wwPDB X-ray Structure Validation Summary Report (i)

Aug 30, 2023 – 05:30 AM EDT

PDB ID	:	3N86
Title	:	Crystal structure of 3-dehydroquinate dehydratase from Mycobacterium tuberculosis in complex with inhibitor 4
Authors	:	Dias, M.V.B.; Snee, W.C.; Bromfield, K.M.; Payne, R.; Palaninathan, S.K.; Ciulli, A.; Howard, N.I.; Abell, C.; Sacchettini, J.C.; Blundell, T.L.
Deposited on	:	2010-05-27
Resolution	:	1.90 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (i)) were used in the production of this report:

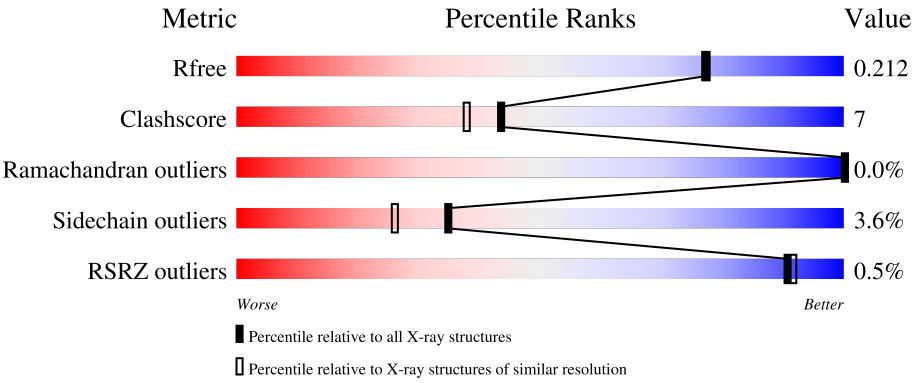
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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 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 <=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	147	86% 8%
1	B	147	85% 8%
1	C	147	85% 10%
1	D	147	84% 11%

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Mol	Chain	Length	Quality of chain
1	E	147	82% 12%
1	F	147	86% 7%
1	G	147	84% 10%
1	H	147	81% 12%
1	I	147	86% 10%
1	J	147	82% 11% 5%
1	K	147	80% 14%
1	L	147	83% 12%
1	M	147	79% 14%
1	N	147	84% 10%
1	O	147	85% 10%
1	P	147	82% 12%
1	Q	147	88% 5%
1	R	147	85% 10%
1	S	147	80% 14%
1	T	147	82% 14%
1	U	147	80% 15%
1	V	147	84% 11%
1	W	147	82% 13%
1	X	147	79% 15%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 30436 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called 3-dehydroquinate dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	B	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	C	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	D	141	Total	C	N	O	S	0	1	0
			1079	679	199	200	1			
1	E	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	F	141	Total	C	N	O	S	0	0	0
			1065	671	193	200	1			
1	G	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	H	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	I	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	J	143	Total	C	N	O	S	0	0	0
			1086	682	198	205	1			
1	K	142	Total	C	N	O	S	0	0	0
			1080	679	197	203	1			
1	L	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	M	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	N	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	O	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	P	142	Total	C	N	O	S	0	0	0
			1080	679	197	203	1			

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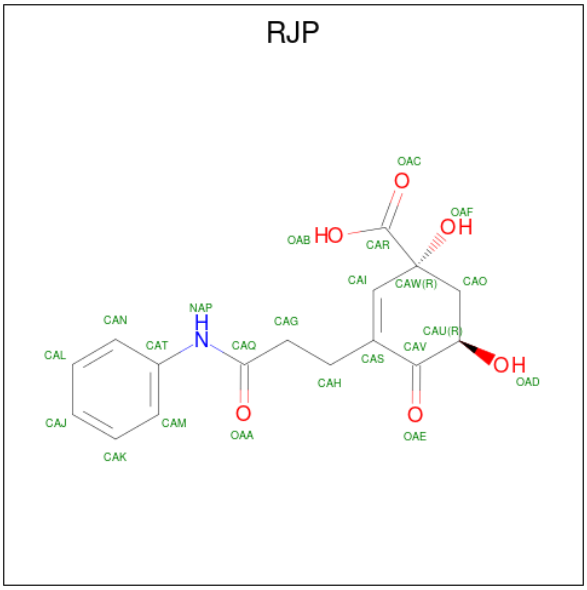
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	R	142	Total	C	N	O	S	0	0	0
			1076	677	197	201	1			
1	S	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	T	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	U	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			
1	V	144	Total	C	N	O	S	0	0	0
			1091	685	199	206	1			
1	W	143	Total	C	N	O	S	0	0	0
			1086	682	198	205	1			
1	X	141	Total	C	N	O	S	0	0	0
			1071	674	196	200	1			

- Molecule 2 is (1R,5R)-1,5-dihydroxy-4-oxo-3-[3-oxo-3-(phenylamino)propyl]cyclohex-2-ene-1-carboxylic acid (three-letter code: RJP) (formula: $C_{16}H_{17}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			23	16	1	6		
2	B	1	Total	C	N	O	0	0
			23	16	1	6		
2	C	1	Total	C	N	O	0	0
			23	16	1	6		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	D	1	Total	C	N	O	0	0
			23	16	1	6		
2	E	1	Total	C	N	O	0	0
			23	16	1	6		
2	F	1	Total	C	N	O	0	0
			23	16	1	6		
2	G	1	Total	C	N	O	0	0
			23	16	1	6		
2	H	1	Total	C	N	O	0	0
			23	16	1	6		
2	I	1	Total	C	N	O	0	0
			23	16	1	6		
2	J	1	Total	C	N	O	0	0
			23	16	1	6		
2	K	1	Total	C	N	O	0	0
			23	16	1	6		
2	L	1	Total	C	N	O	0	0
			23	16	1	6		
2	M	1	Total	C	N	O	0	0
			23	16	1	6		
2	N	1	Total	C	N	O	0	0
			23	16	1	6		
2	O	1	Total	C	N	O	0	0
			23	16	1	6		
2	P	1	Total	C	N	O	0	0
			23	16	1	6		
2	Q	1	Total	C	N	O	0	0
			23	16	1	6		
2	R	1	Total	C	N	O	0	0
			23	16	1	6		
2	S	1	Total	C	N	O	0	0
			23	16	1	6		
2	T	1	Total	C	N	O	0	0
			23	16	1	6		
2	U	1	Total	C	N	O	0	0
			23	16	1	6		
2	V	1	Total	C	N	O	0	0
			23	16	1	6		
2	W	1	Total	C	N	O	0	0
			23	16	1	6		
2	X	1	Total	C	N	O	0	0
			23	16	1	6		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	169	Total O 169 169	0	0
3	B	162	Total O 162 162	0	0
3	C	177	Total O 177 177	0	0
3	D	165	Total O 165 165	0	0
3	E	169	Total O 169 169	0	0
3	F	164	Total O 164 164	0	0
3	G	155	Total O 155 155	0	0
3	H	175	Total O 175 175	0	0
3	I	159	Total O 159 159	0	0
3	J	189	Total O 189 189	0	0
3	K	185	Total O 185 185	0	0
3	L	148	Total O 148 148	0	0
3	M	181	Total O 181 181	0	0
3	N	172	Total O 172 172	0	0
3	O	187	Total O 187 187	0	0
3	P	186	Total O 186 186	0	0
3	Q	151	Total O 151 151	0	0
3	R	175	Total O 175 175	0	0
3	S	184	Total O 184 184	0	0
3	T	204	Total O 204 204	0	0
3	U	150	Total O 150 150	0	0

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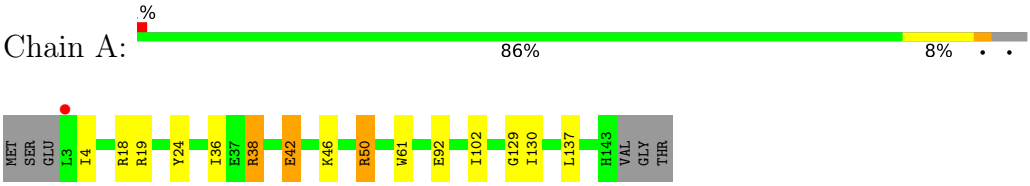
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	V	186	Total O 186 186	0	0
3	W	176	Total O 176 176	0	0
3	X	136	Total O 136 136	0	0

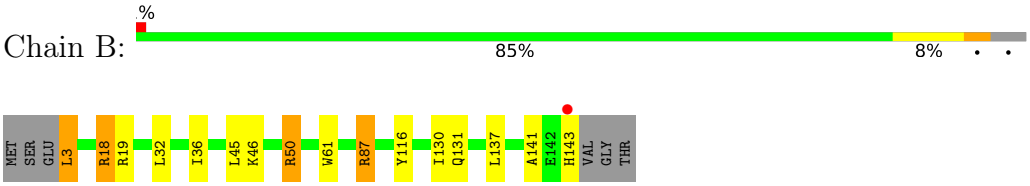
3 Residue-property plots (i)

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.

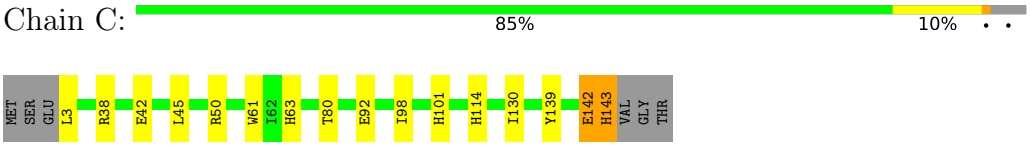
• Molecule 1: 3-dehydroquinate dehydratase



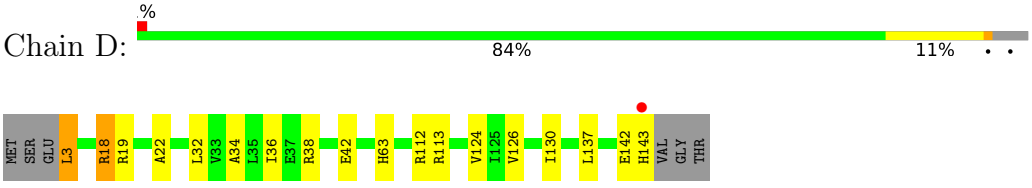
• Molecule 1: 3-dehydroquinate dehydratase



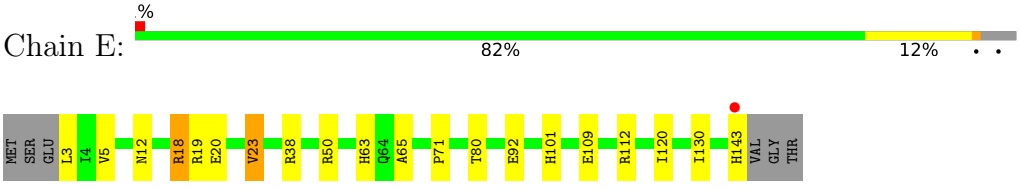
• Molecule 1: 3-dehydroquinate dehydratase



• Molecule 1: 3-dehydroquinate dehydratase

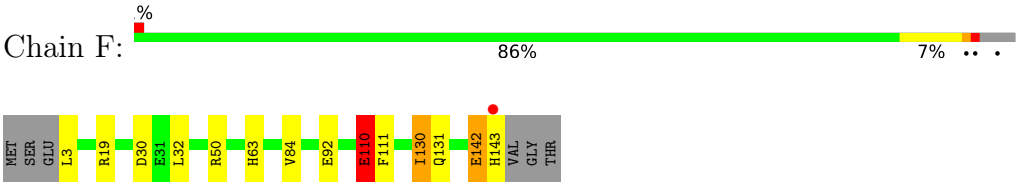


• Molecule 1: 3-dehydroquinate dehydratase



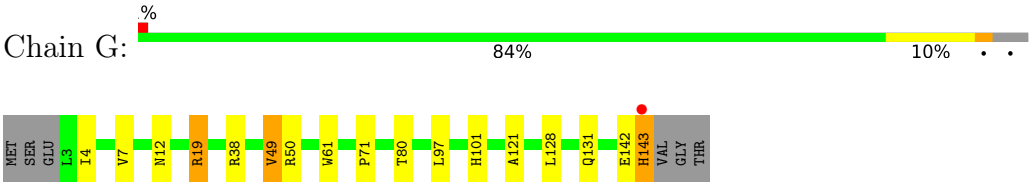
• Molecule 1: 3-dehydroquinate dehydratase

Chain F:



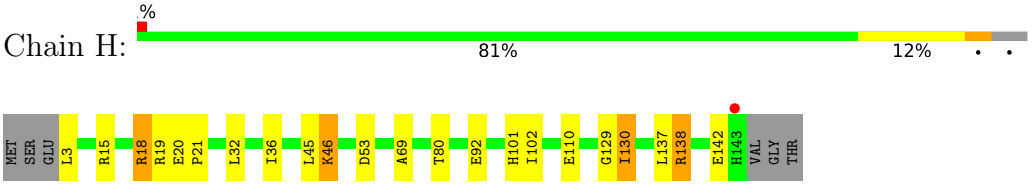
• Molecule 1: 3-dehydroquinate dehydratase

Chain G:



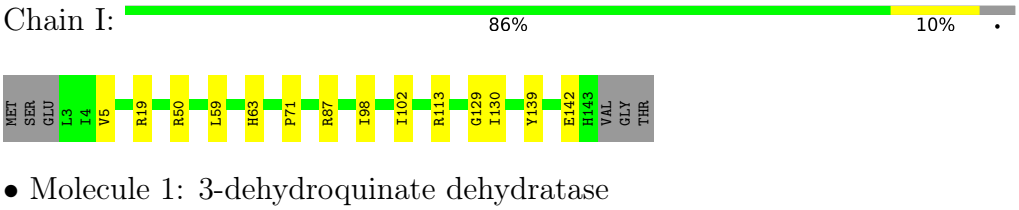
• Molecule 1: 3-dehydroquinate dehydratase

Chain H:



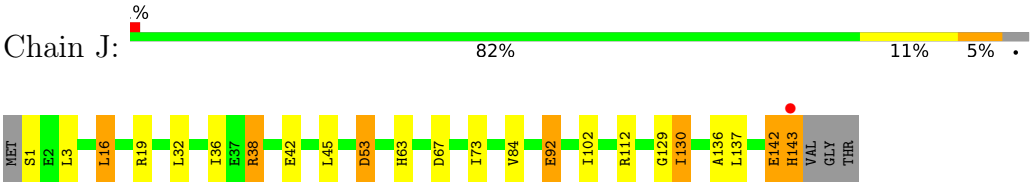
• Molecule 1: 3-dehydroquinate dehydratase

Chain I:



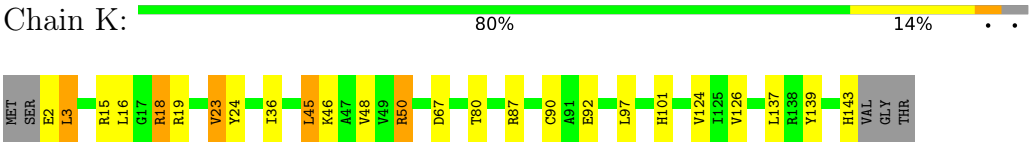
• Molecule 1: 3-dehydroquinate dehydratase

Chain J:



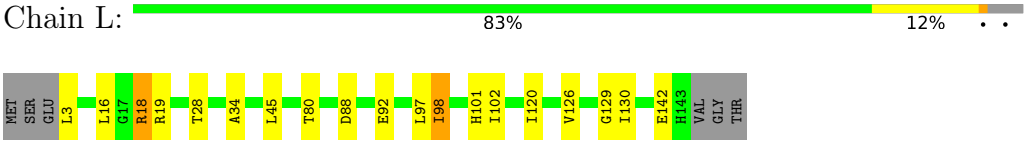
• Molecule 1: 3-dehydroquinate dehydratase

Chain K:



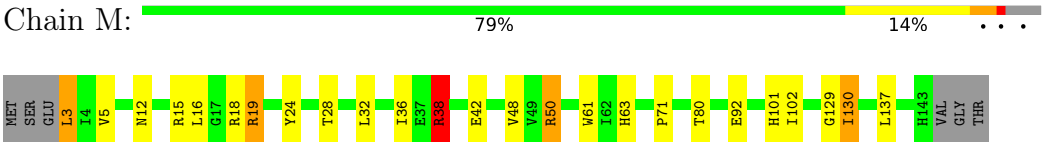
• Molecule 1: 3-dehydroquinate dehydratase

Chain L:



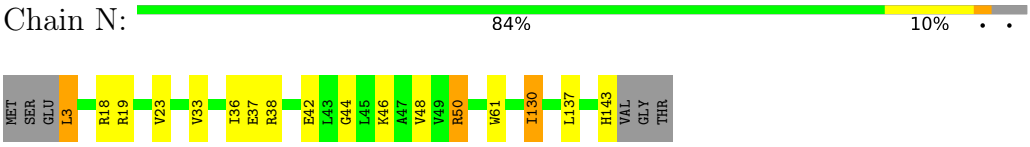
• Molecule 1: 3-dehydroquinate dehydratase

Chain M:



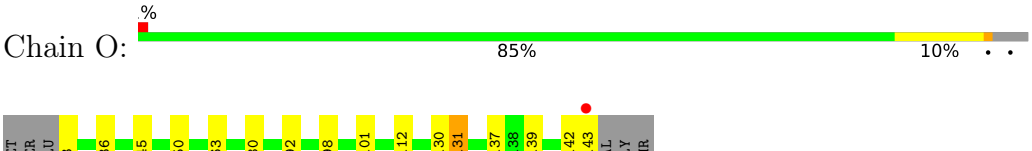
• Molecule 1: 3-dehydroquinate dehydratase

Chain N:



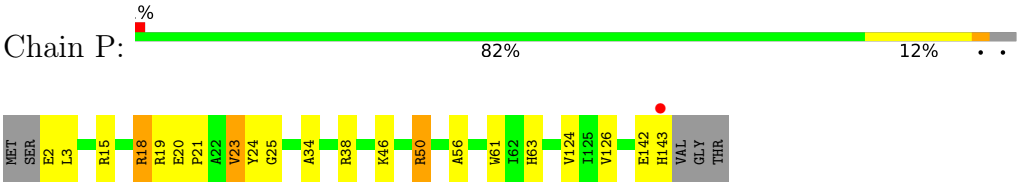
• Molecule 1: 3-dehydroquinate dehydratase

Chain O:



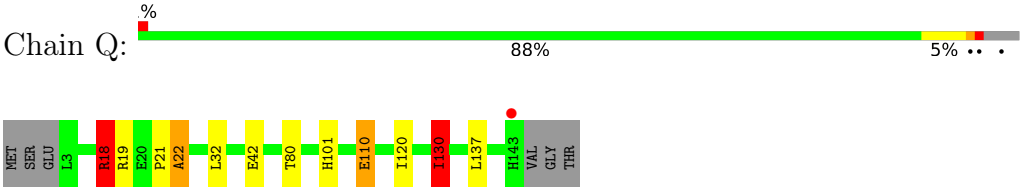
• Molecule 1: 3-dehydroquinate dehydratase

Chain P:



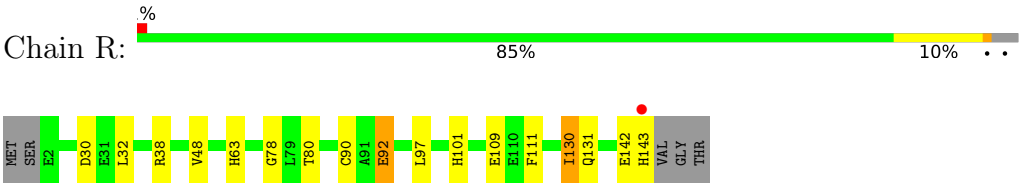
• Molecule 1: 3-dehydroquinate dehydratase

Chain Q:



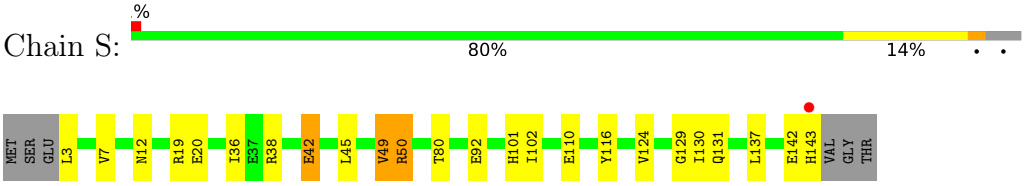
• Molecule 1: 3-dehydroquinate dehydratase

Chain R:



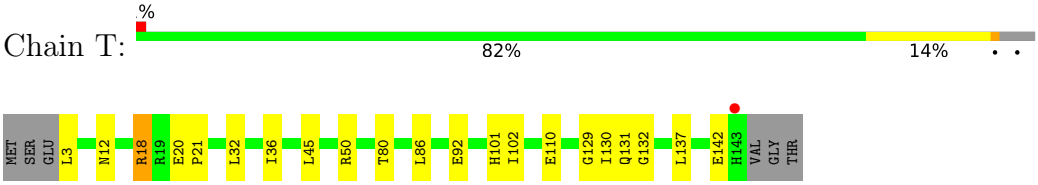
• Molecule 1: 3-dehydroquinate dehydratase

Chain S:



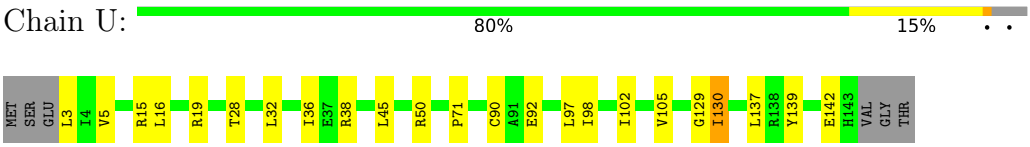
• Molecule 1: 3-dehydroquinate dehydratase

Chain T:



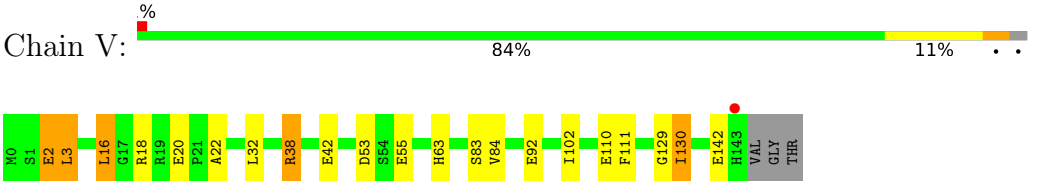
• Molecule 1: 3-dehydroquinate dehydratase

