



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 09:46 pm BST

PDB ID : 6TM4
Title : NatL2 in complex with two molecules of salicylic acid
Authors : Naismith, J.H.; Song, H.
Deposited on : 2019-12-03
Resolution : 1.89 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

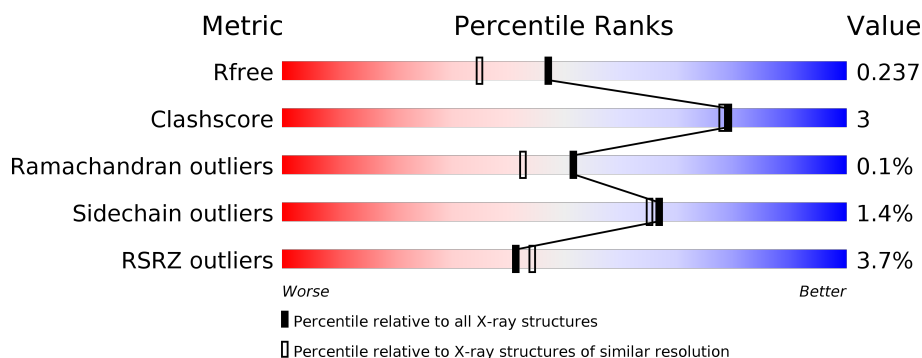
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	AAA	436	<div> <div>5%</div> <div> <div></div> <div>89%</div> <div>9%</div> <div>..</div> </div> </div>
2	BBB	437	<div> <div>3%</div> <div> <div></div> <div>91%</div> <div>7%</div> <div>.</div> </div> </div>

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 6954 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 PaaK-like ligase (AMP-dependent synthetase and ligase).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	430	Total	C	N	O	S	0	0	0
			3354	2120	594	627	13			

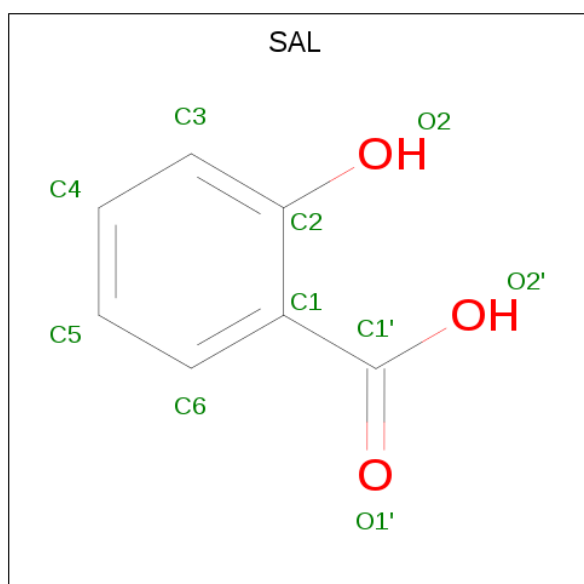
- Molecule 2 is a protein called PaaK-like ligase (AMP-dependent synthetase and ligase).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	BBB	432	Total	C	N	O	S	0	0	0
			3373	2131	599	629	14			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
BBB	0	ALA	-	expression tag	UNP A0A022MRT4

- Molecule 3 is 2-HYDROXYBENZOIC ACID (three-letter code: SAL) (formula: C₇H₆O₃) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	AAA	1	Total C O 10 7 3	0	0
3	AAA	1	Total C O 10 7 3	0	0
3	BBB	1	Total C O 10 7 3	0	0
3	BBB	1	Total C O 10 7 3	0	0

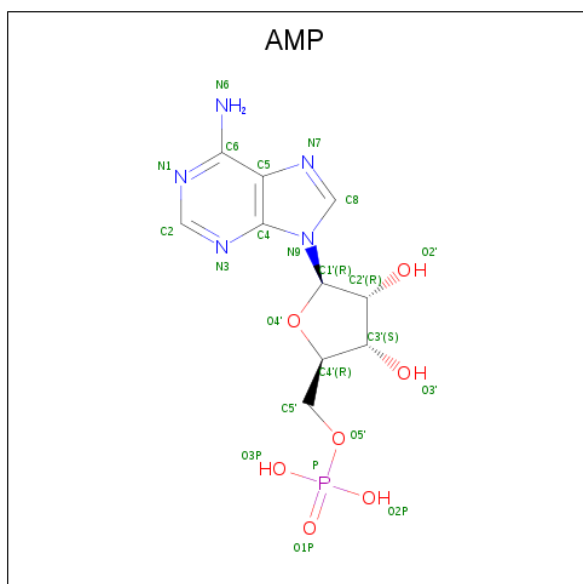
- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	BBB	1	Total Zn 1 1	0	0
4	AAA	1	Total Zn 1 1	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	BBB	1	Total Mg 1 1	0	0
5	AAA	1	Total Mg 1 1	0	0

- Molecule 6 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula: C₁₀H₁₄N₅O₇P) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	BBB	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

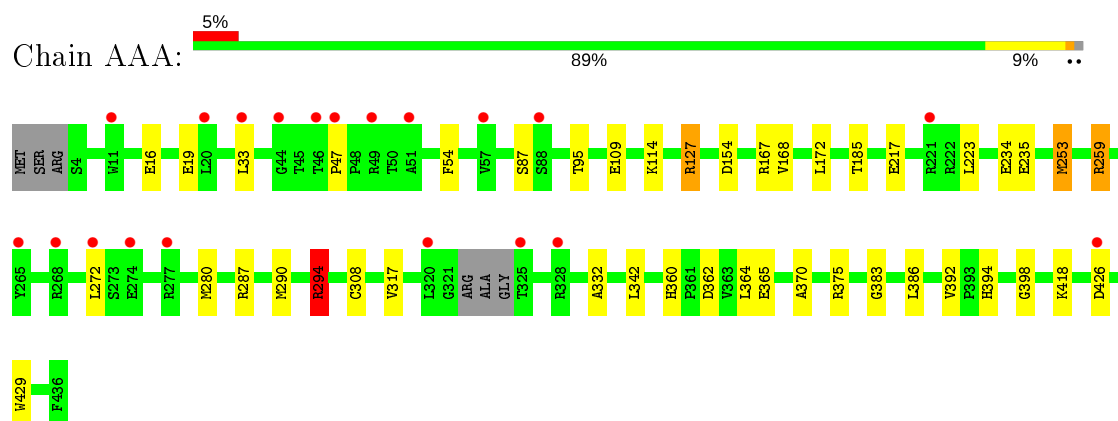
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	AAA	68	Total	O	0	0
			68	68		
7	BBB	92	Total	O	0	0
			92	92		

