



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 03:22 AM BST

PDB ID : 6ZA5
Title : M. tuberculosis salicylate synthase MbtI in complex with salicylate and Mg²⁺
Authors : Mori, M.; Villa, S.; Meneghetti, F.; Bellinzoni, M.
Deposited on : 2020-06-04
Resolution : 2.11 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

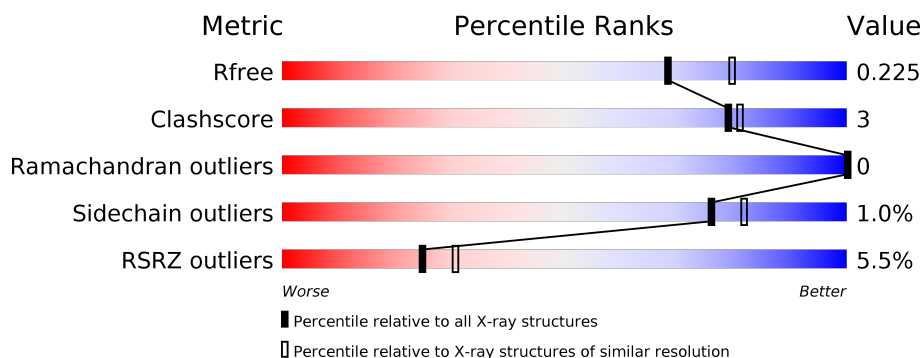
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.11 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	452	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="width: 94%; height: 10px; background-color: green;"></div> <div style="width: 7%; height: 10px; background-color: yellow;"></div> <div style="width: 3%; height: 10px; background-color: grey;"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> % 94% 7% • • </div> </div>
1	B	452	<div> <div style="width: 2%; height: 10px; background-color: red;"></div> <div style="width: 90%; height: 10px; background-color: green;"></div> <div style="width: 7%; height: 10px; background-color: yellow;"></div> <div style="width: 1%; height: 10px; background-color: grey;"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> 2% 90% 7% • </div> </div>
1	C	452	<div> <div style="width: 12%; height: 10px; background-color: red;"></div> <div style="width: 86%; height: 10px; background-color: green;"></div> <div style="width: 10%; height: 10px; background-color: yellow;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> 12% 86% 10% • </div> </div>
1	D	452	<div> <div style="width: 7%; height: 10px; background-color: red;"></div> <div style="width: 91%; height: 10px; background-color: green;"></div> <div style="width: 2%; height: 10px; background-color: yellow;"></div> <div style="width: 1%; height: 10px; background-color: grey;"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> 7% 91% • • • </div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	C	501	-	-	X	-
4	ACT	A	520	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14227 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 Salicylate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	445	Total	C	N	O	S	0	8	0
			3401	2133	612	646	10			
1	B	440	Total	C	N	O	S	0	1	0
			3310	2077	603	620	10			
1	C	435	Total	C	N	O	S	0	6	0
			3245	2042	588	605	10			
1	D	436	Total	C	N	O	S	0	1	0
			3276	2056	594	616	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P9WFX1
A	0	HIS	-	expression tag	UNP P9WFX1
B	-1	GLY	-	expression tag	UNP P9WFX1
B	0	HIS	-	expression tag	UNP P9WFX1
C	-1	GLY	-	expression tag	UNP P9WFX1
C	0	HIS	-	expression tag	UNP P9WFX1
D	-1	GLY	-	expression tag	UNP P9WFX1
D	0	HIS	-	expression tag	UNP P9WFX1

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

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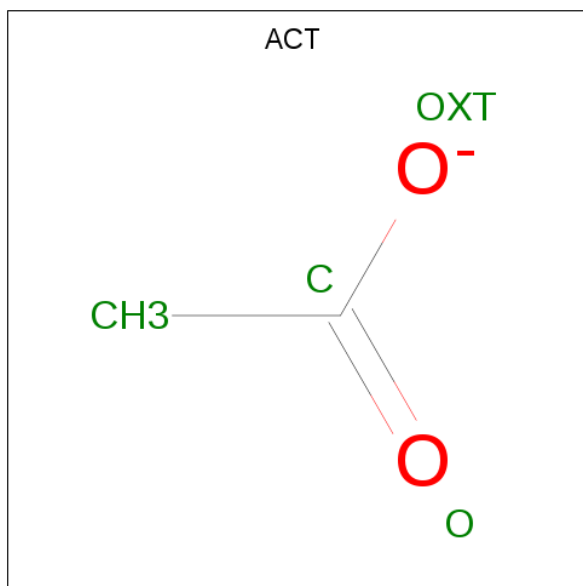
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0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

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by author).

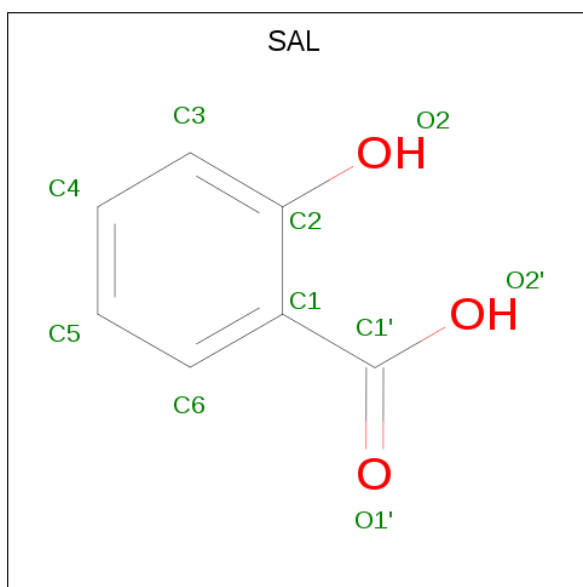
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	A	2	Total	Mg	0	0
			2	2		
3	D	1	Total	Mg	0	0
			1	1		

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is 2-HYDROXYBENZOIC ACID (three-letter code: SAL) (formula: $C_7H_6O_3$) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			10	7	3		
5	D	1	Total	C	O	0	0
			10	7	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	331	Total	O	0	0
			331	331		
6	B	214	Total	O	0	0
			214	214		
6	C	123	Total	O	0	0
			123	123		
6	D	129	Total	O	0	0
			129	129		

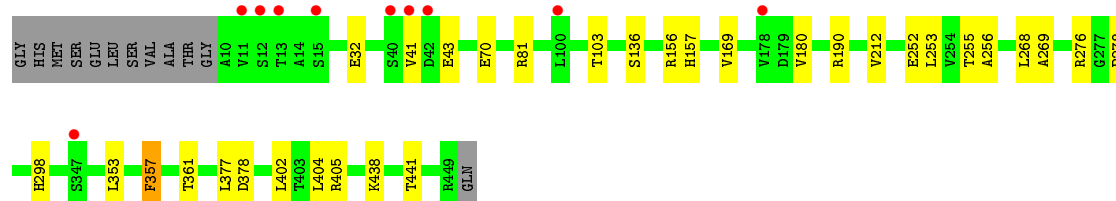
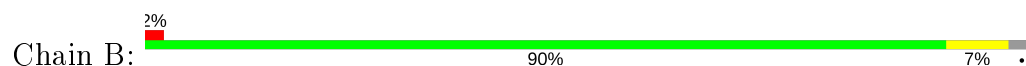
3 Residue-property plots [i](#)

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.

