



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 01:36 PM UTC

PDB ID : 3RV9 / pdb_00003rv9
Title : Structure of a M. tuberculosis Salicylate Synthase, MbtI, in Complex with an Inhibitor with Ethyl 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 : 2.14 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

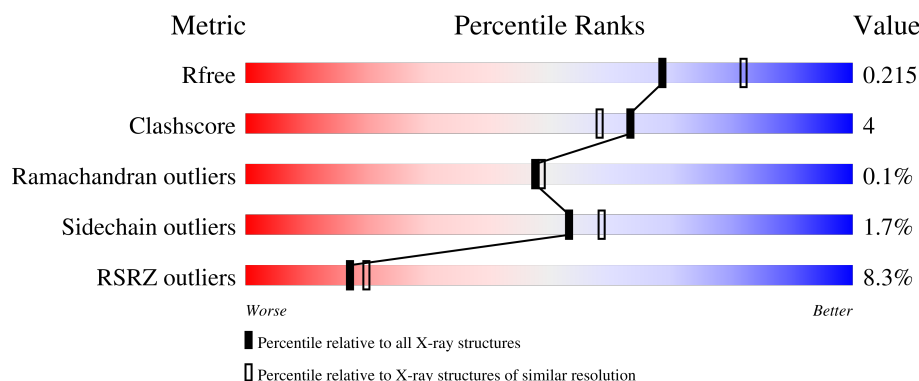
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.14 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	3689 (2.16-2.12)
Clashscore	190562	3812 (2.16-2.12)
Ramachandran outliers	187476	3773 (2.16-2.12)
Sidechain outliers	187428	3772 (2.16-2.12)
RSRZ outliers	180081	3691 (2.16-2.12)

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	7% 90% • 7%
1	B	450	6% 86% 7% • 7%
1	C	450	6% 82% 9% • 8%
1	D	450	12% 83% 10% • 6%

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	RVD	C	451	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13135 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.

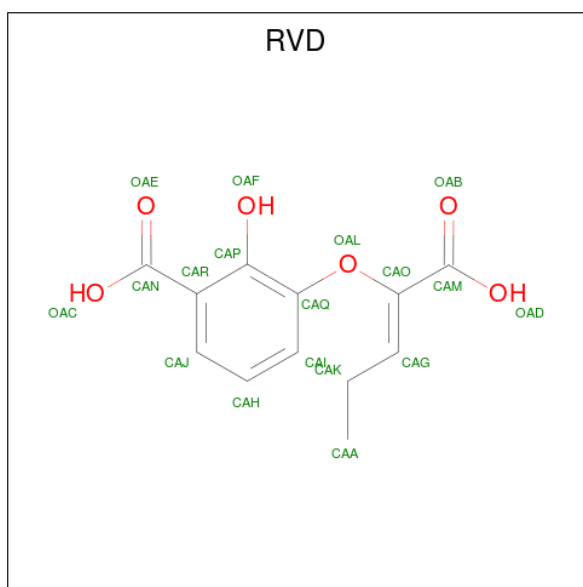
- Molecule 1 is a protein called Isochorismate synthase/isochorismate-pyruvate lyase mbtI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	420	Total	C	N	O	S	0	2	0
			3189	2001	570	608	10			
1	B	420	Total	C	N	O	S	5	2	0
			3188	2004	570	604	10			
1	C	413	Total	C	N	O	S	0	0	0
			3127	1960	557	600	10			
1	D	425	Total	C	N	O	S	0	0	0
			3184	1999	566	609	10			

There are 4 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
C	1	VAL	-	expression tag	UNP Q7D785
D	1	VAL	-	expression tag	UNP Q7D785

- Molecule 2 is 3-([(1Z)-1-carboxybut-1-en-1-yl]oxy)-2-hydroxybenzoic acid (CCD ID: RVD) (formula: C₁₂H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			18	12	6		
2	B	1	Total	C	O	0	0
			18	12	6		
2	C	1	Total	C	O	0	0
			18	12	6		
2	D	1	Total	C	O	0	0
			18	12	6		

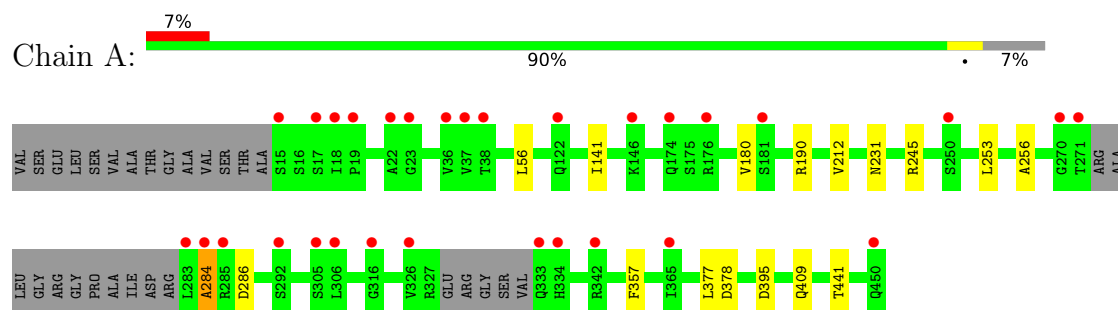
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	111	Total	O	0	0
			111	111		
3	B	102	Total	O	0	0
			102	102		
3	C	102	Total	O	0	0
			102	102		
3	D	60	Total	O	0	0
			60	60		

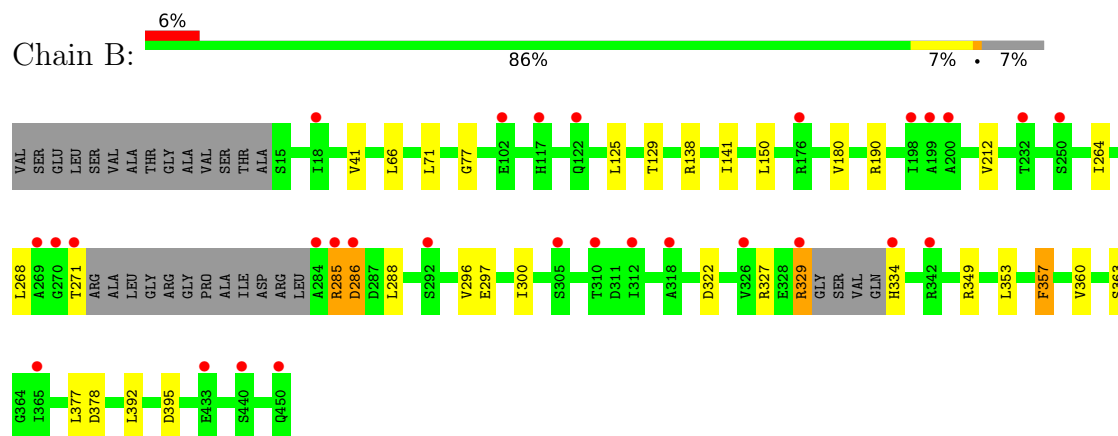
3 Residue-property plots [i](#)

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.