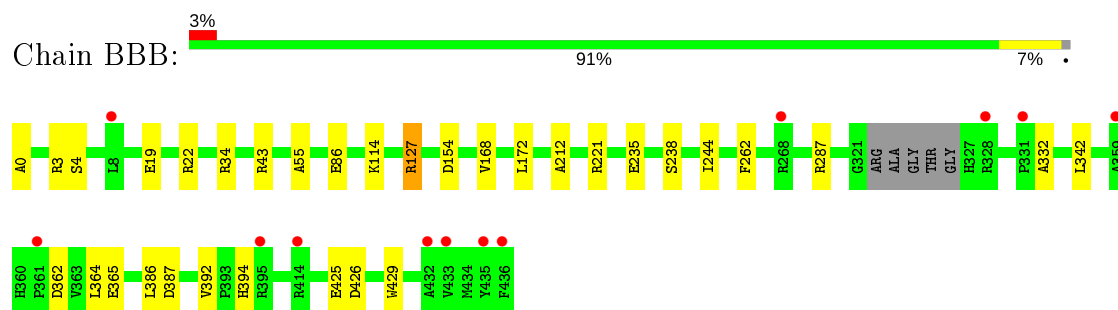
3 Residue-property plots [i](#)

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

- Molecule 1: PaaK-like ligase (AMP-dependent synthetase and ligase)



- Molecule 2: PaaK-like ligase (AMP-dependent synthetase and ligase)



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	107.46Å 139.24Å 129.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	71.13 – 1.89 71.13 – 1.89	Depositor EDS
% Data completeness (in resolution range)	99.9 (71.13-1.89) 99.9 (71.13-1.89)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.198 , 0.230 0.209 , 0.237	Depositor DCC
R_{free} test set	3811 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.0	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	0.96	EDS
Total number of atoms	6954	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.22% 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: AMP, ZN, MG, 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	AAA	0.87	8/3437 (0.2%)	0.87	4/4686 (0.1%)
2	BBB	0.88	6/3456 (0.2%)	0.89	2/4710 (0.0%)
All	All	0.87	14/6893 (0.2%)	0.88	6/9396 (0.1%)

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	BBB	19	GLU	CD-OE2	15.32	1.42	1.25
1	AAA	19	GLU	CD-OE2	12.55	1.39	1.25
2	BBB	19	GLU	CD-OE1	12.53	1.39	1.25
1	AAA	362	ASP	CG-OD1	10.46	1.49	1.25
1	AAA	383	GLY	C-O	9.64	1.39	1.23
1	AAA	19	GLU	CD-OE1	9.10	1.35	1.25
2	BBB	365	GLU	CD-OE2	7.87	1.34	1.25
2	BBB	55	ALA	C-O	6.64	1.35	1.23
1	AAA	308	CYS	C-O	6.55	1.35	1.23
1	AAA	16	GLU	CD-OE1	6.09	1.32	1.25
1	AAA	16	GLU	CD-OE2	5.94	1.32	1.25
2	BBB	362	ASP	CG-OD1	5.85	1.38	1.25
1	AAA	365	GLU	CG-CD	5.42	1.60	1.51
2	BBB	22	ARG	C-O	5.36	1.33	1.23

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AAA	362	ASP	CB-CG-OD1	-8.28	110.85	118.30
1	AAA	294	ARG	CG-CD-NE	8.14	128.90	111.80
2	BBB	34	ARG	NE-CZ-NH1	6.70	123.65	120.30
2	BBB	387	ASP	CB-CG-OD2	-6.03	112.87	118.30
1	AAA	19	GLU	OE1-CD-OE2	5.45	129.84	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AAA	259	ARG	NE-CZ-NH2	-5.05	117.77	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	AAA	3354	0	3298	22	0
2	BBB	3373	0	3323	18	0
3	AAA	20	0	9	1	0
3	BBB	20	0	8	2	0
4	AAA	1	0	0	0	0
4	BBB	1	0	0	0	0
5	AAA	1	0	0	0	0
5	BBB	1	0	0	0	0
6	BBB	23	0	12	3	0
7	AAA	68	0	0	0	0
7	BBB	92	0	0	0	0
All	All	6954	0	6650	38	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 (38) 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:AAA:109:GLU:OE1	1:AAA:259:ARG:NH2	2.16	0.78
1:AAA:54:PHE:CZ	1:AAA:294:ARG:HD2	2.32	0.65
1:AAA:33:LEU:HD11	1:AAA:47:PRO:HB2	1.77	0.64
1:AAA:287:ARG:HD3	1:AAA:290:MET:O	1.99	0.62
1:AAA:342:LEU:HD13	1:AAA:386:LEU:HD12	1.86	0.57
2:BBB:332:ALA:HB3	2:BBB:392:VAL:HG13	1.91	0.53
2:BBB:114:LYS:NZ	2:BBB:235:GLU:OE1	2.39	0.53
1:AAA:114:LYS:NZ	1:AAA:235:GLU:OE1	2.34	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BBB:168:VAL:HG13	2:BBB:172:LEU:HD12	1.92	0.52
1:AAA:418:LYS:HD3	2:BBB:212:ALA:HB1	1.92	0.51
1:AAA:332:ALA:HB3	1:AAA:392:VAL:CG1	2.41	0.51
2:BBB:127:ARG:HD3	2:BBB:154:ASP:O	2.10	0.51
1:AAA:418:LYS:HD2	2:BBB:212:ALA:O	2.12	0.50
1:AAA:370:ALA:O	1:AAA:375:ARG:NH1	2.45	0.49
1:AAA:360:HIS:ND1	2:BBB:425:GLU:HG3	2.28	0.48
6:BBB:501:AMP:H5'2	3:BBB:503:SAL:O1'	2.13	0.47
1:AAA:168:VAL:HG13	1:AAA:172:LEU:HD12	1.97	0.47
1:AAA:332:ALA:HB3	1:AAA:392:VAL:HG13	1.96	0.47
1:AAA:234:GLU:CD	1:AAA:253:MET:HE1	2.35	0.47
1:AAA:280:MET:CE	1:AAA:317:VAL:HG21	2.44	0.47
2:BBB:212:ALA:HB2	3:BBB:502:SAL:C4	2.45	0.47
2:BBB:332:ALA:HB3	2:BBB:392:VAL:CG1	2.45	0.46
2:BBB:244:ILE:HB	2:BBB:262:PHE:CE1	2.52	0.45
1:AAA:375:ARG:HD2	1:AAA:398:GLY:C	2.38	0.44
2:BBB:364:LEU:O	2:BBB:394:HIS:HA	2.18	0.44
2:BBB:238:SER:HA	6:BBB:501:AMP:H5'1	2.00	0.44
2:BBB:287:ARG:HG3	2:BBB:287:ARG:O	2.17	0.44
2:BBB:221:ARG:HH21	2:BBB:221:ARG:HG2	1.83	0.43
1:AAA:127:ARG:HD3	1:AAA:154:ASP:O	2.19	0.43
3:AAA:501:SAL:O1'	3:AAA:501:SAL:O2	2.31	0.43
1:AAA:185:THR:HG23	1:AAA:223:LEU:CD2	2.49	0.42
2:BBB:0:ALA:HB1	2:BBB:3:ARG:O	2.19	0.42
1:AAA:167:ARG:NH2	2:BBB:86:GLU:OE1	2.53	0.42
1:AAA:87:SER:HB2	1:AAA:95:THR:HB	2.00	0.42
1:AAA:272:LEU:HD13	1:AAA:294:ARG:NH1	2.35	0.41
2:BBB:342:LEU:HD13	2:BBB:386:LEU:HD12	2.02	0.40
1:AAA:364:LEU:O	1:AAA:394:HIS:HA	2.20	0.40
2:BBB:212:ALA:HA	6:BBB:501:AMP:C8	2.56	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	AAA	426/436 (98%)	411 (96%)	15 (4%)	0	100	100
2	BBB	428/437 (98%)	415 (97%)	12 (3%)	1 (0%)	47	38
All	All	854/873 (98%)	826 (97%)	27 (3%)	1 (0%)	51	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	BBB	43	ARG