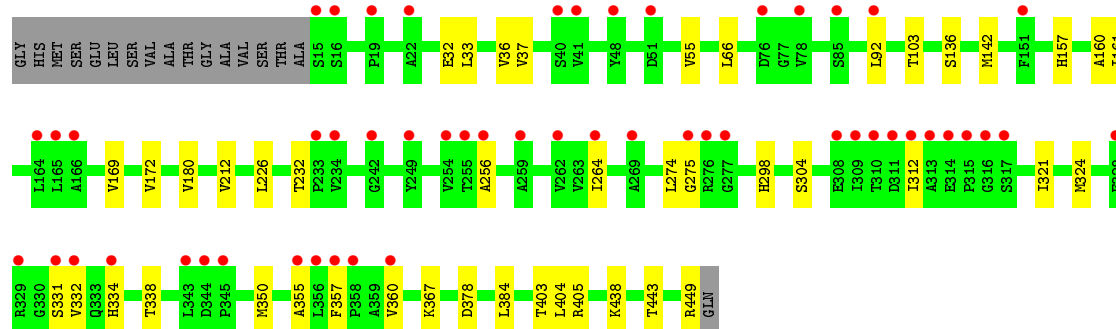
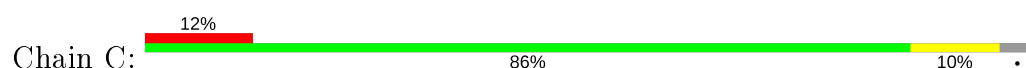
• Molecule 1: Salicylate synthase



• Molecule 1: Salicylate synthase

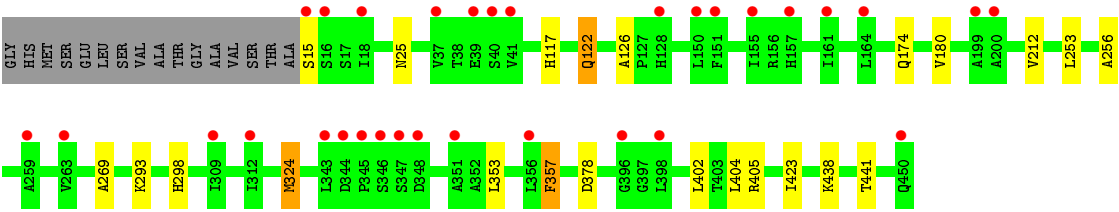


• Molecule 1: Salicylate synthase



• Molecule 1: Salicylate synthase





4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	193.97Å 193.97Å 257.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	154.83 – 2.11 154.83 – 2.11	Depositor EDS
% Data completeness (in resolution range)	93.9 (154.83-2.11) 93.9 (154.83-2.11)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.10Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.195 , 0.217 0.201 , 0.225	Depositor DCC
R_{free} test set	6600 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	43.2	Xtriage
Anisotropy	0.015	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 58.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14227	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 2.05% 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: MG, ACT, SO4, SAL

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	0.44	0/3480	0.59	0/4731
1	B	0.42	0/3371	0.57	0/4584
1	C	0.43	0/3309	0.59	0/4511
1	D	0.37	0/3337	0.56	0/4539
All	All	0.42	0/13497	0.58	0/18365