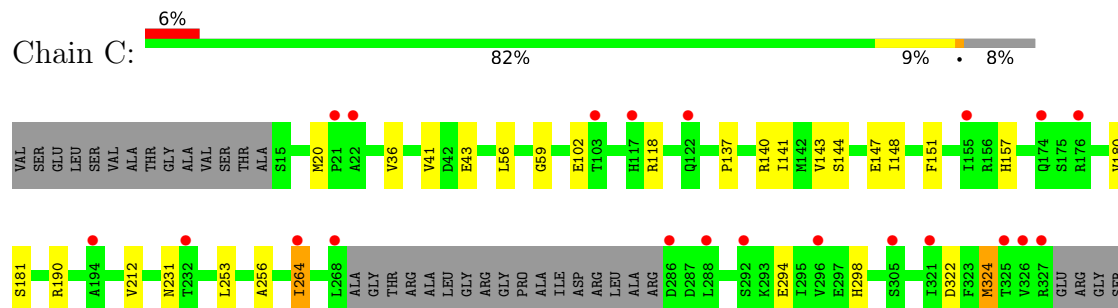
- Molecule 1: Isochorismate synthase/isochorismate-pyruvate lyase mbtI

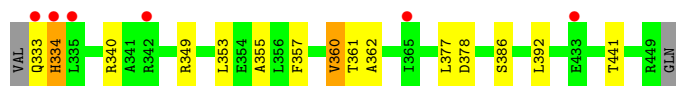


- Molecule 1: Isochorismate synthase/isochorismate-pyruvate lyase mbtI

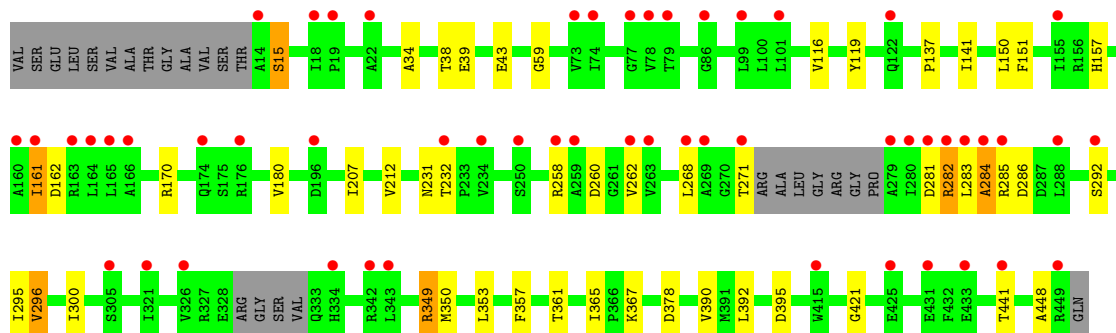
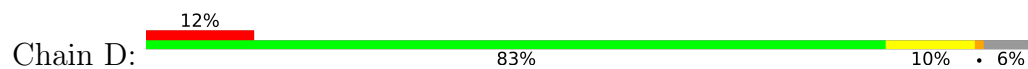


- Molecule 1: Isochorismate synthase/isochorismate-pyruvate lyase mbtI





- 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	87.27Å 115.47Å 95.47Å 90.00° 91.26° 90.00°	Depositor
Resolution (Å)	19.84 – 2.14 19.84 – 2.15	Depositor EDS
% Data completeness (in resolution range)	(Not available) (19.84-2.14) 98.4 (19.84-2.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 2.15Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, R_{free}	0.207 , 0.229 (Not available) , 0.215	Depositor DCC
R_{free} test set	5102 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	36.1	Xtriage
Anisotropy	0.081	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.39$	Xtriage
Estimated twinning fraction	0.006 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13135	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% 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: RVD

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	0.84	1/3246 (0.0%)	1.25	3/4411 (0.1%)
1	B	0.82	2/3249 (0.1%)	1.27	6/4411 (0.1%)
1	C	0.84	1/3181 (0.0%)	1.25	7/4325 (0.2%)
1	D	0.87	2/3239 (0.1%)	1.25	7/4405 (0.2%)
All	All	0.84	6/12915 (0.0%)	1.26	23/17552 (0.1%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	363	SER	CA-C	6.29	1.56	1.52
1	D	232	THR	N-CA	-5.30	1.41	1.46
1	D	295	ILE	C-O	-5.23	1.18	1.24
1	C	324	MET	SD-CE	-5.15	1.66	1.79
1	B	357	PHE	C-O	-5.12	1.17	1.24
1	A	357	PHE	C-O	-5.08	1.17	1.24

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	329[A]	ARG	CB-CA-C	10.86	130.74	110.10
1	B	329[B]	ARG	CB-CA-C	10.86	130.74	110.10
1	C	36	VAL	N-CA-C	8.23	118.27	110.53
1	D	232	THR	CB-CA-C	7.02	121.45	110.71
1	C	360	VAL	CA-C-N	6.12	129.31	120.38
1	C	360	VAL	C-N-CA	6.12	129.31	120.38
1	C	181	SER	N-CA-C	6.02	117.84	111.28
1	C	334	HIS	N-CA-C	6.01	118.56	108.23
1	C	102	GLU	CB-CG-CD	5.85	122.55	112.60
1	C	147	GLU	N-CA-C	5.59	117.47	109.14
1	D	349	ARG	N-CA-C	5.44	118.71	111.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	284	ALA	CA-C-N	5.32	131.27	121.70
1	A	284	ALA	C-N-CA	5.32	131.27	121.70
1	D	395	ASP	CA-CB-CG	5.28	117.88	112.60
1	B	395	ASP	CA-CB-CG	5.24	117.83	112.60
1	D	448	ALA	CA-C-N	5.23	131.12	121.70
1	D	448	ALA	C-N-CA	5.23	131.12	121.70
1	B	360	VAL	CA-C-N	5.18	127.95	120.38
1	B	360	VAL	C-N-CA	5.18	127.95	120.38
1	D	390	VAL	N-CA-C	5.13	116.75	108.85
1	A	395	ASP	CA-CB-CG	5.12	117.72	112.60
1	D	15	SER	N-CA-C	5.11	117.14	109.23
1	B	286	ASP	N-CA-C	5.00	117.21	110.06

