



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 19, 2026 – 11:49 AM UTC

PDB ID : 3RV6 / pdb_00003rv6
Title : Structure of a M. tuberculosis Salicylate Synthase, MbtI, in Complex with an Inhibitor with Phenyl R-Group
Authors : Chi, G.; Bulloch, E.M.M.; Manos-Turvey, A.; Payne, R.J.; Lott, J.S.; TB Structural Genomics Consortium (TBSGC)
Deposited on : 2011-05-06
Resolution : 2.04 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

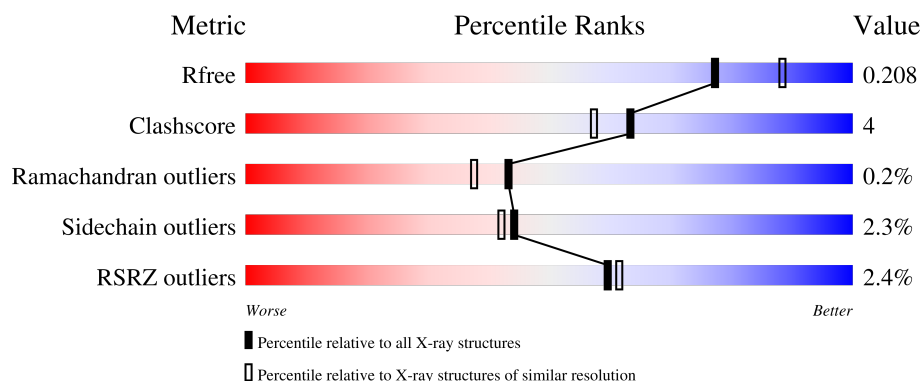
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

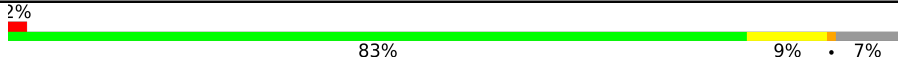
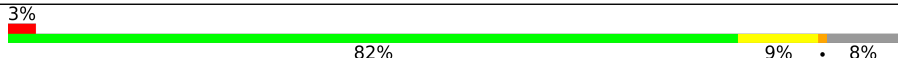
The reported resolution of this entry is 2.04 Å.

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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	450	
1	B	450	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	VAE	A	451	-	X	-	-
5	RVA	A	454	-	X	-	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7015 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.

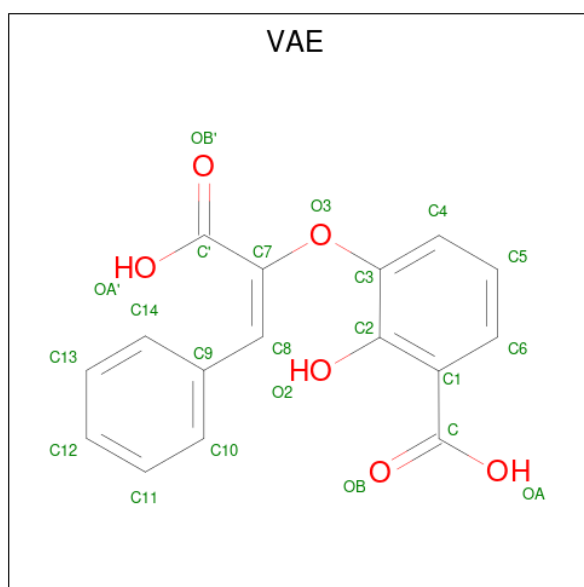
- Molecule 1 is a protein called Isochorismate synthase/ischorismate-pyruvate lyase mbtI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	417	Total	C	N	O	S	0	2	0
			3158	1983	566	599	10			
1	B	413	Total	C	N	O	S	0	1	0
			3122	1962	557	593	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	VAL	-	expression tag	UNP Q7D785
B	1	VAL	-	expression tag	UNP Q7D785

- Molecule 2 is 3-[[*(E)*-1-carboxy-2-phenylethenyl]oxy]-2-hydroxybenzoic acid (CCD ID: VAE) (formula: C₁₆H₁₂O₆).

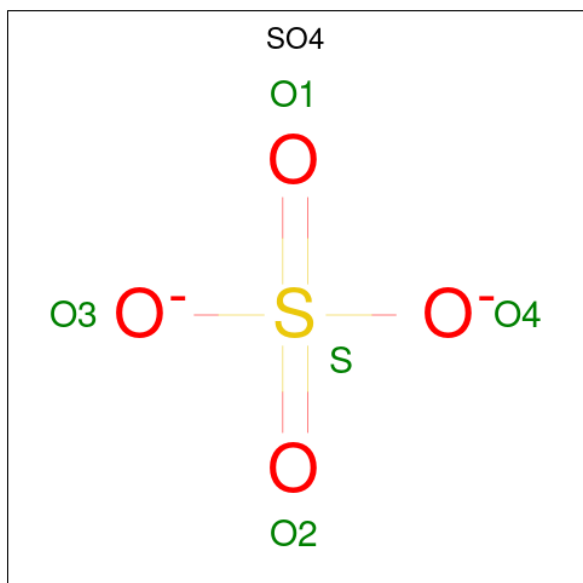


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			22	16	6		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

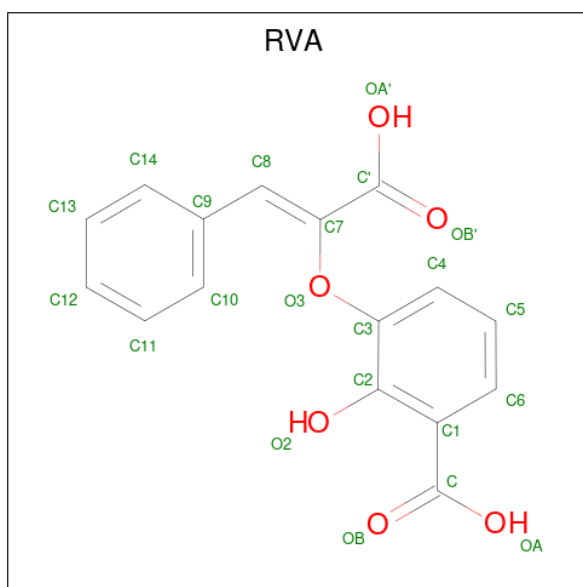
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 5 is 3-{[(Z)-1-carboxy-2-phenylethenyl]oxy}-2-hydroxybenzoic acid (CCD ID: RVA) (formula: C₁₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total 22	C 16	O 6	0	0
5	A	1	Total 22	C 16	O 6	0	0
5	B	1	Total 22	C 16	O 6	0	0
5	B	1	Total 22	C 16	O 6	0	0

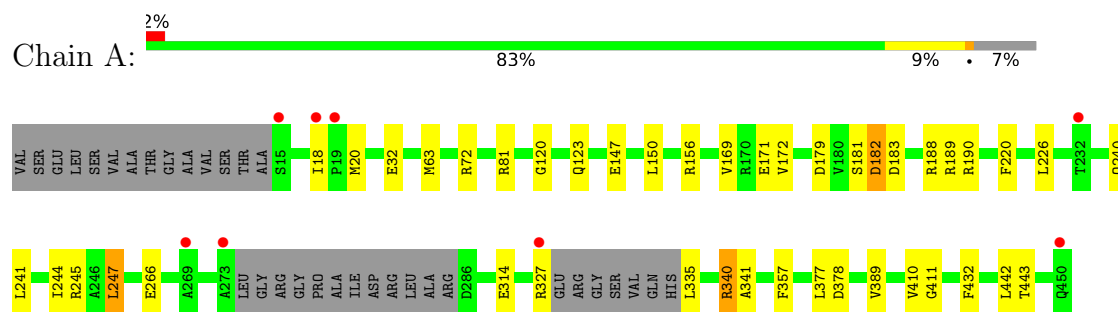
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	328	Total O 328 328	0	0
6	B	290	Total O 290 290	0	0