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	3401	0	3420	13	0
1	B	3310	0	3301	16	0
1	C	3245	0	3174	34	0
1	D	3276	0	3252	13	0
2	A	85	0	0	0	0
2	B	25	0	0	1	0
2	C	20	0	0	2	0
2	D	40	0	0	1	0
3	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
3	D	1	0	0	0	0
4	A	4	0	3	3	0
5	B	10	0	4	0	0
5	D	10	0	4	0	0
6	A	331	0	0	1	0
6	B	214	0	0	0	0
6	C	123	0	0	2	0
6	D	129	0	0	0	0
All	All	14227	0	13158	75	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 (75) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:298:HIS:CD2	1:D:324:MET:HG2	2.24	0.73
1:C:332:VAL:HG23	1:C:334:HIS:HE1	1.55	0.72
1:C:312:ILE:CG1	1:C:355:ALA:HB2	2.21	0.70
1:C:298:HIS:CD2	1:C:324:MET:HG2	2.28	0.69
1:C:312:ILE:HG12	1:C:355:ALA:HB2	1.79	0.64
1:C:332:VAL:HG23	1:C:334:HIS:CE1	2.33	0.62
1:C:304:SER:HB3	1:C:360:VAL:HG21	1.82	0.62
1:B:268:LEU:H	1:B:361:THR:HG21	1.66	0.59
1:C:36:VAL:HG12	1:C:160:ALA:HB1	1.84	0.58
1:C:312:ILE:HG13	1:C:355:ALA:HB2	1.84	0.58
1:A:103:THR:HG21	1:A:136:SER:HB2	1.86	0.58
1:C:37:VAL:HG13	1:C:157:HIS:HD2	1.69	0.57
1:C:274:LEU:HD13	1:C:275:GLY:N	2.19	0.57
1:C:232:THR:CB	1:C:331:SER:H	2.17	0.57
1:B:438:LYS:HD3	2:B:503:SO4:O4	2.07	0.55
1:C:37:VAL:HG13	1:C:157:HIS:CD2	2.42	0.54
1:A:405:ARG:HH21	4:A:520:ACT:H3	1.73	0.54
1:B:103:THR:HG21	1:B:136:SER:HB2	1.90	0.54
1:B:190:ARG:HD2	1:B:377:LEU:O	2.08	0.52
1:A:174:GLN:H	1:A:174:GLN:CD	2.12	0.52
1:C:321:ILE:HG13	1:C:338:THR:HB	1.90	0.52
1:B:353:LEU:O	1:B:357:PHE:HB2	2.10	0.51
1:D:293:LYS:HG2	1:D:423:ILE:HD11	1.90	0.51
1:C:384:LEU:HG	1:C:403:THR:HG21	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:353:LEU:O	1:D:357:PHE:HB2	2.11	0.50
1:C:33:LEU:HD21	1:C:161:ILE:HG12	1.93	0.50
1:C:55:VAL:HG22	1:C:142:MET:HG2	1.94	0.49
1:C:404:LEU:HB3	2:C:501:SO4:O4	2.11	0.49
1:C:103:THR:HG21	1:C:136:SER:HB2	1.92	0.49
1:D:402:LEU:HG	1:D:404:LEU:HG	1.94	0.49
1:C:438:LYS:HD3	2:C:501:SO4:O3	2.13	0.48
1:B:402:LEU:HG	1:B:404:LEU:HG	1.95	0.48
1:D:438:LYS:HD3	2:D:504:SO4:O2	2.13	0.48
1:D:174:GLN:CD	1:D:174:GLN:H	2.15	0.48
1:A:260:ASP:O	1:A:342:ARG:NH1	2.48	0.47
1:D:180:VAL:HG12	1:D:212:VAL:HG11	1.95	0.47
1:A:180:VAL:HG12	1:A:212:VAL:HG11	1.97	0.46
1:A:404:LEU:HD13	4:A:520:ACT:H3	1.96	0.46
1:C:274:LEU:HG	6:C:659:HOH:O	2.15	0.46
1:C:180:VAL:HG11	1:C:443:THR:HG21	1.97	0.46
1:C:274:LEU:HD13	1:C:275:GLY:H	1.80	0.46
1:C:180:VAL:HG12	1:C:212:VAL:HG11	1.98	0.45
1:D:253:LEU:HD13	1:D:256:ALA:HB2	1.98	0.45
1:B:180:VAL:HG12	1:B:212:VAL:HG11	1.97	0.45
1:B:405:ARG:HB3	1:B:441:THR:HG21	1.99	0.45
1:D:405:ARG:HB3	1:D:441:THR:HG21	1.98	0.45
1:C:66:LEU:HD22	1:C:92:LEU:HB2	1.98	0.45
1:B:41:VAL:HG12	1:B:43:GLU:HB3	1.97	0.44
1:A:405:ARG:HB3	1:A:441:THR:HG21	1.99	0.44
1:D:298:HIS:HD2	1:D:324:MET:HG2	1.80	0.44
1:C:312:ILE:HG13	1:C:355:ALA:CB	2.46	0.44
1:C:332:VAL:CG2	1:C:334:HIS:CE1	2.99	0.44
1:C:32:GLU:HG3	1:C:169:VAL:HB	1.98	0.44
1:C:92:LEU:HD23	1:C:350:MET:HE1	2.00	0.43
1:B:278:PRO:HG3	1:D:126:ALA:HB3	2.00	0.43
1:C:256:ALA:HB3	1:C:264:ILE:HG13	1.99	0.43
1:A:191:VAL:HG22	1:A:422:ILE:HD11	2.01	0.43
1:B:253:LEU:HD13	1:B:256:ALA:HB2	1.99	0.43
1:D:269:ALA:HB3	1:D:298:HIS:HA	2.01	0.43
1:A:32:GLU:HG3	1:A:169:VAL:HB	2.01	0.43
1:C:92:LEU:HD23	1:C:350:MET:CE	2.50	0.42
1:A:405:ARG:HH21	4:A:520:ACT:CH3	2.31	0.42
1:A:237:PHE:HE1	6:A:619:HOH:O	2.03	0.42
1:B:269:ALA:HB3	1:B:298:HIS:HA	2.02	0.42
1:C:360:VAL:HB	1:C:367:LYS:HD2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:GLU:HG3	1:B:169:VAL:HB	2.02	0.41
1:C:312:ILE:CG1	1:C:355:ALA:CB	2.97	0.41
1:A:402:LEU:HG	1:A:404:LEU:HG	2.01	0.41
1:B:70:GLU:OE1	1:B:81:ARG:NH2	2.54	0.41
1:B:156:ARG:NH1	1:B:157:HIS:CE1	2.89	0.41
1:D:117:HIS:HA	1:D:122:GLN:HE21	1.85	0.41
1:B:252:GLU:HB3	1:B:268:LEU:HD22	2.02	0.40
1:C:172:VAL:HG11	1:C:226:LEU:HB2	2.03	0.40
1:A:64:VAL:HG23	1:A:99:LEU:HD11	2.04	0.40
1:C:405:ARG:HB3	6:C:624:HOH:O	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	451/452 (100%)	445 (99%)	6 (1%)	0	100	100
1	B	439/452 (97%)	434 (99%)	5 (1%)	0	100	100
1	C	439/452 (97%)	427 (97%)	12 (3%)	0	100	100
1	D	435/452 (96%)	430 (99%)	5 (1%)	0	100	100
All	All	1764/1808 (98%)	1736 (98%)	28 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	356/359 (99%)	355 (100%)	1 (0%)	92	95
1	B	337/359 (94%)	333 (99%)	4 (1%)	71	77
1	C	321/359 (89%)	318 (99%)	3 (1%)	78	84
1	D	333/359 (93%)	327 (98%)	6 (2%)	59	65
All	All	1347/1436 (94%)	1333 (99%)	14 (1%)	76	82

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	378	ASP
1	B	255	THR
1	B	276	ARG
1	B	357	PHE
1	B	378	ASP
1	C	357	PHE
1	C	378	ASP
1	C	449	ARG
1	D	15	SER
1	D	25	ASN
1	D	122	GLN
1	D	324	MET
1	D	357	PHE
1	D	378	ASP

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

Mol	Chain	Res	Type
1	C	157	HIS
1	C	334	HIS
1	D	122	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 41 ligands modelled in this entry, 4 are monoatomic - leaving 37 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	$\# Z > 2$
2	SO4	C	501	-	4,4,4	0.47	0	6,6,6	0.74	0
2	SO4	D	505	-	4,4,4	0.19	0	6,6,6	0.39	0
2	SO4	A	512	-	4,4,4	0.15	0	6,6,6	0.07	0
2	SO4	A	502	-	4,4,4	0.16	0	6,6,6	0.08	0
5	SAL	B	502	3	8,10,10	1.38	1 (12%)	9,13,13	0.46	0
2	SO4	B	506	-	4,4,4	0.17	0	6,6,6	0.04	0
2	SO4	B	507	-	4,4,4	0.16	0	6,6,6	0.08	0
2	SO4	D	504	-	4,4,4	0.11	0	6,6,6	0.30	0
2	SO4	A	506	-	4,4,4	0.16	0	6,6,6	0.09	0
2	SO4	D	507	-	4,4,4	0.13	0	6,6,6	0.07	0
2	SO4	B	504	-	4,4,4	0.16	0	6,6,6	0.36	0
2	SO4	A	501	3	4,4,4	0.17	0	6,6,6	0.51	0
2	SO4	A	508	-	4,4,4	0.13	0	6,6,6	0.36	0
2	SO4	D	506	-	4,4,4	0.16	0	6,6,6	0.15	0
2	SO4	A	513	-	4,4,4	0.13	0	6,6,6	0.08	0
2	SO4	A	503	-	4,4,4	0.31	0	6,6,6	0.38	0
2	SO4	A	507	-	4,4,4	0.45	0	6,6,6	0.49	0
2	SO4	D	509	-	4,4,4	0.15	0	6,6,6	0.05	0
2	SO4	A	505	-	4,4,4	0.18	0	6,6,6	0.36	0
2	SO4	A	515	-	4,4,4	0.17	0	6,6,6	0.26	0
2	SO4	A	509	-	4,4,4	0.15	0	6,6,6	0.12	0
2	SO4	D	503	-	4,4,4	0.29	0	6,6,6	0.19	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	D	508	-	4,4,4	0.14	0	6,6,6	0.06	0
2	SO4	A	511	-	4,4,4	0.22	0	6,6,6	0.22	0
2	SO4	D	510	-	4,4,4	0.13	0	6,6,6	0.07	0
2	SO4	A	516	-	4,4,4	0.15	0	6,6,6	0.07	0
2	SO4	C	504	-	4,4,4	0.14	0	6,6,6	0.08	0
2	SO4	A	517	-	4,4,4	0.15	0	6,6,6	0.04	0
5	SAL	D	502	3	8,10,10	1.32	1 (12%)	9,13,13	0.43	0
2	SO4	A	504	-	4,4,4	0.53	0	6,6,6	0.23	0
2	SO4	C	503	-	4,4,4	0.16	0	6,6,6	0.39	0
2	SO4	A	514	-	4,4,4	0.16	0	6,6,6	0.10	0
2	SO4	B	503	-	4,4,4	0.41	0	6,6,6	0.44	0
2	SO4	A	510	-	4,4,4	0.10	0	6,6,6	0.09	0
2	SO4	C	502	-	4,4,4	0.11	0	6,6,6	0.11	0
4	ACT	A	520	-	1,3,3	5.13	1 (100%)	0,3,3	0.00	-
2	SO4	B	505	-	4,4,4	0.12	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	SAL	B	502	3	-	0/0/4/4	0/1/1/1
5	SAL	D	502	3	-	0/0/4/4	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	520	ACT	CH3-C	5.13	1.55	1.48
5	B	502	SAL	C1-C1'	3.81	1.51	1.47
5	D	502	SAL	C1-C1'	3.64	1.51	1.47