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	AAA	346/350 (99%)	340 (98%)	6 (2%)	60	57
2	BBB	348/350 (99%)	344 (99%)	4 (1%)	73	73
All	All	694/700 (99%)	684 (99%)	10 (1%)	67	65

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	127	ARG
1	AAA	217	GLU
1	AAA	253	MET
1	AAA	294	ARG
1	AAA	426	ASP
1	AAA	429	TRP
2	BBB	4	SER
2	BBB	127	ARG
2	BBB	426	ASP
2	BBB	429	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no

such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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$
6	AMP	BBB	501	-	22,25,25	0.59	0	25,38,38	1.14	3 (12%)
3	SAL	AAA	501	-	8,10,10	1.71	1 (12%)	9,13,13	1.36	1 (11%)
3	SAL	BBB	502	-	8,10,10	1.19	1 (12%)	9,13,13	1.12	1 (11%)
3	SAL	BBB	503	-	8,10,10	1.37	2 (25%)	9,13,13	2.05	3 (33%)
3	SAL	AAA	502	-	8,10,10	1.61	2 (25%)	9,13,13	1.02	1 (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
6	AMP	BBB	501	-	-	2/6/26/26	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SAL	AAA	501	-	-	0/0/4/4	0/1/1/1
3	SAL	BBB	502	-	-	0/0/4/4	0/1/1/1
3	SAL	BBB	503	-	-	0/0/4/4	0/1/1/1
3	SAL	AAA	502	-	-	0/0/4/4	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	AAA	501	SAL	C1-C1'	3.98	1.51	1.47
3	AAA	502	SAL	C1-C1'	3.54	1.50	1.47
3	BBB	503	SAL	C1-C2	2.53	1.47	1.40
3	BBB	502	SAL	C1-C2	2.36	1.47	1.40
3	AAA	502	SAL	C1-C2	2.31	1.47	1.40
3	BBB	503	SAL	O2-C2	-2.10	1.32	1.36

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	BBB	503	SAL	C4-C3-C2	-3.70	115.36	120.05
3	BBB	503	SAL	C5-C6-C1	-2.91	115.42	120.33
3	BBB	503	SAL	C4-C5-C6	2.87	124.56	120.19
3	AAA	501	SAL	C6-C1-C1'	2.86	124.46	120.20
6	BBB	501	AMP	C5-C6-N6	2.83	124.65	120.35
6	BBB	501	AMP	O3P-P-O5'	-2.76	99.38	106.73
3	BBB	502	SAL	C6-C1-C2	2.30	121.74	117.45
3	AAA	502	SAL	C6-C1-C1'	2.05	123.26	120.20
6	BBB	501	AMP	P-O5'-C5'	2.03	123.89	118.30

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	BBB	501	AMP	C3'-C4'-C5'-O5'
6	BBB	501	AMP	O4'-C4'-C5'-O5'

There are no ring outliers.

4 monomers are involved in 5 short contacts:

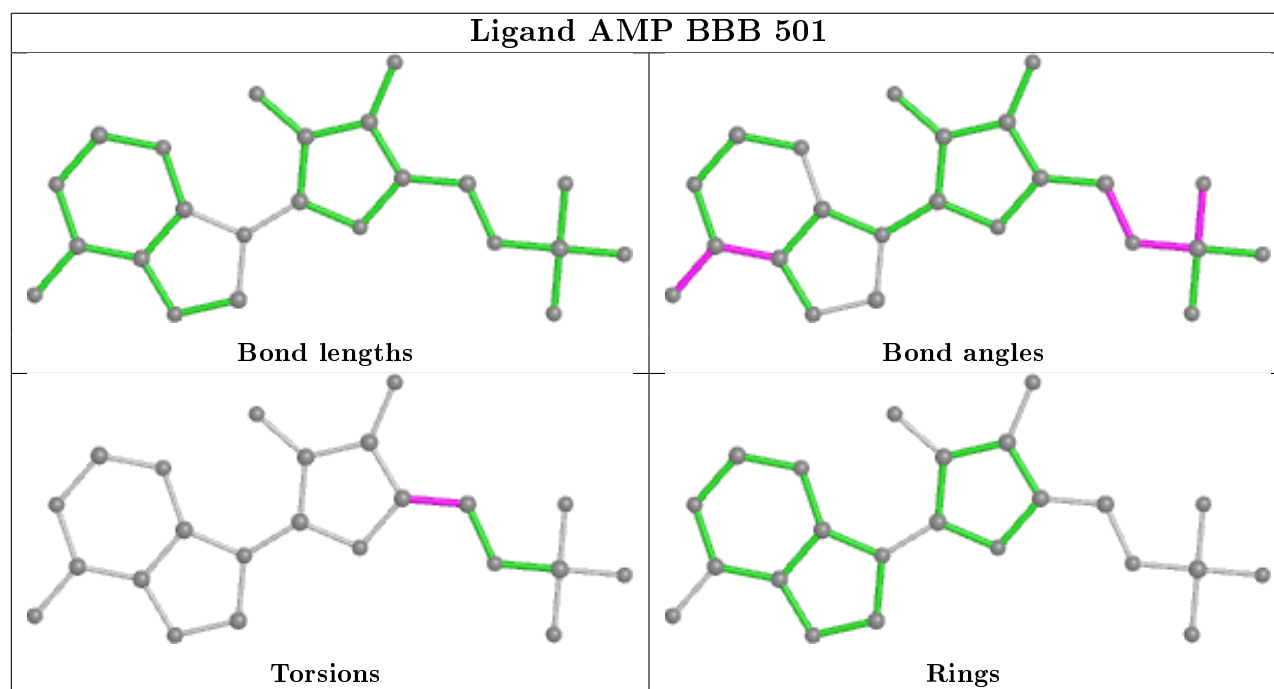
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	BBB	501	AMP	3	0
3	AAA	501	SAL	1	0

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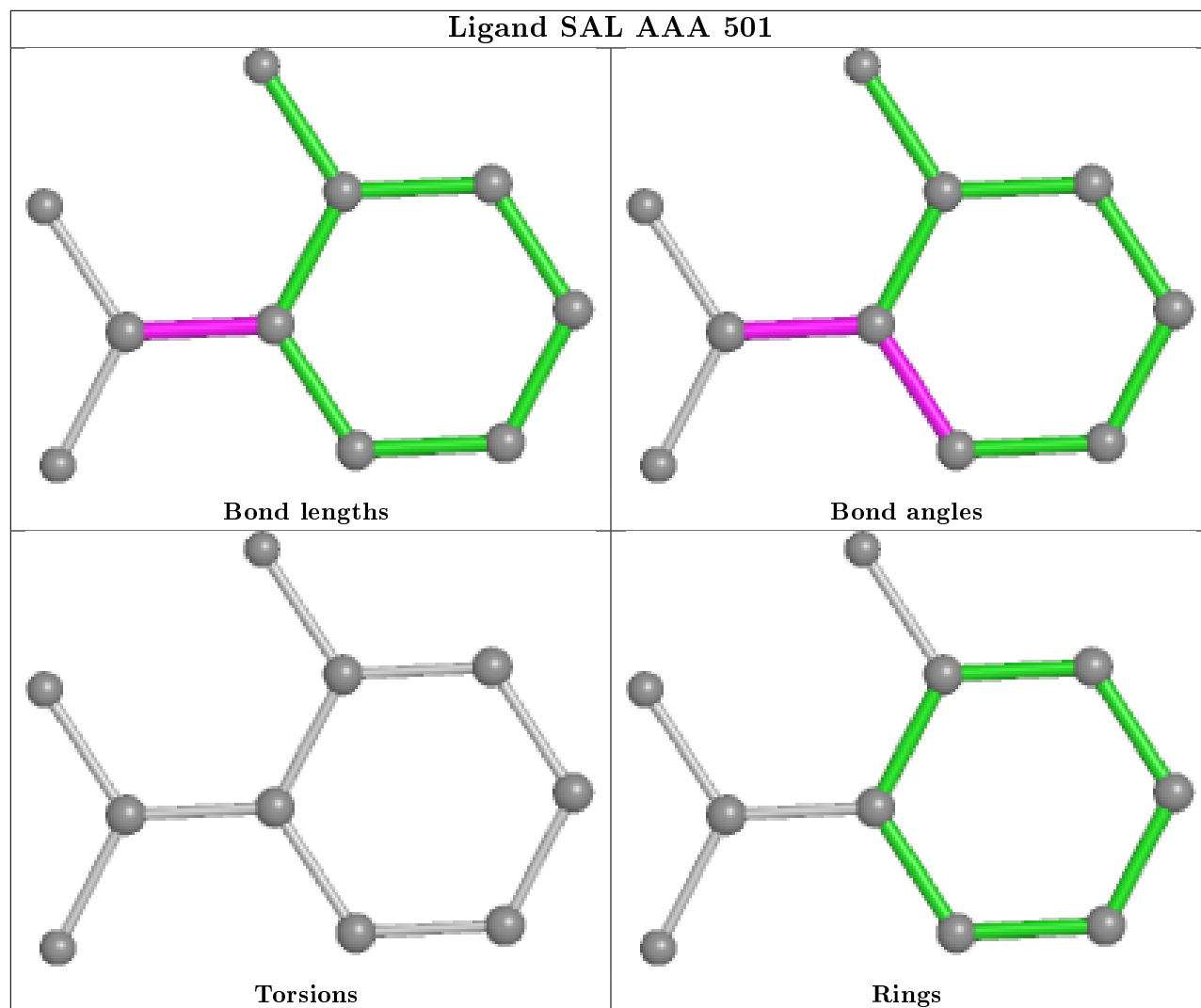
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	BBB	502	SAL	1	0
3	BBB	503	SAL	1	0