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	3189	0	3153	10	0
1	B	3188	0	3164	27	0
1	C	3127	0	3080	25	0
1	D	3184	0	3115	36	0
2	A	18	0	9	0	0
2	B	18	0	9	0	0
2	C	18	0	9	7	0
2	D	18	0	9	3	0
3	A	111	0	0	0	0
3	B	102	0	0	0	0
3	C	102	0	0	0	0
3	D	60	0	0	0	0
All	All	13135	0	12548	100	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 (100) 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:284:ALA:N	1:D:285:ARG:HB3	1.42	1.35
1:D:284:ALA:HA	1:D:286:ASP:N	1.44	1.31
1:D:284:ALA:H	1:D:285:ARG:CB	1.51	1.23
1:B:285:ARG:O	1:B:285:ARG:HD3	1.45	1.15
1:D:284:ALA:CA	1:D:286:ASP:H	1.60	1.12
1:D:284:ALA:HA	1:D:286:ASP:H	0.82	0.96
1:C:361:THR:HG21	2:C:451:RVD:HAAA	1.48	0.95
1:A:284:ALA:HA	1:A:286:ASP:H	1.31	0.93
1:B:285:ARG:HD3	1:B:285:ARG:C	1.92	0.92
1:C:361:THR:HG21	2:C:451:RVD:CAA	2.00	0.90
1:B:285:ARG:NH2	1:B:288:LEU:HB2	1.86	0.88
1:D:260:ASP:OD1	1:D:262:VAL:HG12	1.75	0.87
1:C:361:THR:CG2	2:C:451:RVD:HAAA	2.10	0.80
1:D:283:LEU:O	1:D:284:ALA:CB	2.30	0.80
1:B:77:GLY:HA3	1:D:170:ARG:HD3	1.63	0.79
1:B:285:ARG:HH22	1:B:288:LEU:CB	1.96	0.79
1:B:285:ARG:C	1:B:285:ARG:CD	2.53	0.78
1:C:20:MET:HE3	1:C:144:SER:O	1.85	0.76
1:B:285:ARG:O	1:B:285:ARG:CD	2.30	0.76
1:D:284:ALA:CA	1:D:286:ASP:N	2.32	0.74
1:B:285:ARG:NH2	1:B:288:LEU:H	1.86	0.73
1:C:20:MET:HE1	1:C:143:VAL:HG13	1.72	0.72
1:D:157:HIS:O	1:D:161:ILE:HG23	1.90	0.71
1:B:285:ARG:HH22	1:B:288:LEU:HB2	1.50	0.69
1:B:285:ARG:NH2	1:B:288:LEU:CB	2.52	0.68
1:D:283:LEU:O	1:D:284:ALA:HB3	1.95	0.66
1:D:281:ASP:O	1:D:282:ARG:HB2	1.98	0.62
1:D:141:ILE:HD12	1:D:150:LEU:HD22	1.82	0.62
1:B:141:ILE:HD12	1:B:150:LEU:HD22	1.82	0.61
1:B:285:ARG:O	1:B:285:ARG:NH1	2.34	0.60
2:D:451:RVD:HAI	2:D:451:RVD:CAM	2.33	0.58
1:C:361:THR:CG2	2:C:451:RVD:CAA	2.75	0.57
1:A:231:ASN:OD1	1:A:441:THR:HG23	2.05	0.57
1:D:284:ALA:N	1:D:286:ASP:H	2.04	0.55
1:A:284:ALA:HA	1:A:286:ASP:N	2.11	0.55
1:D:284:ALA:N	1:D:285:ARG:CB	2.30	0.55
1:B:66:LEU:HD13	1:B:71:LEU:HD12	1.89	0.55
1:D:207:ILE:HG13	1:D:421:GLY:HA2	1.89	0.54
1:C:353:LEU:O	1:C:357:PHE:HB2	2.09	0.52
1:B:285:ARG:HH21	1:B:288:LEU:H	1.57	0.52
1:C:41:VAL:HG21	1:C:157:HIS:CE1	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:180:VAL:HG12	1:B:212:VAL:HG11	1.92	0.52
1:D:284:ALA:H	1:D:285:ARG:HB3	0.57	0.51
1:C:231:ASN:OD1	1:C:441:THR:HG23	2.11	0.51
1:A:180:VAL:HG12	1:A:212:VAL:HG11	1.93	0.50
1:B:190:ARG:HD2	1:B:377:LEU:O	2.10	0.50
1:C:20:MET:HE2	1:C:148:ILE:HG13	1.93	0.50
1:D:34:ALA:O	1:D:38:THR:OG1	2.26	0.50
1:C:43:GLU:HB2	1:C:59:GLY:HA2	1.94	0.50
1:B:322:ASP:HB3	1:B:327:ARG:HH22	1.77	0.49
1:D:292:SER:O	1:D:296:VAL:HG12	2.12	0.49
1:B:296:VAL:O	1:B:300:ILE:HG12	2.12	0.48
1:A:56:LEU:HD23	1:A:141:ILE:HD12	1.94	0.48
1:C:180:VAL:HG12	1:C:212:VAL:HG11	1.95	0.48
1:D:180:VAL:HG12	1:D:212:VAL:HG11	1.95	0.48
1:D:258:ARG:HG3	1:D:262:VAL:HG13	1.94	0.47
1:D:43:GLU:OE2	1:D:157:HIS:NE2	2.44	0.47
1:C:190:ARG:HD2	1:C:377:LEU:O	2.14	0.47
1:D:284:ALA:CA	1:D:285:ARG:HB3	2.35	0.47
1:C:361:THR:HG21	2:C:451:RVD:HAAB	1.91	0.47
1:C:56:LEU:HD23	1:C:141:ILE:HD12	1.96	0.46
1:A:245[B]:ARG:HB2	1:A:409:GLN:HB3	1.97	0.46
1:C:264:ILE:CD1	1:C:340:ARG:HG3	2.45	0.46
1:C:298:HIS:HB3	1:C:324:MET:HE2	1.98	0.46
2:C:451:RVD:CAM	2:C:451:RVD:HAI	2.45	0.46
1:D:161:ILE:HG13	1:D:162:ASP:N	2.31	0.46
1:B:285:ARG:HH22	1:B:288:LEU:HB3	1.76	0.46
1:B:353:LEU:O	1:B:357:PHE:HB2	2.16	0.46
1:A:245[A]:ARG:HB2	1:A:409:GLN:HB3	1.98	0.45
1:C:118:ARG:NH1	1:C:355:ALA:O	2.49	0.45
1:C:59:GLY:O	1:C:137:PRO:HA	2.16	0.45
1:C:20:MET:CE	1:C:148:ILE:HG13	2.47	0.45
1:C:362:ALA:HB1	1:C:386:SER:HB2	1.99	0.45
1:D:15:SER:HB3	1:D:151:PHE:CE1	2.52	0.45
1:C:333:GLN:HG3	1:C:334:HIS:H	1.81	0.45
1:A:190:ARG:HD2	1:A:377:LEU:O	2.17	0.44
2:C:451:RVD:CAM	2:C:451:RVD:CAI	2.93	0.44
1:D:300:ILE:HG21	1:D:365:ILE:HG21	1.99	0.44
1:B:285:ARG:NH2	1:B:288:LEU:N	2.59	0.44
2:D:451:RVD:CAM	2:D:451:RVD:CAI	2.87	0.44
1:B:322:ASP:HB3	1:B:327:ARG:NH2	2.32	0.44
1:D:282:ARG:HB3	1:D:283:LEU:H	1.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:283:LEU:O	1:D:284:ALA:HB2	2.12	0.44
1:D:284:ALA:CA	1:D:285:ARG:CB	2.95	0.44
1:B:285:ARG:HH21	1:B:288:LEU:N	2.14	0.43
1:D:119:TYR:OH	1:D:367:LYS:HD2	2.18	0.43
1:D:361:THR:HG21	2:D:451:RVD:HAK	2.00	0.43
1:D:353:LEU:O	1:D:357:PHE:HB2	2.19	0.43
1:B:285:ARG:CZ	1:B:286:ASP:HA	2.49	0.42
1:B:138:ARG:NH2	1:D:39:GLU:OE2	2.46	0.42
1:C:140:ARG:HB2	1:C:151:PHE:HB2	2.02	0.41
1:B:125:LEU:HD22	1:B:129:THR:HG21	2.03	0.41
1:D:231:ASN:OD1	1:D:441:THR:HG23	2.20	0.41
1:A:253:LEU:HD21	1:A:256:ALA:HB2	2.03	0.41
1:D:349:ARG:HD2	1:D:392:LEU:HD23	2.03	0.41
1:A:245[B]:ARG:NH1	1:A:409:GLN:OE1	2.52	0.41
1:B:349:ARG:HD2	1:B:392:LEU:HD23	2.03	0.40
1:C:253:LEU:HD21	1:C:256:ALA:HB2	2.04	0.40
1:C:349:ARG:HD2	1:C:392:LEU:HD23	2.03	0.40
1:D:59:GLY:O	1:D:137:PRO:HA	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	416/450 (92%)	411 (99%)	5 (1%)	0	100	100
1	B	415/450 (92%)	410 (99%)	5 (1%)	0	100	100
1	C	407/450 (90%)	403 (99%)	4 (1%)	0	100	100
1	D	419/450 (93%)	411 (98%)	6 (1%)	2 (0%)	24	19
All	All	1657/1800 (92%)	1635 (99%)	20 (1%)	2 (0%)	48	49