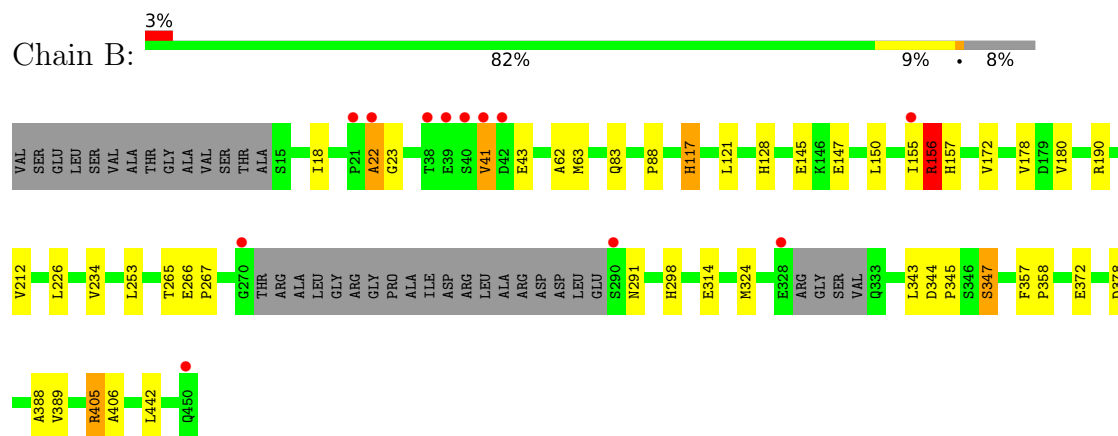
3 Residue-property plots [i](#)

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

- Molecule 1: Isochorismate synthase/isochorismate-pyruvate lyase mbtI



- Molecule 1: Isochorismate synthase/isochorismate-pyruvate lyase mbtI



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.75Å 91.53Å 96.47Å 90.00° 104.83° 90.00°	Depositor
Resolution (Å)	16.72 – 2.04 16.72 – 2.04	Depositor EDS
% Data completeness (in resolution range)	(Not available) (16.72-2.04) 96.6 (16.72-2.04)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.03Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, R_{free}	0.171 , 0.205 0.176 , 0.208	Depositor DCC
R_{free} test set	2762 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	24.3	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 60.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.018 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7015	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% 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: RVA, VAE, MG, SO4

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	0.97	4/3212 (0.1%)	1.15	8/4365 (0.2%)
1	B	0.96	5/3177 (0.2%)	1.16	8/4319 (0.2%)
All	All	0.97	9/6389 (0.1%)	1.15	16/8684 (0.2%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	443	THR	C-O	-7.47	1.17	1.24
1	A	20	MET	SD-CE	-5.77	1.65	1.79
1	B	117	HIS	C-O	-5.36	1.17	1.24
1	A	266	GLU	C-O	-5.25	1.18	1.24
1	B	63	MET	SD-CE	-5.25	1.66	1.79
1	B	62	ALA	CA-CB	-5.23	1.44	1.53
1	B	128	HIS	C-O	-5.15	1.16	1.23
1	B	62	ALA	C-O	-5.07	1.17	1.23
1	A	247	LEU	C-O	-5.05	1.17	1.23

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	157	HIS	N-CA-C	8.93	120.94	111.03
1	B	147	GLU	N-CA-C	8.31	119.59	108.38
1	A	147	GLU	N-CA-C	6.37	116.97	108.38
1	B	291	ASN	CA-C-N	6.02	128.34	120.28
1	B	291	ASN	C-N-CA	6.02	128.34	120.28
1	A	181	SER	CA-C-N	5.85	129.59	120.75
1	A	181	SER	C-N-CA	5.85	129.59	120.75
1	B	406	ALA	N-CA-C	5.43	117.06	108.96
1	A	182	ASP	N-CA-CB	-5.34	102.19	110.04
1	A	188	ARG	CA-C-N	5.32	127.42	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	188	ARG	C-N-CA	5.32	127.42	120.28
1	A	389	VAL	N-CA-C	-5.32	100.72	108.17
1	A	171	GLU	N-CA-C	-5.28	102.47	110.28
1	B	83	GLN	N-CA-C	-5.19	102.79	110.48
1	B	156	ARG	CB-CA-C	5.13	120.63	110.42
1	B	389	VAL	N-CA-C	-5.08	101.05	108.17

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3158	0	3136	20	0
1	B	3122	0	3079	30	0
2	A	22	0	10	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	5	0	0	0	0
5	A	44	0	19	5	0
5	B	44	0	20	0	0
6	A	328	0	0	4	0
6	B	290	0	0	1	0
All	All	7015	0	6264	51	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 4.

All (51) 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:B:22:ALA:CB	1:B:23:GLY:HA3	1.94	0.96
1:A:123:GLN:NE2	6:A:576:HOH:O	2.01	0.91
1:B:22:ALA:HB1	1:B:23:GLY:HA3	1.51	0.88
1:A:340:ARG:C	1:A:340:ARG:HD3	2.12	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:179:ASP:O	6:A:467:HOH:O	2.06	0.74
1:B:253:LEU:HD11	1:B:265:THR:HG22	1.70	0.73
1:B:155:ILE:O	1:B:156:ARG:C	2.32	0.72
5:A:455:RVA:H14	5:A:455:RVA:O3	1.93	0.69
1:B:22:ALA:CB	1:B:23:GLY:CA	2.75	0.63
1:B:22:ALA:HB3	1:B:23:GLY:HA3	1.78	0.63
1:B:298:HIS:HD2	1:B:324:MET:HG2	1.63	0.63
1:A:156:ARG:CB	5:A:455:RVA:OA'	2.48	0.62
1:A:156:ARG:HB2	5:A:455:RVA:OA'	2.01	0.61
1:B:314:GLU:HG3	1:B:345:PRO:HD2	1.84	0.60
1:B:357:PHE:CZ	1:B:388:ALA:HB2	2.37	0.59
1:B:405:ARG:HG2	6:B:740:HOH:O	2.05	0.56
1:B:41:VAL:HB	1:B:43:GLU:HG3	1.90	0.54
1:B:253:LEU:HD11	1:B:265:THR:CG2	2.36	0.53
1:B:88:PRO:HD2	1:B:117:HIS:CE1	2.45	0.52
1:A:182:ASP:HB3	6:A:467:HOH:O	2.09	0.52
1:B:298:HIS:CD2	1:B:324:MET:HG2	2.44	0.52
1:A:32:GLU:HG3	1:A:169:VAL:HB	1.91	0.51
1:A:189:ARG:HD2	6:A:483:HOH:O	2.10	0.51
1:A:18:ILE:HD12	1:A:150:LEU:HD11	1.95	0.49
1:B:266:GLU:HB2	1:B:267:PRO:HA	1.96	0.48
1:B:324:MET:HE2	1:B:324:MET:HB3	1.74	0.47
1:B:357:PHE:HA	1:B:358:PRO:C	2.41	0.46
1:B:18:ILE:HD12	1:B:150:LEU:HD11	1.97	0.46
1:B:405:ARG:HA	1:B:405:ARG:NE	2.32	0.45
1:B:344:ASP:HB3	1:B:347:SER:HB2	1.99	0.45
5:A:455:RVA:O3	5:A:455:RVA:C14	2.61	0.45
1:B:344:ASP:HA	1:B:345:PRO:HD2	1.66	0.45
1:A:340:ARG:HD3	1:A:341:ALA:N	2.32	0.45
1:A:156:ARG:N	5:A:455:RVA:OA'	2.45	0.45
1:B:22:ALA:HB1	1:B:23:GLY:CA	2.35	0.45
1:B:357:PHE:CE1	1:B:388:ALA:HB2	2.51	0.45
1:A:72:ARG:HG2	1:A:81:ARG:HG2	1.99	0.44
1:A:244:ILE:HG12	1:A:410:VAL:HG22	2.00	0.44
1:A:190:ARG:HD2	1:A:377:LEU:O	2.18	0.44
1:B:343:LEU:O	1:B:345:PRO:HD3	2.18	0.43
1:A:120:GLY:HA3	1:B:190:ARG:HG2	2.00	0.43
1:B:180:VAL:HG12	1:B:212:VAL:HG11	2.00	0.43
1:A:63:MET:HE1	1:A:241:LEU:HD11	2.02	0.42
1:A:220:PHE:CD1	1:A:247:LEU:HD23	2.55	0.41
1:A:240:GLN:OE1	1:A:245:ARG:CZ	2.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:234:VAL:HG21	1:B:267:PRO:HD3	2.01	0.41
1:A:172:VAL:HG11	1:A:226:LEU:HB2	2.04	0.40
1:A:183:ASP:HB2	1:A:432:PHE:CE2	2.56	0.40
1:B:405:ARG:HA	1:B:405:ARG:CZ	2.51	0.40
1:B:172:VAL:HG11	1:B:226:LEU:HB2	2.04	0.40
1:B:121:LEU:HD21	1:B:372:GLU:HG2	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	413/450 (92%)	406 (98%)	6 (2%)	1 (0%)	43	37
1	B	408/450 (91%)	402 (98%)	5 (1%)	1 (0%)	43	37
All	All	821/900 (91%)	808 (98%)	11 (1%)	2 (0%)	43	37