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 7 short contacts:

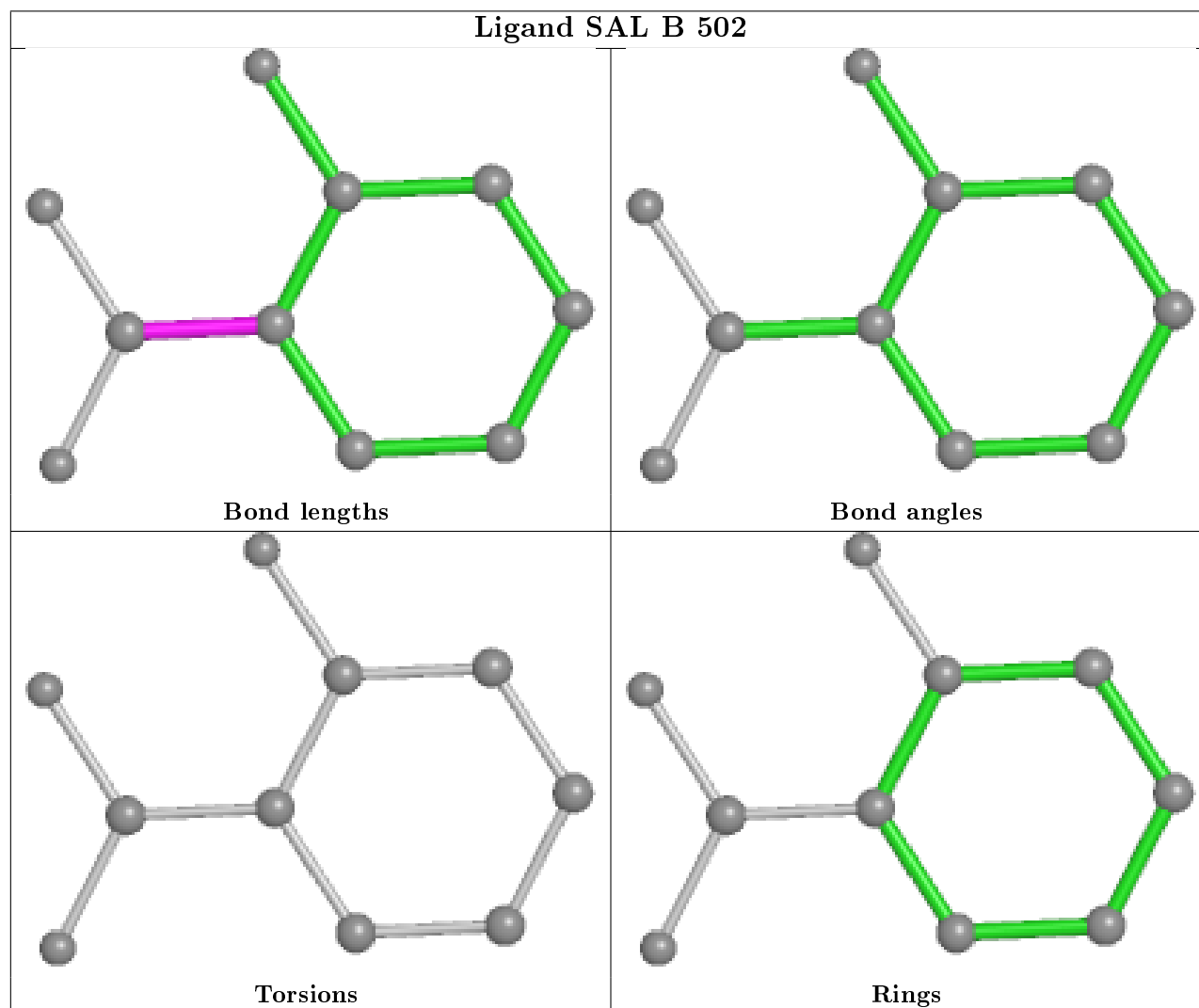
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	501	SO4	2	0