Chain U:



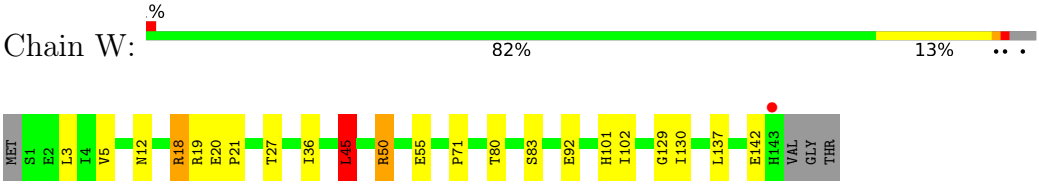
• Molecule 1: 3-dehydroquinate dehydratase

Chain V:



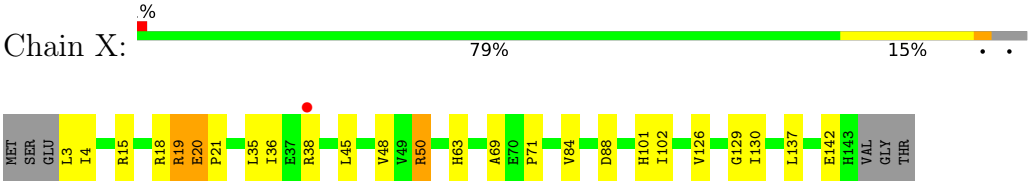
• Molecule 1: 3-dehydroquinate dehydratase

Chain W:



• Molecule 1: 3-dehydroquinate dehydratase

Chain X:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	91.24Å 143.98Å 148.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	56.50 – 1.90 56.52 – 1.90	Depositor EDS
% Data completeness (in resolution range)	92.2 (56.50-1.90) 92.2 (56.52-1.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	2.34 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.155 , 0.212 0.156 , 0.212	Depositor DCC
R_{free} test set	14004 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	16.9	Xtriage
Anisotropy	0.015	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 31.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.009 for -h,l,k 0.009 for -h,-l,-k 0.478 for h,-k,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	30436	wwPDB-VP
Average B, all atoms (Å^2)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.08% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: RJP

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	1.13	1/1089 (0.1%)	0.93	1/1483 (0.1%)
1	B	1.04	0/1089	0.89	1/1483 (0.1%)
1	C	1.10	0/1089	0.92	0/1483
1	D	1.17	0/1100	0.93	4/1497 (0.3%)
1	E	1.13	2/1089 (0.2%)	0.93	2/1483 (0.1%)
1	F	1.23	5/1083 (0.5%)	0.96	2/1476 (0.1%)
1	G	1.02	0/1089	0.99	2/1483 (0.1%)
1	H	1.12	1/1089 (0.1%)	0.92	2/1483 (0.1%)
1	I	1.03	0/1089	0.94	3/1483 (0.2%)
1	J	1.13	2/1104 (0.2%)	0.99	2/1503 (0.1%)
1	K	1.12	1/1098 (0.1%)	1.00	2/1495 (0.1%)
1	L	1.05	0/1089	0.96	2/1483 (0.1%)
1	M	1.07	1/1089 (0.1%)	0.98	3/1483 (0.2%)
1	N	1.05	1/1089 (0.1%)	0.88	0/1483
1	O	1.16	0/1089	0.98	2/1483 (0.1%)
1	P	1.14	1/1098 (0.1%)	0.90	0/1495
1	Q	1.09	1/1089 (0.1%)	0.93	3/1483 (0.2%)
1	R	1.13	5/1094 (0.5%)	0.95	1/1490 (0.1%)
1	S	1.07	4/1089 (0.4%)	0.96	0/1483
1	T	1.10	1/1089 (0.1%)	0.91	2/1483 (0.1%)
1	U	1.04	0/1089	0.91	1/1483 (0.1%)
1	V	1.15	3/1109 (0.3%)	1.03	4/1510 (0.3%)
1	W	1.09	0/1104	0.99	4/1503 (0.3%)
1	X	1.06	2/1089 (0.2%)	0.94	1/1483 (0.1%)
All	All	1.10	31/26214 (0.1%)	0.95	44/35697 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	J	0	1
1	W	0	1
All	All	0	2

The worst 5 of 31 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	110	GLU	CB-CG	12.38	1.75	1.52
1	F	110	GLU	CD-OE2	8.86	1.35	1.25
1	V	92	GLU	CB-CG	-8.40	1.36	1.52
1	F	84	VAL	CB-CG1	6.99	1.67	1.52
1	S	124	VAL	CB-CG2	6.56	1.66	1.52

The worst 5 of 44 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	G	19	ARG	NE-CZ-NH1	-10.78	114.91	120.30
1	X	88	ASP	CB-CG-OD2	9.51	126.86	118.30
1	G	19	ARG	NE-CZ-NH2	8.64	124.62	120.30
1	H	18	ARG	NE-CZ-NH1	8.62	124.61	120.30
1	V	3	LEU	CA-CB-CG	8.45	134.74	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	J	19	ARG	Peptide
1	W	142	GLU	Peptide

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	1071	0	1079	16	0
1	B	1071	0	1079	13	0
1	C	1071	0	1079	13	0
1	D	1079	0	1092	21	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1071	0	1079	12	0
1	F	1065	0	1068	13	0
1	G	1071	0	1079	13	0
1	H	1071	0	1079	18	0
1	I	1071	0	1079	7	0
1	J	1086	0	1093	16	0
1	K	1080	0	1085	25	0
1	L	1071	0	1079	29	0
1	M	1071	0	1079	28	2
1	N	1071	0	1079	10	0
1	O	1071	0	1079	11	0
1	P	1080	0	1085	16	0
1	Q	1071	0	1079	11	0
1	R	1076	0	1081	11	0
1	S	1071	0	1079	14	0
1	T	1071	0	1079	20	0
1	U	1071	0	1079	13	0
1	V	1091	0	1095	12	0
1	W	1086	0	1093	18	0
1	X	1071	0	1079	24	0
2	A	23	0	16	1	0
2	B	23	0	16	1	0
2	C	23	0	16	0	0
2	D	23	0	16	1	0
2	E	23	0	16	1	0
2	F	23	0	16	0	0
2	G	23	0	16	0	0
2	H	23	0	16	0	0
2	I	23	0	16	0	0
2	J	23	0	16	0	0
2	K	23	0	16	0	0
2	L	23	0	16	1	0
2	M	23	0	16	0	0
2	N	23	0	16	0	0
2	O	23	0	16	0	0
2	P	23	0	16	1	0
2	Q	23	0	16	0	0
2	R	23	0	16	0	0
2	S	23	0	16	0	0
2	T	23	0	16	0	0
2	U	23	0	16	1	0
2	V	23	0	16	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	W	23	0	16	0	0
2	X	23	0	16	0	0
3	A	169	0	0	2	0
3	B	162	0	0	4	0
3	C	177	0	0	4	0
3	D	165	0	0	3	3
3	E	169	0	0	3	0
3	F	164	0	0	3	0
3	G	155	0	0	2	0
3	H	175	0	0	2	0
3	I	159	0	0	1	0
3	J	189	0	0	1	0
3	K	185	0	0	5	0
3	L	148	0	0	1	0
3	M	181	0	0	2	0
3	N	172	0	0	2	0
3	O	187	0	0	2	0
3	P	186	0	0	1	0
3	Q	151	0	0	2	2
3	R	175	0	0	2	0
3	S	184	0	0	3	0
3	T	204	0	0	3	1
3	U	150	0	0	3	0
3	V	186	0	0	6	2
3	W	176	0	0	1	0
3	X	136	0	0	3	0
All	All	30436	0	26340	353	5

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 7.

The worst 5 of 353 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:F:110:GLU:CB	1:F:110:GLU:CG	1.75	1.63
1:F:92:GLU:OE1	1:G:19:ARG:HD2	1.27	1.31
1:D:18[B]:ARG:HH21	1:D:18[B]:ARG:CG	1.52	1.17
1:W:18:ARG:HH11	1:W:18:ARG:HG2	1.14	1.13
1:D:18[A]:ARG:HH21	1:D:18[A]:ARG:CG	1.64	1.11

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:3503:HOH:O	3:V:3695:HOH:O[2_546]	1.54	0.66
3:D:3504:HOH:O	3:T:3696:HOH:O[2_546]	1.67	0.53
1:M:38:ARG:NH1	3:Q:3949:HOH:O[2_646]	1.72	0.48
1:M:38:ARG:NE	3:Q:3730:HOH:O[2_646]	1.97	0.23
3:D:3503:HOH:O	3:V:728:HOH:O[2_546]	2.09	0.11

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	139/147 (95%)	135 (97%)	4 (3%)	0	100	100
1	B	139/147 (95%)	136 (98%)	3 (2%)	0	100	100
1	C	139/147 (95%)	137 (99%)	2 (1%)	0	100	100
1	D	140/147 (95%)	137 (98%)	3 (2%)	0	100	100
1	E	139/147 (95%)	137 (99%)	2 (1%)	0	100	100
1	F	139/147 (95%)	136 (98%)	3 (2%)	0	100	100
1	G	139/147 (95%)	135 (97%)	4 (3%)	0	100	100
1	H	139/147 (95%)	136 (98%)	3 (2%)	0	100	100
1	I	139/147 (95%)	137 (99%)	2 (1%)	0	100	100
1	J	141/147 (96%)	136 (96%)	5 (4%)	0	100	100
1	K	140/147 (95%)	137 (98%)	3 (2%)	0	100	100
1	L	139/147 (95%)	134 (96%)	5 (4%)	0	100	100
1	M	139/147 (95%)	133 (96%)	6 (4%)	0	100	100
1	N	139/147 (95%)	135 (97%)	4 (3%)	0	100	100
1	O	139/147 (95%)	137 (99%)	2 (1%)	0	100	100
1	P	140/147 (95%)	138 (99%)	2 (1%)	0	100	100
1	Q	139/147 (95%)	135 (97%)	3 (2%)	1 (1%)	22	12

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	R	140/147 (95%)	137 (98%)	3 (2%)	0	100	100
1	S	139/147 (95%)	136 (98%)	3 (2%)	0	100	100
1	T	139/147 (95%)	135 (97%)	4 (3%)	0	100	100
1	U	139/147 (95%)	136 (98%)	3 (2%)	0	100	100
1	V	142/147 (97%)	139 (98%)	3 (2%)	0	100	100
1	W	141/147 (96%)	138 (98%)	3 (2%)	0	100	100
1	X	139/147 (95%)	135 (97%)	4 (3%)	0	100	100
All	All	3347/3528 (95%)	3267 (98%)	79 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	Q	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	110/115 (96%)	106 (96%)	4 (4%)	35	26
1	B	110/115 (96%)	106 (96%)	4 (4%)	35	26
1	C	110/115 (96%)	106 (96%)	4 (4%)	35	26
1	D	111/115 (96%)	108 (97%)	3 (3%)	44	38
1	E	110/115 (96%)	105 (96%)	5 (4%)	27	18
1	F	109/115 (95%)	106 (97%)	3 (3%)	43	36
1	G	110/115 (96%)	106 (96%)	4 (4%)	35	26
1	H	110/115 (96%)	105 (96%)	5 (4%)	27	18
1	I	110/115 (96%)	108 (98%)	2 (2%)	59	55
1	J	112/115 (97%)	105 (94%)	7 (6%)	18	8
1	K	111/115 (96%)	106 (96%)	5 (4%)	27	18

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	L	110/115 (96%)	106 (96%)	4 (4%)	35 26
1	M	110/115 (96%)	105 (96%)	5 (4%)	27 18
1	N	110/115 (96%)	106 (96%)	4 (4%)	35 26
1	O	110/115 (96%)	108 (98%)	2 (2%)	59 55
1	P	111/115 (96%)	105 (95%)	6 (5%)	22 13
1	Q	110/115 (96%)	107 (97%)	3 (3%)	44 38
1	R	110/115 (96%)	109 (99%)	1 (1%)	78 79
1	S	110/115 (96%)	106 (96%)	4 (4%)	35 26
1	T	110/115 (96%)	108 (98%)	2 (2%)	59 55
1	U	110/115 (96%)	107 (97%)	3 (3%)	44 38
1	V	112/115 (97%)	105 (94%)	7 (6%)	18 8
1	W	112/115 (97%)	109 (97%)	3 (3%)	44 38
1	X	110/115 (96%)	105 (96%)	5 (4%)	27 18
All	All	2648/2760 (96%)	2553 (96%)	95 (4%)	35 26

5 of 95 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	N	50	ARG
1	S	49	VAL
1	O	130	ILE
1	P	142	GLU
1	T	18	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
1	R	131	GLN
1	X	63	HIS
1	V	63	HIS
1	I	131	GLN
1	R	63	HIS

5.3.3 RNA

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

24 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	RJP	B	147	-	22,24,24	2.49	4 (18%)	23,34,34	1.17	3 (13%)
2	RJP	W	147	-	22,24,24	2.59	4 (18%)	23,34,34	1.37	4 (17%)
2	RJP	D	147	-	22,24,24	2.60	4 (18%)	23,34,34	1.55	4 (17%)
2	RJP	J	147	-	22,24,24	2.10	4 (18%)	23,34,34	1.30	4 (17%)
2	RJP	S	147	-	22,24,24	2.49	4 (18%)	23,34,34	1.38	4 (17%)
2	RJP	H	147	-	22,24,24	2.85	5 (22%)	23,34,34	1.52	4 (17%)
2	RJP	A	147	-	22,24,24	2.61	5 (22%)	23,34,34	1.70	5 (21%)
2	RJP	M	147	-	22,24,24	2.68	4 (18%)	23,34,34	1.63	5 (21%)
2	RJP	C	147	-	22,24,24	2.44	5 (22%)	23,34,34	1.50	3 (13%)
2	RJP	L	147	-	22,24,24	2.38	4 (18%)	23,34,34	1.05	2 (8%)
2	RJP	K	147	-	22,24,24	2.47	4 (18%)	23,34,34	1.31	3 (13%)
2	RJP	X	147	-	22,24,24	2.81	4 (18%)	23,34,34	1.08	2 (8%)
2	RJP	U	147	-	22,24,24	2.66	5 (22%)	23,34,34	1.09	0
2	RJP	O	147	-	22,24,24	2.35	5 (22%)	23,34,34	1.24	2 (8%)
2	RJP	N	147	-	22,24,24	2.66	4 (18%)	23,34,34	1.35	4 (17%)
2	RJP	I	147	-	22,24,24	2.55	3 (13%)	23,34,34	1.11	2 (8%)
2	RJP	V	147	-	22,24,24	2.07	7 (31%)	23,34,34	1.56	3 (13%)
2	RJP	P	147	-	22,24,24	2.54	6 (27%)	23,34,34	1.52	4 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	RJP	G	147	-	22,24,24	2.58	4 (18%)	23,34,34	1.66	4 (17%)
2	RJP	R	147	-	22,24,24	2.40	5 (22%)	23,34,34	1.27	2 (8%)
2	RJP	F	147	-	22,24,24	2.59	6 (27%)	23,34,34	1.29	3 (13%)
2	RJP	T	147	-	22,24,24	2.66	4 (18%)	23,34,34	1.55	4 (17%)
2	RJP	Q	147	-	22,24,24	2.45	5 (22%)	23,34,34	1.26	2 (8%)
2	RJP	E	147	-	22,24,24	2.67	5 (22%)	23,34,34	0.97	1 (4%)

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	RJP	B	147	-	-	2/15/33/33	0/2/2/2
2	RJP	W	147	-	-	3/15/33/33	0/2/2/2
2	RJP	D	147	-	-	1/15/33/33	0/2/2/2
2	RJP	J	147	-	-	2/15/33/33	0/2/2/2
2	RJP	S	147	-	-	1/15/33/33	0/2/2/2
2	RJP	H	147	-	-	3/15/33/33	0/2/2/2
2	RJP	A	147	-	-	2/15/33/33	0/2/2/2
2	RJP	M	147	-	-	2/15/33/33	0/2/2/2
2	RJP	C	147	-	-	2/15/33/33	0/2/2/2
2	RJP	L	147	-	-	2/15/33/33	0/2/2/2
2	RJP	K	147	-	-	2/15/33/33	0/2/2/2
2	RJP	X	147	-	-	3/15/33/33	0/2/2/2
2	RJP	U	147	-	-	3/15/33/33	0/2/2/2
2	RJP	O	147	-	-	1/15/33/33	0/2/2/2
2	RJP	N	147	-	-	2/15/33/33	0/2/2/2
2	RJP	I	147	-	-	2/15/33/33	0/2/2/2
2	RJP	V	147	-	-	1/15/33/33	0/2/2/2
2	RJP	P	147	-	-	2/15/33/33	0/2/2/2
2	RJP	G	147	-	-	3/15/33/33	0/2/2/2
2	RJP	R	147	-	-	2/15/33/33	0/2/2/2
2	RJP	F	147	-	-	1/15/33/33	0/2/2/2
2	RJP	T	147	-	-	1/15/33/33	0/2/2/2
2	RJP	Q	147	-	-	2/15/33/33	0/2/2/2
2	RJP	E	147	-	-	1/15/33/33	0/2/2/2

The worst 5 of 110 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	U	147	RJP	OAE-CAV	8.95	1.37	1.22
2	M	147	RJP	OAE-CAV	8.68	1.37	1.22
2	H	147	RJP	OAE-CAV	8.60	1.37	1.22
2	I	147	RJP	OAE-CAV	8.40	1.37	1.22
2	D	147	RJP	OAE-CAV	8.21	1.36	1.22

The worst 5 of 74 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	G	147	RJP	CAH-CAG-CAQ	-4.09	105.38	112.56
2	V	147	RJP	CAH-CAG-CAQ	-4.09	105.39	112.56
2	D	147	RJP	CAH-CAG-CAQ	-3.83	105.84	112.56
2	C	147	RJP	CAG-CAQ-NAP	-3.81	107.88	114.59
2	V	147	RJP	OAA-CAQ-NAP	3.72	130.41	123.63

There are no chirality outliers.

5 of 46 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	147	RJP	OAB-CAR-CAW-OAF
2	B	147	RJP	OAB-CAR-CAW-OAF
2	B	147	RJP	OAC-CAR-CAW-CAO
2	C	147	RJP	OAC-CAR-CAW-CAI
2	D	147	RJP	OAB-CAR-CAW-OAF

There are no ring outliers.

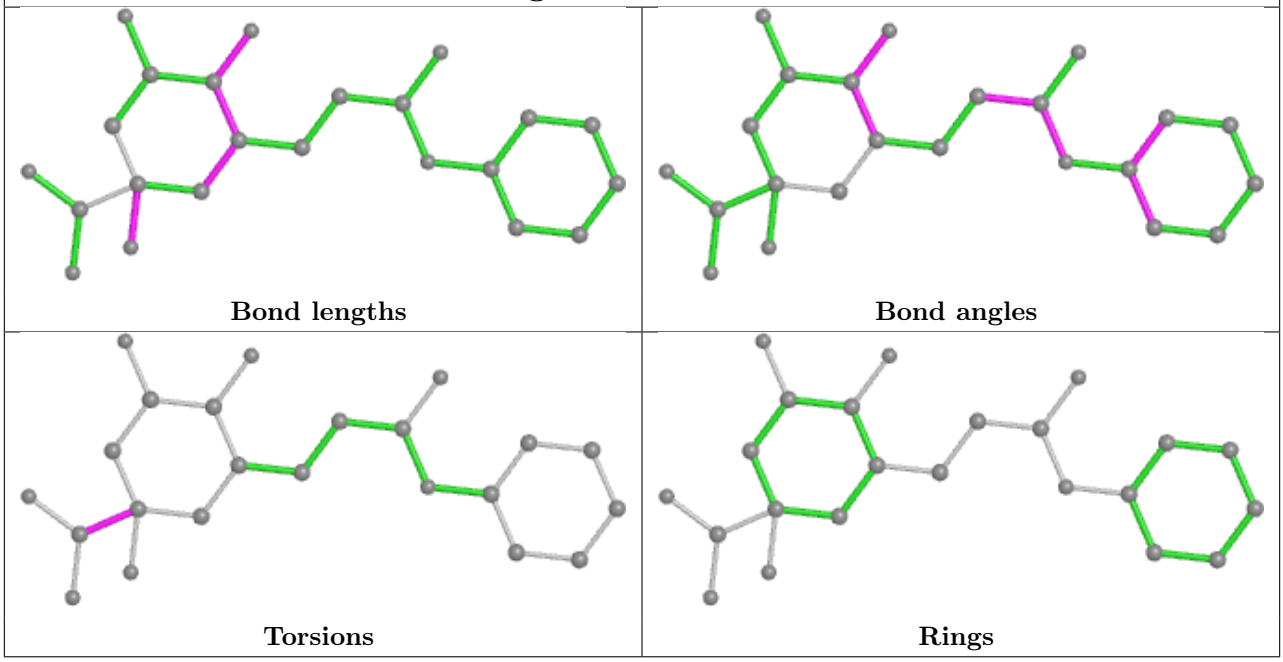
8 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	147	RJP	1	0
2	D	147	RJP	1	0
2	A	147	RJP	1	0
2	L	147	RJP	1	0
2	U	147	RJP	1	0
2	V	147	RJP	1	0
2	P	147	RJP	1	0
2	E	147	RJP	1	0

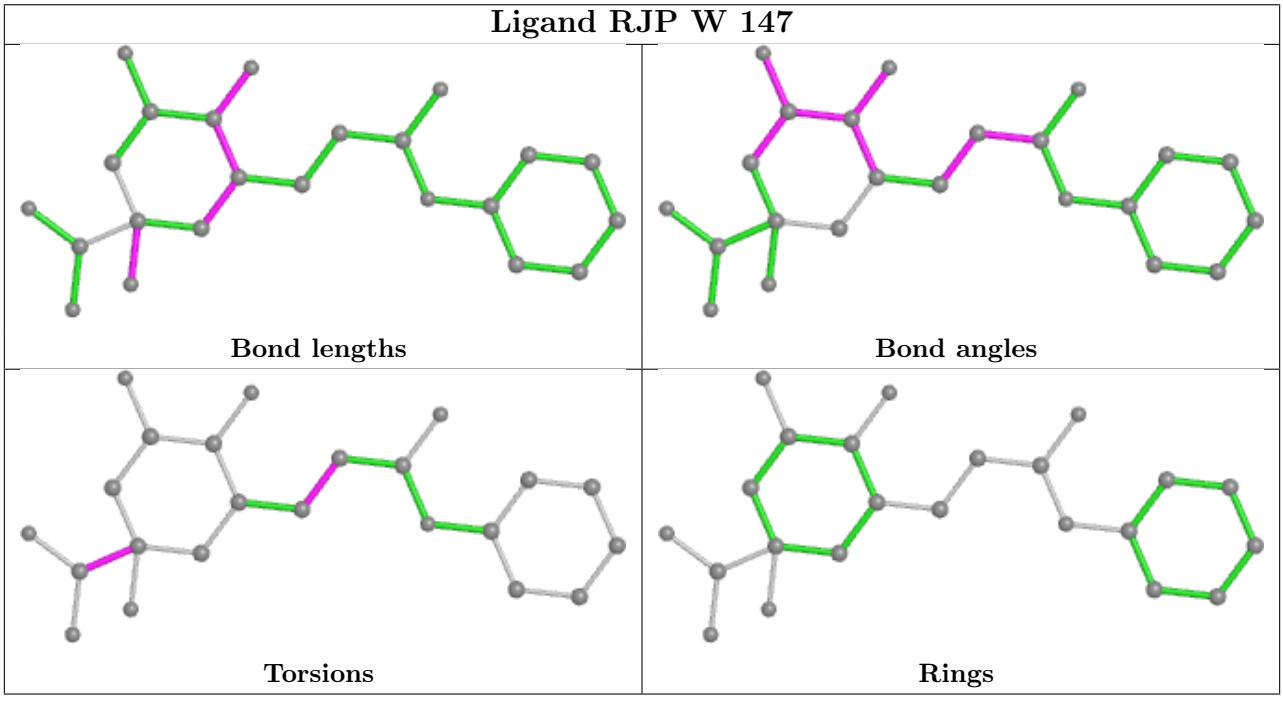
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

also be included. For torsion angles, if less then 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.