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	411	GLY
1	B	22	ALA

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	324/358 (90%)	317 (98%)	7 (2%)	45	44
1	B	319/358 (89%)	311 (98%)	8 (2%)	42	39
All	All	643/716 (90%)	628 (98%)	15 (2%)	44	42

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	314	GLU
1	A	327	ARG
1	A	335	LEU
1	A	340	ARG
1	A	357	PHE
1	A	378	ASP
1	A	442	LEU
1	B	41	VAL
1	B	145	GLU
1	B	156	ARG
1	B	178	VAL
1	B	347	SER
1	B	378	ASP
1	B	405	ARG
1	B	442	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	105	GLN
1	A	128	HIS
1	B	117	HIS
1	B	334	HIS
1	B	409	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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
5	RVA	A	455	-	23,23,23	2.31	7 (30%)	30,31,31	1.77	6 (20%)
5	RVA	B	451	-	23,23,23	4.40	17 (73%)	30,31,31	1.84	8 (26%)
4	SO4	A	453	-	4,4,4	0.35	0	6,6,6	0.17	0
5	RVA	A	454	-	23,23,23	4.56	18 (78%)	30,31,31	2.18	14 (46%)
2	VAE	A	451	-	23,23,23	4.48	16 (69%)	30,31,31	2.24	9 (30%)
5	RVA	B	453	-	23,23,23	4.42	16 (69%)	30,31,31	1.74	7 (23%)

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	RVA	A	455	-	-	2/16/16/16	0/2/2/2
5	RVA	B	451	-	-	3/16/16/16	0/2/2/2
5	RVA	A	454	-	-	2/16/16/16	0/2/2/2
2	VAE	A	451	-	-	5/16/16/16	0/2/2/2
5	RVA	B	453	-	-	0/16/16/16	0/2/2/2

All (74) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	454	RVA	C13-C14	9.73	1.55	1.38
2	A	451	VAE	C13-C14	9.02	1.54	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	451	RVA	C13-C14	8.79	1.53	1.38
5	A	454	RVA	C10-C9	8.65	1.57	1.39
5	B	453	RVA	C13-C14	8.57	1.53	1.38
2	A	451	VAE	C5-C6	-7.49	1.26	1.38
2	A	451	VAE	C10-C9	7.42	1.54	1.39
5	B	451	RVA	C10-C9	7.07	1.53	1.39
2	A	451	VAE	C8-C7	-7.00	1.25	1.34
5	B	451	RVA	O2-C2	6.97	1.52	1.36
5	B	453	RVA	C10-C9	6.97	1.53	1.39
5	B	453	RVA	C5-C6	-6.48	1.27	1.38
5	B	453	RVA	C8-C7	-6.25	1.26	1.34
5	A	454	RVA	C5-C6	-6.11	1.28	1.38
5	B	451	RVA	C3-C2	5.88	1.48	1.40
2	A	451	VAE	O2-C2	5.76	1.50	1.36
5	A	454	RVA	C3-C2	5.76	1.48	1.40
5	B	453	RVA	O2-C2	5.69	1.49	1.36
5	B	451	RVA	C8-C7	-5.68	1.27	1.34
5	B	453	RVA	C6-C1	5.66	1.48	1.39
5	A	454	RVA	C6-C1	5.61	1.48	1.39
5	B	453	RVA	OB'-C'	-5.59	1.08	1.22
5	B	451	RVA	C5-C6	-5.48	1.29	1.38
2	A	451	VAE	OB'-C'	-5.46	1.09	1.22
5	A	454	RVA	C13-C12	5.38	1.50	1.38
2	A	451	VAE	C13-C12	5.29	1.49	1.38
5	B	453	RVA	C13-C12	5.28	1.49	1.38
5	A	455	RVA	C7-C'	5.24	1.60	1.48
5	B	451	RVA	C5-C4	5.13	1.47	1.38
5	B	451	RVA	C6-C1	5.09	1.47	1.39
5	A	454	RVA	O3-C7	5.08	1.49	1.38
5	B	453	RVA	C3-C2	5.04	1.47	1.40
2	A	451	VAE	O3-C7	4.99	1.49	1.38
5	B	451	RVA	O3-C7	4.89	1.49	1.38
5	B	453	RVA	O3-C7	4.79	1.49	1.38
5	B	453	RVA	OB-C	4.64	1.36	1.22
5	B	451	RVA	OB'-C'	-4.56	1.11	1.22
5	A	454	RVA	C5-C4	4.50	1.46	1.38
5	A	455	RVA	OA'-C'	-4.43	1.18	1.30
5	B	451	RVA	C7-C'	4.35	1.58	1.48
5	A	454	RVA	OB-C	4.35	1.35	1.22
5	A	454	RVA	C9-C8	4.24	1.54	1.46
5	A	454	RVA	C7-C'	4.13	1.58	1.48
5	A	454	RVA	OB'-C'	-4.12	1.12	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	451	VAE	C7-C'	4.11	1.58	1.48
2	A	451	VAE	C3-C2	4.10	1.46	1.40
5	B	453	RVA	C5-C4	4.07	1.45	1.38
5	A	455	RVA	C9-C8	-4.05	1.39	1.46
5	A	454	RVA	O2-C2	4.03	1.46	1.36
2	A	451	VAE	C5-C4	4.01	1.45	1.38
5	B	451	RVA	C13-C12	4.01	1.47	1.38
5	A	455	RVA	OA-C	-3.89	1.19	1.30
5	A	455	RVA	O3-C3	-3.72	1.33	1.41
2	A	451	VAE	C6-C1	3.39	1.45	1.39
5	A	455	RVA	C8-C7	3.28	1.37	1.34
2	A	451	VAE	C9-C8	3.21	1.52	1.46
5	B	451	RVA	C1-C	-3.19	1.42	1.49
5	A	454	RVA	C8-C7	-3.17	1.30	1.34
2	A	451	VAE	C4-C3	-3.10	1.33	1.39
5	B	451	RVA	C9-C8	3.09	1.52	1.46
5	B	453	RVA	OA'-C'	-2.88	1.22	1.30
5	A	454	RVA	OA'-C'	-2.77	1.23	1.30
5	A	454	RVA	C4-C3	-2.77	1.33	1.39
5	B	453	RVA	C7-C'	2.75	1.55	1.48
5	B	453	RVA	OA-C	-2.71	1.22	1.30
5	A	454	RVA	C14-C9	2.55	1.44	1.39
5	B	451	RVA	OA'-C'	-2.54	1.23	1.30
5	B	451	RVA	C14-C9	2.50	1.44	1.39
2	A	451	VAE	OA'-C'	-2.45	1.23	1.30
5	A	454	RVA	OA-C	-2.42	1.23	1.30
5	A	455	RVA	C1-C	2.41	1.54	1.49
5	B	453	RVA	C14-C9	2.28	1.44	1.39
2	A	451	VAE	C1-C	-2.19	1.44	1.49
5	B	451	RVA	C1-C2	2.17	1.45	1.41