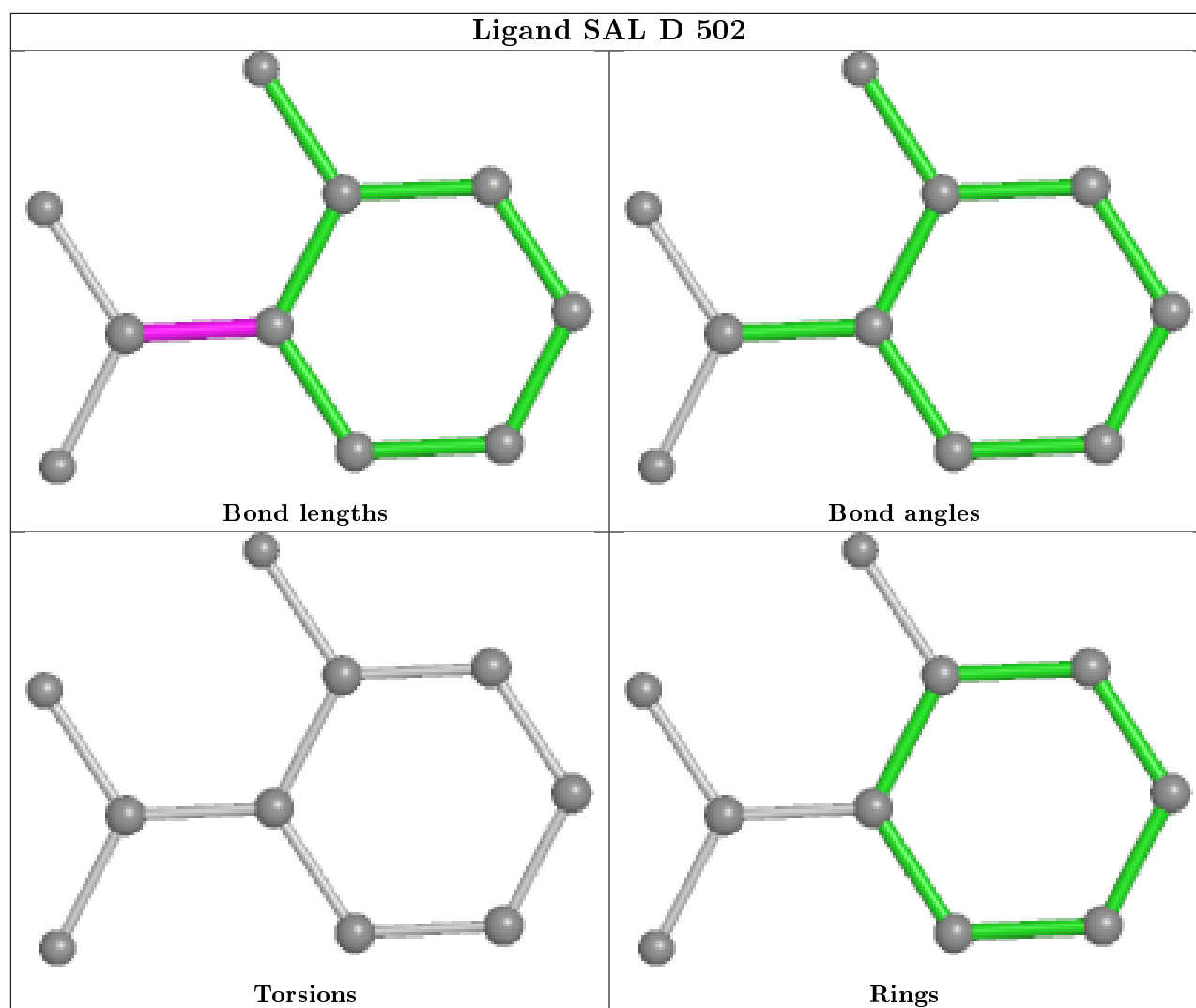
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	504	SO4	1	0
2	B	503	SO4	1	0
4	A	520	ACT	3	0

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.





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	445/452 (98%)	0.27	3 (0%) 87 89	28, 38, 56, 79	0
1	B	440/452 (97%)	0.18	10 (2%) 60 65	32, 45, 63, 84	0
1	C	435/452 (96%)	0.87	53 (12%) 4 5	33, 60, 90, 101	0
1	D	436/452 (96%)	0.55	31 (7%) 16 20	36, 59, 85, 97	0
All	All	1756/1808 (97%)	0.46	97 (5%) 25 31	28, 48, 84, 101	0

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	355	ALA	8.0
1	C	234	VAL	6.4
1	C	312	ILE	5.8
1	D	347	SER	5.6
1	C	310	THR	5.3
1	C	332	VAL	5.1
1	C	40	SER	4.8
1	D	18	ILE	4.7
1	B	41	VAL	4.7
1	C	41	VAL	4.5
1	C	315[A]	PRO	4.4
1	C	275	GLY	4.4
1	C	15	SER	4.4
1	C	22	ALA	4.3
1	C	254	VAL	4.3
1	D	200	ALA	4.3
1	B	40	SER	4.3
1	D	396	GLY	4.1
1	D	15	SER	4.0
1	C	76	ASP	3.9
1	D	344	ASP	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	329	ARG	3.9
1	C	356	LEU	3.8
1	C	345	PRO	3.6
1	D	346	SER	3.4
1	C	343	LEU	3.4
1	C	48	TYR	3.3
1	D	151	PHE	3.3
1	C	269	ALA	3.3
1	C	233	PRO	3.2
1	C	334	HIS	3.2
1	C	78	VAL	3.2
1	C	316[A]	GLY	3.2
1	C	309	ILE	3.1
1	C	16	SER	3.1
1	D	37	VAL	3.1
1	D	150	LEU	3.0
1	D	199	ALA	3.0
1	B	15	SER	3.0
1	D	161	ILE	2.9
1	C	262	VAL	2.9
1	C	264	ILE	2.9
1	D	40	SER	2.9
1	D	351	ALA	2.9
1	C	344	ASP	2.8
1	A	40	SER	2.8
1	D	16	SER	2.8
1	C	51	ASP	2.7
1	D	343	LEU	2.7
1	A	37	VAL	2.7
1	D	259	ALA	2.6
1	C	151	PHE	2.6
1	C	314[A]	GLU	2.6
1	C	331	SER	2.6
1	D	157	HIS	2.6
1	C	92	LEU	2.6
1	D	348	ASP	2.5
1	D	155	ILE	2.5
1	B	13	THR	2.5
1	B	11	VAL	2.4
1	C	357	PHE	2.4
1	D	39	GLU	2.4
1	C	164	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	263	VAL	2.3
1	C	256	ALA	2.3
1	D	345	PRO	2.3
1	C	360	VAL	2.3
1	C	166	ALA	2.3
1	C	313[A]	ALA	2.3
1	C	328	GLU	2.3
1	B	347	SER	2.3
1	D	41	VAL	2.3
1	C	277	GLY	2.3
1	C	308	GLU	2.3
1	C	85	SER	2.2
1	C	317	SER	2.2
1	D	398	LEU	2.2
1	D	450	GLN	2.2
1	C	19	PRO	2.2
1	C	311	ASP	2.2
1	D	312	ILE	2.2
1	C	242	GLY	2.2
1	C	165	LEU	2.2
1	D	164	LEU	2.1
1	D	356	LEU	2.1
1	B	100	LEU	2.1
1	C	255	THR	2.1
1	C	259	ALA	2.1
1	C	358	PRO	2.1
1	D	128	HIS	2.1
1	B	12	SER	2.1
1	C	276	ARG	2.1
1	B	178	VAL	2.1
1	C	249	TYR	2.0
1	B	42	ASP	2.0
1	D	309	ILE	2.0
1	A	229	ARG	2.0