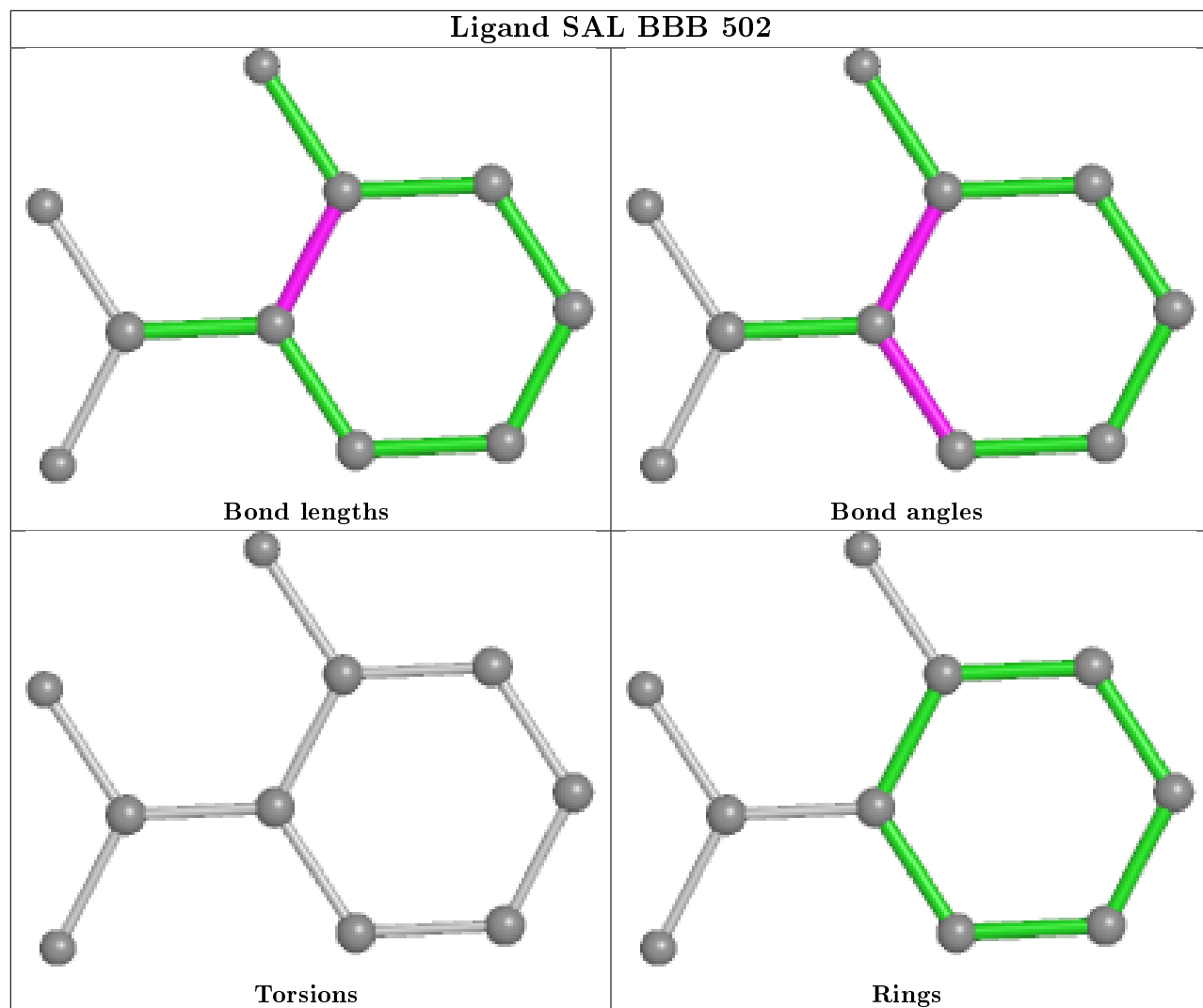
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