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	282	ARG
1	D	284	ALA

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	327/358 (91%)	326 (100%)	1 (0%)	86	90
1	B	327/358 (91%)	317 (97%)	10 (3%)	35	35
1	C	323/358 (90%)	318 (98%)	5 (2%)	57	63
1	D	320/358 (89%)	313 (98%)	7 (2%)	45	49
All	All	1297/1432 (91%)	1274 (98%)	23 (2%)	53	57

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	378	ASP
1	B	41	VAL
1	B	264	ILE
1	B	268	LEU
1	B	271	THR
1	B	285	ARG
1	B	297	GLU
1	B	329[A]	ARG
1	B	329[B]	ARG
1	B	334	HIS
1	B	378	ASP
1	C	264	ILE
1	C	294	GLU
1	C	322	ASP
1	C	360	VAL
1	C	378	ASP
1	D	116	VAL
1	D	161	ILE
1	D	268	LEU
1	D	271	THR

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Mol	Chain	Res	Type
1	D	296	VAL
1	D	350	MET
1	D	378	ASP

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	82	GLN
1	B	409	GLN
1	C	291	ASN
1	D	291	ASN
1	D	409	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	$\# Z > 2$
2	RVD	B	451	-	18,18,18	2.24	7 (38%)	23,24,24	1.31	4 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	RVD	C	451	-	18,18,18	2.44	7 (38%)	23,24,24	1.16	2 (8%)
2	RVD	A	451	-	18,18,18	2.67	9 (50%)	23,24,24	1.95	6 (26%)
2	RVD	D	451	-	18,18,18	2.37	6 (33%)	23,24,24	1.84	4 (17%)

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
2	RVD	B	451	-	-	3/15/15/15	0/1/1/1
2	RVD	C	451	-	-	4/15/15/15	0/1/1/1
2	RVD	A	451	-	-	3/15/15/15	0/1/1/1
2	RVD	D	451	-	-	4/15/15/15	0/1/1/1

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	451	RVD	OAL-CAQ	-6.06	1.29	1.41
2	B	451	RVD	CAK-CAG	-6.01	1.36	1.50
2	D	451	RVD	CAK-CAG	-5.92	1.37	1.50
2	C	451	RVD	CAK-CAG	-5.54	1.37	1.50
2	A	451	RVD	CAK-CAG	-4.87	1.39	1.50
2	C	451	RVD	OAL-CAQ	-4.65	1.31	1.41
2	B	451	RVD	OAL-CAQ	-3.96	1.33	1.41
2	A	451	RVD	OAL-CAO	-3.79	1.30	1.38
2	D	451	RVD	OAL-CAQ	-3.77	1.33	1.41
2	C	451	RVD	CAQ-CAP	-3.40	1.35	1.40
2	D	451	RVD	CAG-CAO	-3.31	1.30	1.33
2	C	451	RVD	OAD-CAM	-3.22	1.21	1.30
2	A	451	RVD	OAD-CAM	-3.18	1.21	1.30
2	D	451	RVD	CAR-CAP	-3.00	1.36	1.41
2	C	451	RVD	OAC-CAN	-2.99	1.21	1.30
2	D	451	RVD	OAC-CAN	-2.95	1.21	1.30
2	A	451	RVD	CAR-CAP	-2.93	1.36	1.41
2	D	451	RVD	OAD-CAM	-2.91	1.22	1.30
2	B	451	RVD	CAQ-CAP	-2.69	1.36	1.40
2	B	451	RVD	CAO-CAM	2.67	1.54	1.48
2	A	451	RVD	CAQ-CAP	-2.60	1.36	1.40
2	A	451	RVD	OAC-CAN	-2.59	1.22	1.30
2	A	451	RVD	CAG-CAO	-2.57	1.31	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	451	RVD	CAR-CAP	-2.52	1.37	1.41
2	B	451	RVD	CAR-CAP	-2.35	1.37	1.41
2	B	451	RVD	OAD-CAM	-2.27	1.24	1.30
2	B	451	RVD	OAC-CAN	-2.10	1.24	1.30
2	C	451	RVD	OAF-CAP	-2.03	1.32	1.36
2	A	451	RVD	OAF-CAP	-2.00	1.32	1.36

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	451	RVD	CAK-CAG-CAO	-5.21	118.37	126.33
2	D	451	RVD	CAK-CAG-CAO	-5.07	118.59	126.33
2	D	451	RVD	CAQ-OAL-CAO	-4.22	111.54	118.49
2	A	451	RVD	OAL-CAO-CAG	-3.52	111.60	118.86
2	A	451	RVD	CAQ-OAL-CAO	-3.16	113.28	118.49
2	B	451	RVD	OAL-CAO-CAM	2.79	122.14	114.54
2	A	451	RVD	CAI-CAQ-CAP	2.79	122.56	119.98
2	C	451	RVD	OAL-CAO-CAM	2.72	121.94	114.54
2	B	451	RVD	OAE-CAN-CAR	-2.63	115.69	121.97
2	D	451	RVD	OAL-CAO-CAM	2.52	121.40	114.54
2	B	451	RVD	OAD-CAM-CAO	2.38	119.46	114.06
2	A	451	RVD	OAE-CAN-CAR	-2.35	116.35	121.97
2	C	451	RVD	CAK-CAG-CAO	-2.31	122.81	126.33
2	A	451	RVD	OAD-CAM-CAO	2.15	118.94	114.06
2	D	451	RVD	OAD-CAM-CAO	2.03	118.66	114.06
2	B	451	RVD	CAK-CAG-CAO	-2.01	123.26	126.33

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	451	RVD	CAK-CAG-CAO-CAM
2	A	451	RVD	OAB-CAM-CAO-CAG
2	B	451	RVD	OAB-CAM-CAO-OAL
2	B	451	RVD	OAD-CAM-CAO-OAL
2	C	451	RVD	OAB-CAM-CAO-OAL
2	C	451	RVD	OAD-CAM-CAO-OAL
2	A	451	RVD	OAD-CAM-CAO-CAG
2	C	451	RVD	OAB-CAM-CAO-CAG
2	C	451	RVD	OAD-CAM-CAO-CAG
2	D	451	RVD	OAB-CAM-CAO-CAG
2	D	451	RVD	OAD-CAM-CAO-CAG

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Mol	Chain	Res	Type	Atoms
2	D	451	RVD	OAB-CAM-CAO-OAL
2	D	451	RVD	OAD-CAM-CAO-OAL
2	B	451	RVD	OAD-CAM-CAO-CAG