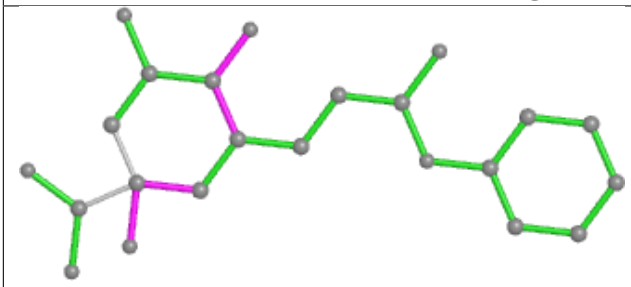
Ligand RJP B 147



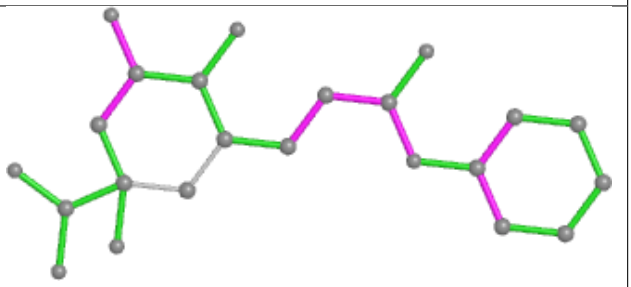
Ligand RJP W 147



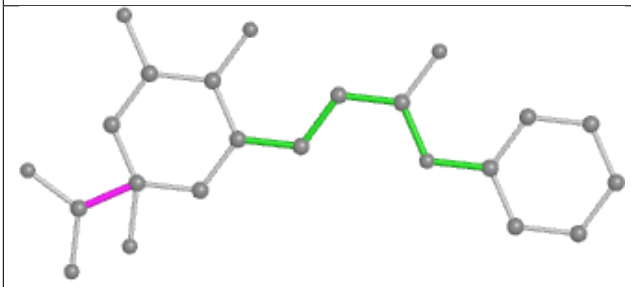
Ligand RJP D 147



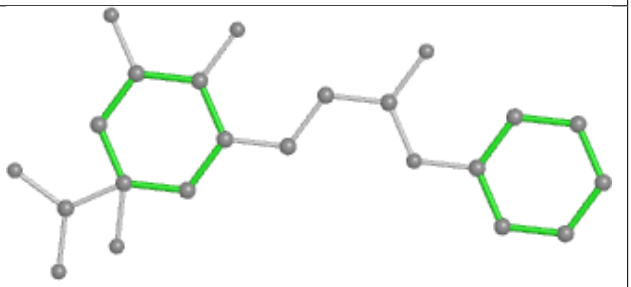
Bond lengths



Bond angles

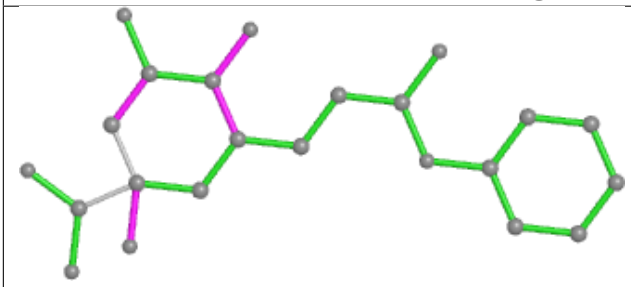


Torsions



Rings

Ligand RJP J 147



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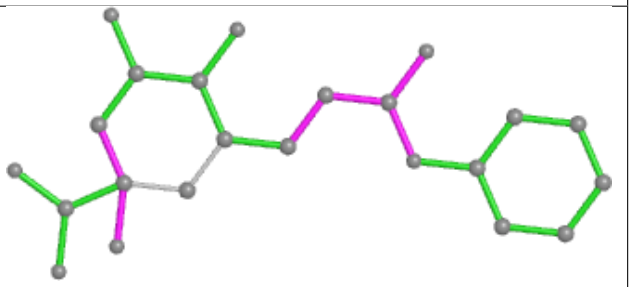
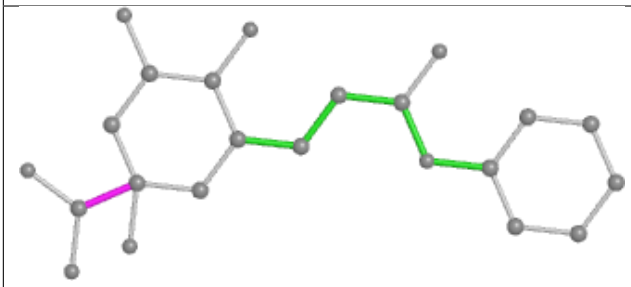


Bond angles



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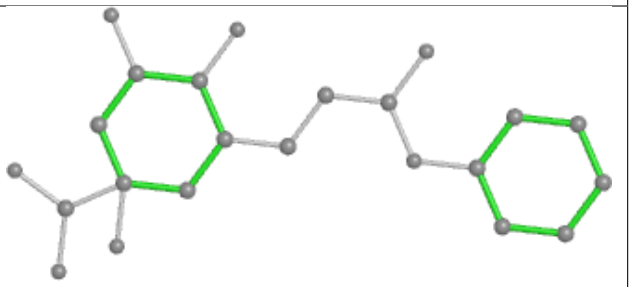
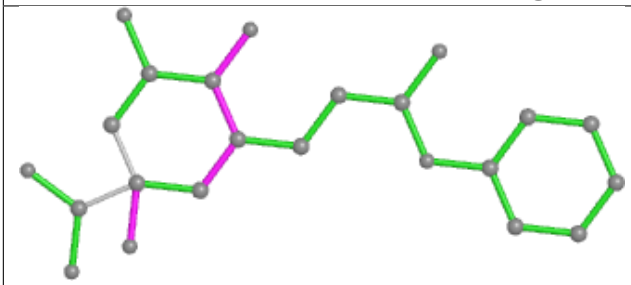


Rings

Ligand RJP S 147



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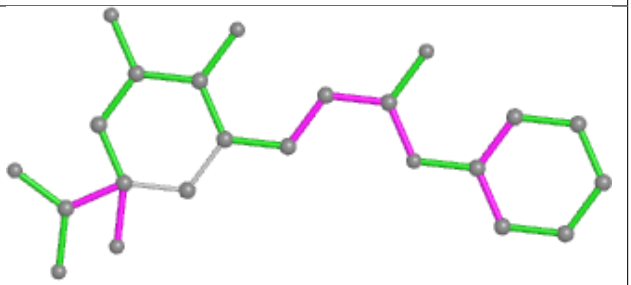
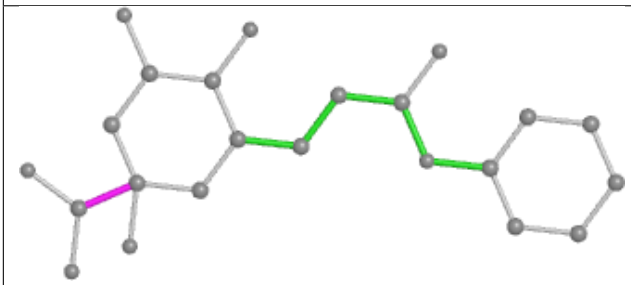
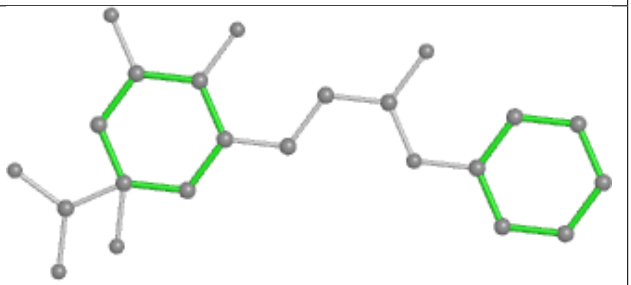


Bond angles

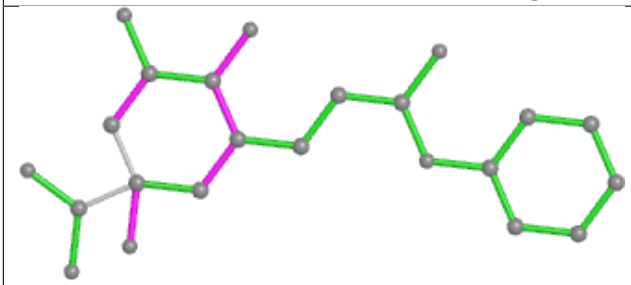


Torsions



Rings

Ligand RJP H 147



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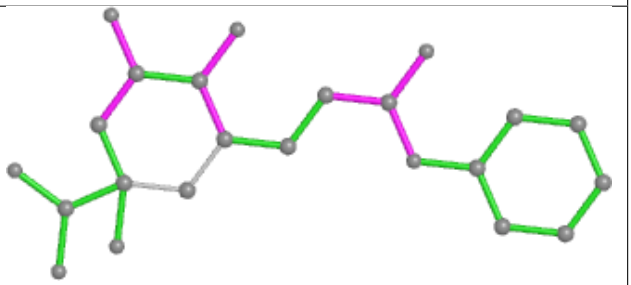
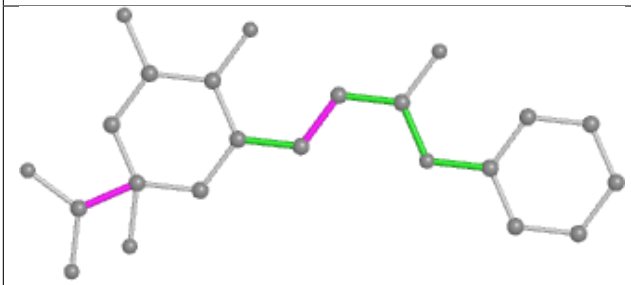


Bond angles



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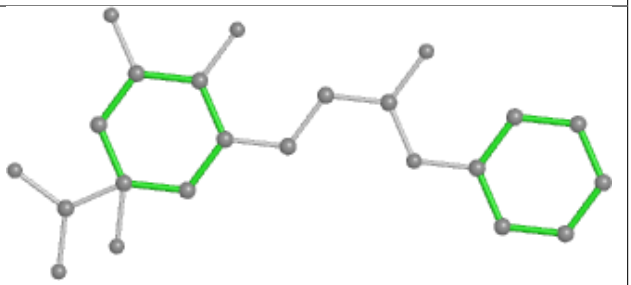
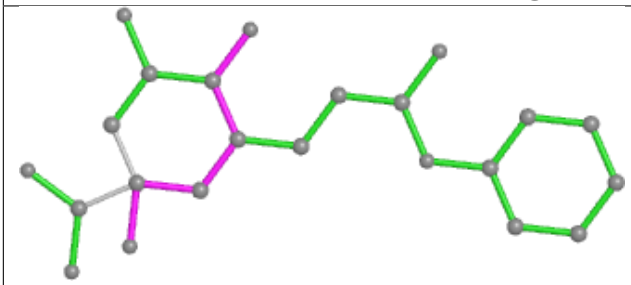


Rings

Ligand RJP A 147



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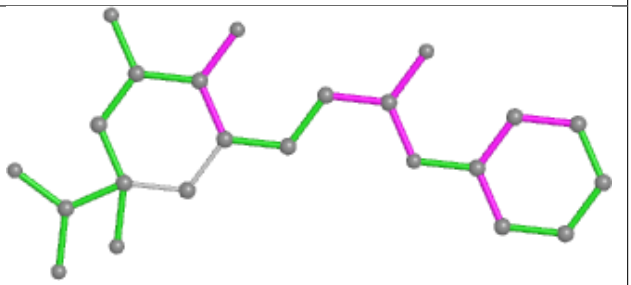
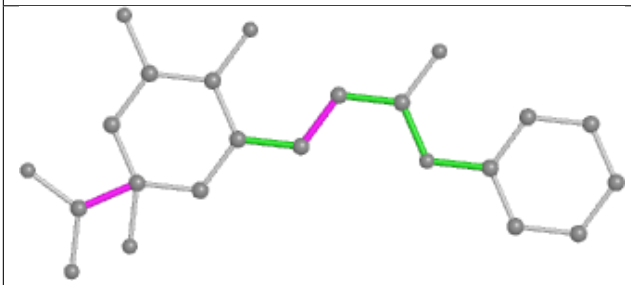


Bond angles



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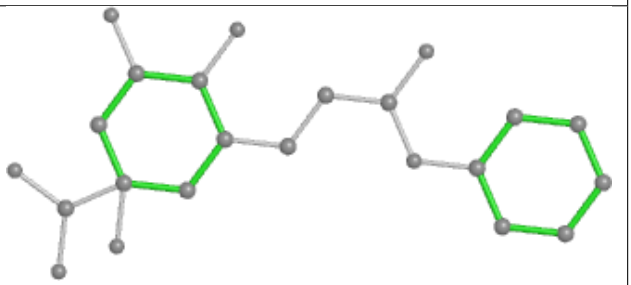
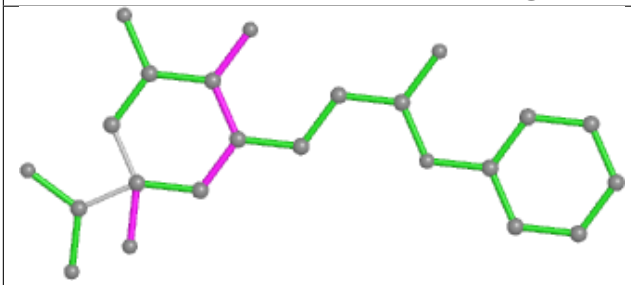


Rings

Ligand RJP M 147



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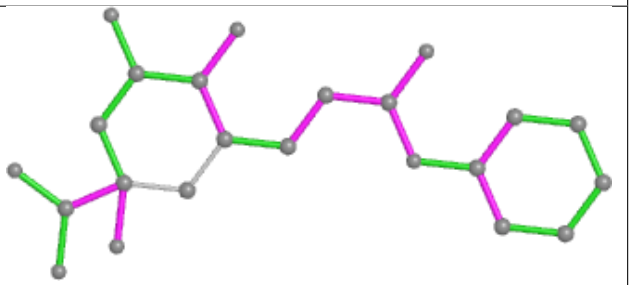

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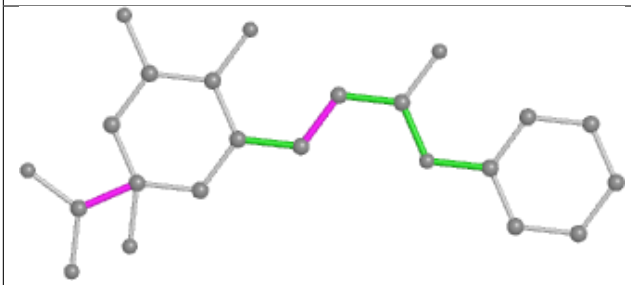

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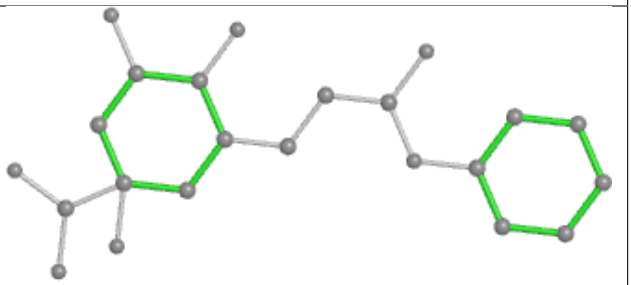
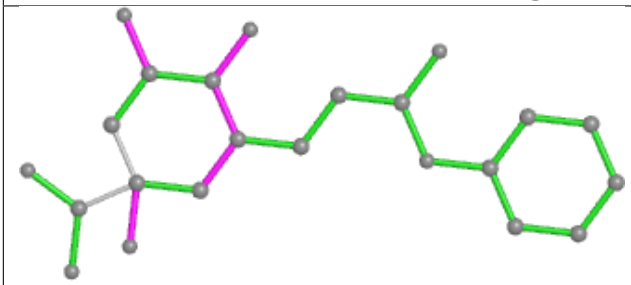


Rings

Ligand RJP C 147



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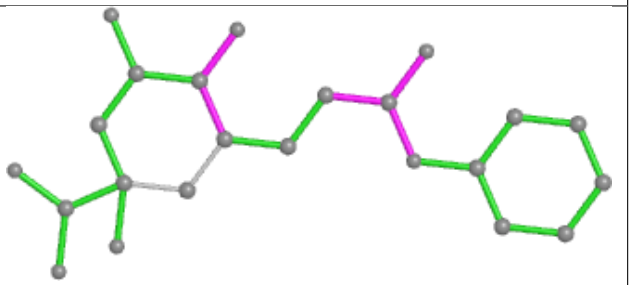

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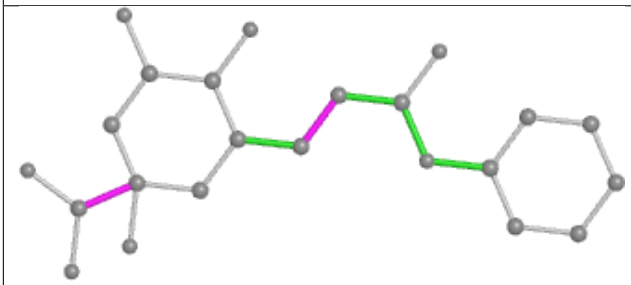

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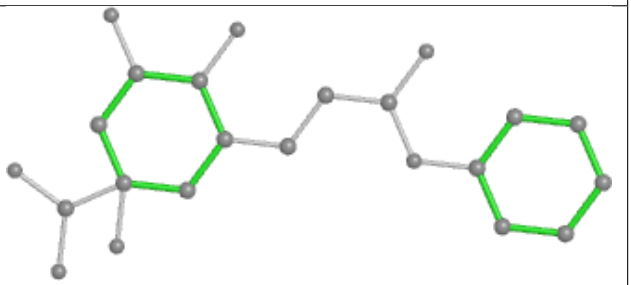
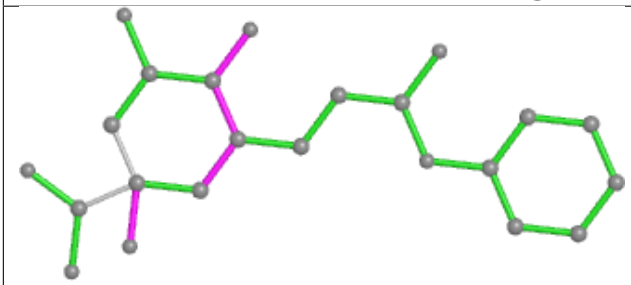


Rings

Ligand RJP L 147



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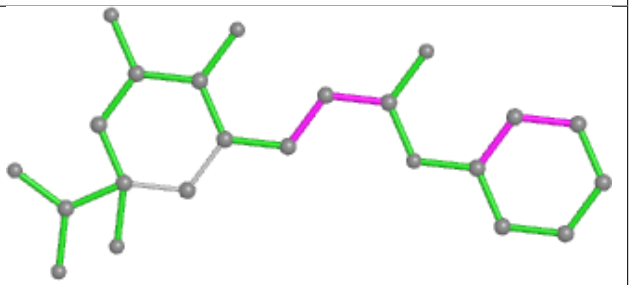
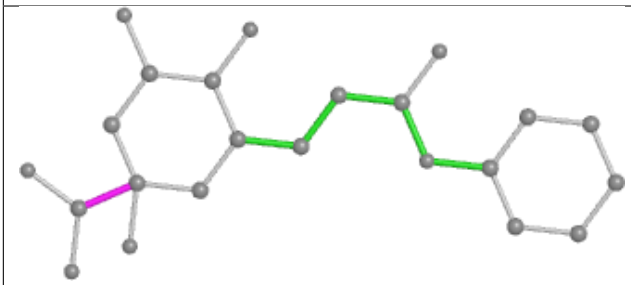