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	451	VAE	O3-C3-C4	6.29	134.03	118.89
5	A	454	RVA	O3-C3-C4	5.26	131.55	118.89
5	B	451	RVA	O3-C3-C4	5.19	131.38	118.89
5	B	453	RVA	O3-C3-C4	5.08	131.12	118.89
5	A	455	RVA	O3-C7-C'	5.01	128.19	114.54
2	A	451	VAE	C12-C13-C14	-4.22	115.03	120.24
5	A	454	RVA	C11-C10-C9	4.10	125.44	120.64
5	B	451	RVA	C12-C13-C14	-3.99	115.31	120.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	451	VAE	C13-C12-C11	3.97	125.31	119.87
5	A	455	RVA	C2-C1-C	3.96	124.28	119.86
2	A	451	VAE	O3-C3-C2	-3.94	109.55	117.50
2	A	451	VAE	C4-C3-C2	-3.86	116.41	119.98
5	B	453	RVA	C12-C13-C14	-3.66	115.72	120.24
5	B	451	RVA	C13-C12-C11	3.60	124.79	119.87
2	A	451	VAE	C11-C10-C9	3.53	124.78	120.64
5	A	454	RVA	C13-C12-C11	3.48	124.64	119.87
5	A	454	RVA	O3-C3-C2	-3.19	111.05	117.50
5	B	451	RVA	C4-C3-C2	-3.11	117.10	119.98
5	B	451	RVA	O3-C3-C2	-2.99	111.46	117.50
5	B	453	RVA	O3-C3-C2	-2.94	111.57	117.50
5	A	454	RVA	C12-C13-C14	-2.92	116.64	120.24
5	B	453	RVA	C13-C12-C11	2.90	123.84	119.87
5	B	453	RVA	C4-C3-C2	-2.88	117.31	119.98
5	A	454	RVA	C2-C1-C	2.83	123.03	119.86
5	A	454	RVA	C6-C5-C4	2.80	123.84	120.24
5	A	454	RVA	C4-C3-C2	-2.78	117.40	119.98
5	A	455	RVA	C9-C8-C7	-2.68	126.43	130.09
5	B	453	RVA	C11-C10-C9	2.62	123.71	120.64
5	A	454	RVA	OB-C-C1	-2.59	115.78	121.97
5	A	455	RVA	C8-C7-C'	-2.49	115.78	122.83
5	A	455	RVA	OA'-C'-OB'	-2.44	118.09	123.90
5	B	451	RVA	C3-O3-C7	-2.43	114.48	118.49
5	A	455	RVA	C3-O3-C7	2.38	122.40	118.49
5	B	453	RVA	OB-C-C1	-2.37	116.31	121.97
5	B	451	RVA	C6-C5-C4	2.34	123.25	120.24
5	A	454	RVA	C8-C7-C'	2.33	129.43	122.83
5	A	454	RVA	C1-C2-C3	2.28	121.42	119.56
2	A	451	VAE	C6-C5-C4	2.22	123.10	120.24
2	A	451	VAE	C9-C8-C7	2.19	133.09	130.09
5	A	454	RVA	C3-O3-C7	2.11	121.97	118.49
5	A	454	RVA	OA-C-C1	2.07	121.17	115.28
2	A	451	VAE	OB-C-C1	-2.02	117.16	121.97
5	B	451	RVA	C9-C8-C7	2.01	132.84	130.09
5	A	454	RVA	C6-C1-C2	-2.01	116.82	118.76

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	451	VAE	OA'-C'-C7-O3

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Mol	Chain	Res	Type	Atoms
2	A	451	VAE	OB'-C'-C7-O3
5	A	454	RVA	OA'-C'-C7-O3
5	B	451	RVA	OA'-C'-C7-O3
5	B	451	RVA	C7-C8-C9-C10
2	A	451	VAE	OA'-C'-C7-C8
2	A	451	VAE	OB'-C'-C7-C8
5	A	454	RVA	OB'-C'-C7-O3
5	A	455	RVA	C2-C3-O3-C7
2	A	451	VAE	C8-C7-O3-C3
5	A	455	RVA	C8-C7-O3-C3
5	B	451	RVA	C7-C8-C9-C14

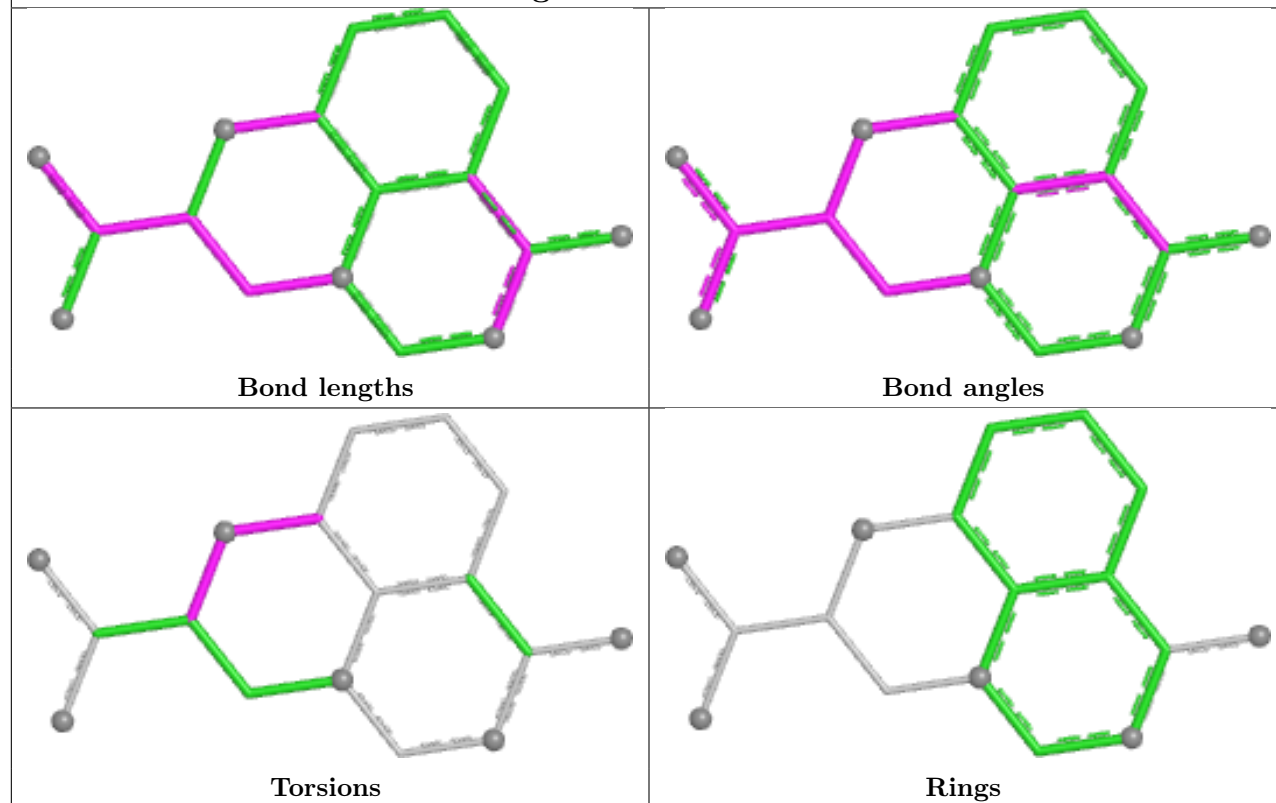
There are no ring outliers.