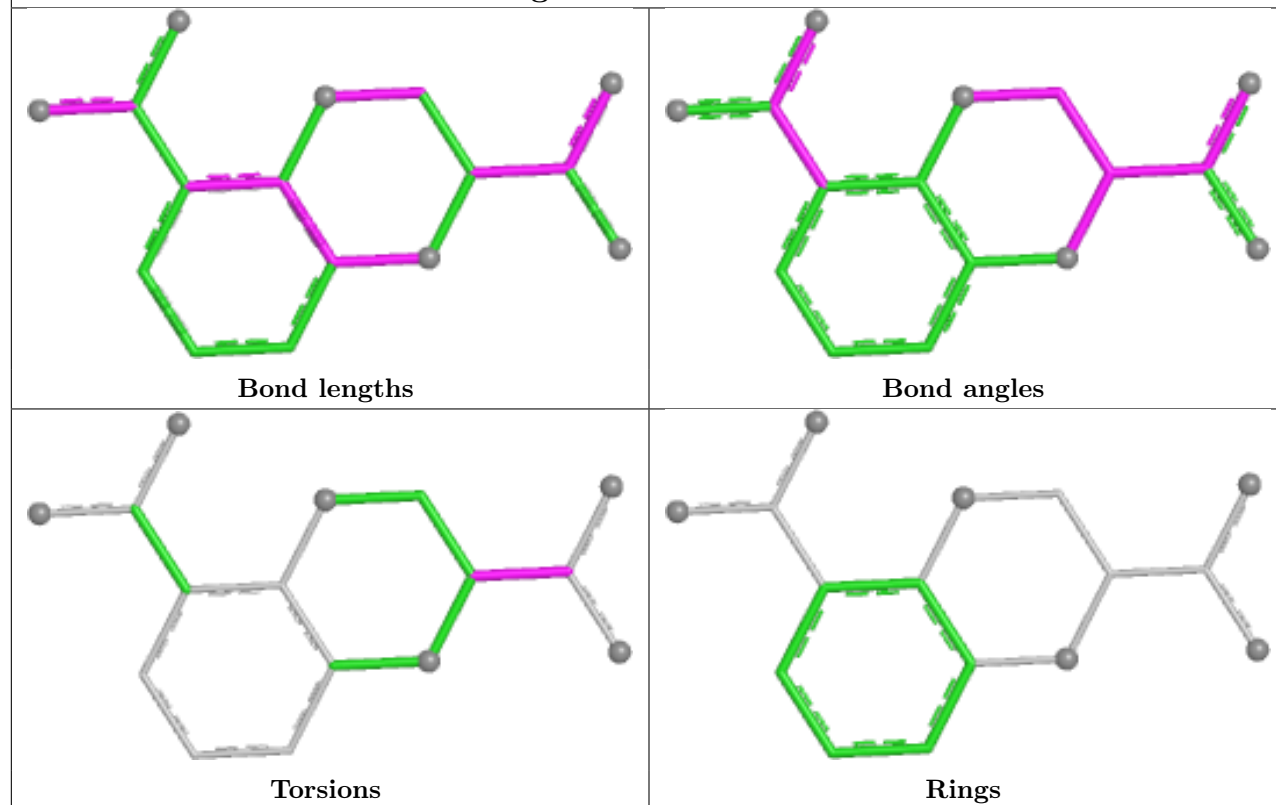
There are no ring outliers.

2 monomers are involved in 10 short contacts:

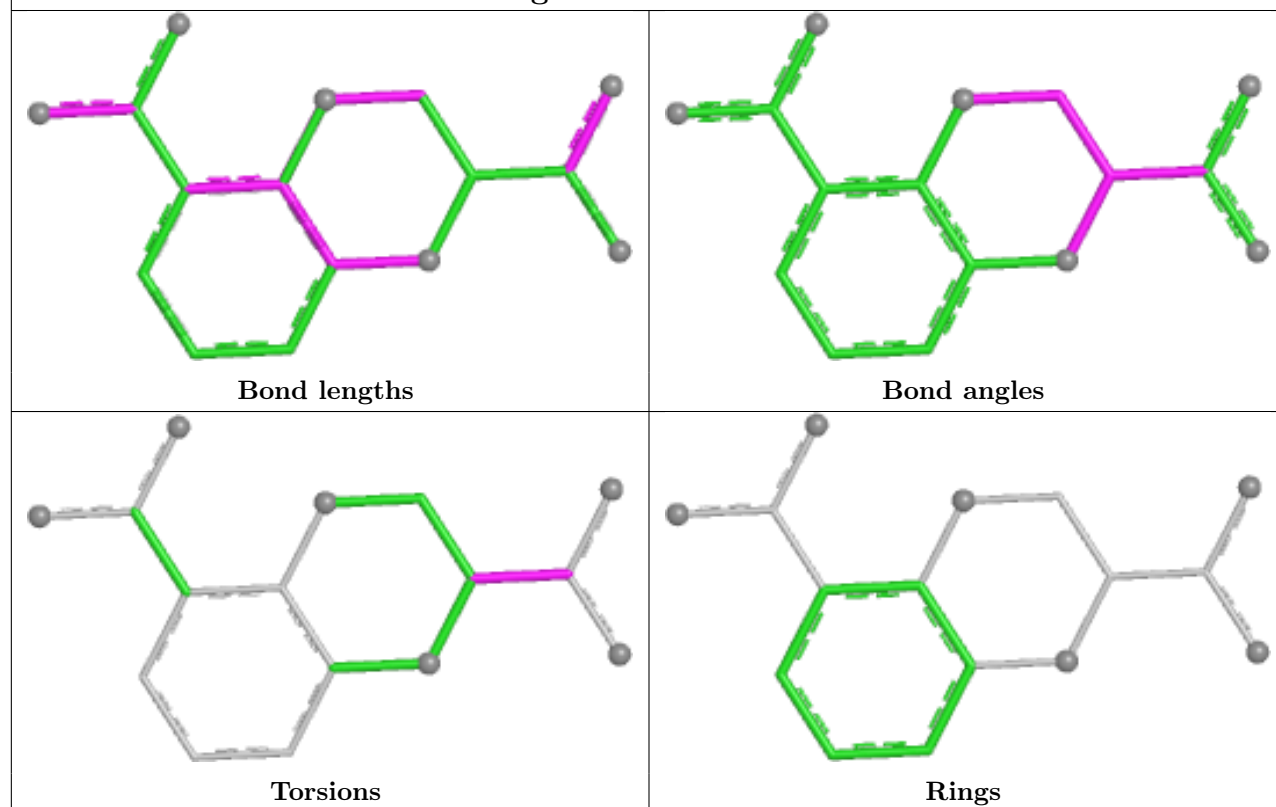
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	451	RVD	7	0
2	D	451	RVD	3	0