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

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

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	A	511	5/5	0.73	0.26	105,105,106,106	0
2	SO4	A	513	5/5	0.77	0.21	143,143,143,143	0
2	SO4	D	509	5/5	0.79	0.12	125,125,125,125	0
2	SO4	B	507	5/5	0.82	0.18	127,127,127,127	0
2	SO4	C	502	5/5	0.83	0.14	121,121,122,122	0
2	SO4	D	510	5/5	0.84	0.37	157,157,157,157	0
2	SO4	D	508	5/5	0.84	0.25	144,144,144,144	0
2	SO4	D	507	5/5	0.85	0.14	121,121,121,121	0
2	SO4	A	517	5/5	0.87	0.14	136,136,136,136	0
2	SO4	A	509	5/5	0.87	0.17	116,116,116,116	0
2	SO4	B	505	5/5	0.89	0.11	105,106,106,106	0
5	SAL	D	502	10/10	0.90	0.25	52,52,53,55	0
2	SO4	B	506	5/5	0.90	0.11	147,147,147,147	0
2	SO4	A	510	5/5	0.91	0.19	119,119,119,119	0
5	SAL	B	502	10/10	0.92	0.21	56,59,59,60	0
2	SO4	A	516	5/5	0.92	0.08	113,113,113,113	0
2	SO4	A	505	5/5	0.92	0.17	64,65,65,66	5
2	SO4	A	504	5/5	0.92	0.14	54,54,54,54	5
2	SO4	A	512	5/5	0.92	0.08	128,128,128,128	0
2	SO4	A	508	5/5	0.92	0.34	84,85,85,85	0
4	ACT	A	520	4/4	0.92	0.19	35,37,38,39	0
2	SO4	A	502	5/5	0.92	0.11	109,109,109,109	0
3	MG	D	501	1/1	0.93	0.10	50,50,50,50	0
2	SO4	D	506	5/5	0.94	0.22	88,89,89,89	0
2	SO4	A	506	5/5	0.94	0.13	97,97,97,97	0
2	SO4	A	507	5/5	0.94	0.27	59,59,60,60	5
2	SO4	A	514	5/5	0.94	0.21	105,105,105,105	0
2	SO4	C	503	5/5	0.95	0.22	36,36,37,38	5
3	MG	B	501	1/1	0.95	0.06	55,55,55,55	0
2	SO4	A	515	5/5	0.95	0.16	86,86,86,86	0
2	SO4	A	503	5/5	0.96	0.14	55,56,57,57	0
2	SO4	C	501	5/5	0.96	0.18	44,45,46,46	5

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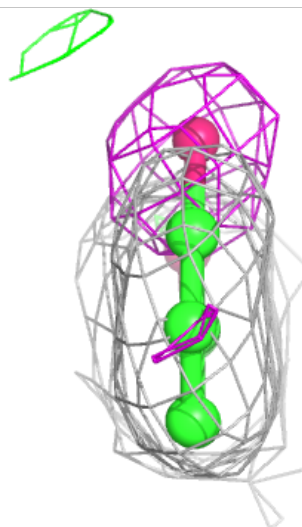
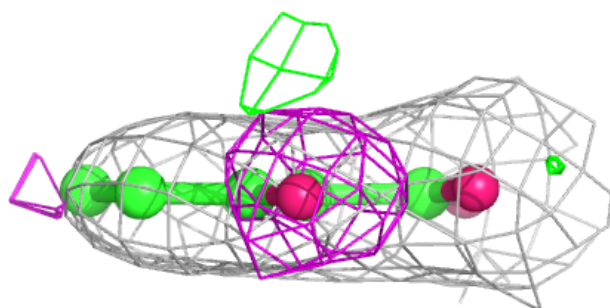
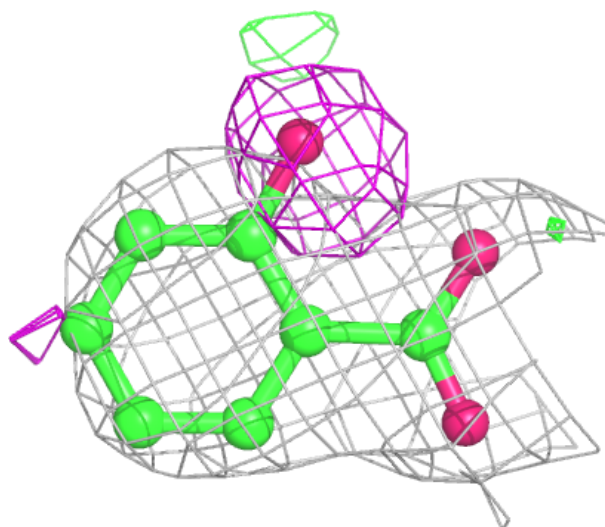
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	SO4	C	504	5/5	0.96	0.12	114,114,114,114	0
2	SO4	A	501	5/5	0.97	0.14	37,39,39,39	5
2	SO4	D	504	5/5	0.98	0.16	57,57,58,58	0
3	MG	A	519	1/1	0.98	0.15	36,36,36,36	0
2	SO4	B	504	5/5	0.98	0.13	68,68,69,69	0
2	SO4	D	505	5/5	0.98	0.13	51,51,52,52	5
2	SO4	B	503	5/5	0.99	0.18	45,45,45,45	5
2	SO4	D	503	5/5	0.99	0.15	58,58,59,59	0
3	MG	A	518	1/1	0.99	0.06	44,44,44,44	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.

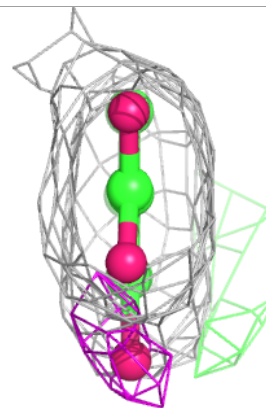
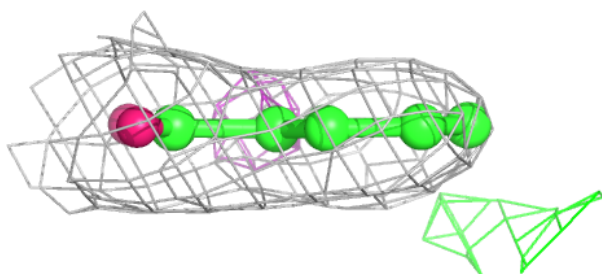
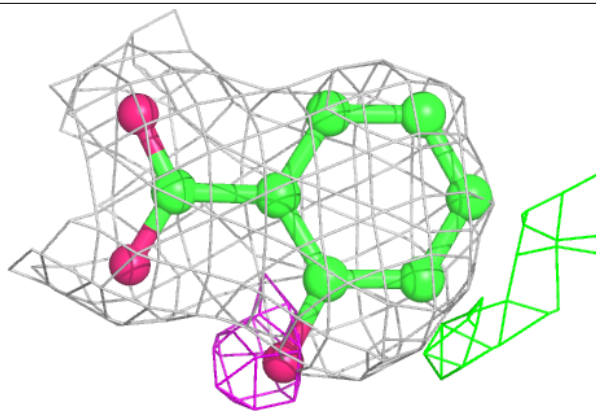
Electron density around SAL D 502:

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



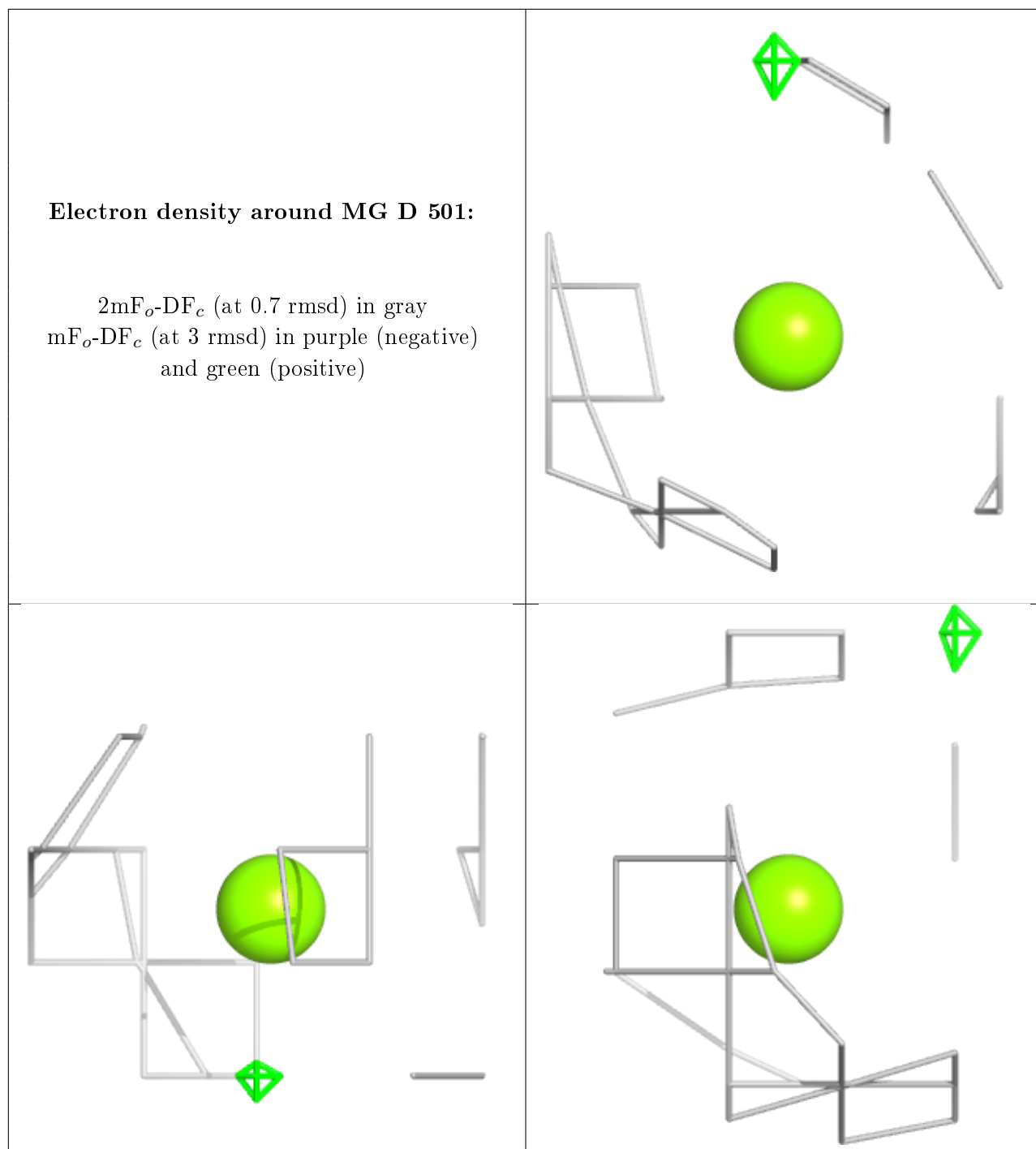
Electron density around SAL B 502:

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



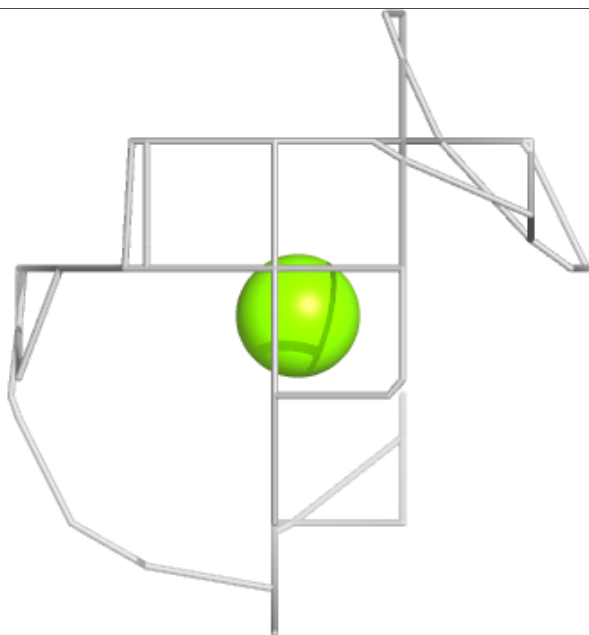
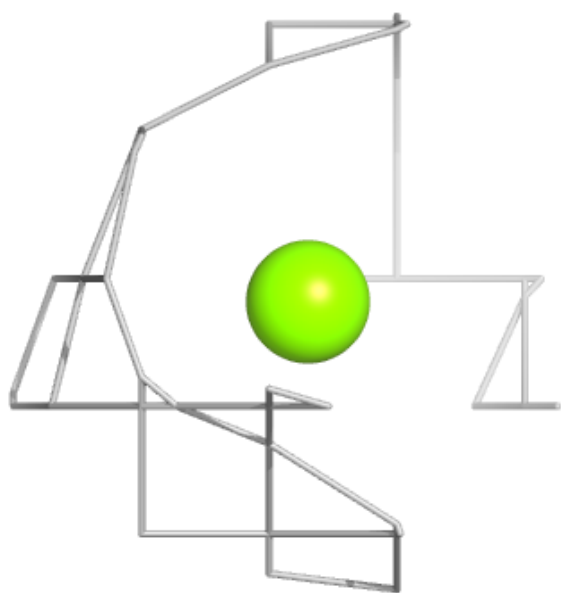
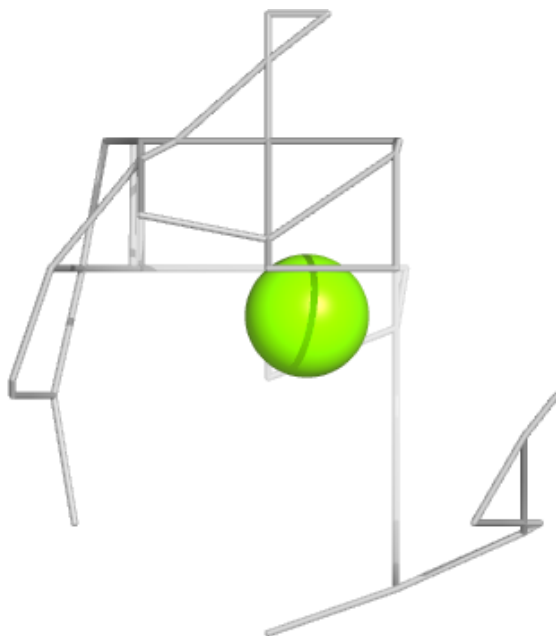
Electron density around MG D 501:

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



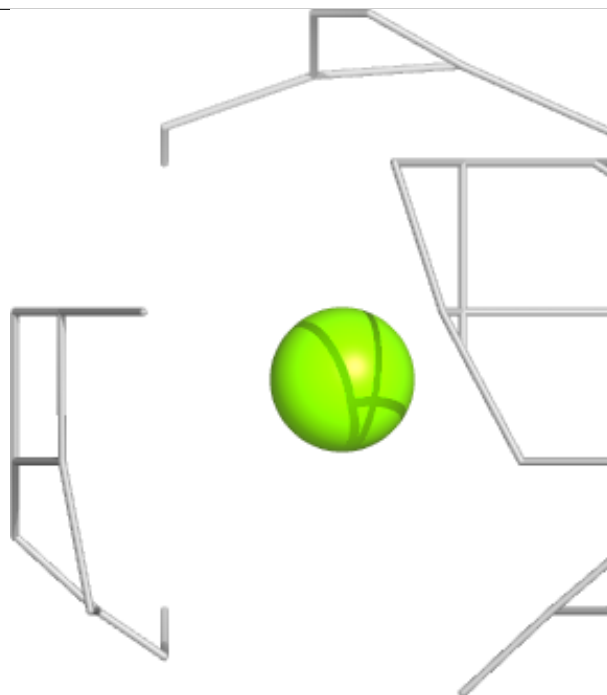
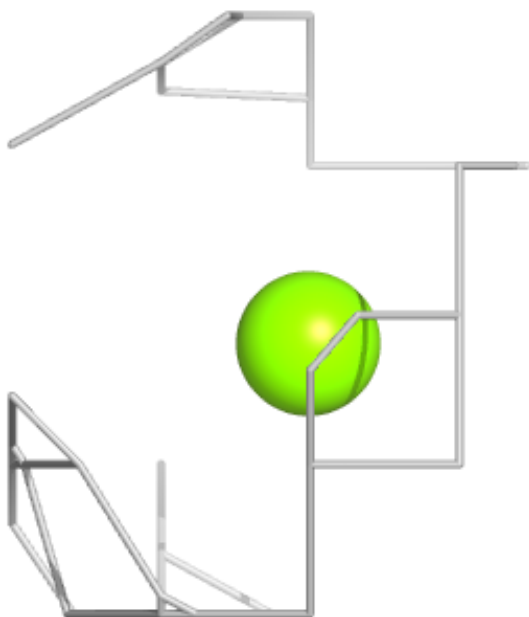
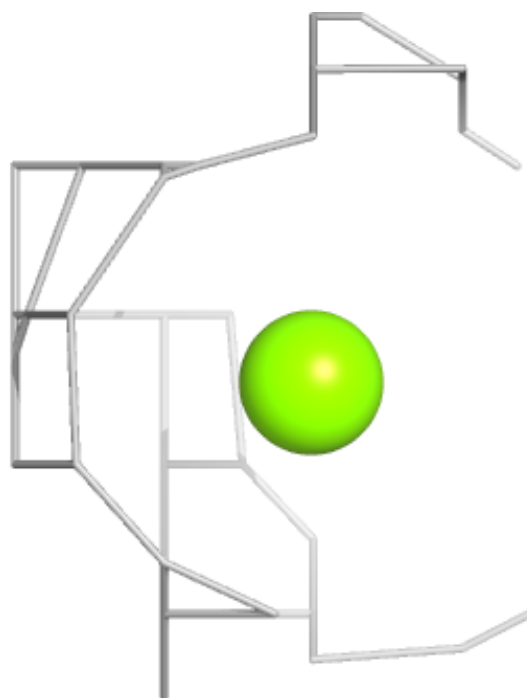
Electron density around MG B 501:

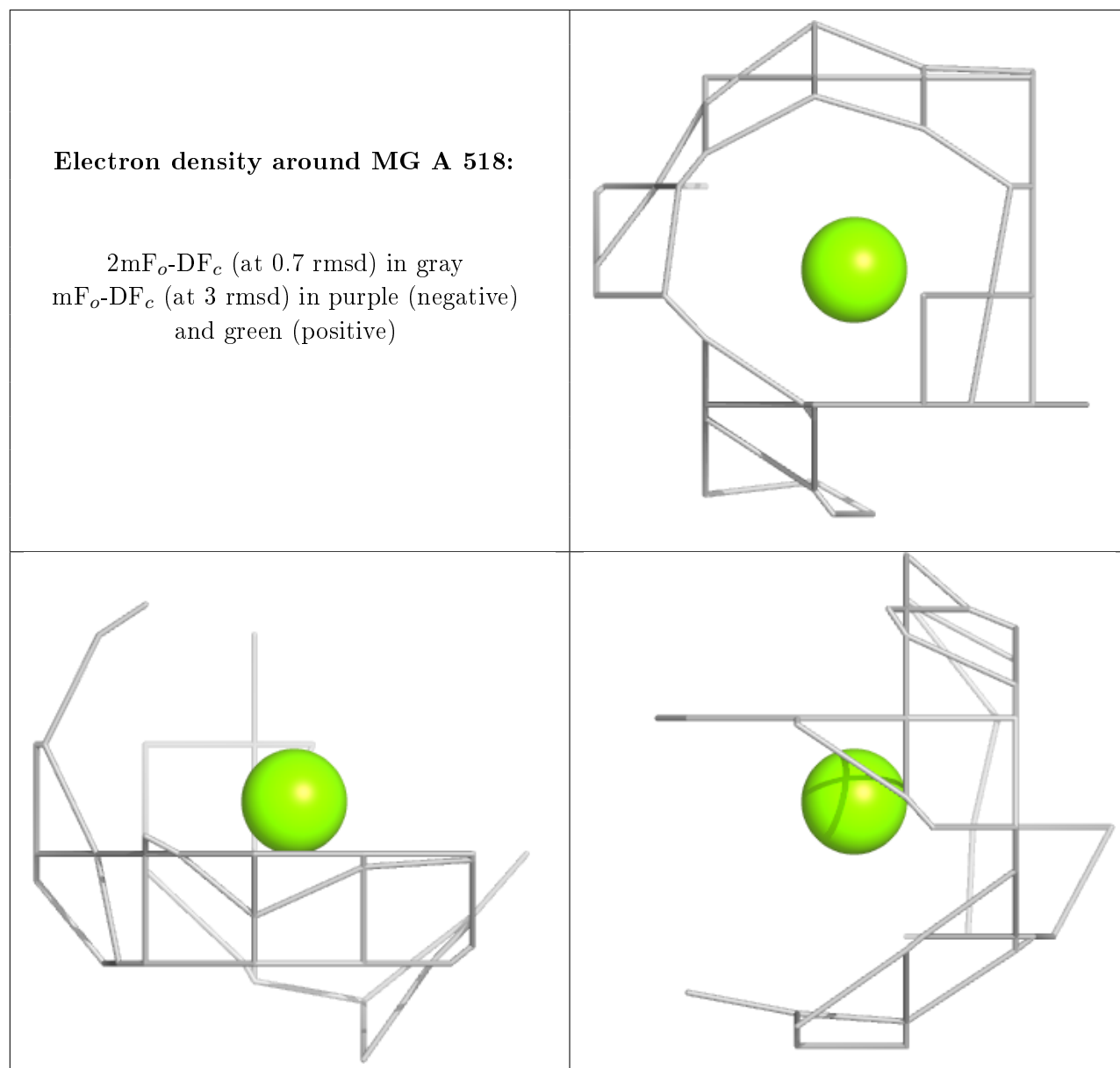
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around MG A 519:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.