1 monomer is involved in 5 short contacts:

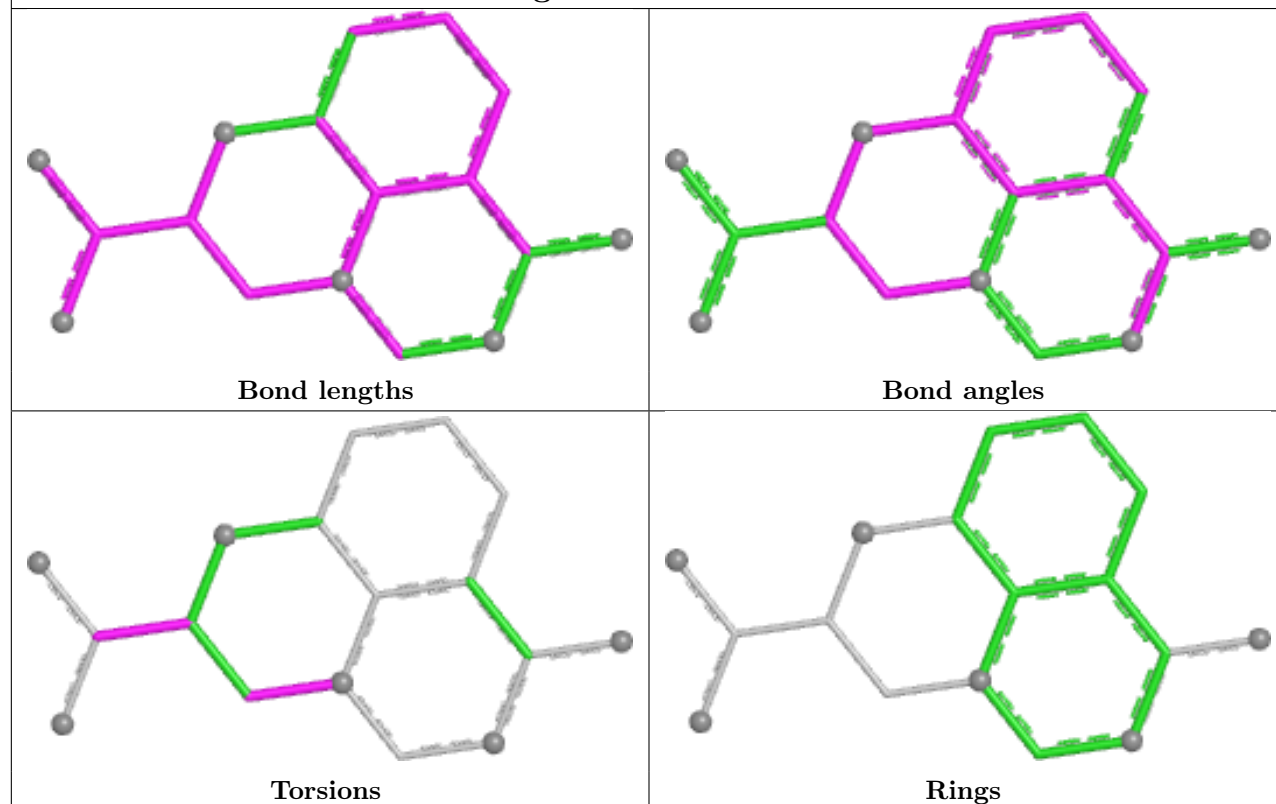
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	455	RVA	5	0

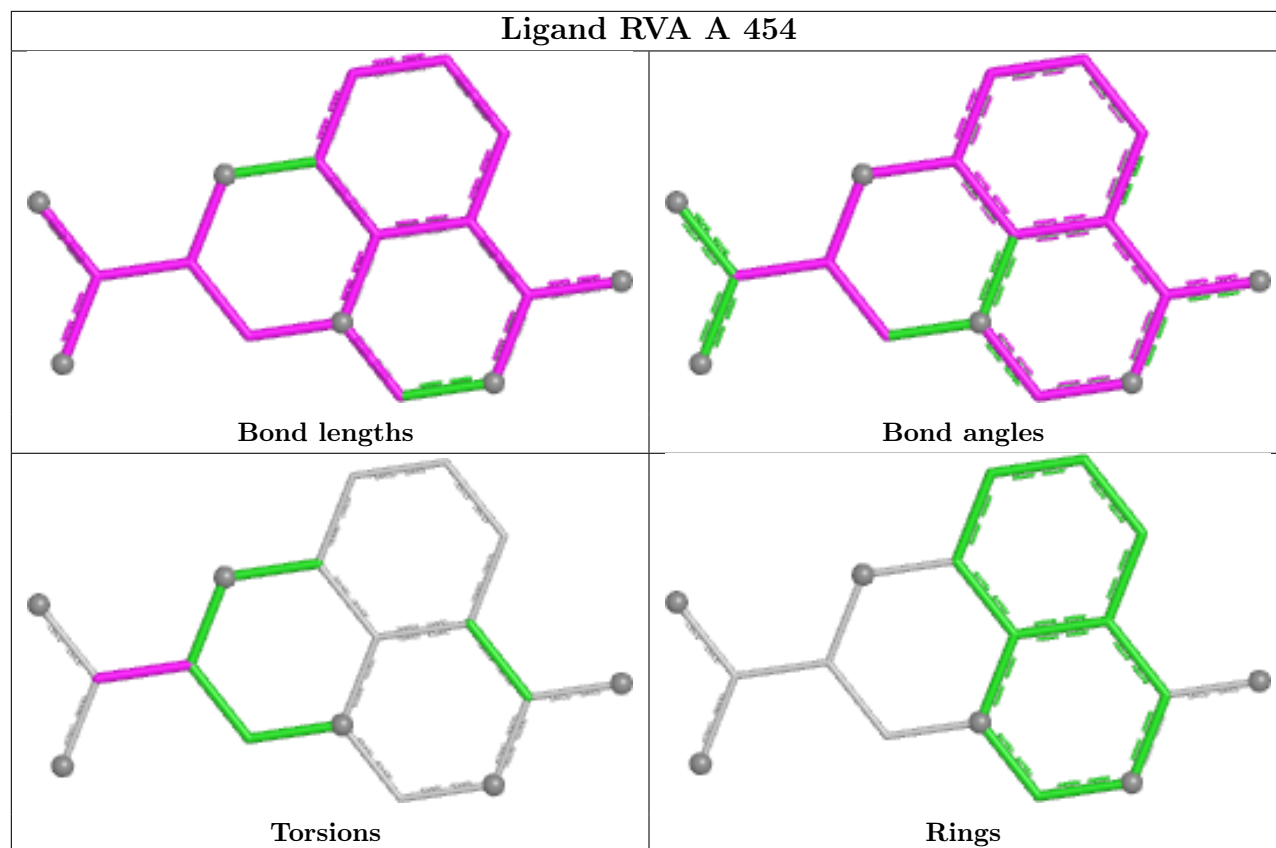
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 RVA A 455