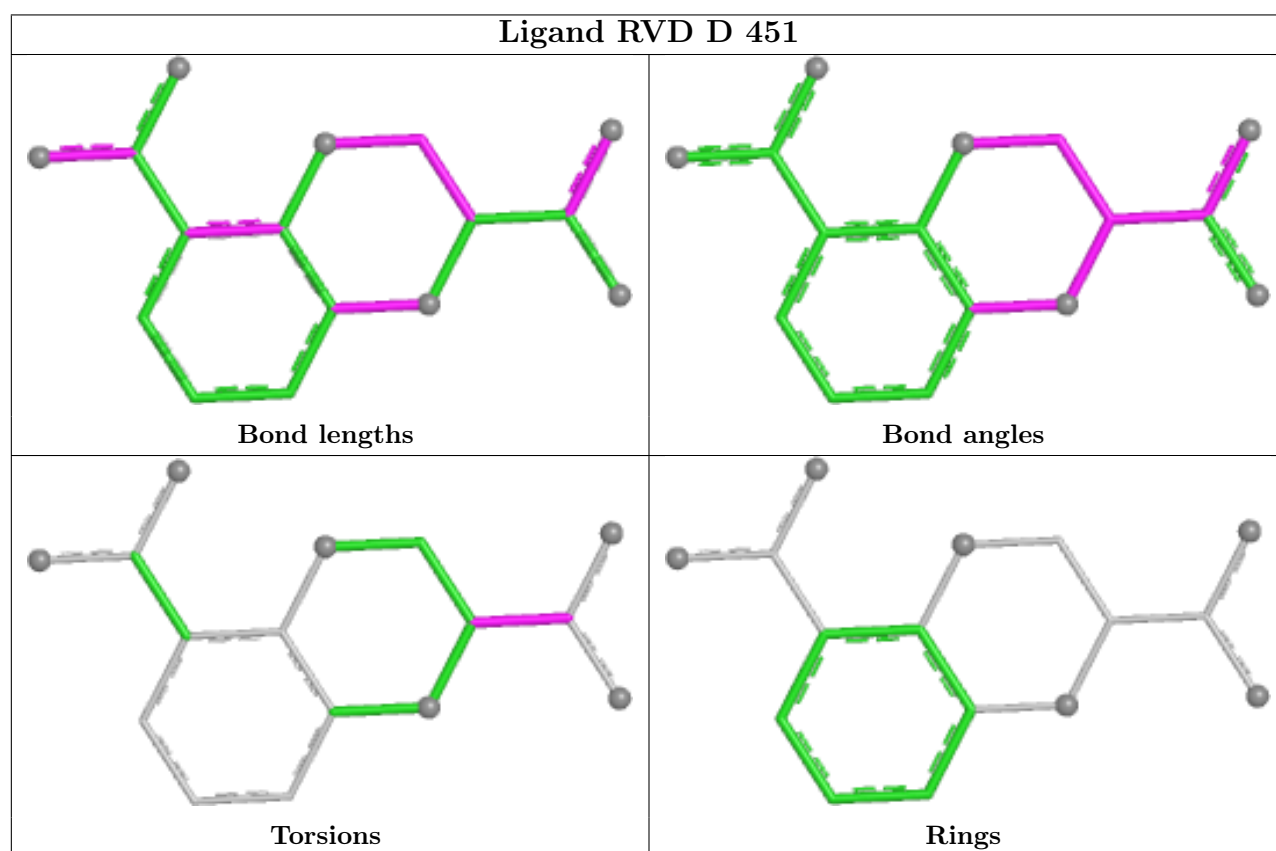
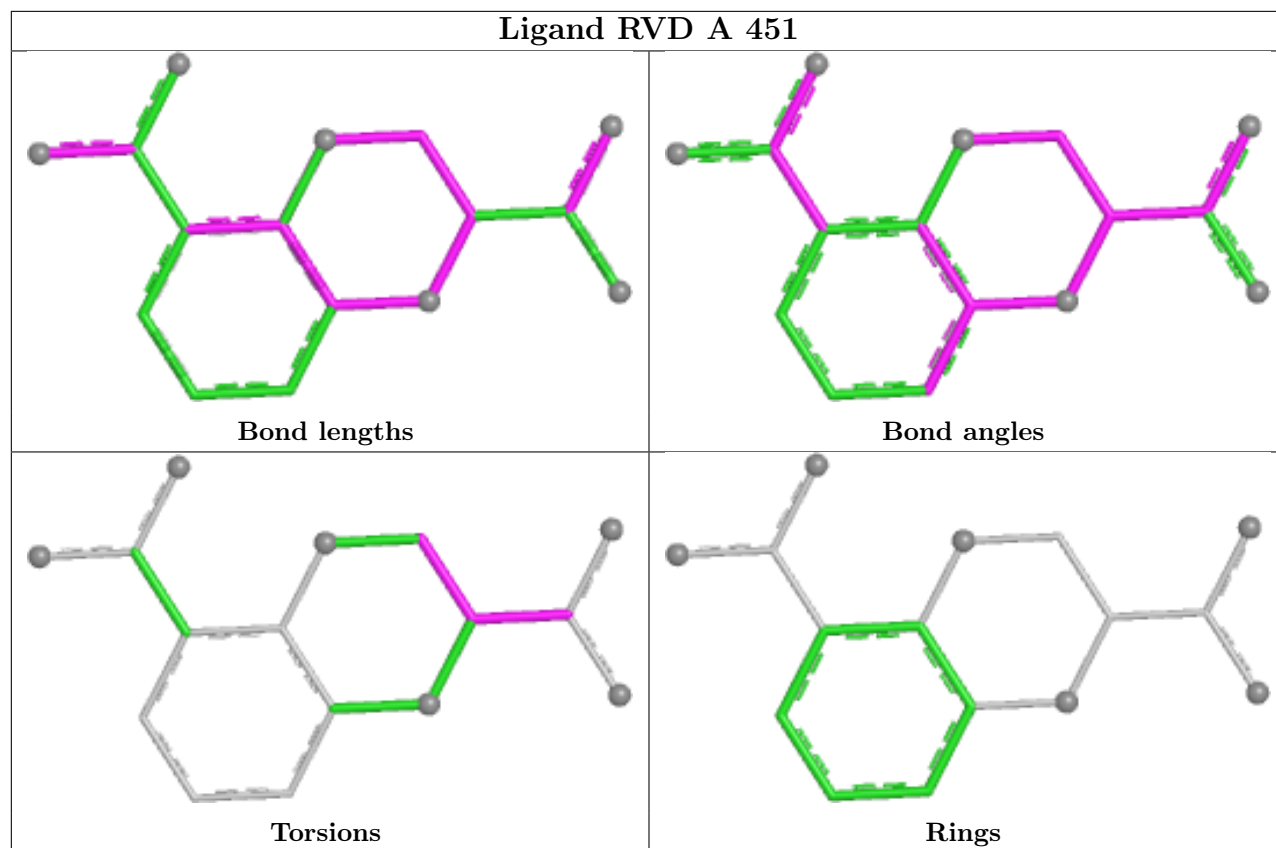
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

Ligand RVD B 451



Ligand RVD C 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	420/450 (93%)	0.40	30 (7%)	22	25	10, 37, 62, 90	10 (2%)
1	B	419/450 (93%)	0.43	29 (6%)	23	26	14, 37, 61, 82	10 (2%)
1	C	413/450 (91%)	0.45	27 (6%)	25	29	13, 38, 60, 87	8 (1%)
1	D	425/450 (94%)	0.90	54 (12%)	8	9	16, 48, 76, 91	8 (1%)
All	All	1677/1800 (93%)	0.55	140 (8%)	17	20	10, 40, 68, 91	36 (2%)

All (140) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	342	ARG	4.7
1	B	271	THR	4.5
1	D	279	ALA	4.4
1	D	326	VAL	4.4
1	D	282	ARG	4.4
1	A	283	LEU	4.4
1	A	271	THR	4.2
1	A	285[A]	ARG	4.1
1	D	176	ARG	4.1
1	D	78	VAL	4.1
1	B	176	ARG	4.0
1	B	329[A]	ARG	4.0
1	A	305	SER	3.9
1	C	292	SER	3.9
1	A	176	ARG	3.9
1	D	14	ALA	3.9
1	D	271	THR	3.9
1	D	73	VAL	3.8
1	B	292	SER	3.8
1	C	288	LEU	3.8
1	C	176	ARG	3.7