Bond angles



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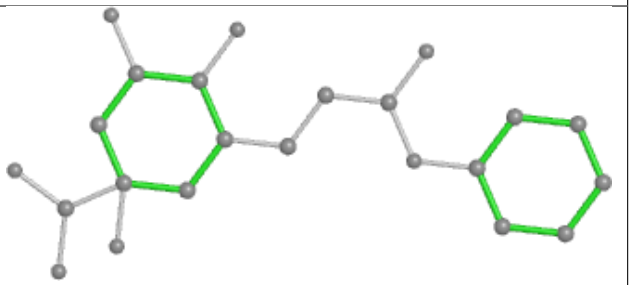
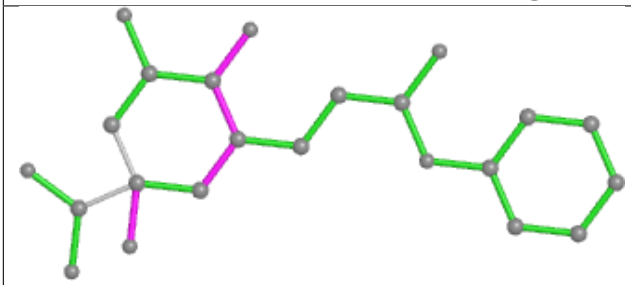


Rings

Ligand RJP K 147



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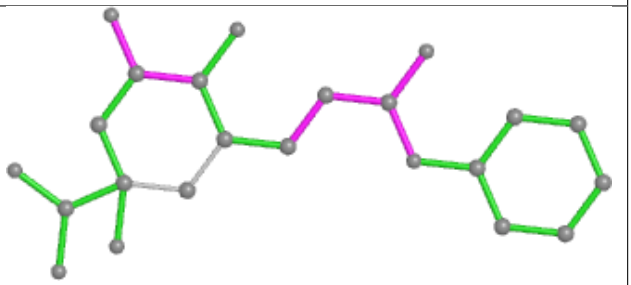
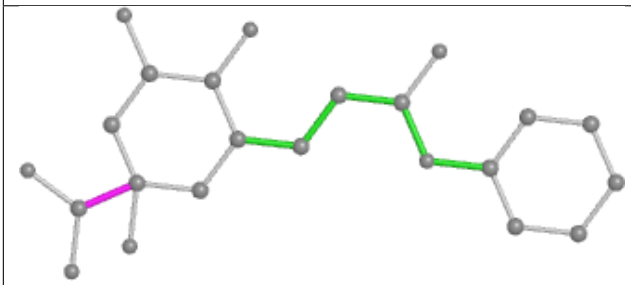
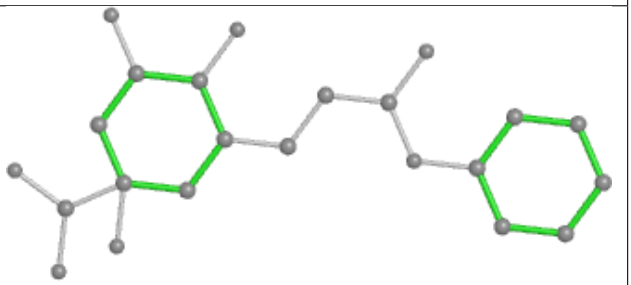


Bond angles

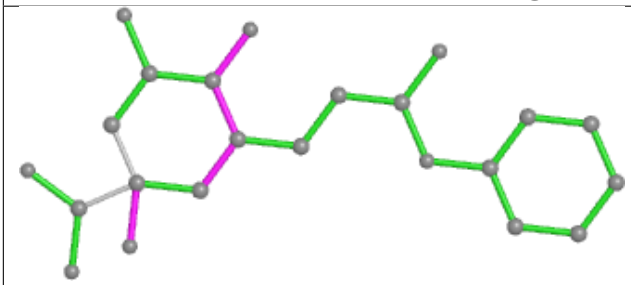


Torsions



Rings

Ligand RJP X 147



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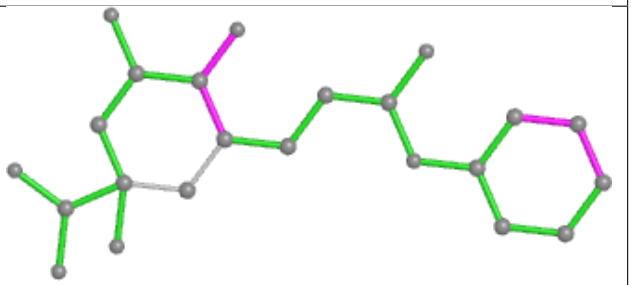
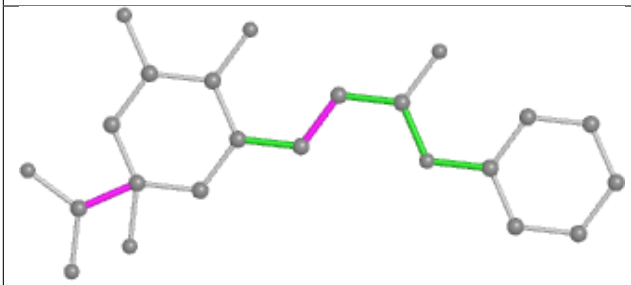


Bond angles



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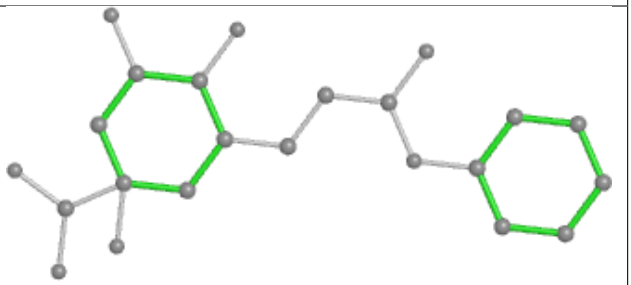
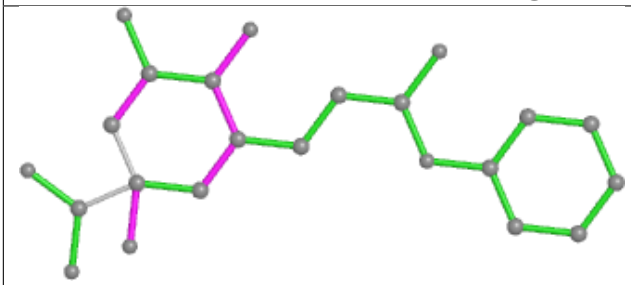


Rings

Ligand RJP U 147



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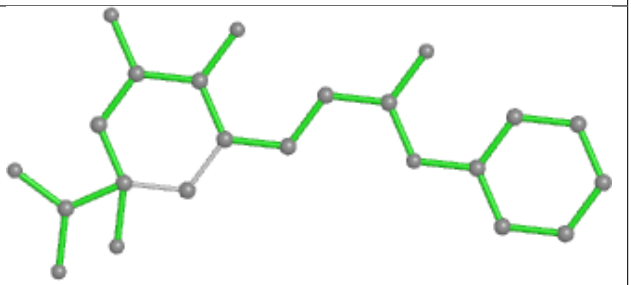
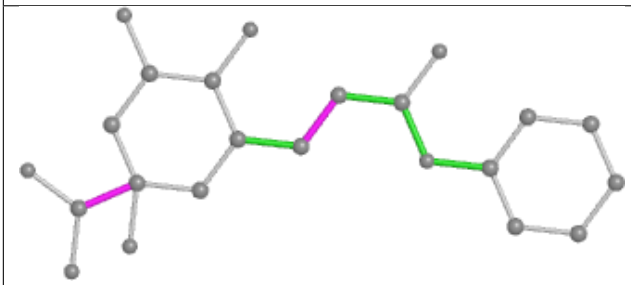


Bond angles



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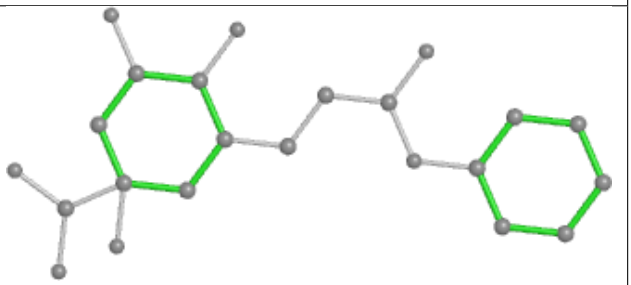
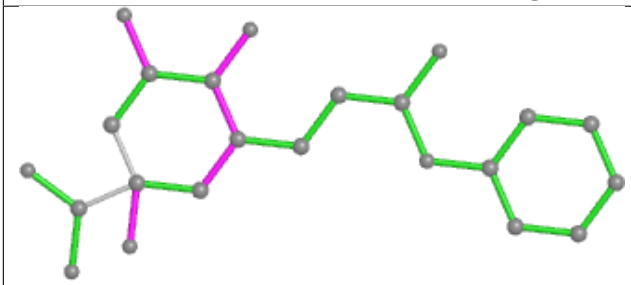


Rings

Ligand RJP O 147



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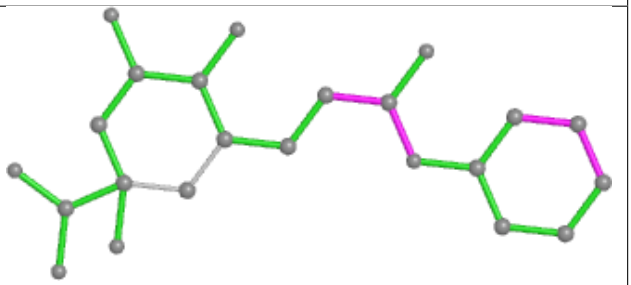
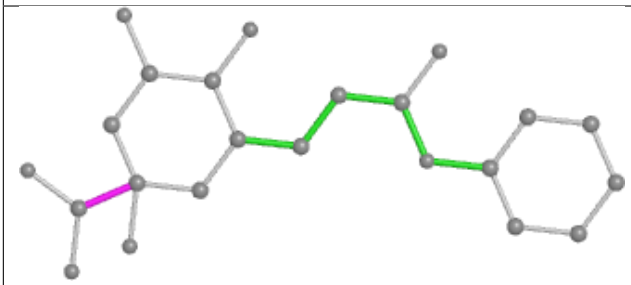


Bond angles



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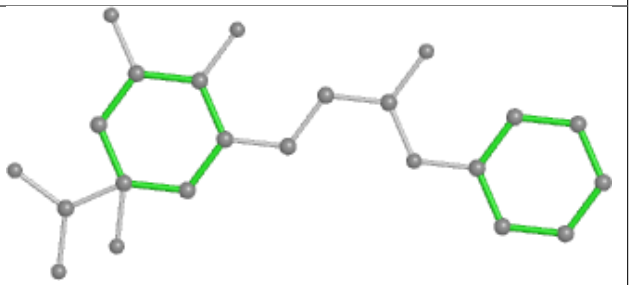
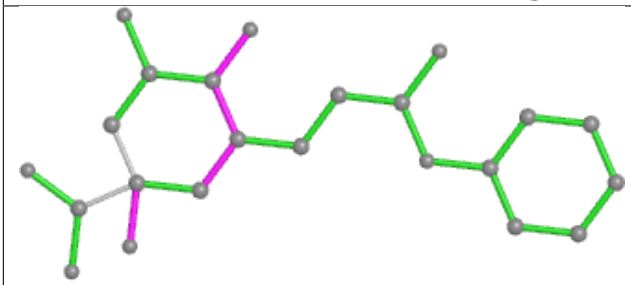


Rings

Ligand RJP N 147



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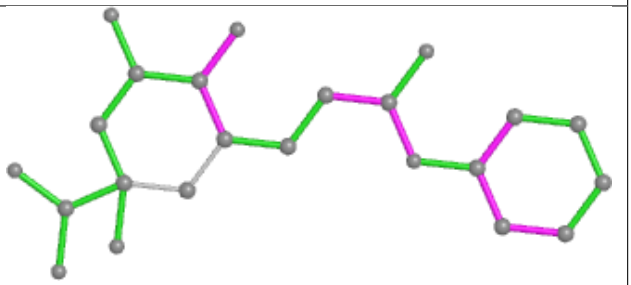
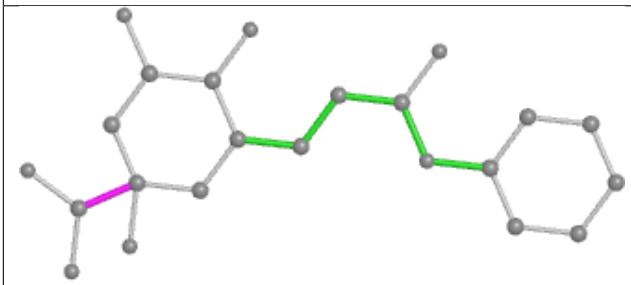


Bond angles



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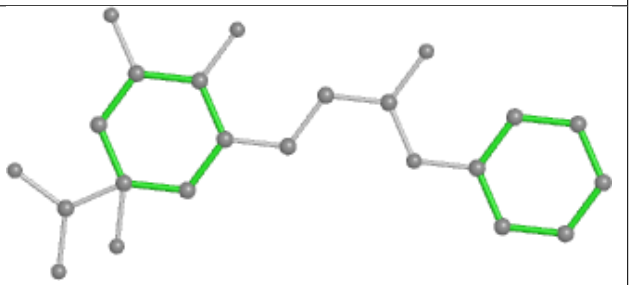
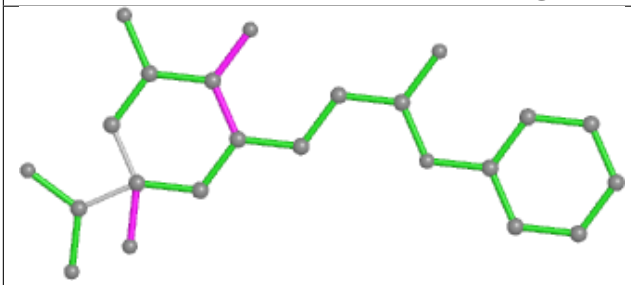


Rings

Ligand RJP I 147



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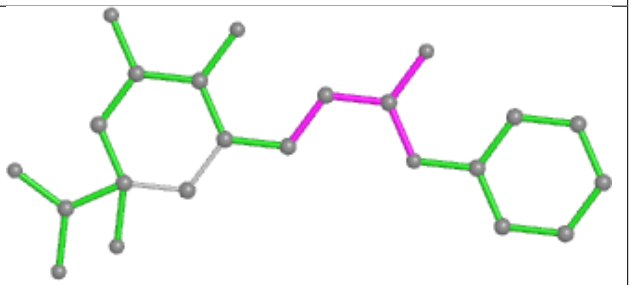
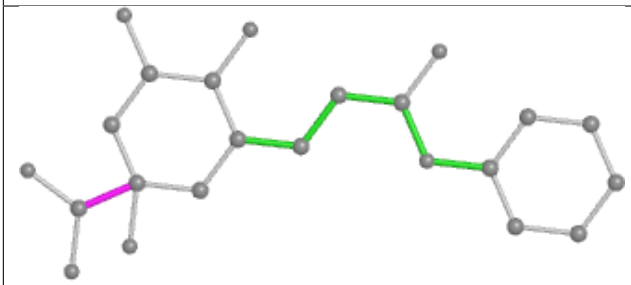
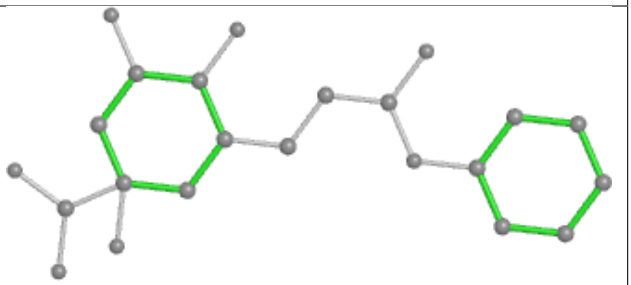