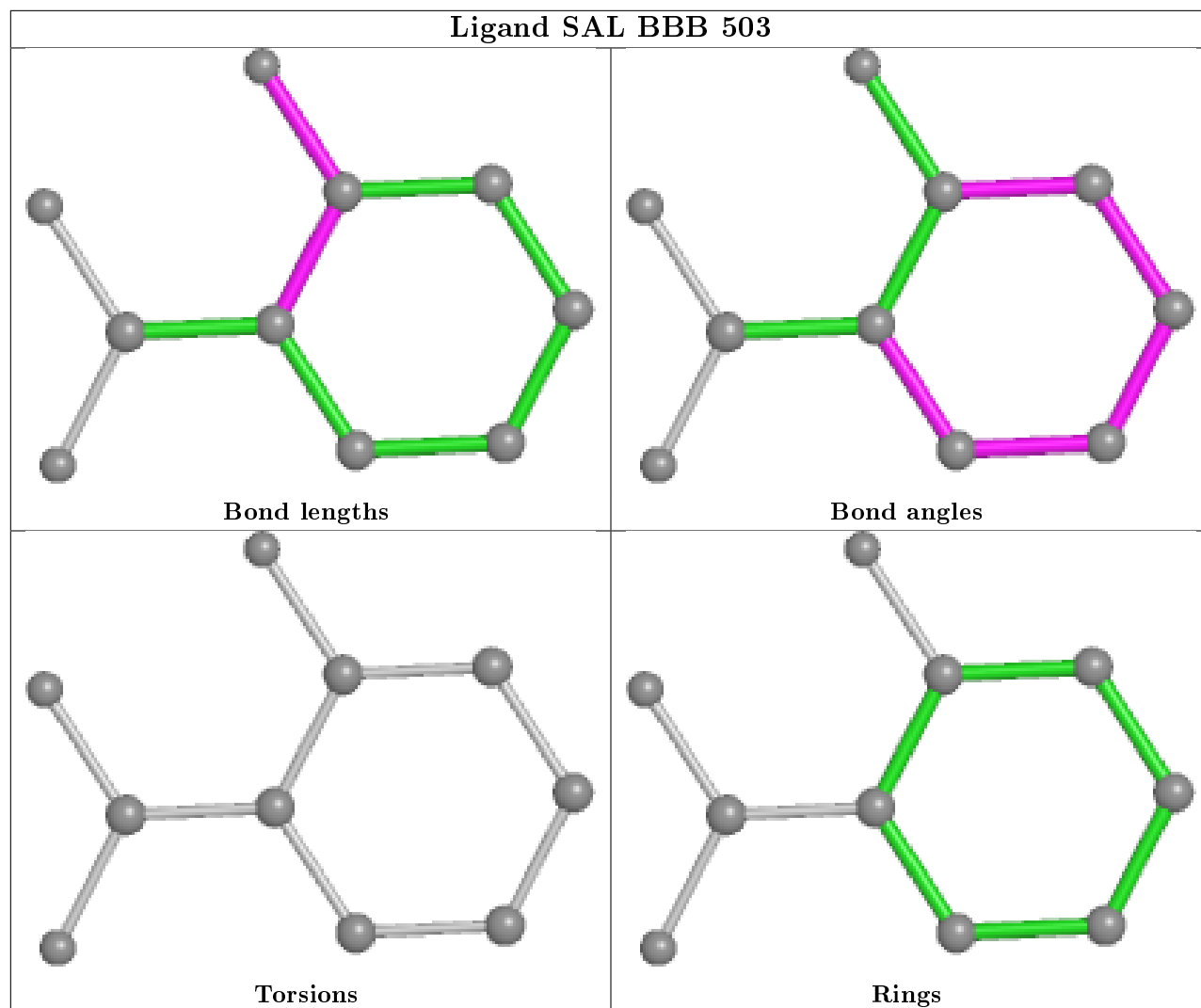
Ligand SAL AAA 501

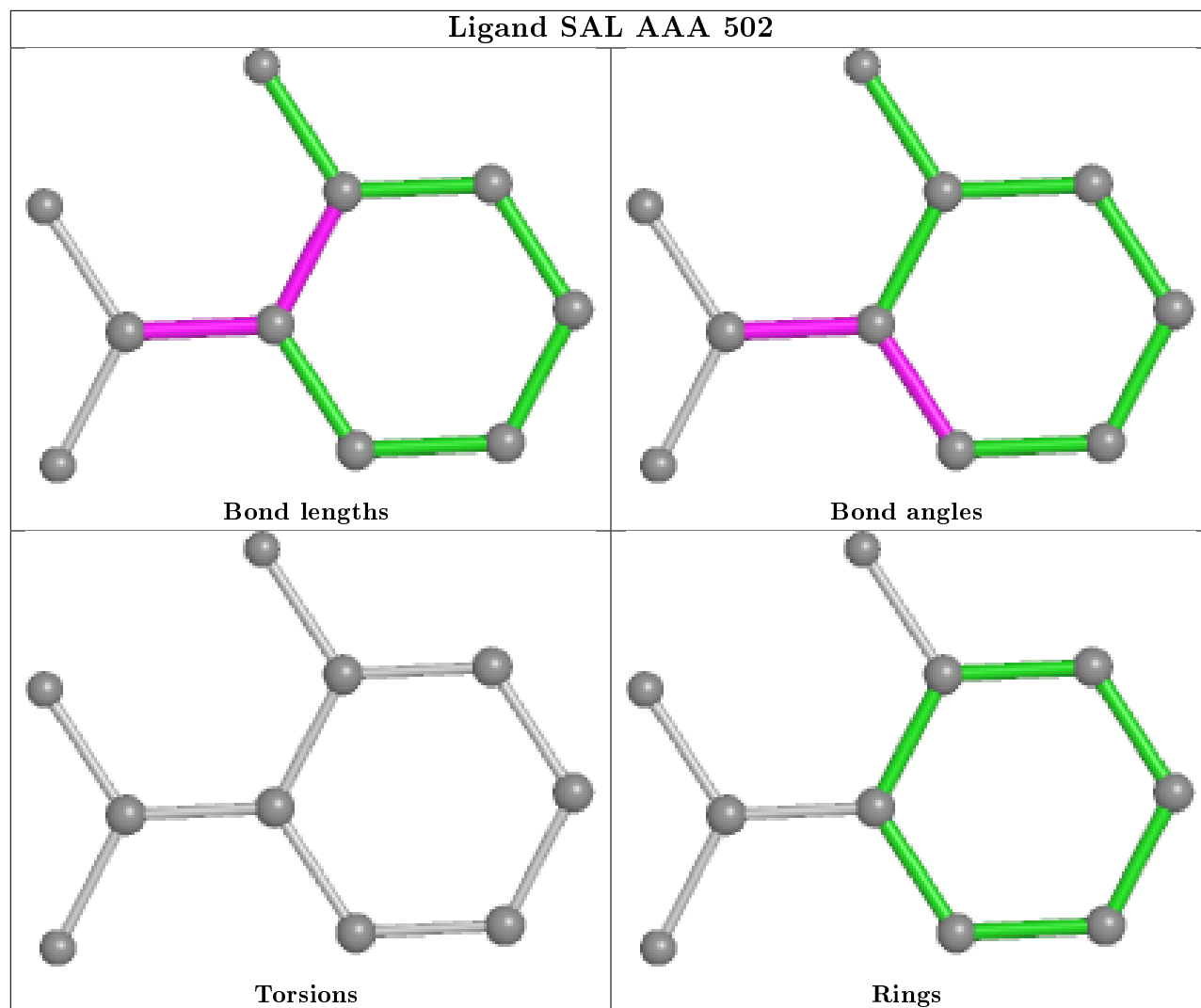


Ligand SAL BBB 502



Ligand SAL BBB 503





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	AAA	430/436 (98%)	0.28	20 (4%) 31 34	27, 52, 95, 115	0
2	BBB	432/437 (98%)	0.07	12 (2%) 53 56	26, 42, 74, 118	0
All	All	862/873 (98%)	0.17	32 (3%) 41 44	26, 46, 91, 118	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	BBB	331	PRO	5.2
2	BBB	436	PHE	4.7
1	AAA	268	ARG	4.5
1	AAA	320	LEU	4.2
2	BBB	433	VAL	4.2
1	AAA	49	ARG	4.1
1	AAA	88	SER	3.3
1	AAA	274	GLU	3.0
2	BBB	435	TYR	2.9
2	BBB	268	ARG	2.9
1	AAA	20	LEU	2.8
1	AAA	11	TRP	2.7
1	AAA	46	THR	2.7
2	BBB	361	PRO	2.6
1	AAA	325	THR	2.6
1	AAA	265	TYR	2.6
1	AAA	33	LEU	2.5
2	BBB	432	ALA	2.5
2	BBB	8	LEU	2.5
1	AAA	44	GLY	2.3
1	AAA	328	ARG	2.3
1	AAA	57	VAL	2.2
2	BBB	359	ALA	2.2
1	AAA	47	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	AAA	277	ARG	2.2
1	AAA	272	LEU	2.2
2	BBB	395	ARG	2.1
2	BBB	328	ARG	2.1
1	AAA	51	ALA	2.0
1	AAA	221	ARG	2.0
2	BBB	414	ARG	2.0
1	AAA	426	ASP	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