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Mol	Chain	Res	Type	RSRZ
1	D	280	ILE	3.7
1	B	269	ALA	3.6
1	C	342	ARG	3.5
1	D	86	GLY	3.5
1	D	283	LEU	3.4
1	C	335	LEU	3.4
1	D	288	LEU	3.4
1	A	18	ILE	3.3
1	D	343	LEU	3.3
1	D	415	TRP	3.2
1	A	334	HIS	3.2
1	D	74	ILE	3.2
1	B	270	GLY	3.2
1	D	305	SER	3.2
1	C	333	GLN	3.2
1	C	232	THR	3.2
1	D	284	ALA	3.1
1	B	342	ARG	3.1
1	C	433	GLU	3.1
1	D	174	GLN	3.1
1	D	292	SER	3.1
1	C	286	ASP	3.1
1	D	263	VAL	3.1
1	D	79	THR	3.0
1	D	433	GLU	3.0
1	A	37	VAL	3.0
1	C	155	ILE	3.0
1	B	305	SER	3.0
1	B	285	ARG	3.0
1	C	103	THR	3.0
1	C	174	GLN	2.9
1	C	326	VAL	2.9
1	B	433	GLU	2.9
1	A	365	ILE	2.9
1	C	334	HIS	2.9
1	D	281	ASP	2.9
1	A	326	VAL	2.8
1	A	15	SER	2.8
1	B	334	HIS	2.8
1	B	284	ALA	2.8
1	A	36	VAL	2.8
1	D	196	ASP	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	22	ALA	2.7
1	A	333	GLN	2.7
1	B	450	GLN	2.7
1	D	19	PRO	2.7
1	B	310	THR	2.7
1	B	250	SER	2.7
1	D	18	ILE	2.7
1	A	250	SER	2.6
1	C	21	PRO	2.6
1	B	326	VAL	2.6
1	D	165	LEU	2.6
1	A	19	PRO	2.6
1	A	270	GLY	2.5
1	D	232	THR	2.5
1	C	321	ILE	2.5
1	D	321	ILE	2.5
1	B	286	ASP	2.5
1	A	174	GLN	2.5
1	D	101	LEU	2.5
1	B	232	THR	2.5
1	B	200	ALA	2.5
1	A	284	ALA	2.4
1	D	285	ARG	2.4
1	B	102	GLU	2.4
1	C	305	SER	2.4
1	D	425	GLU	2.4
1	B	198	ILE	2.4
1	C	264	ILE	2.4
1	D	262	VAL	2.4
1	A	122	GLN	2.3
1	D	164	LEU	2.3
1	A	146	LYS	2.3
1	A	342	ARG	2.3
1	C	327	ARG	2.3
1	B	440	SER	2.3
1	D	77	GLY	2.3
1	A	292	SER	2.3
1	B	122	GLN	2.3
1	C	325	THR	2.3
1	B	117	HIS	2.3
1	C	117	HIS	2.3
1	B	365	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	122	GLN	2.3
1	D	122	GLN	2.3
1	C	296	VAL	2.3
1	D	431	GLU	2.2
1	B	18	ILE	2.2
1	D	155	ILE	2.2
1	D	161	ILE	2.2
1	D	160	ALA	2.2
1	D	166	ALA	2.2
1	D	334	HIS	2.2
1	D	259	ALA	2.2
1	D	269	ALA	2.2
1	D	268	LEU	2.2
1	D	234	VAL	2.2
1	B	199	ALA	2.2
1	A	17	SER	2.2
1	D	258	ARG	2.1
1	B	318	ALA	2.1
1	C	22	ALA	2.1
1	D	250	SER	2.1
1	A	38	THR	2.1
1	D	163	ARG	2.1
1	D	449	ARG	2.1
1	A	22	ALA	2.1
1	C	194	ALA	2.1
1	A	306	LEU	2.1
1	A	450	GLN	2.1
1	D	441	THR	2.1
1	A	23	GLY	2.1
1	A	316	GLY	2.1
1	C	365	ILE	2.1
1	C	268	LEU	2.0
1	D	99	LEU	2.0
1	B	312	ILE	2.0
1	A	181	SER	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 [i](#)

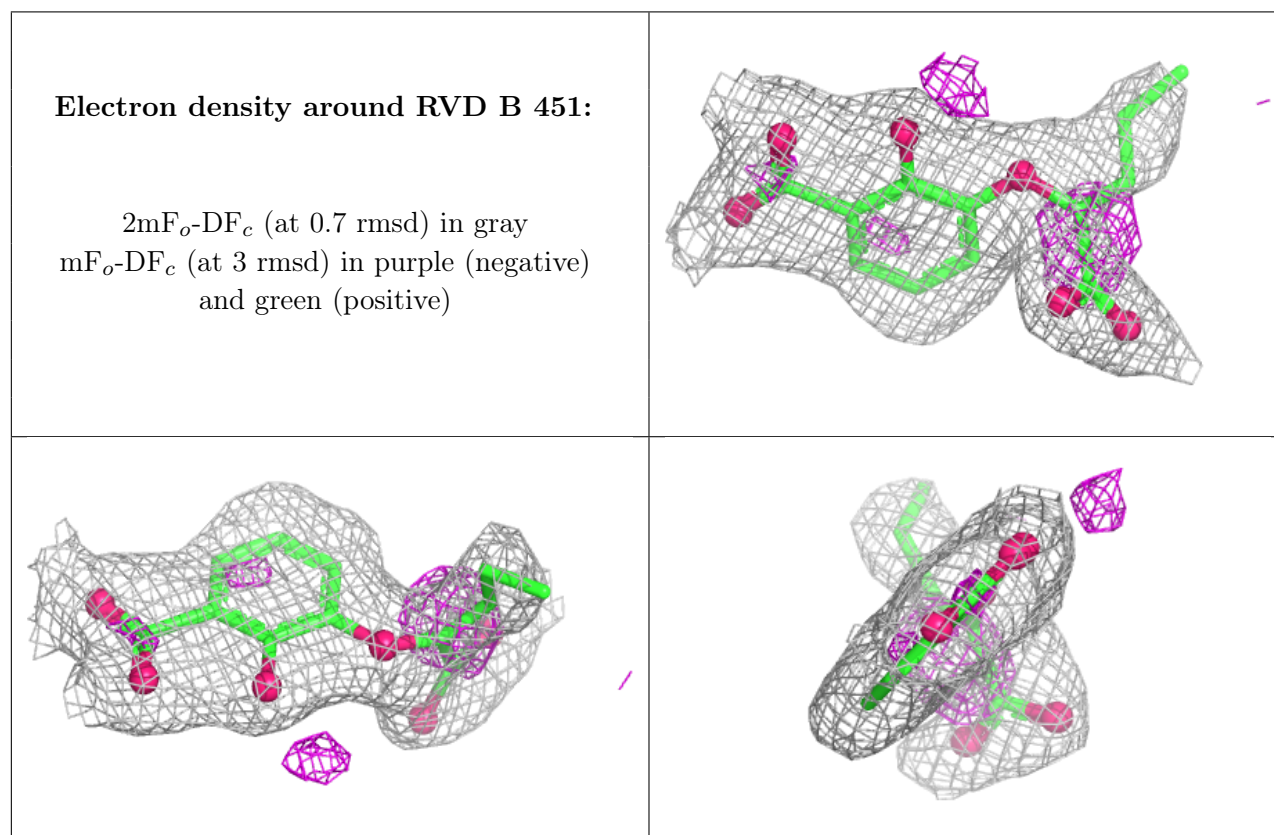
There are no oligosaccharides in this entry.