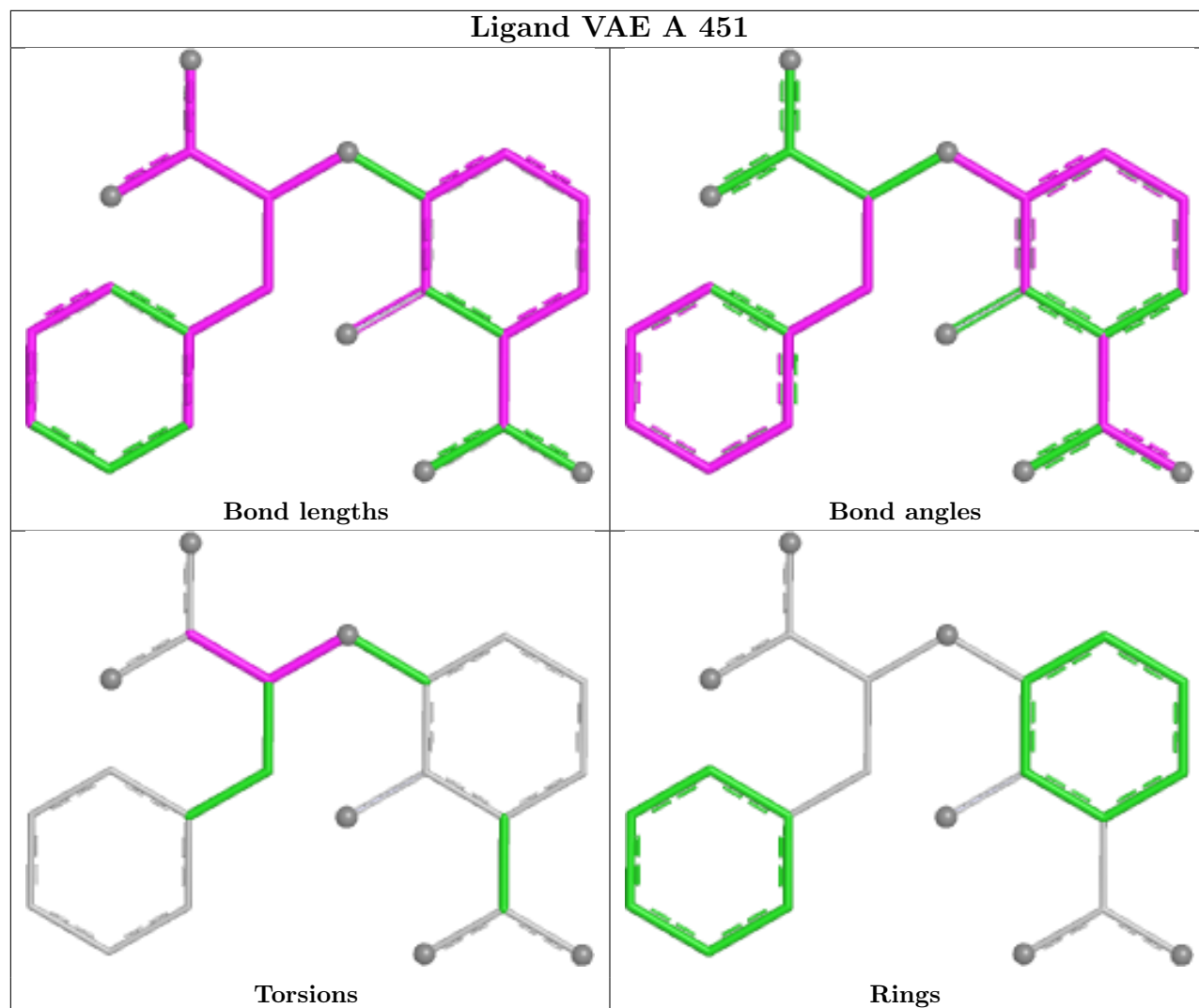


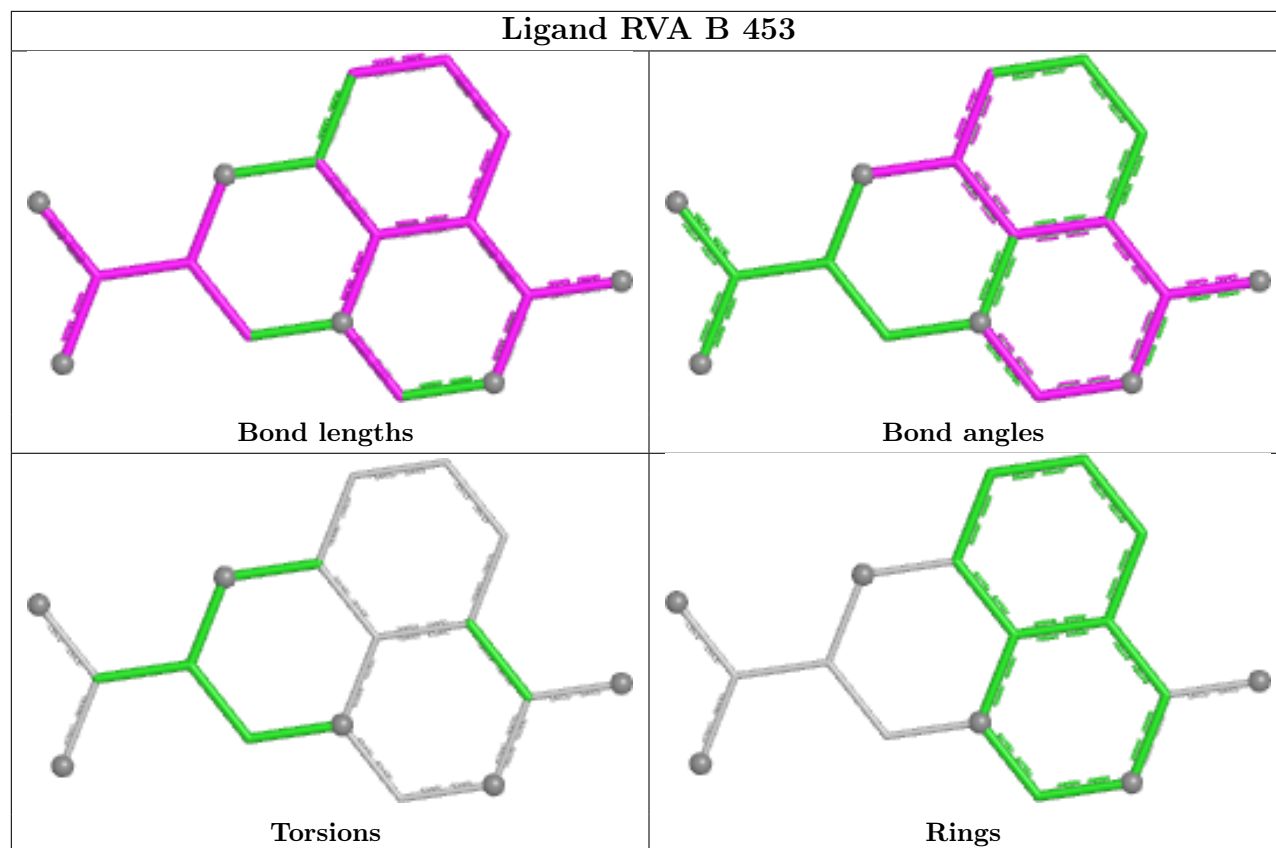
Ligand RVA B 451





Ligand VAE A 451





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	417/450 (92%)	-0.35	8 (1%) 66 68	12, 25, 46, 67	3 (0%)
1	B	413/450 (91%)	-0.21	12 (2%) 53 55	14, 26, 54, 63	3 (0%)
All	All	830/900 (92%)	-0.28	20 (2%) 59 61	12, 25, 51, 67	6 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	450	GLN	3.8
1	B	328	GLU	3.0
1	B	290	SER	2.9
1	B	41	VAL	2.8
1	B	270	GLY	2.8
1	B	40	SER	2.7
1	A	19	PRO	2.6
1	A	273	ALA	2.6
1	A	232	THR	2.6
1	A	15	SER	2.6
1	B	21	PRO	2.5
1	B	22	ALA	2.5
1	A	269	ALA	2.4
1	B	38	THR	2.3
1	A	327	ARG	2.3
1	A	18	ILE	2.2
1	B	39	GLU	2.1
1	B	155	ILE	2.1
1	A	450	GLN	2.1
