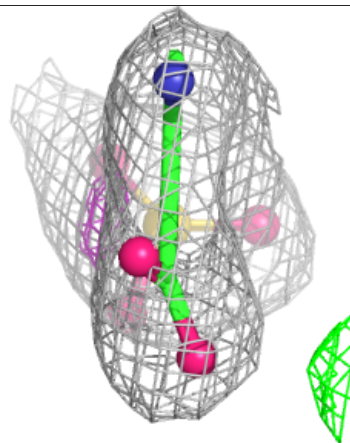
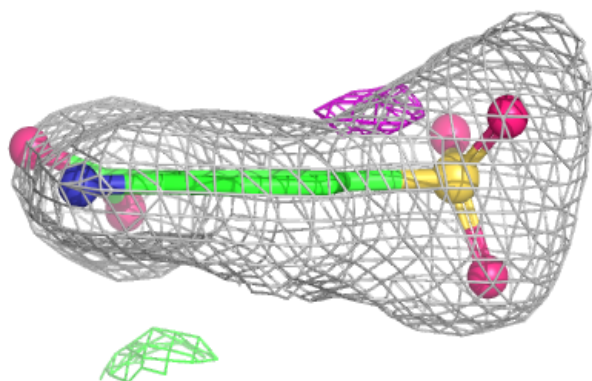
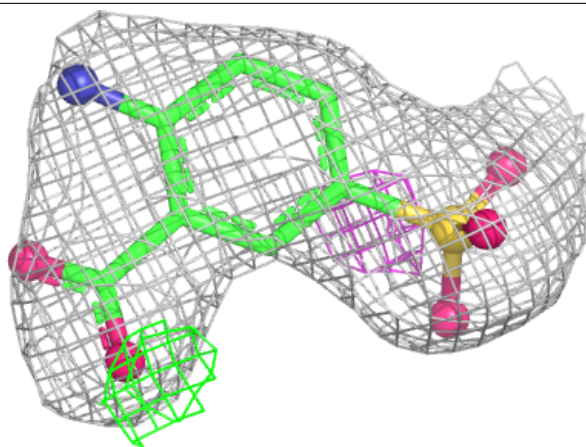
Electron density around VXZ A 809:

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



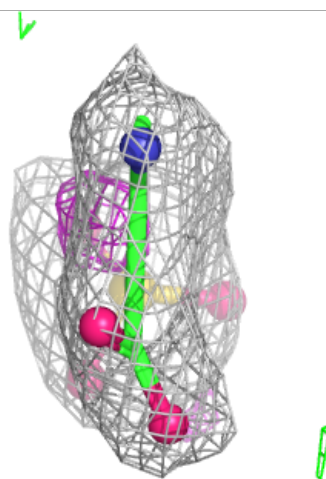
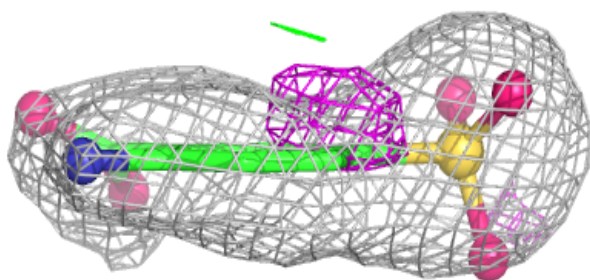
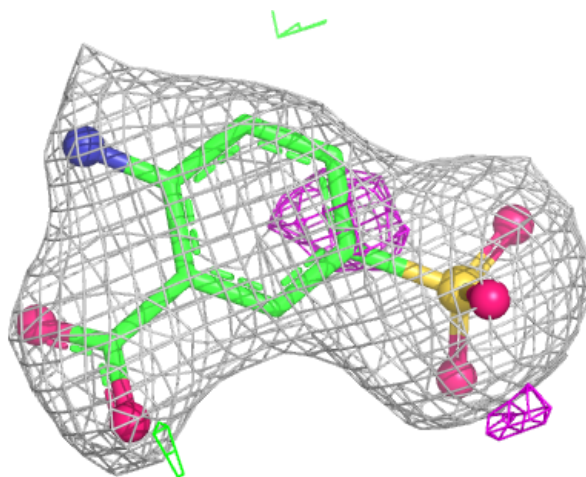
Electron density around VXZ C 505:

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



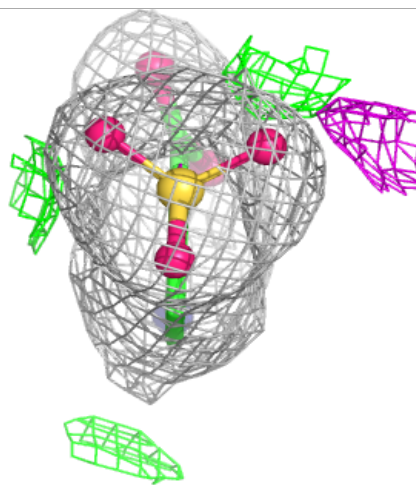
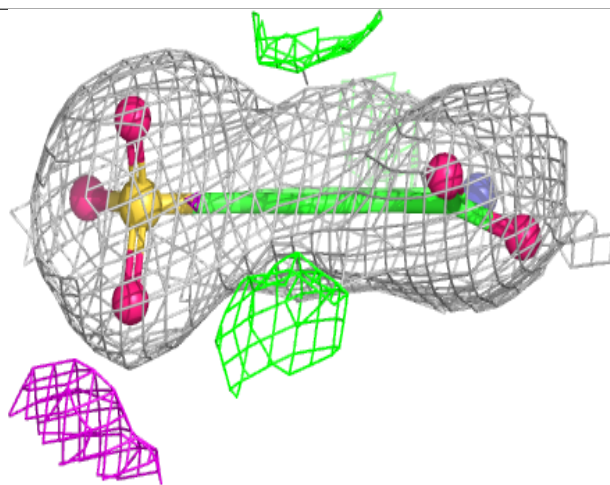
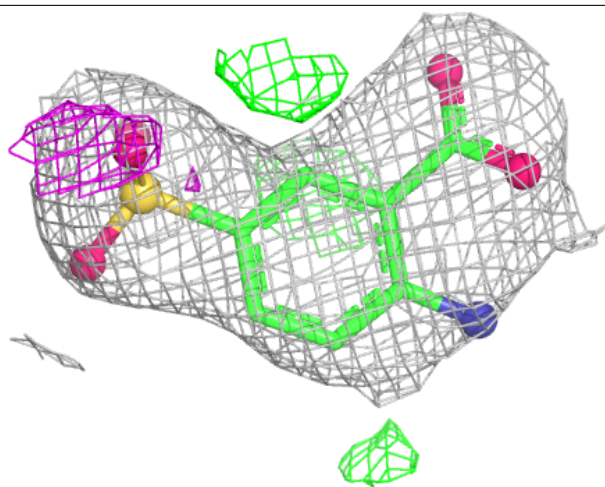
Electron density around VXZ D 506:

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



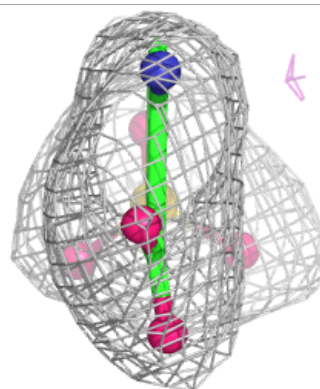
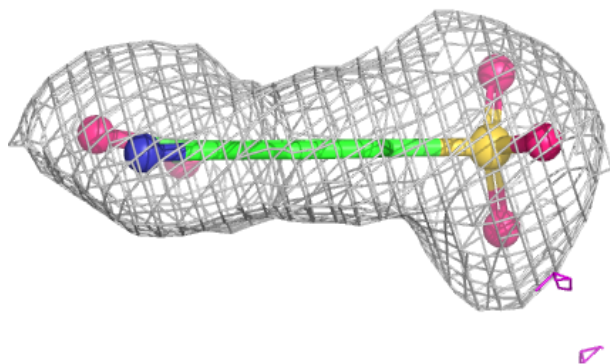
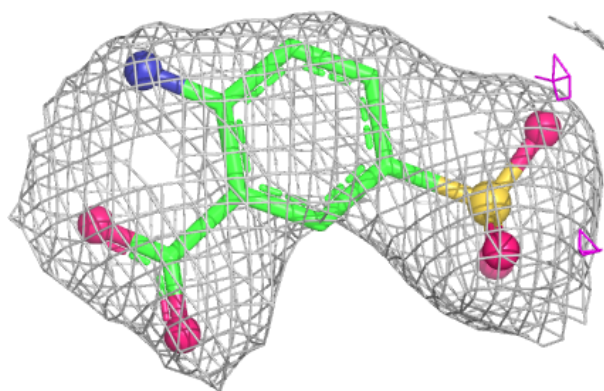
Electron density around VXZ B 807:

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

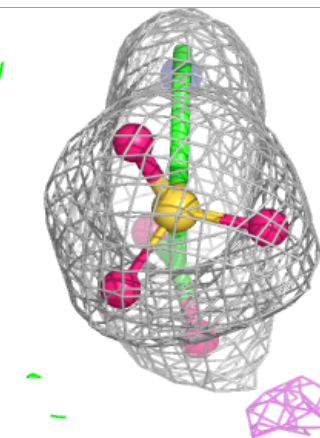
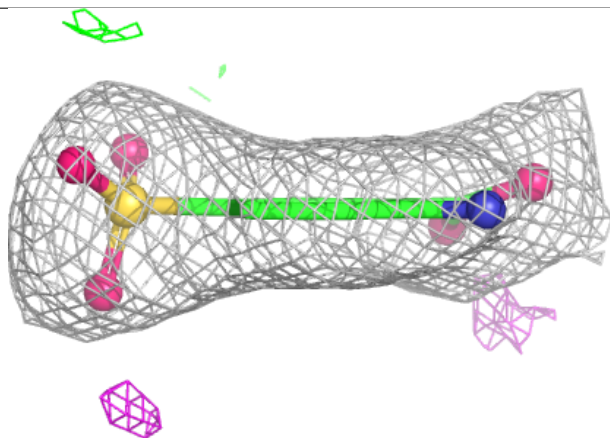
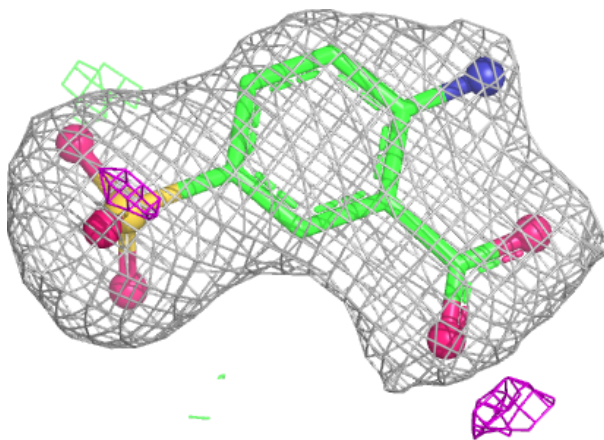