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.

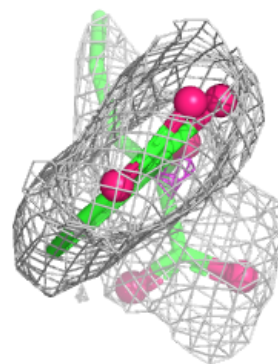
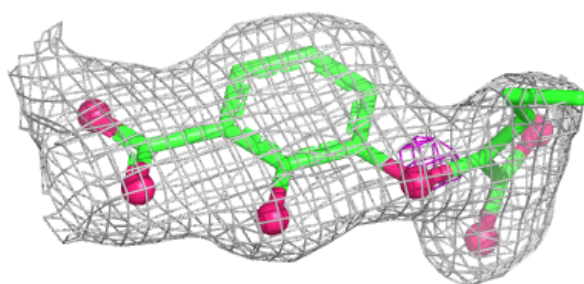
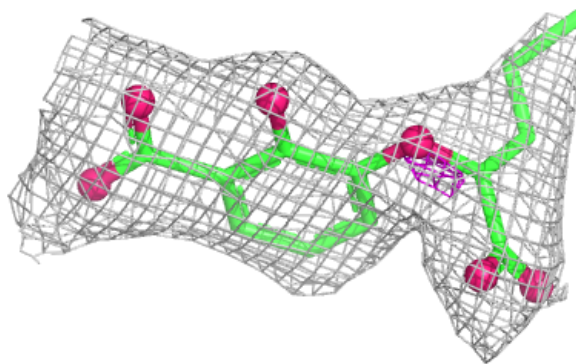
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	RVD	B	451	18/18	0.84	0.10	33,38,51,51	0
2	RVD	D	451	18/18	0.88	0.10	37,46,62,64	0
2	RVD	C	451	18/18	0.90	0.08	32,34,50,51	0
2	RVD	A	451	18/18	0.90	0.08	23,33,45,47	0

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.

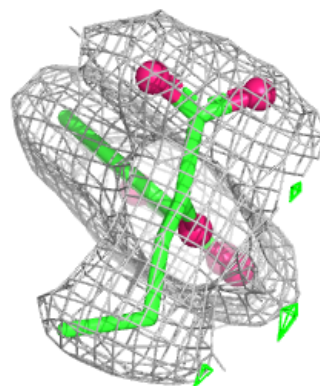
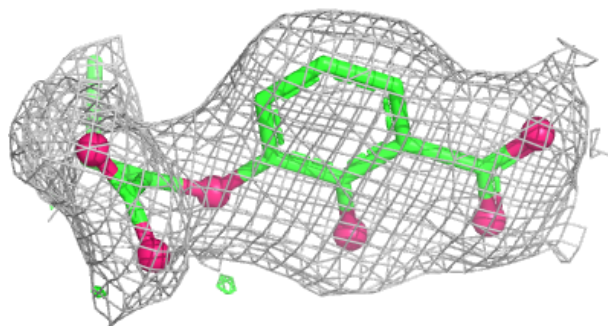
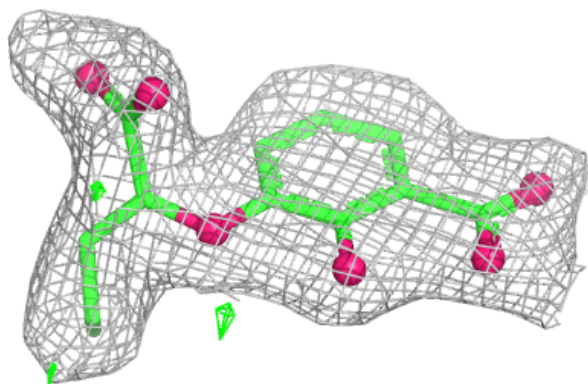


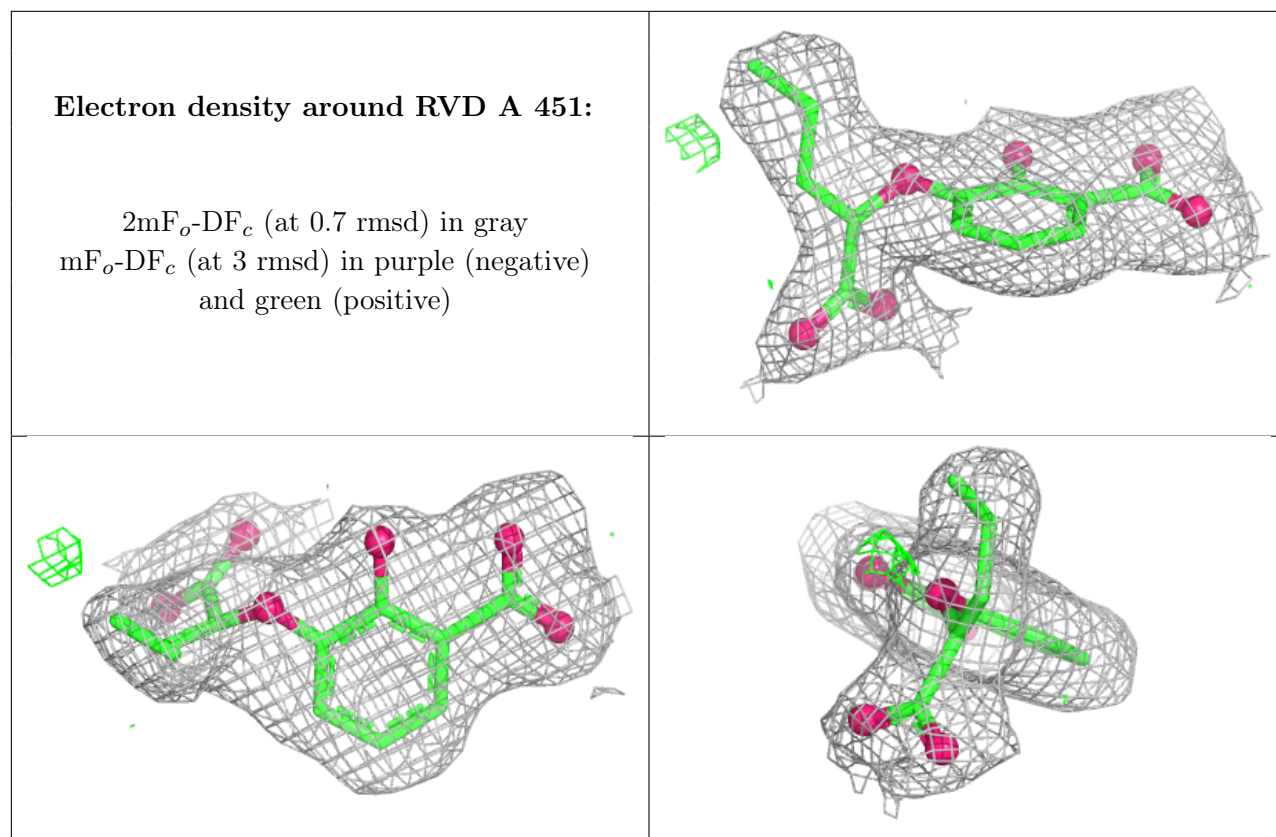
Electron density around RVD D 451:

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

**Electron density around RVD C 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.