Bond angles

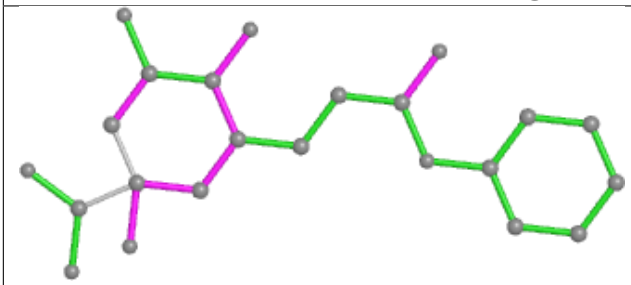


Torsions

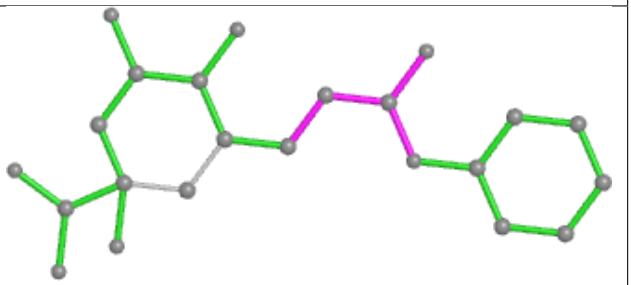


Rings

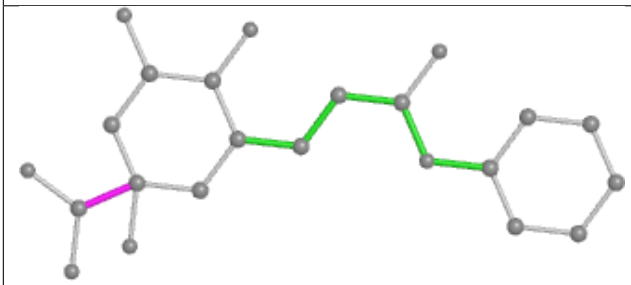
Ligand RJP V 147



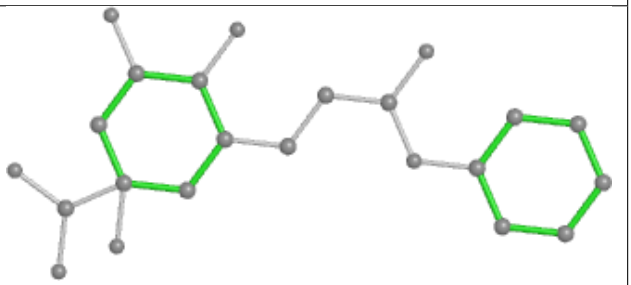
Bond lengths



Bond angles

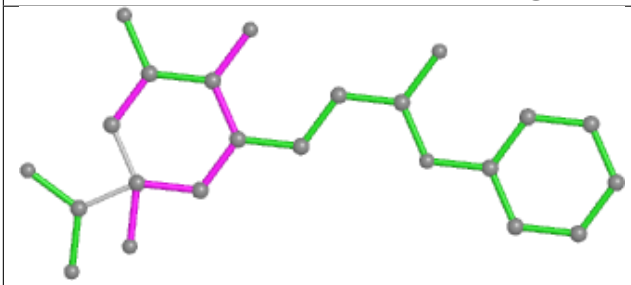


Torsions

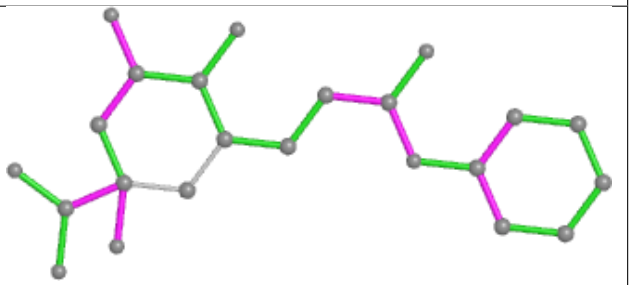


Rings

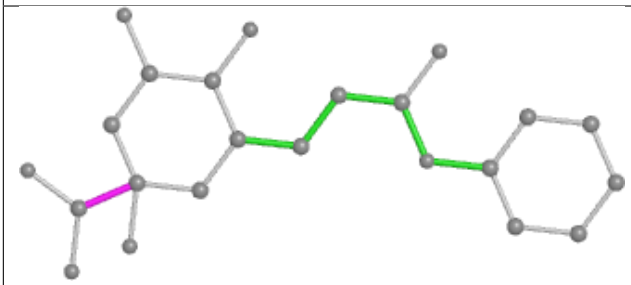
Ligand RJP P 147



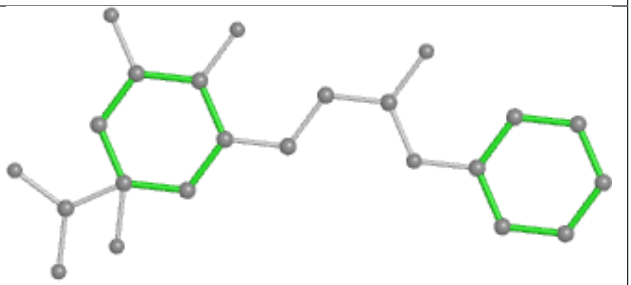
Bond lengths



Bond angles

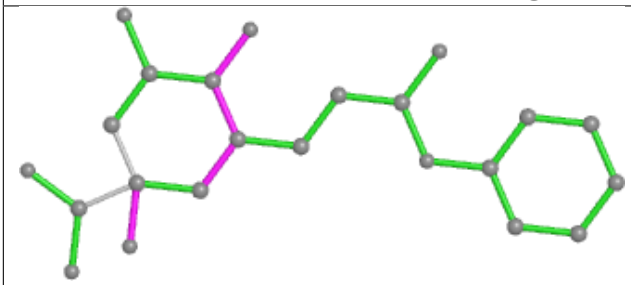


Torsions

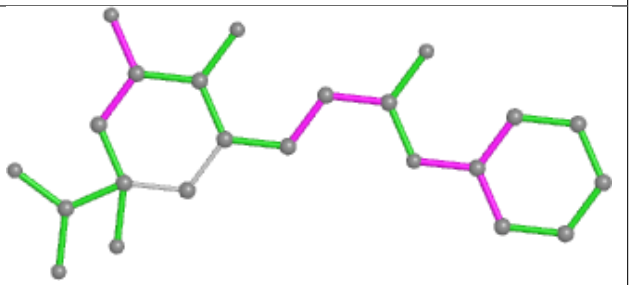


Rings

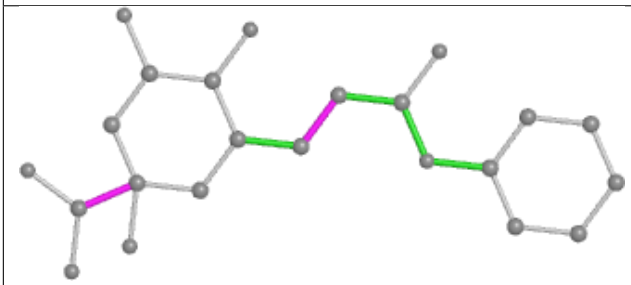
Ligand RJP G 147



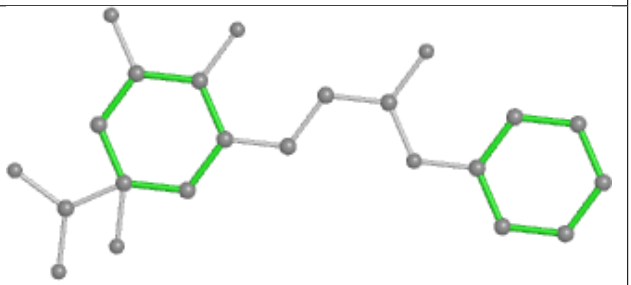
Bond lengths



Bond angles

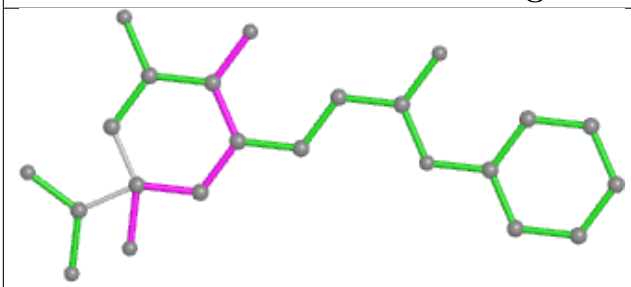


Torsions

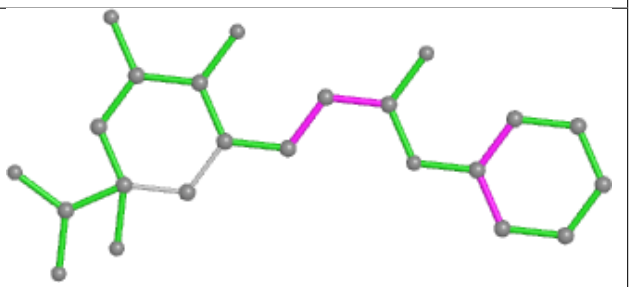


Rings

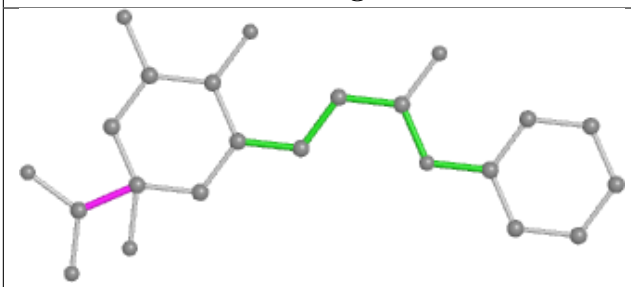
Ligand RJP R 147



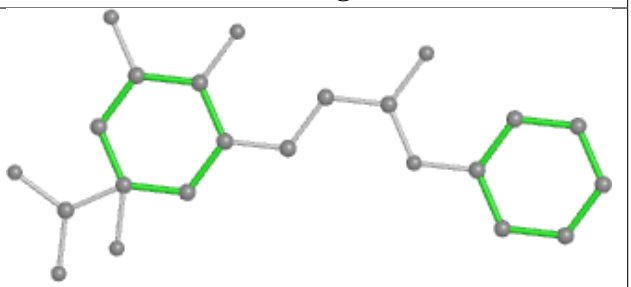
Bond lengths



Bond angles

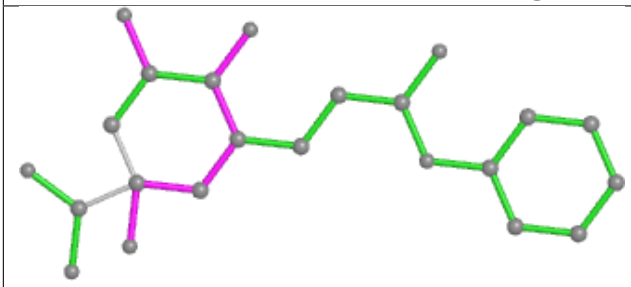


Torsions

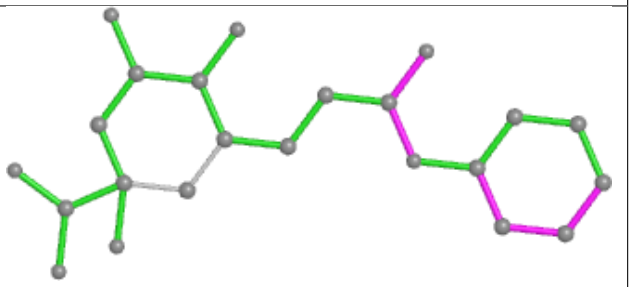


Rings

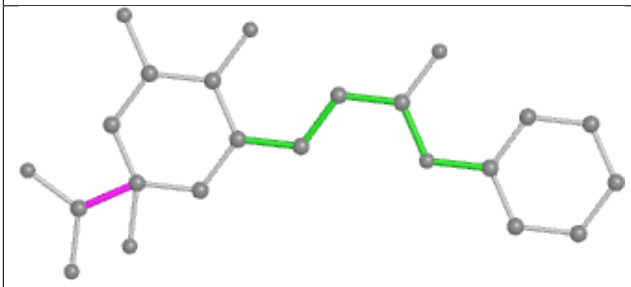
Ligand RJP F 147



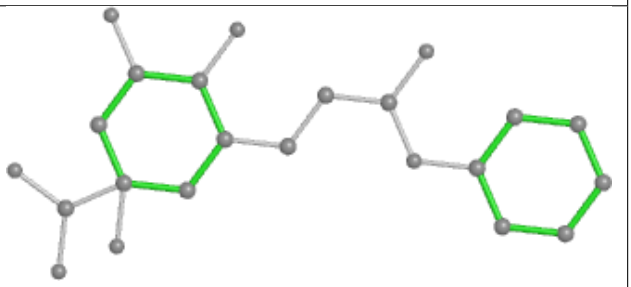
Bond lengths



Bond angles

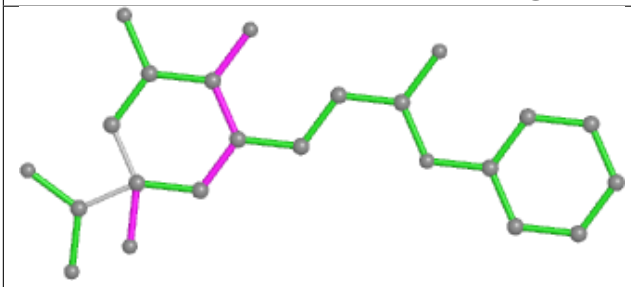


Torsions

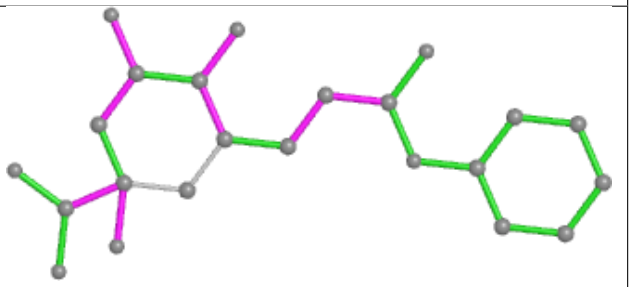


Rings

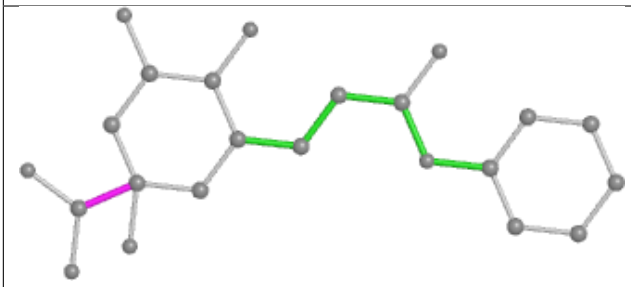
Ligand RJP T 147



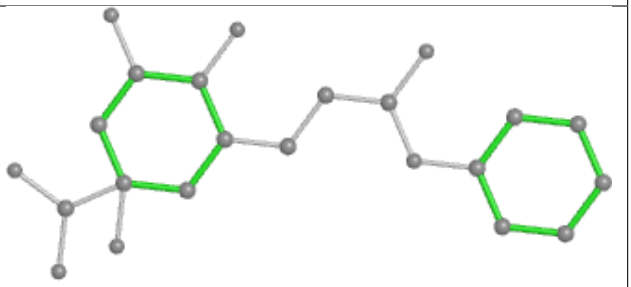
Bond lengths



Bond angles

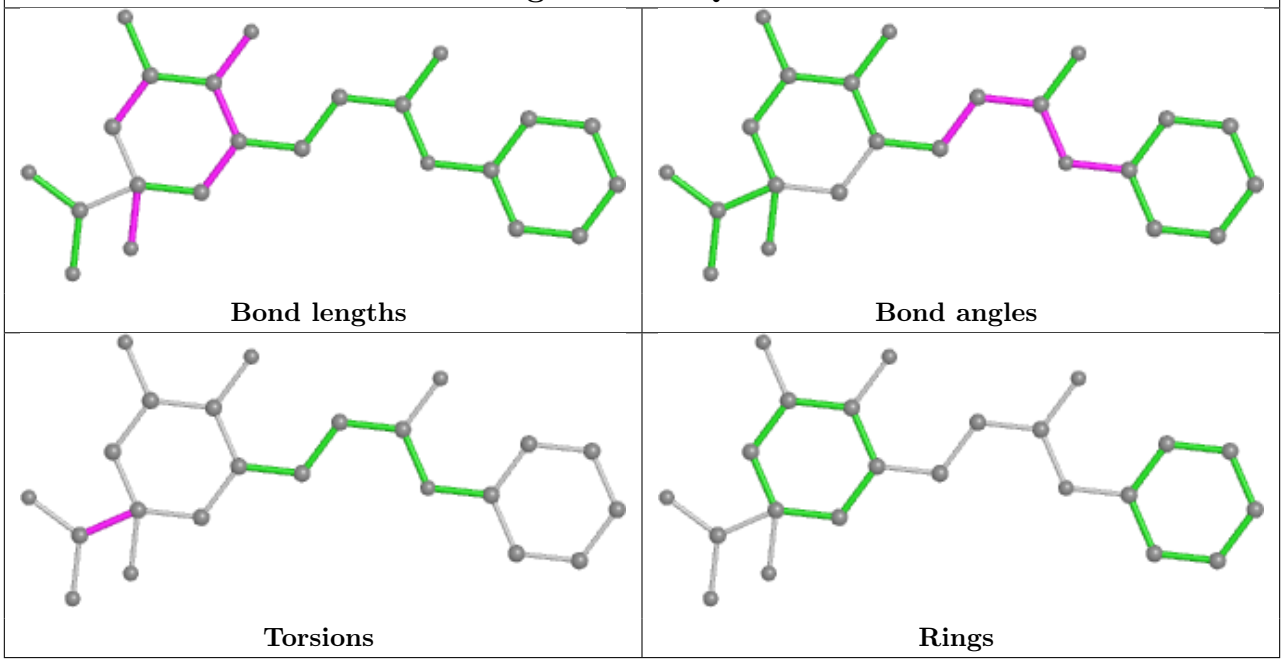


Torsions

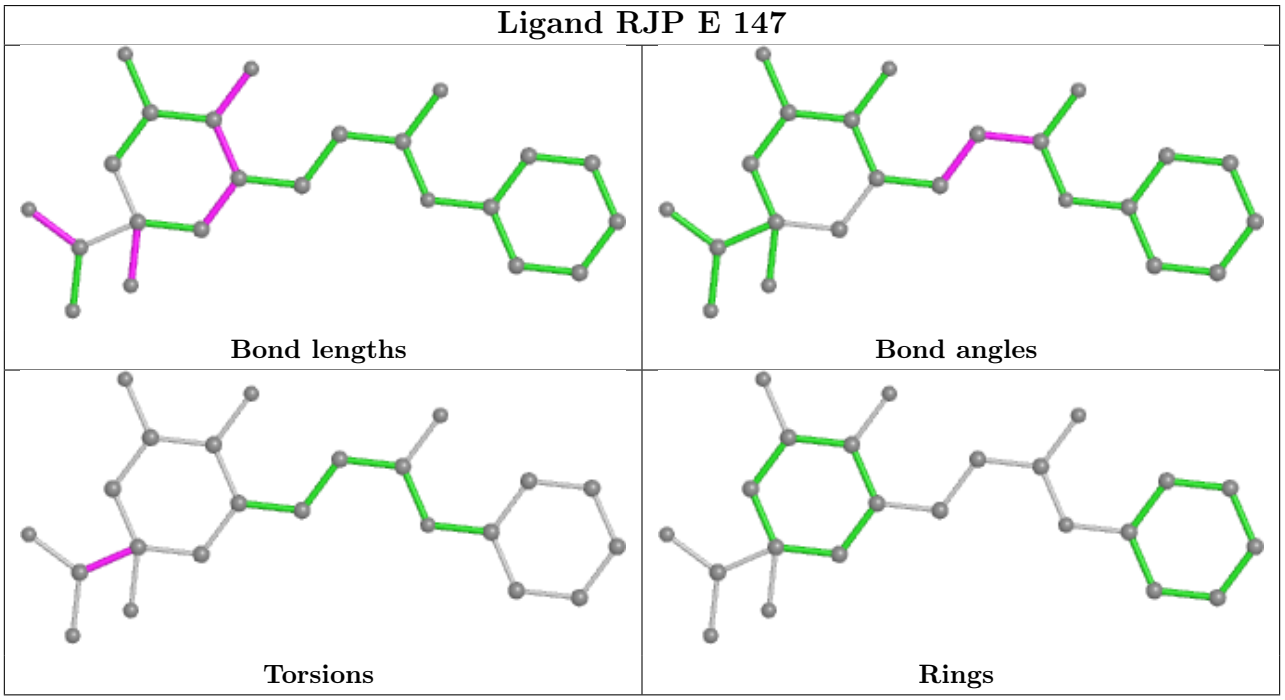


Rings

Ligand RJP Q 147



Ligand RJP E 147



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, 95^{th} 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	141/147 (95%)	-0.71	1 (0%) 87 88	11, 16, 29, 49	0
1	B	141/147 (95%)	-0.69	1 (0%) 87 88	7, 13, 34, 52	0
1	C	141/147 (95%)	-0.72	0 100 100	9, 14, 29, 48	0
1	D	141/147 (95%)	-0.69	1 (0%) 87 88	9, 13, 34, 46	0
1	E	141/147 (95%)	-0.65	1 (0%) 87 88	9, 14, 33, 51	0
1	F	141/147 (95%)	-0.68	1 (0%) 87 88	8, 14, 28, 53	0
1	G	141/147 (95%)	-0.67	1 (0%) 87 88	10, 16, 30, 53	1 (0%)
1	H	141/147 (95%)	-0.70	1 (0%) 87 88	9, 14, 29, 56	0
1	I	141/147 (95%)	-0.62	0 100 100	9, 14, 38, 49	0
1	J	143/147 (97%)	-0.67	1 (0%) 87 88	9, 13, 32, 51	0
1	K	142/147 (96%)	-0.72	0 100 100	10, 14, 29, 44	0
1	L	141/147 (95%)	-0.63	0 100 100	10, 16, 36, 53	0
1	M	141/147 (95%)	-0.69	0 100 100	10, 16, 29, 46	0
1	N	141/147 (95%)	-0.68	0 100 100	8, 13, 35, 48	0
1	O	141/147 (95%)	-0.69	1 (0%) 87 88	9, 13, 28, 47	0
1	P	142/147 (96%)	-0.69	1 (0%) 87 88	8, 13, 31, 49	0
1	Q	141/147 (95%)	-0.68	1 (0%) 87 88	9, 14, 33, 53	0
1	R	142/147 (96%)	-0.66	1 (0%) 87 88	9, 14, 29, 54	0
1	S	141/147 (95%)	-0.69	1 (0%) 87 88	10, 15, 28, 51	1 (0%)
1	T	141/147 (95%)	-0.72	1 (0%) 87 88	9, 13, 29, 54	0
1	U	141/147 (95%)	-0.60	0 100 100	9, 15, 36, 52	0
1	V	144/147 (97%)	-0.62	1 (0%) 87 88	8, 14, 32, 51	0
1	W	143/147 (97%)	-0.69	1 (0%) 87 88	10, 14, 28, 47	0
1	X	141/147 (95%)	-0.56	1 (0%) 87 88	9, 17, 36, 56	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å^2)	Q<0.9
All	All	3394/3528 (96%)	-0.67	17 (0%)	91 92	7, 14, 34, 56	2 (0%)

The worst 5 of 17 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	S	143	HIS	3.9
1	R	143	HIS	3.8
1	H	143	HIS	3.7
1	F	143	HIS	3.5
1	Q	143	HIS	3.4

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 monosaccharides 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(Å^2)	Q<0.9
2	RJP	F	147	23/23	0.94	0.10	10,14,21,27	0
2	RJP	Q	147	23/23	0.94	0.10	10,17,23,28	0
2	RJP	W	147	23/23	0.94	0.10	11,18,25,27	0
2	RJP	M	147	23/23	0.95	0.08	12,17,23,25	0
2	RJP	U	147	23/23	0.95	0.12	10,17,25,30	0
2	RJP	P	147	23/23	0.95	0.10	11,18,25,28	0
2	RJP	H	147	23/23	0.96	0.10	9,13,17,21	0
2	RJP	I	147	23/23	0.96	0.09	10,18,24,27	0
2	RJP	J	147	23/23	0.96	0.08	10,18,26,27	0
2	RJP	L	147	23/23	0.96	0.09	12,18,23,26	0
2	RJP	B	147	23/23	0.96	0.09	8,15,25,28	0
2	RJP	N	147	23/23	0.96	0.09	9,14,26,28	0
2	RJP	O	147	23/23	0.96	0.09	10,14,17,23	0

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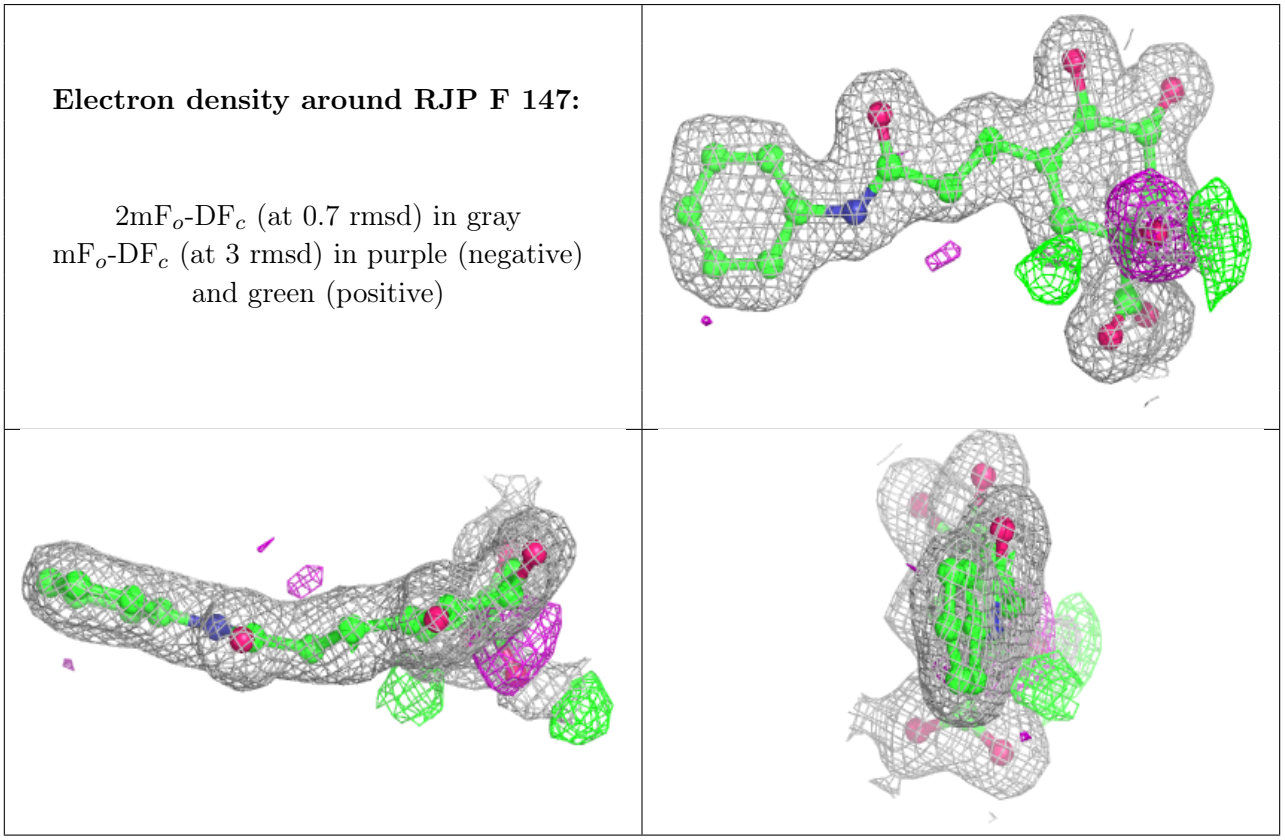
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	RJP	C	147	23/23	0.96	0.08	8,13,19,20	0
2	RJP	D	147	23/23	0.96	0.09	12,16,26,27	0
2	RJP	R	147	23/23	0.96	0.10	9,15,18,25	0
2	RJP	S	147	23/23	0.96	0.09	11,16,21,27	0
2	RJP	T	147	23/23	0.96	0.08	8,13,17,26	0
2	RJP	A	147	23/23	0.96	0.09	13,16,26,26	0
2	RJP	V	147	23/23	0.96	0.09	9,17,29,30	0
2	RJP	G	147	23/23	0.96	0.08	11,16,21,26	0
2	RJP	X	147	23/23	0.96	0.10	11,17,24,24	0
2	RJP	K	147	23/23	0.97	0.08	10,16,24,26	0
2	RJP	E	147	23/23	0.97	0.08	9,17,25,26	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.