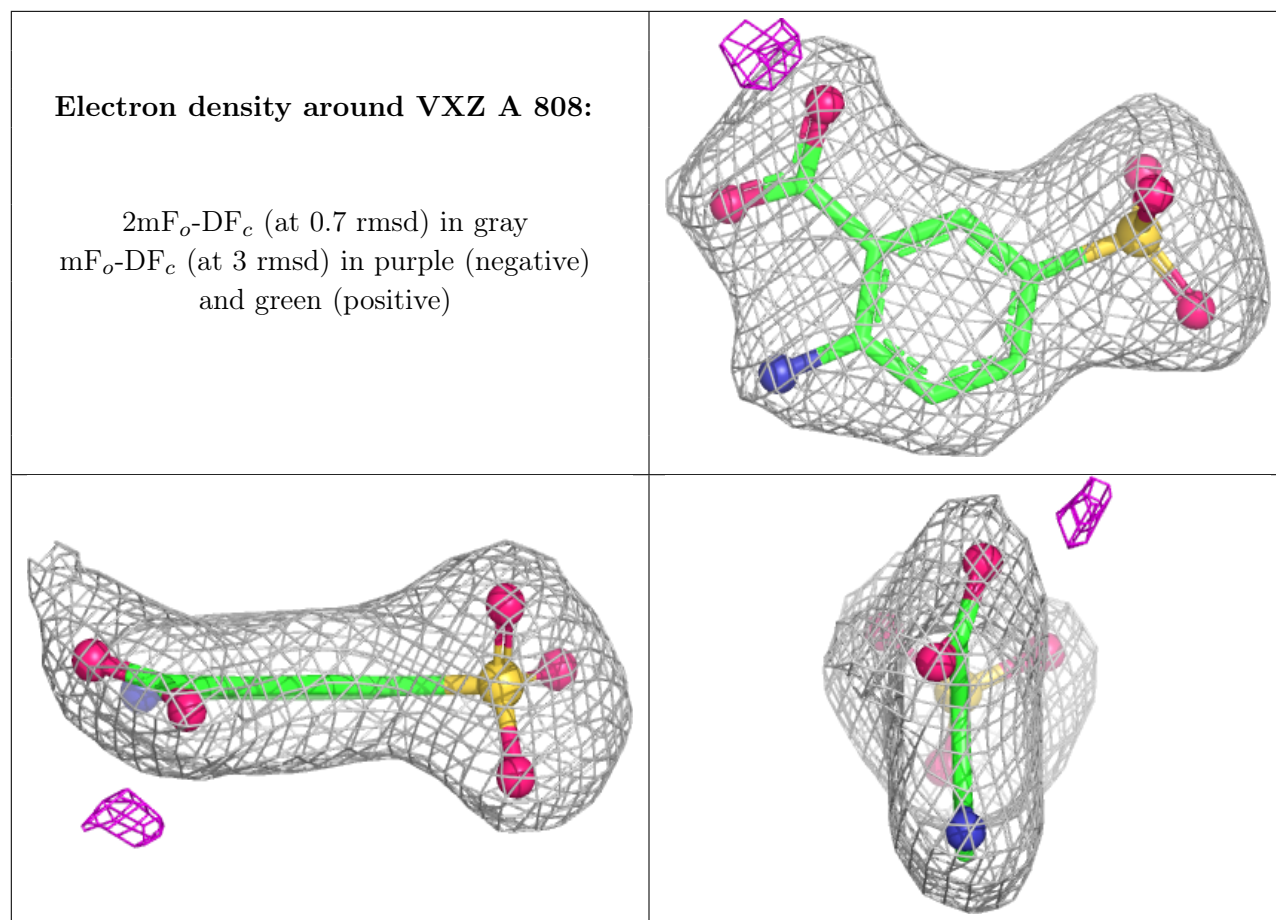
Electron density around VXZ A 807:

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

**Electron density around VXZ B 806:**

$2mF_o-DF_c$ (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.