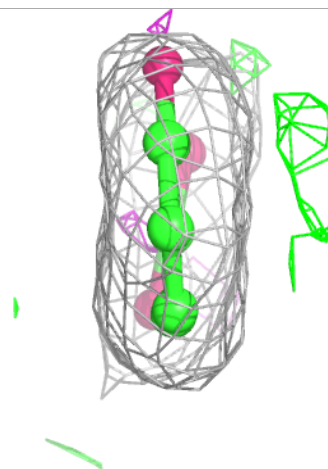
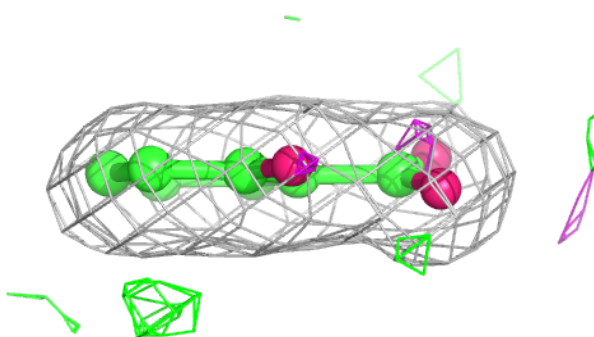
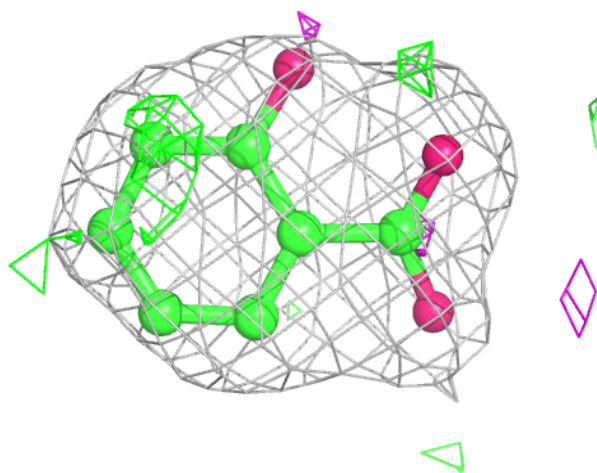
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	MG	BBB	505	1/1	0.80	0.14	36,36,36,36	0
5	MG	AAA	504	1/1	0.84	0.09	45,45,45,45	0
3	SAL	BBB	503	10/10	0.88	0.15	41,43,50,51	0
6	AMP	BBB	501	23/23	0.92	0.16	40,45,54,64	0
3	SAL	AAA	502	10/10	0.93	0.14	53,53,59,65	0
4	ZN	AAA	503	1/1	0.94	0.07	64,64,64,64	0
3	SAL	BBB	502	10/10	0.96	0.14	39,48,55,55	0
3	SAL	AAA	501	10/10	0.97	0.17	42,46,51,52	0
4	ZN	BBB	504	1/1	0.99	0.10	39,39,39,39	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.

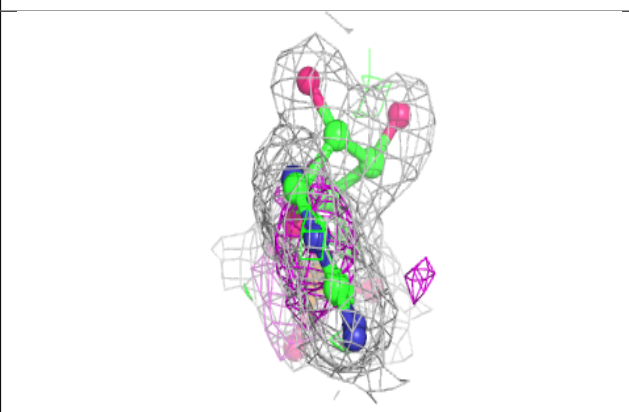
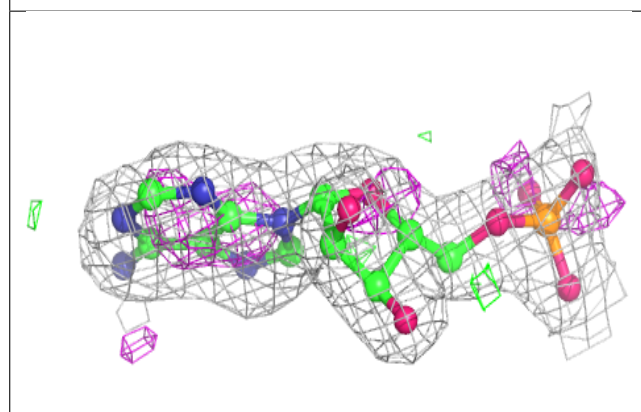
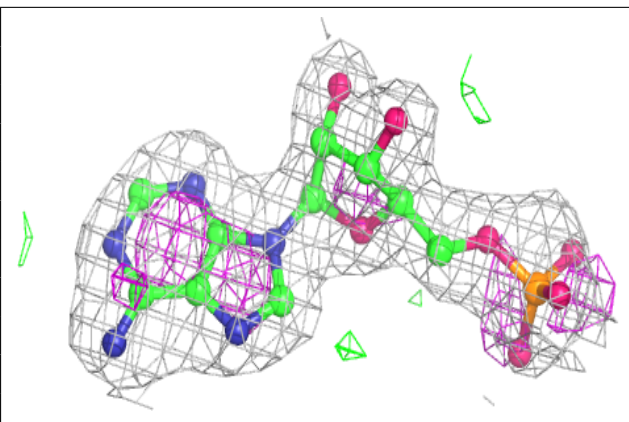
Electron density around SAL BBB 503:

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



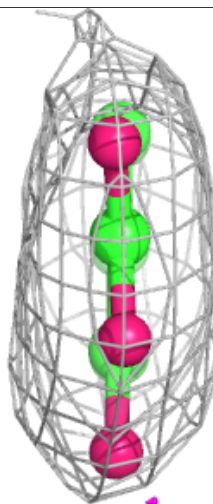
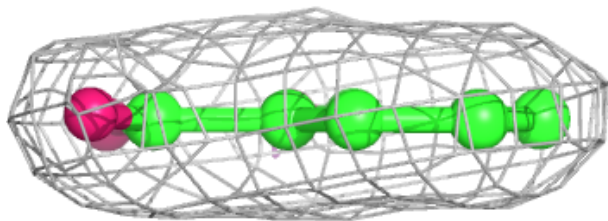
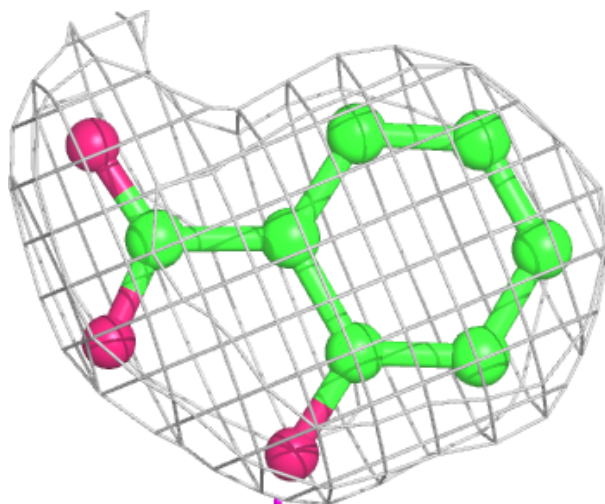
Electron density around AMP BBB 501:

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



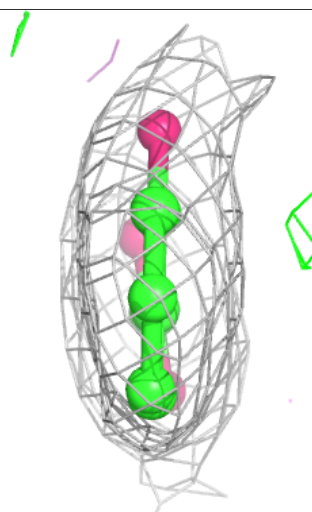
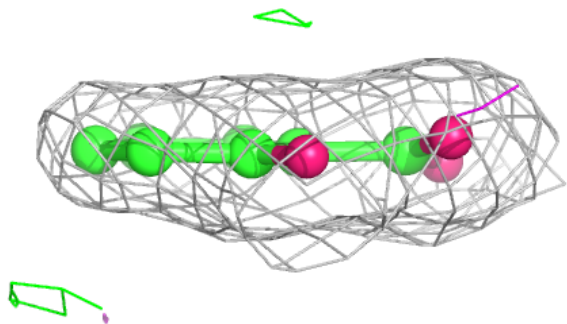
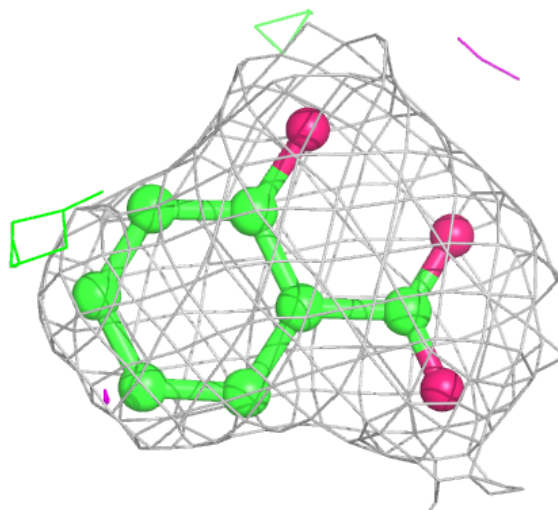
Electron density around SAL AAA 502:

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



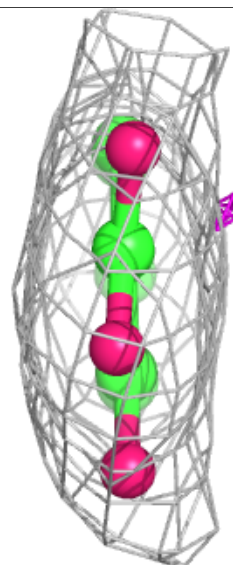
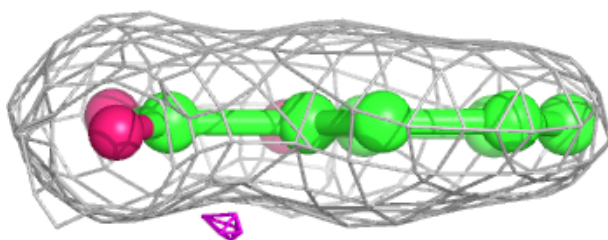
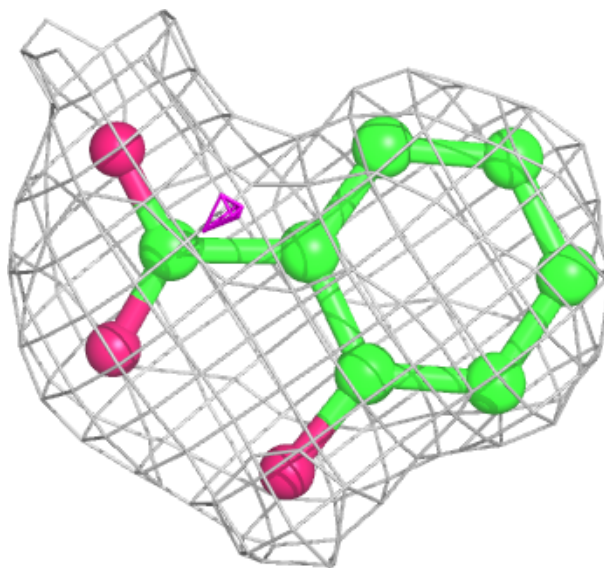
Electron density around SAL BBB 502:

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



Electron density around SAL AAA 501:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.