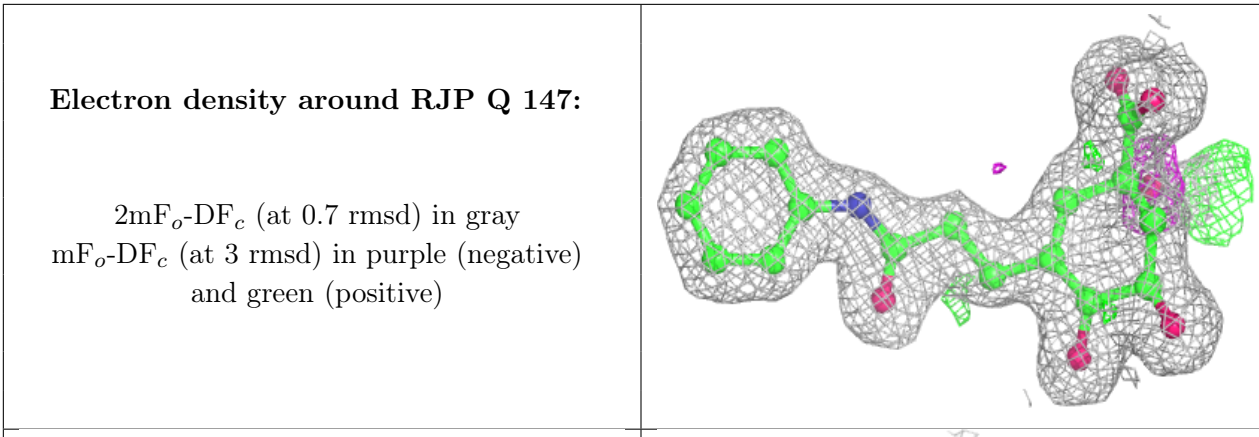
Electron density around RJP F 147:

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



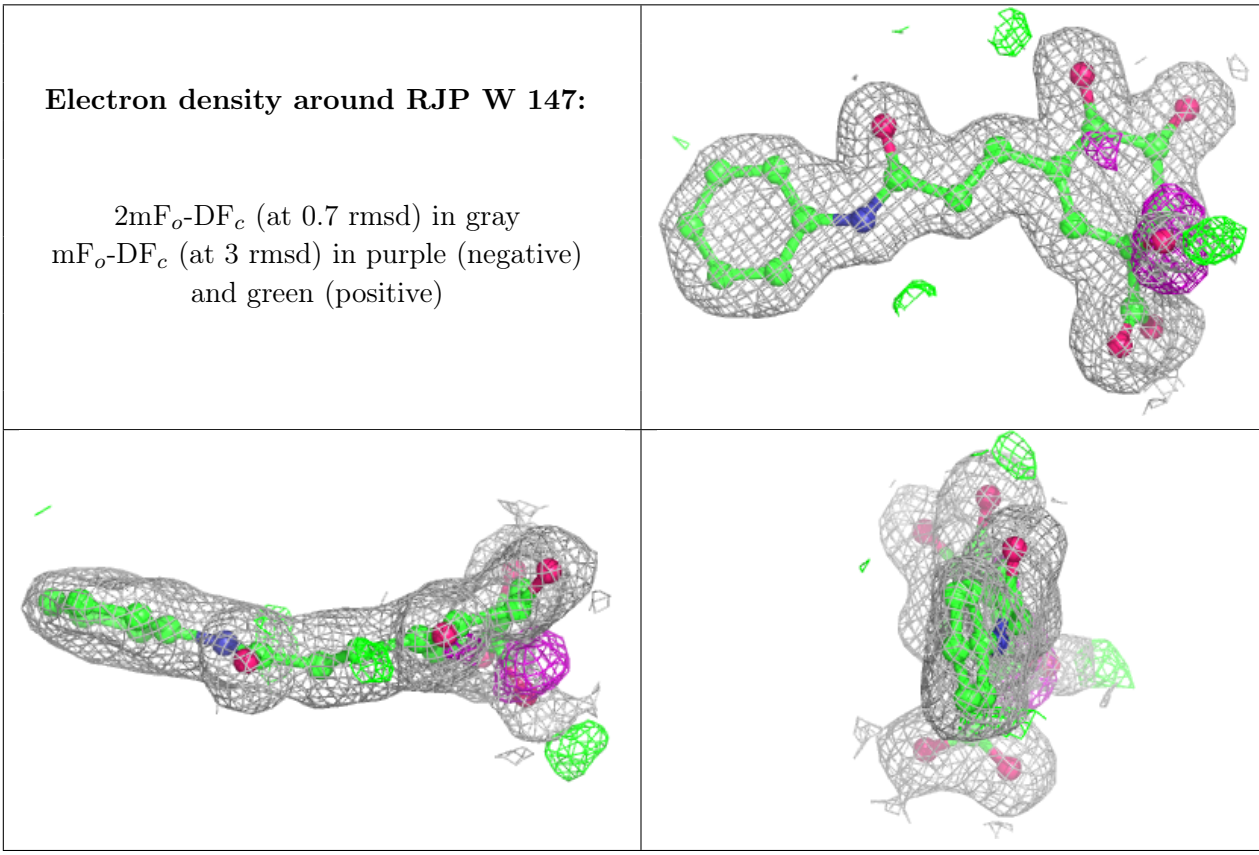
Electron density around RJP Q 147:

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



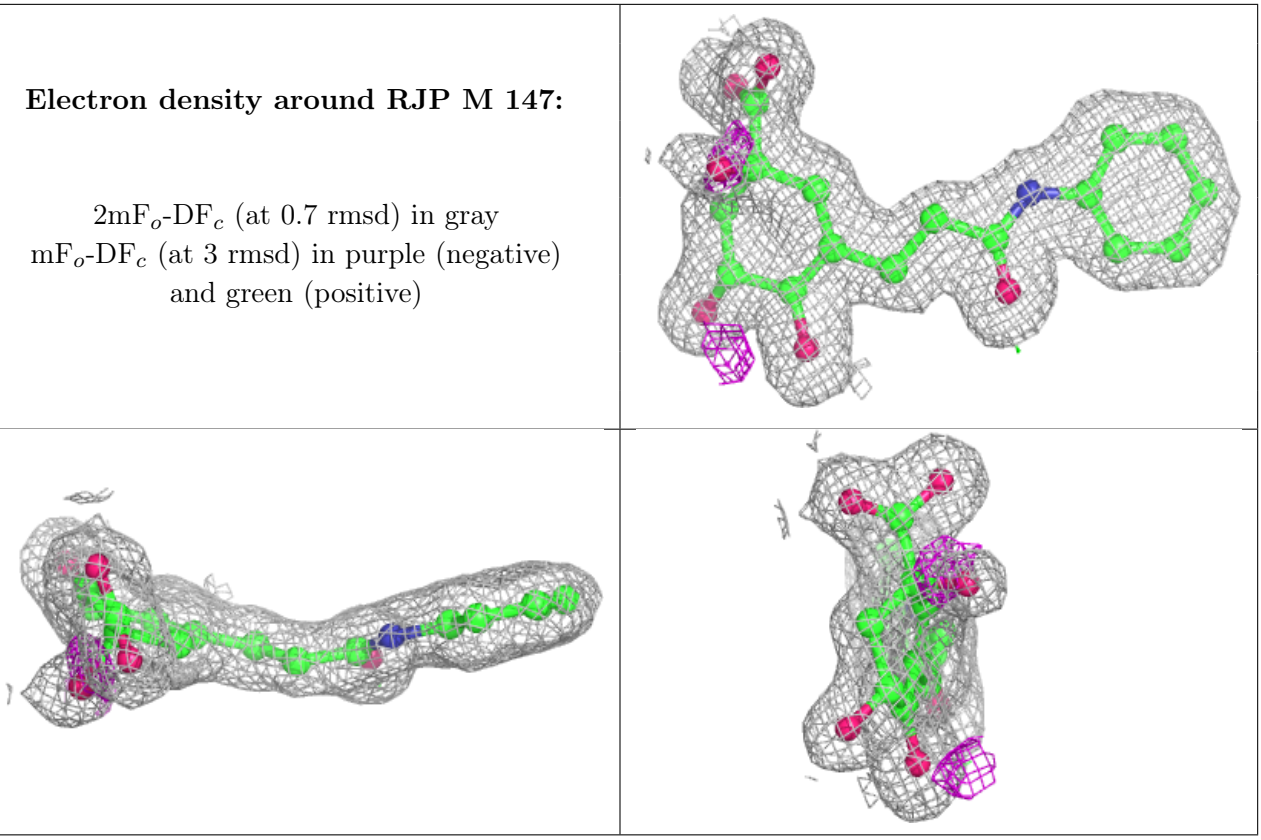
Electron density around RJP W 147:

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



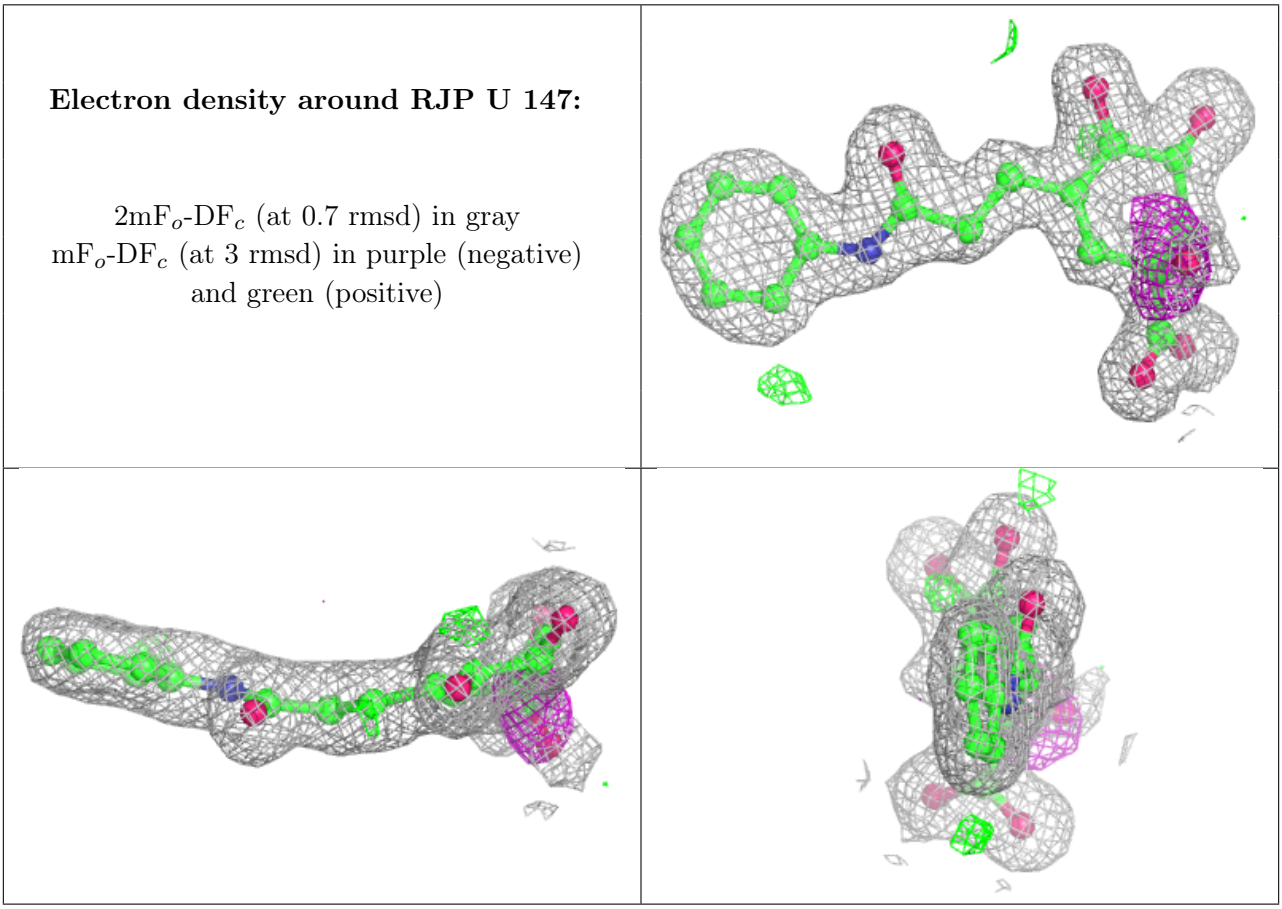
Electron density around RJP M 147:

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



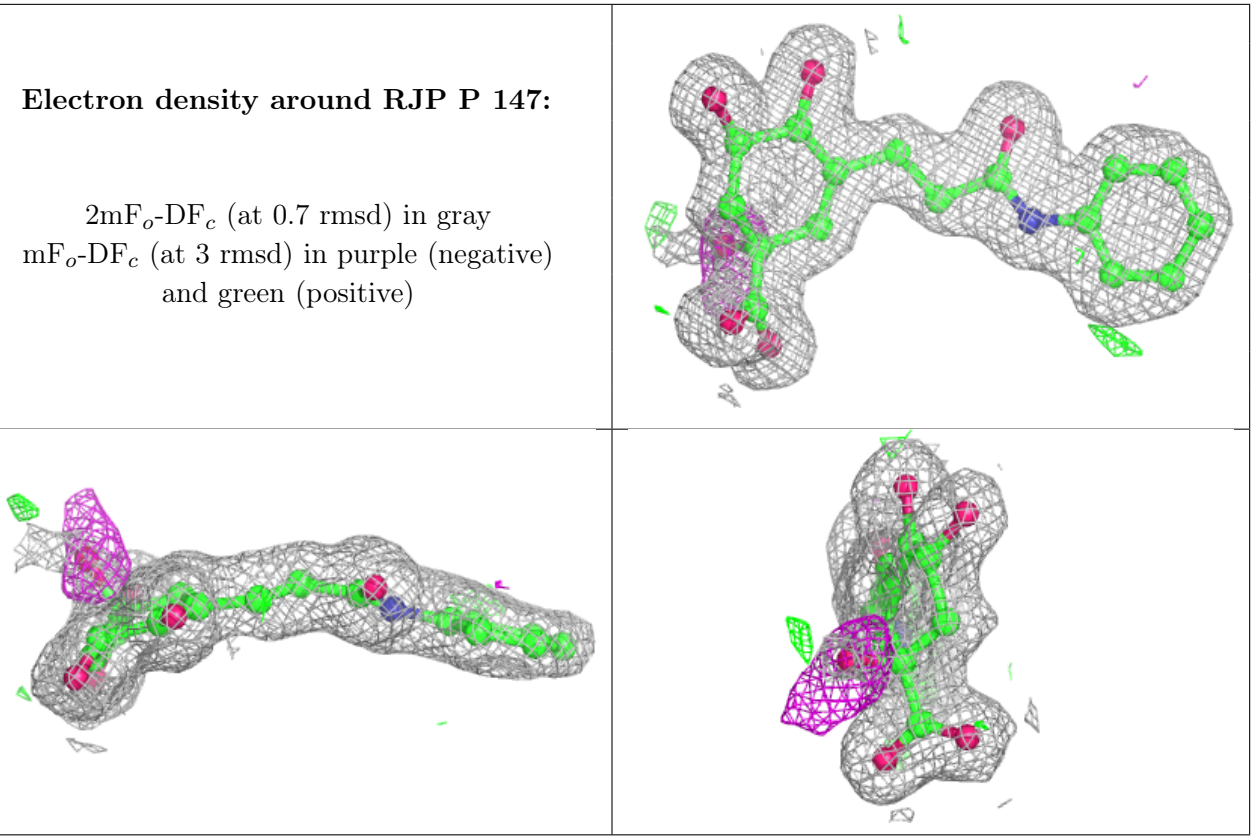
Electron density around RJP U 147:

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



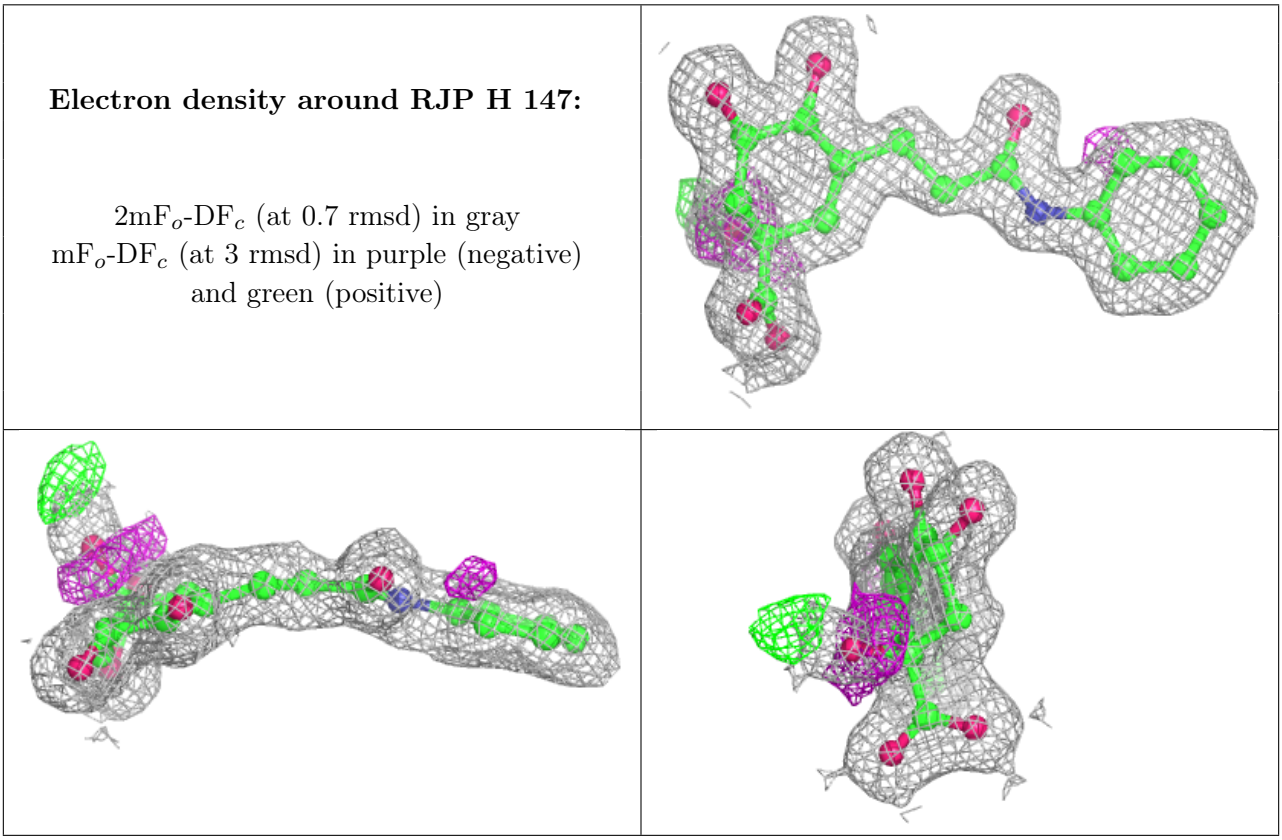
Electron density around RJP P 147:

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



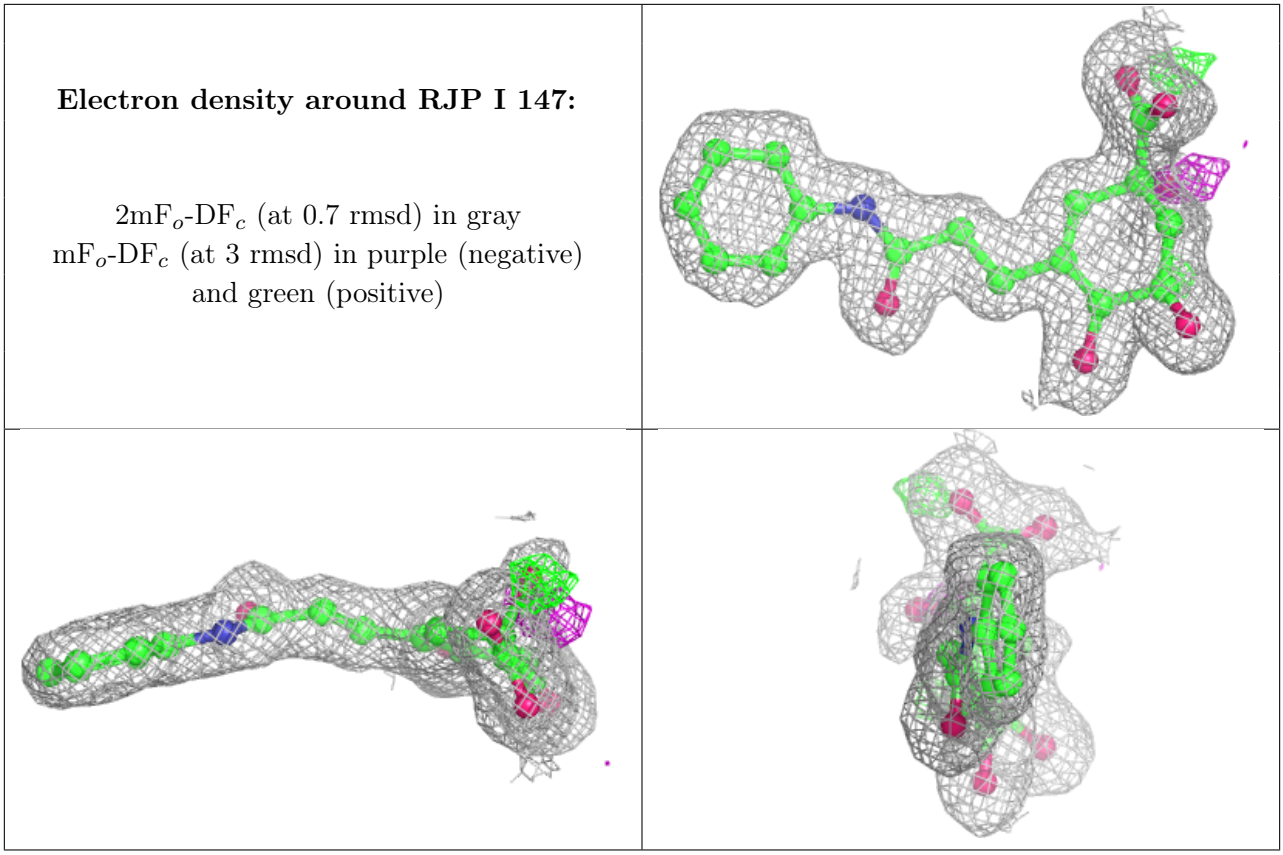
Electron density around RJP H 147:

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



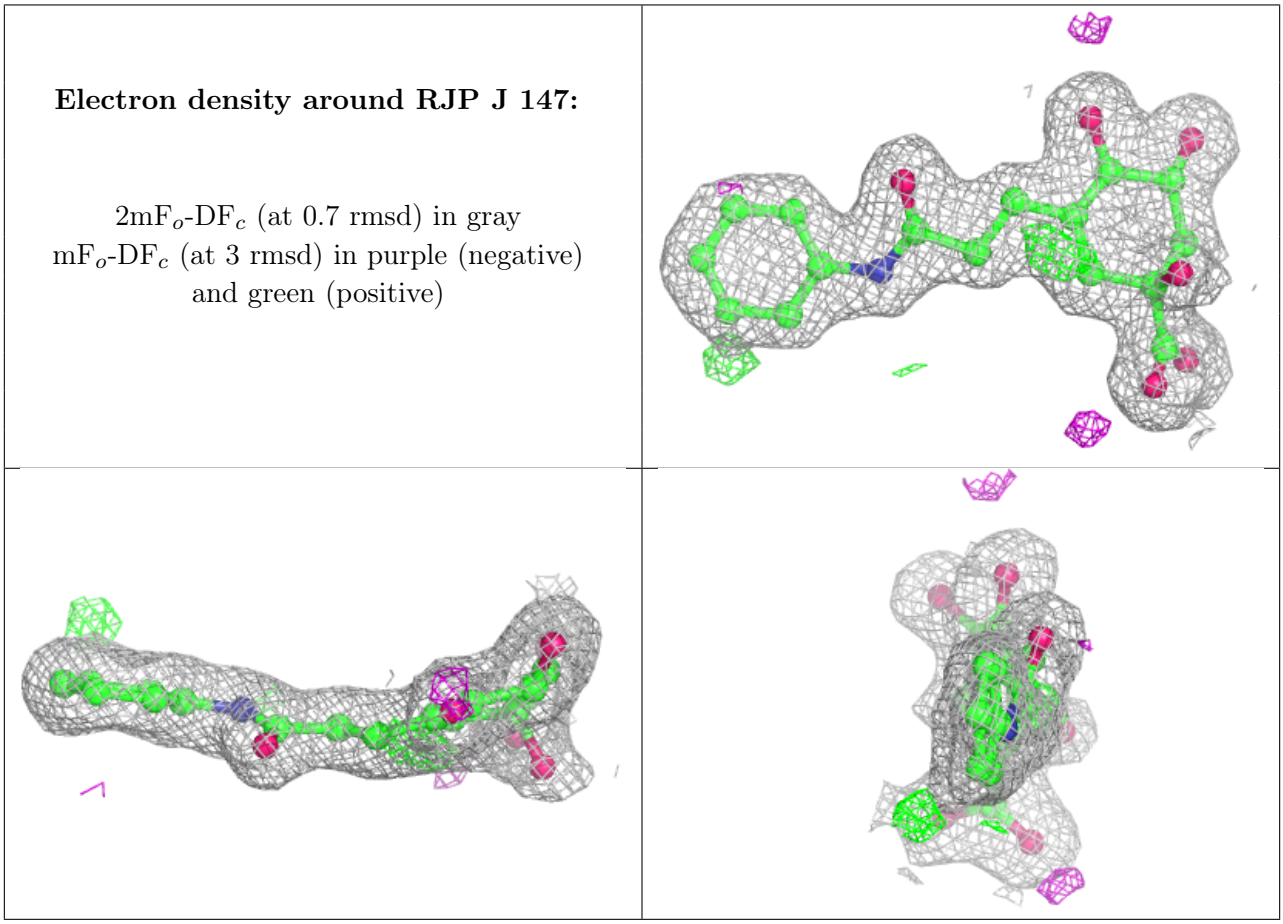
Electron density around RJP I 147:

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



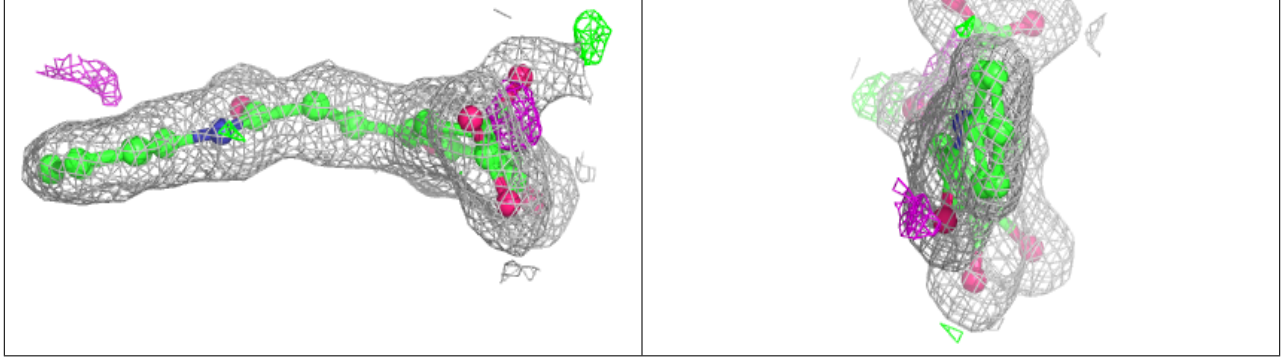
Electron density around RJP J 147:

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



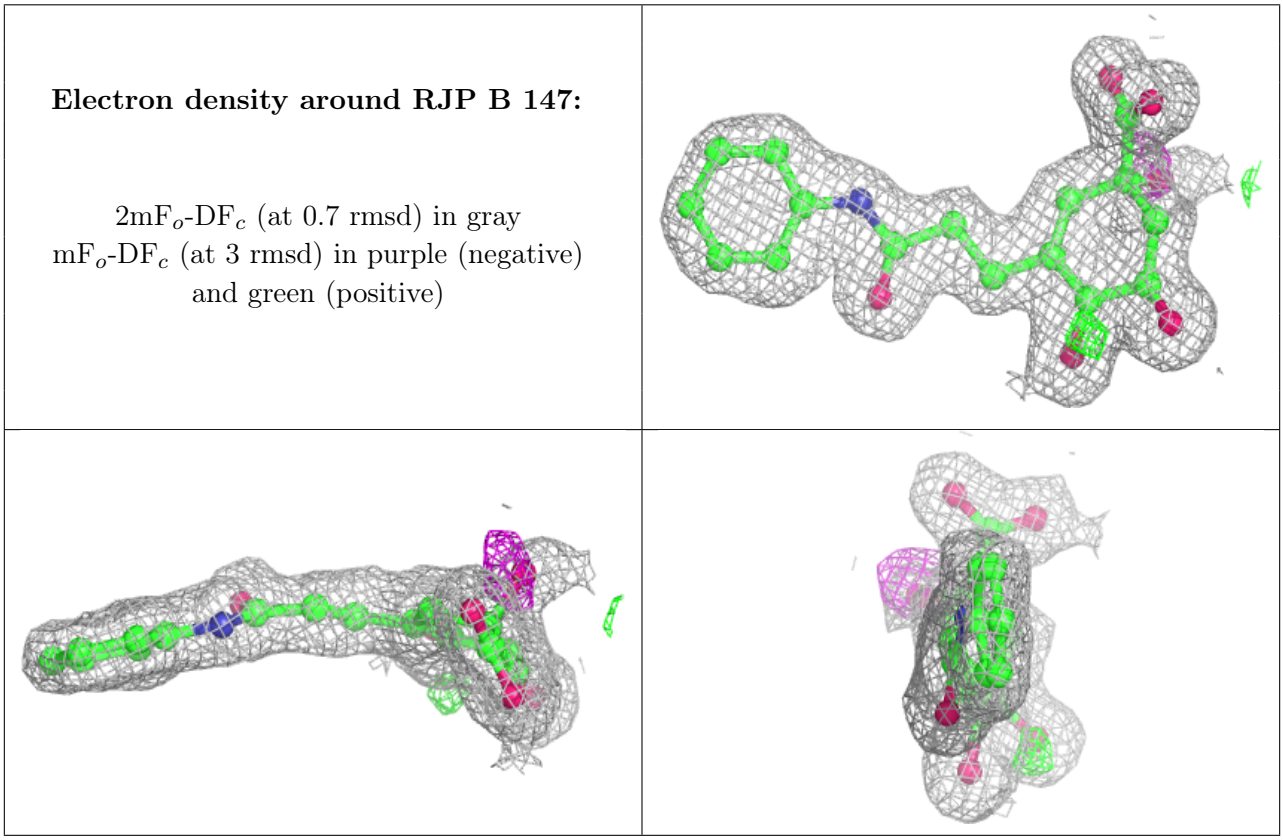
Electron density around RJP L 147:

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



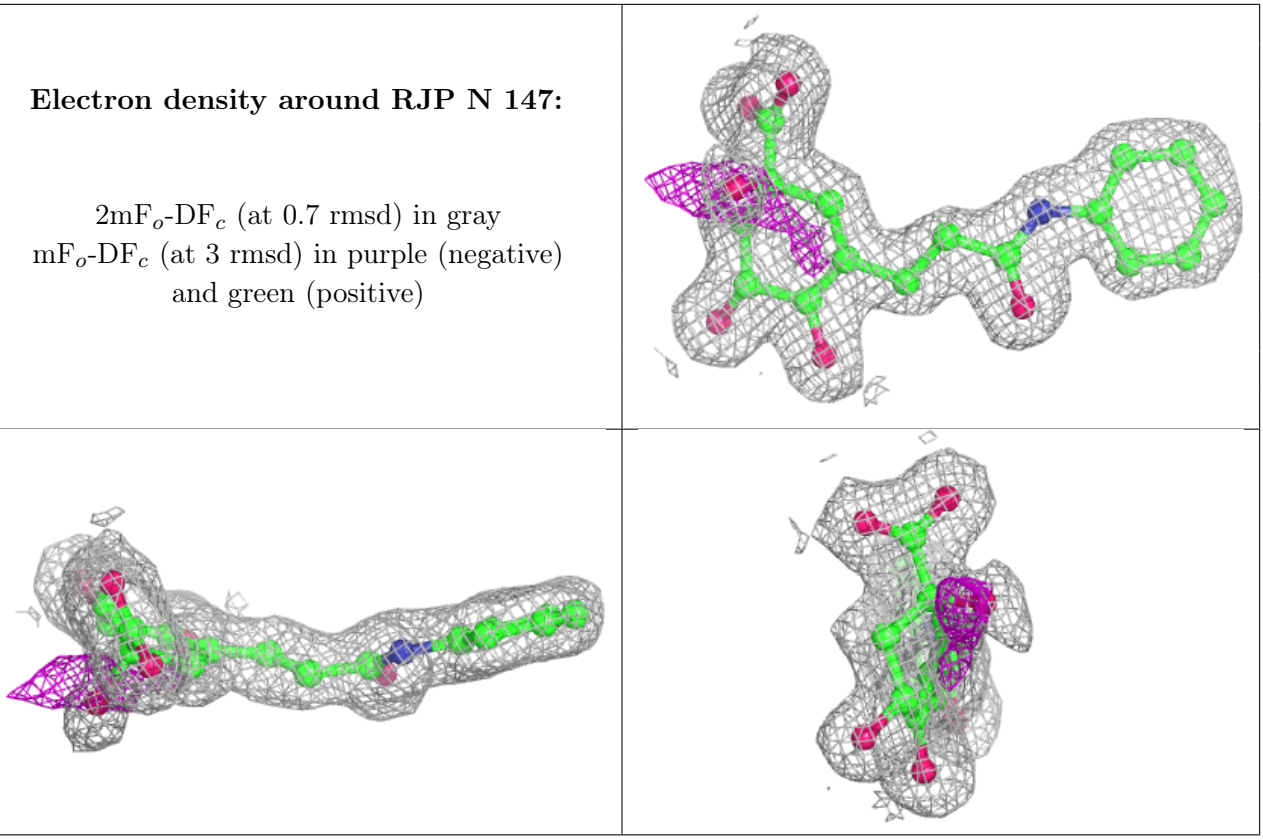
Electron density around RJP B 147:

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



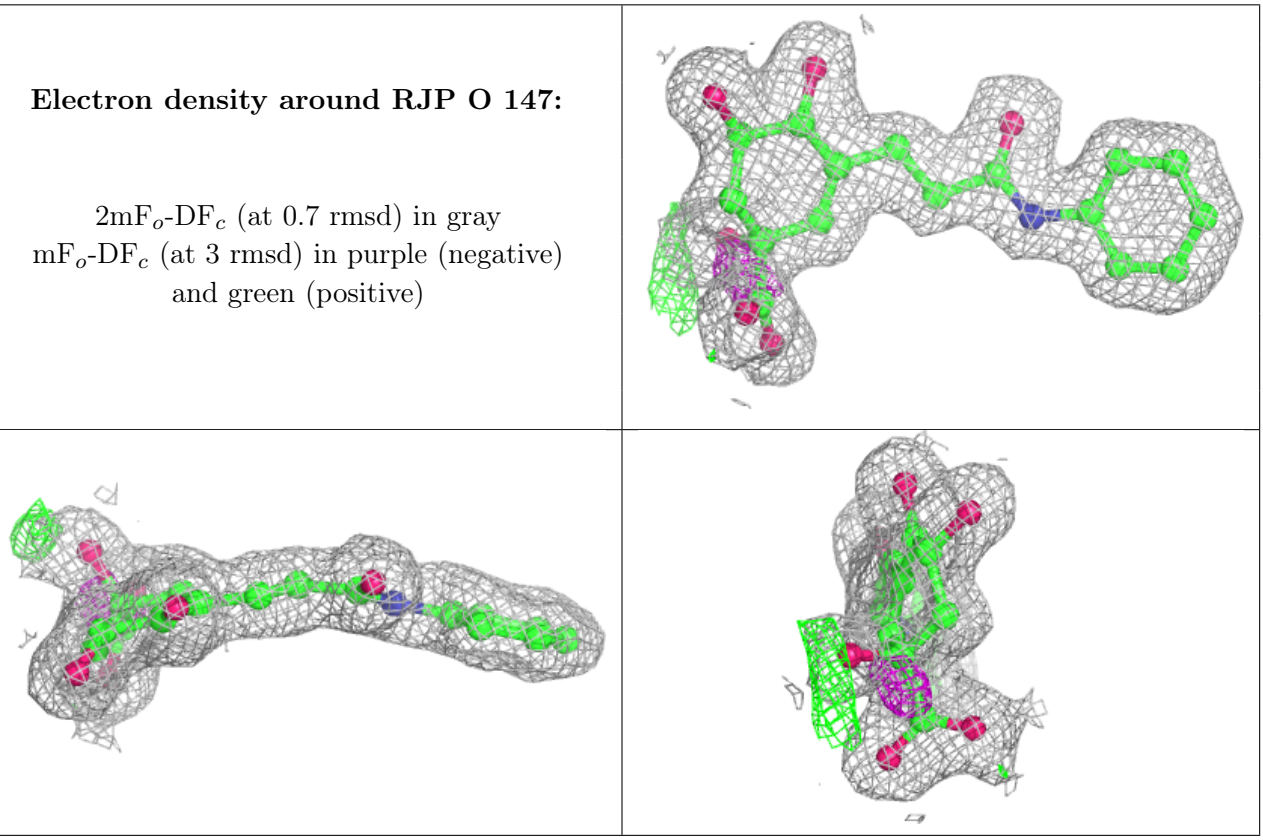
Electron density around RJP N 147:

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



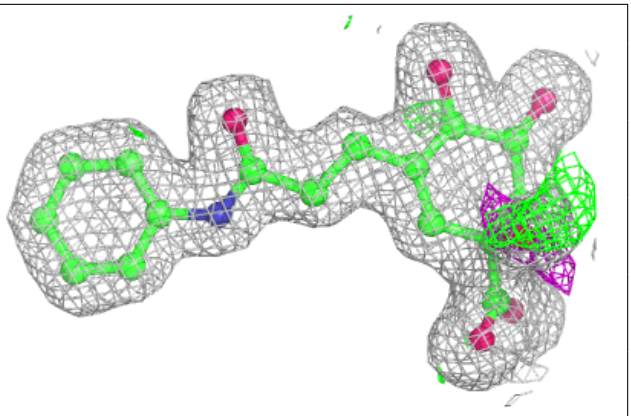
Electron density around RJP O 147:

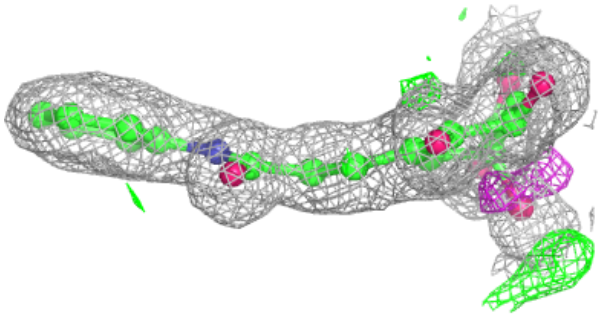
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

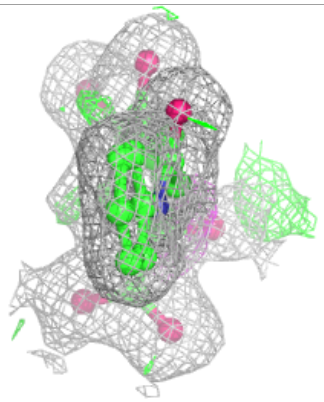


Electron density around RJP C 147:

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

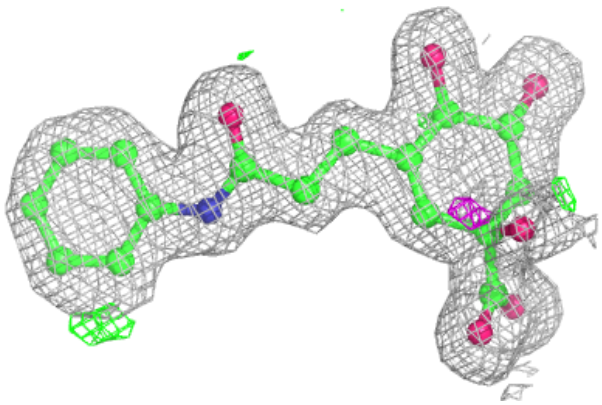


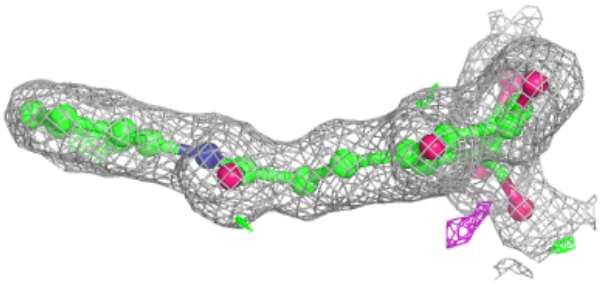


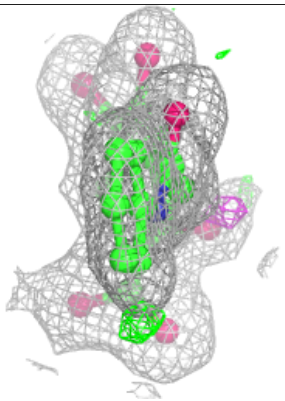


Electron density around RJP D 147:

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

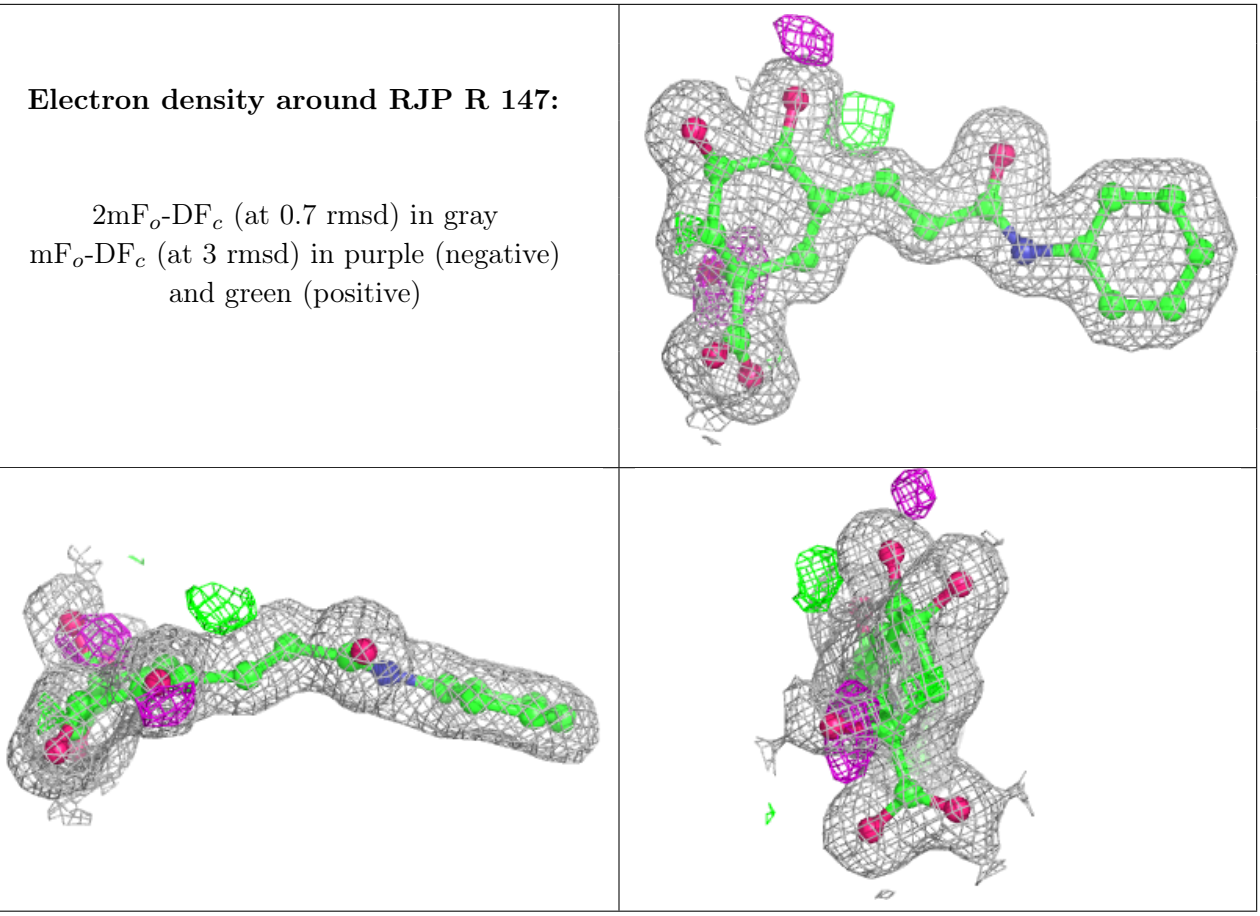






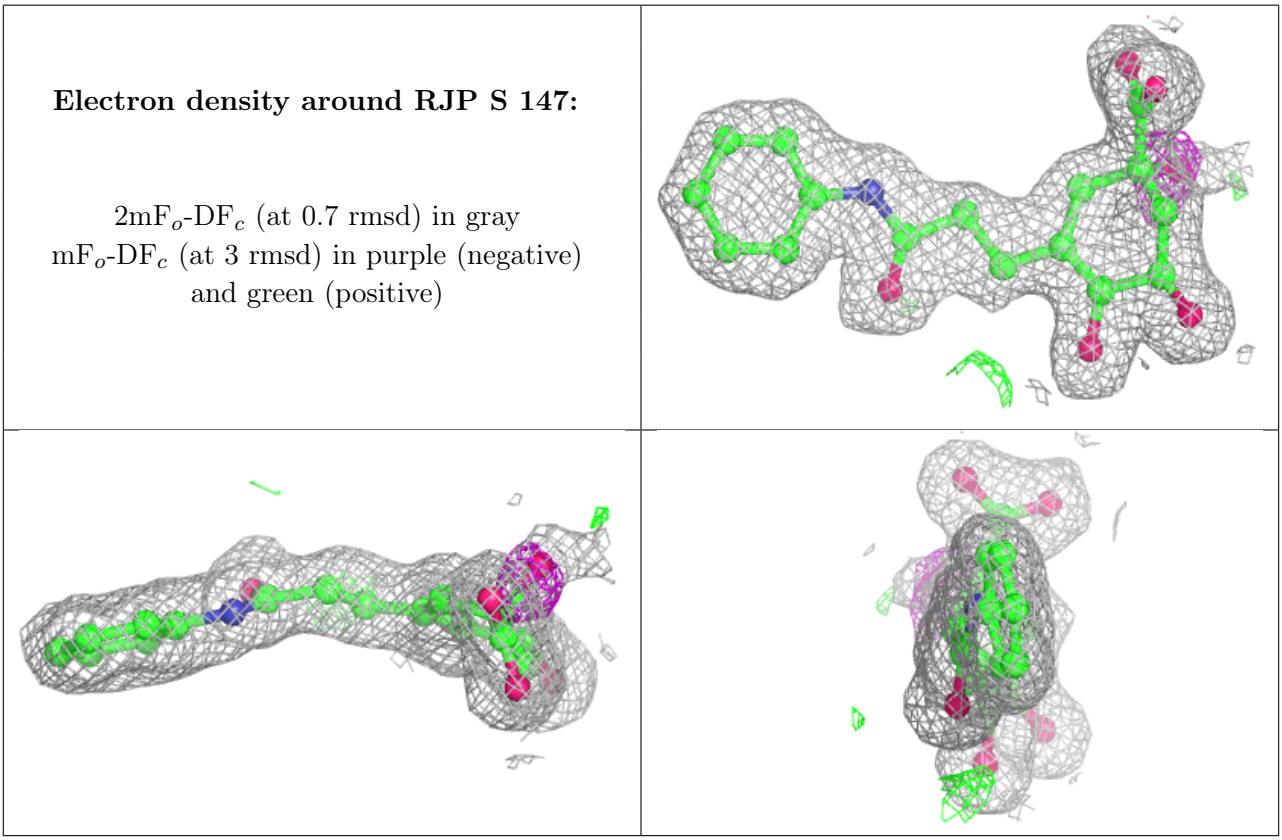
Electron density around RJP R 147:

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



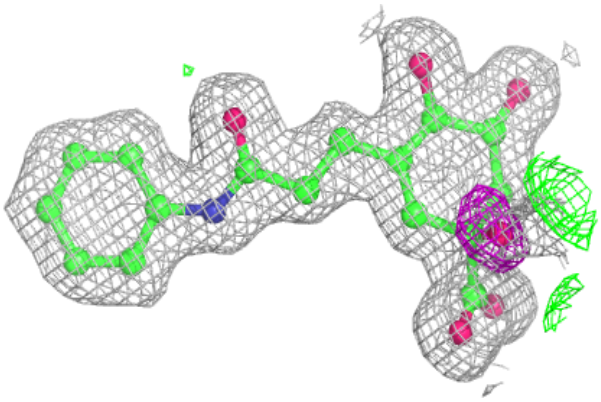
Electron density around RJP S 147:

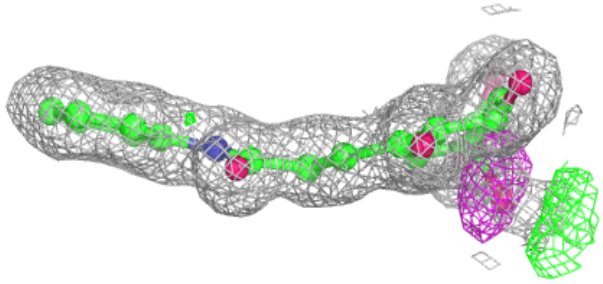
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

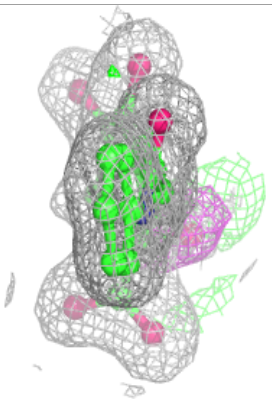


Electron density around RJP T 147:

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

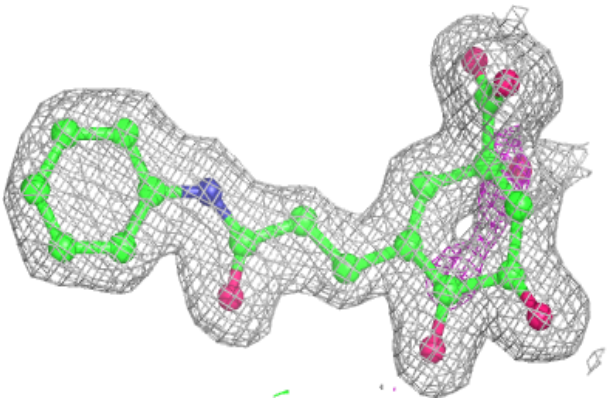


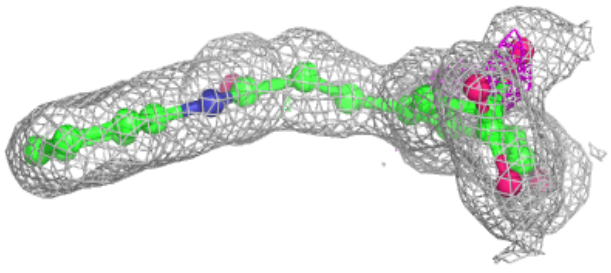


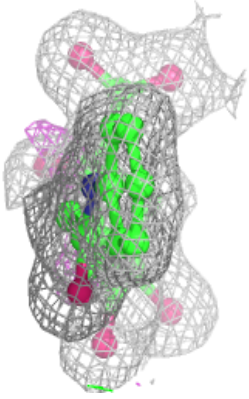


Electron density around RJP A 147:

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

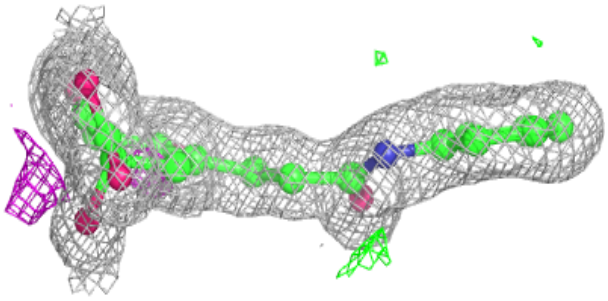


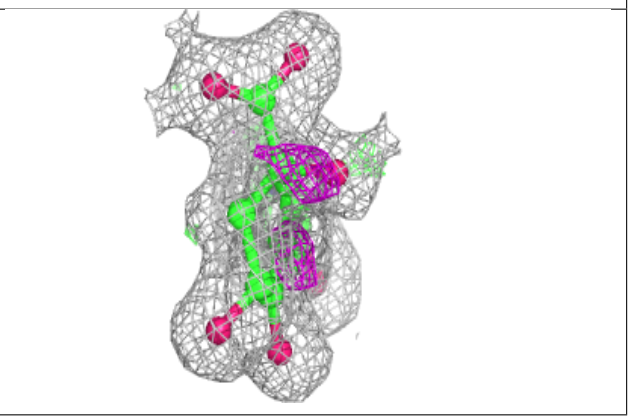




Electron density around RJP V 147:

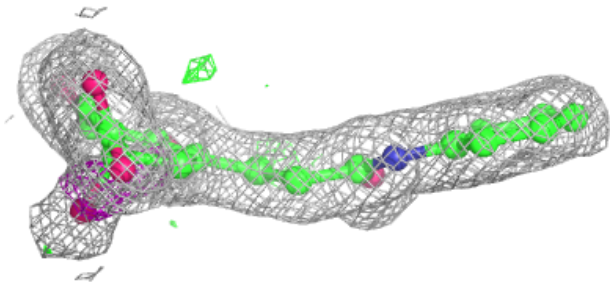
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

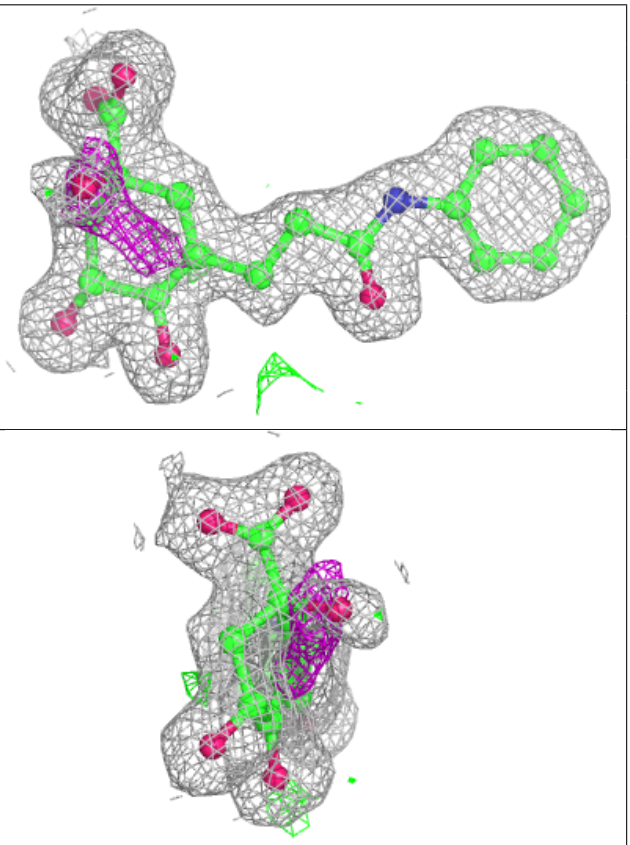




Electron density around RJP G 147:

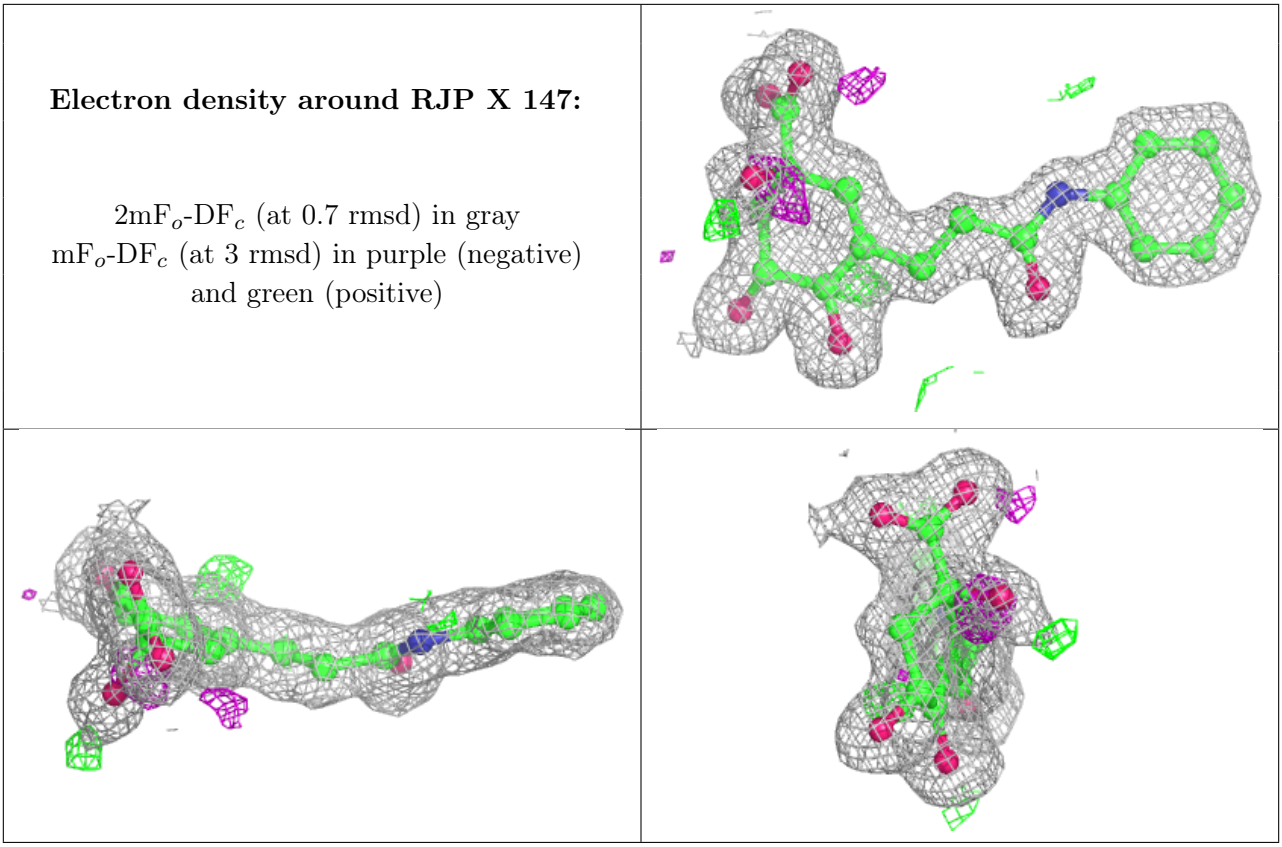
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)





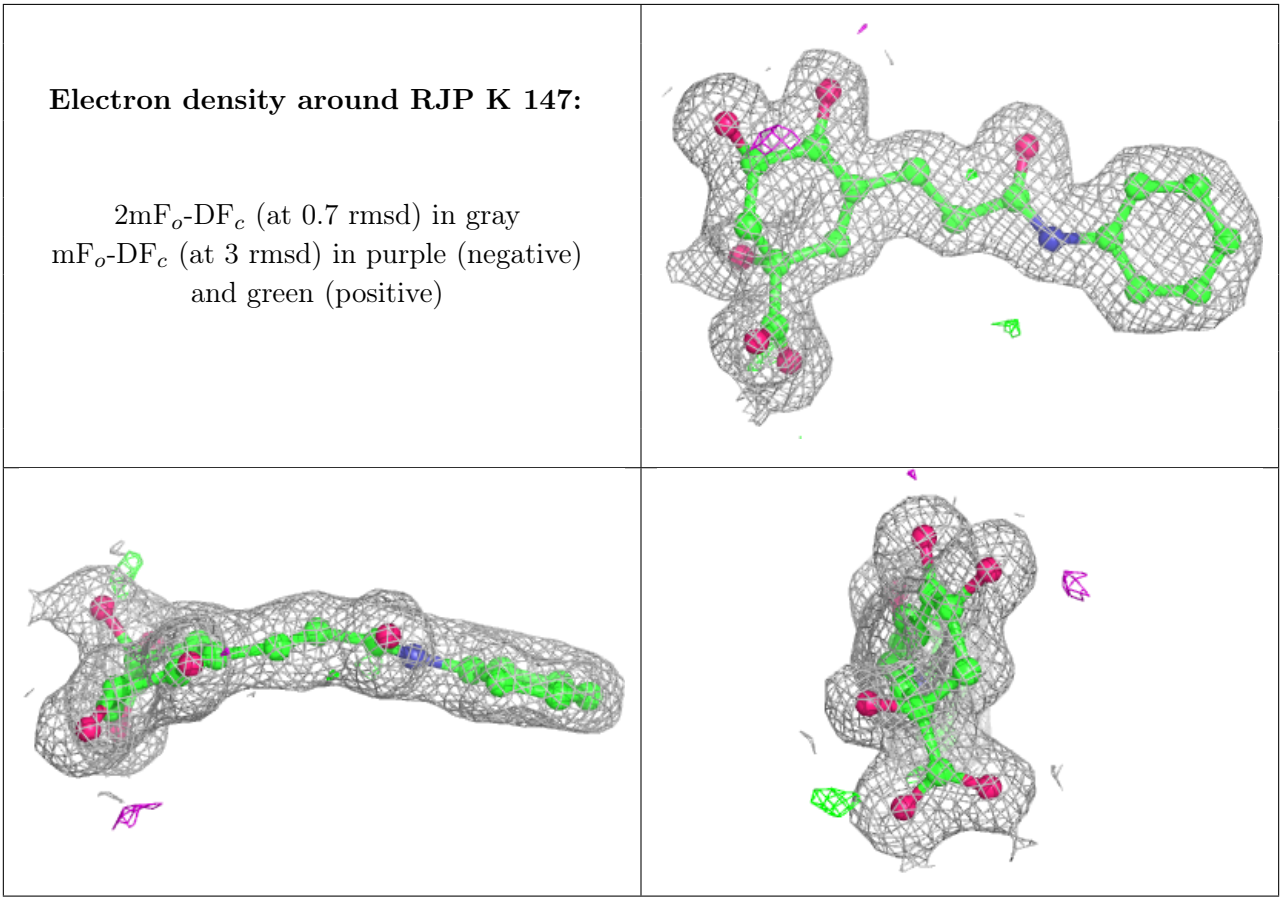
Electron density around RJP X 147:

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



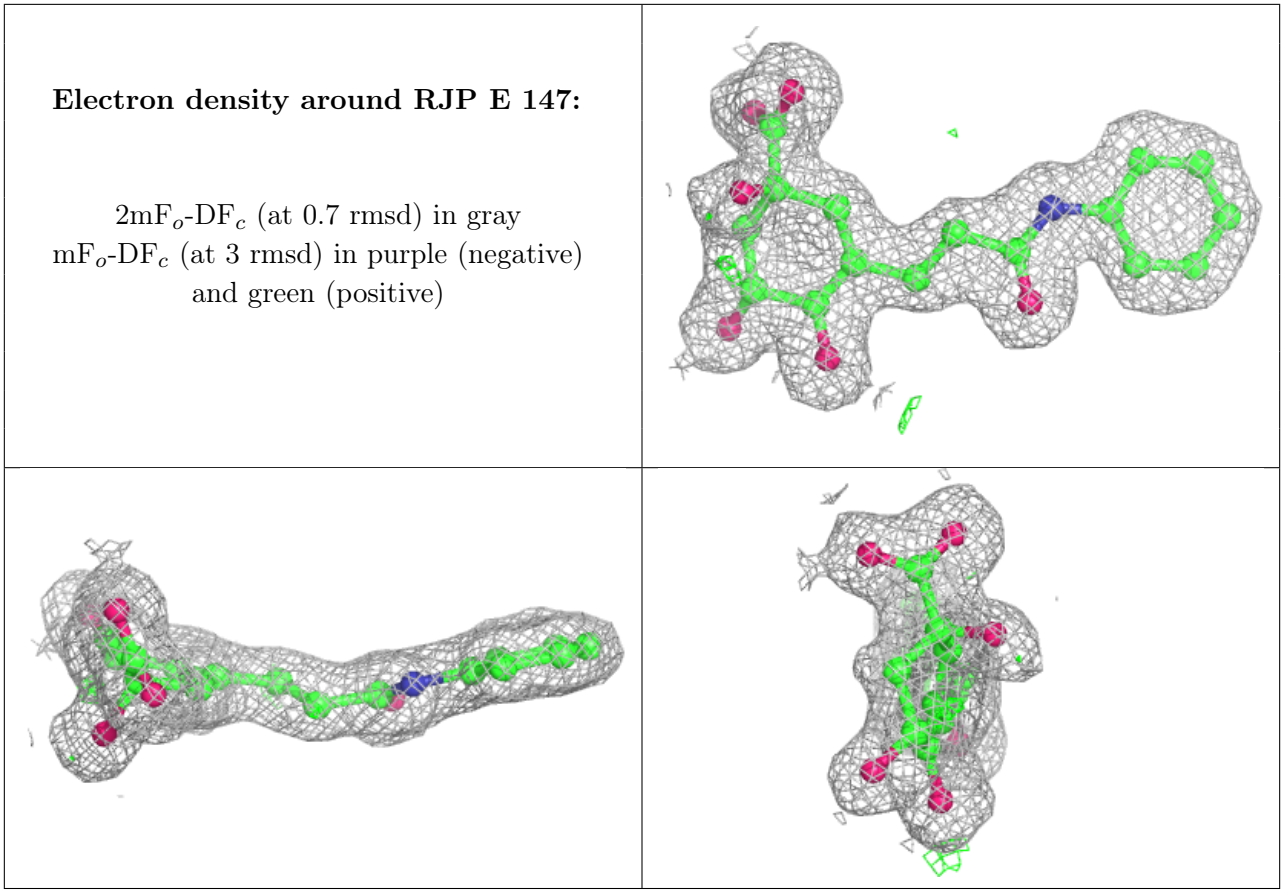
Electron density around RJP K 147:

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



Electron density around RJP E 147:

$2mF_o$-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.