1	B	42	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

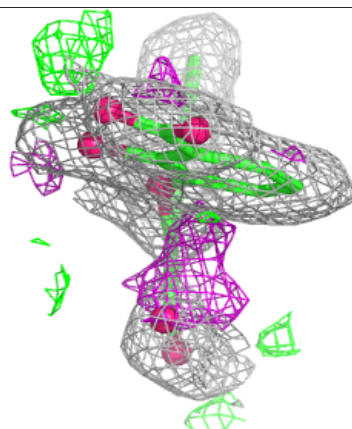
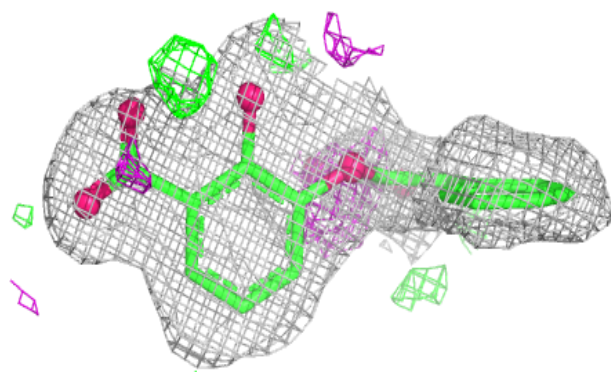
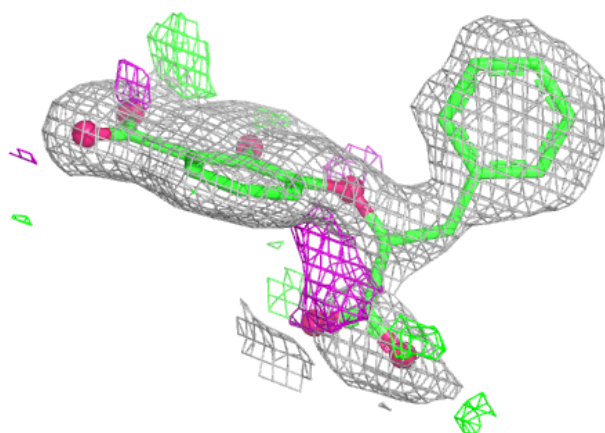
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	RVA	A	455	22/22	0.82	0.11	29,35,45,47	0
3	MG	B	452	1/1	0.88	0.12	54,54,54,54	0
5	RVA	A	454	22/22	0.89	0.08	27,41,43,47	0
4	SO4	A	453	5/5	0.89	0.13	62,66,67,68	0
5	RVA	B	453	22/22	0.93	0.07	31,34,42,43	0
5	RVA	B	451	22/22	0.96	0.05	18,23,36,40	0
2	VAE	A	451	22/22	0.97	0.05	14,23,26,26	0
3	MG	A	452	1/1	0.99	0.02	27,27,27,27	0

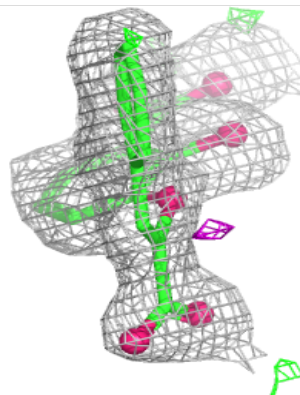
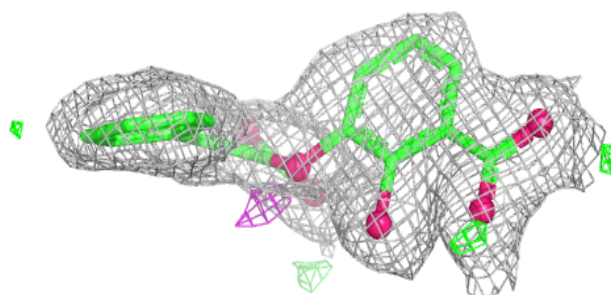
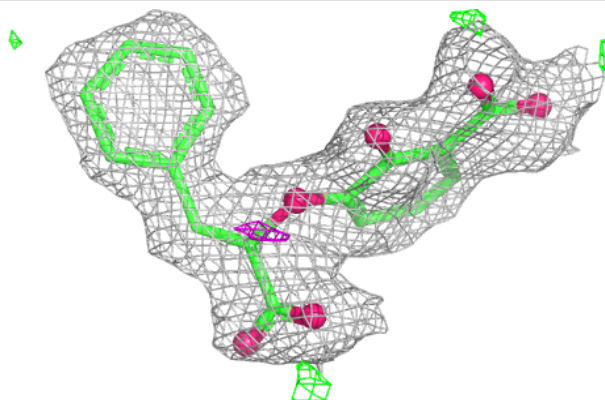
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around RVA A 455:

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

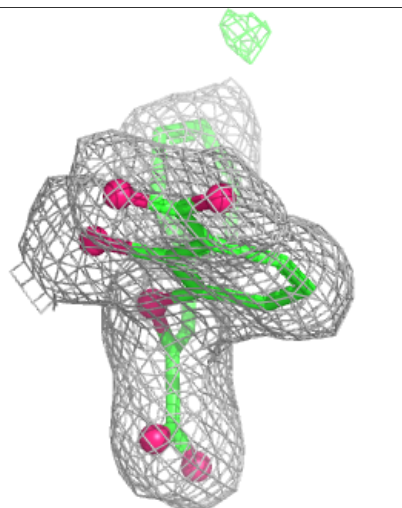
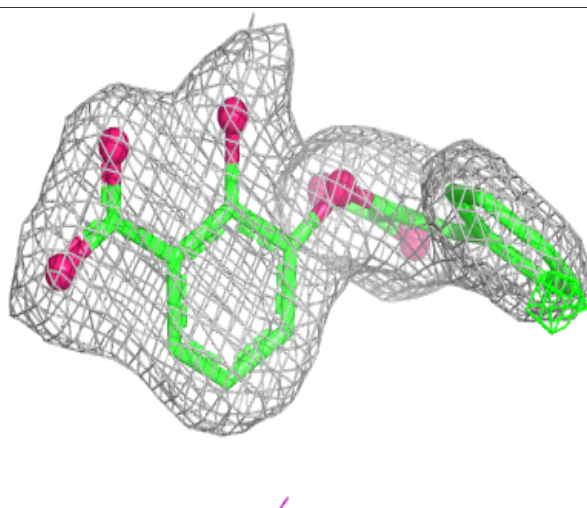
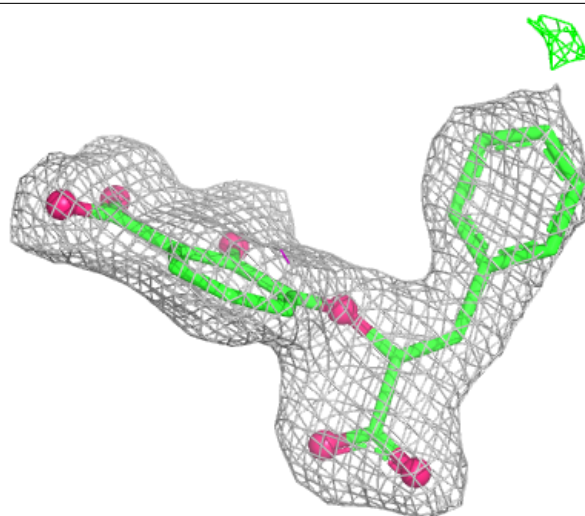
**Electron density around RVA A 454:**

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



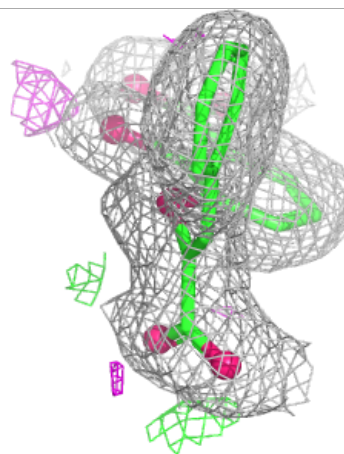
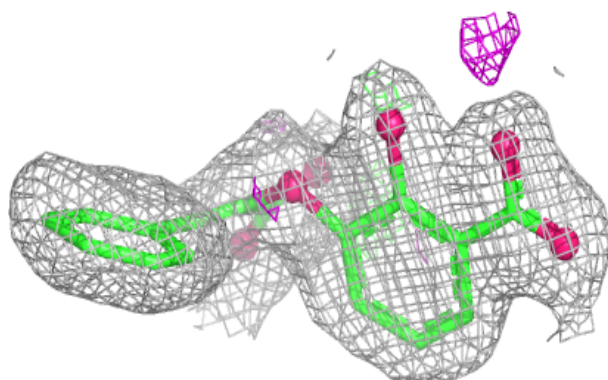
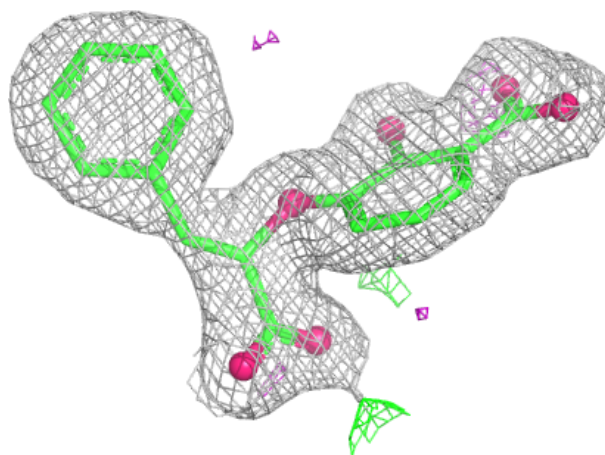
Electron density around RVA B 453:

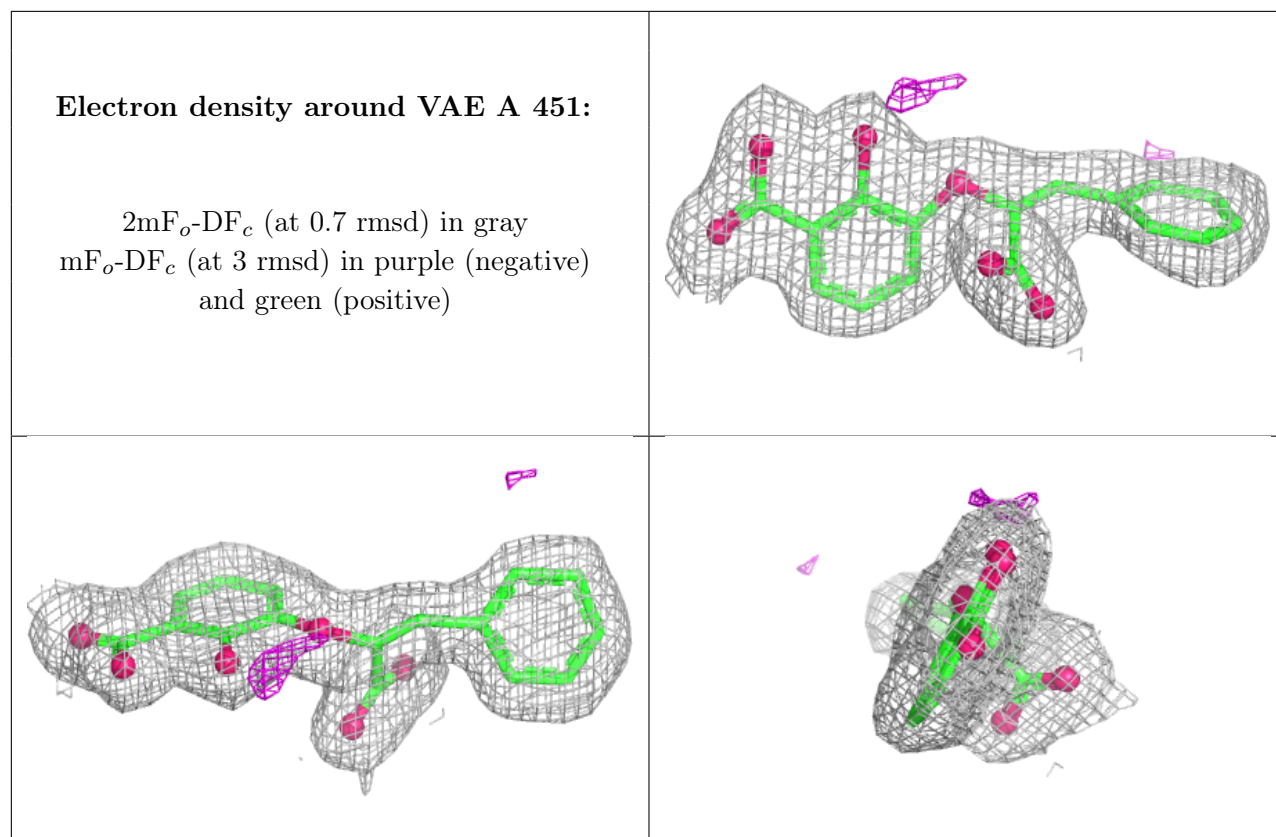
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around RVA B 451:

$2mF_o-DF_c$ (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.