



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 5, 2022 – 09:14 pm GMT

PDB ID : 6ZK3  
Title : Plant nucleoside hydrolase - ZmNRh2b in complex with ribose  
Authors : Morera, S.; Vigouroux, A.; Kopecny, D.  
Deposited on : 2020-06-29  
Resolution : 1.75 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](mailto: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.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.24
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.24

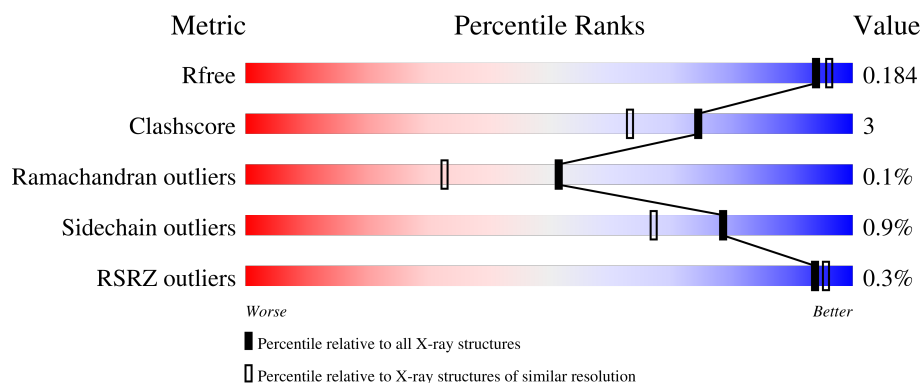
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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  $\geq 3$ , 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	343	
1	B	343	
1	C	343	
1	D	343	
1	E	343	

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Mol	Chain	Length	Quality of chain
1	F	343	 84% 8% 8%

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
4	EDO	E	404	-	-	X	-
4	EDO	F	404	-	-	X	-
5	PGE	B	407	-	-	X	-
5	PGE	E	408	-	-	X	-
6	PEG	F	409	-	-	X	-

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 16641 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 Pyrimidine-specific ribonucleoside hydrolase rihA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	S	0	1	0
			2435	1564	411	450	10			
1	B	318	Total	C	N	O	S	0	1	0
			2428	1561	409	448	10			
1	C	318	Total	C	N	O	S	0	0	0
			2421	1556	408	447	10			
1	D	317	Total	C	N	O	S	0	3	0
			2428	1561	408	447	12			
1	E	316	Total	C	N	O	S	0	1	0
			2405	1546	403	445	11			
1	F	317	Total	C	N	O	S	0	3	0
			2428	1560	409	448	11			

There are 108 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	MET	-	initiating methionine	UNP B6THD4
A	-16	GLY	-	expression tag	UNP B6THD4
A	-15	SER	-	expression tag	UNP B6THD4
A	-14	SER	-	expression tag	UNP B6THD4
A	-13	HIS	-	expression tag	UNP B6THD4
A	-12	HIS	-	expression tag	UNP B6THD4
A	-11	HIS	-	expression tag	UNP B6THD4
A	-10	HIS	-	expression tag	UNP B6THD4
A	-9	HIS	-	expression tag	UNP B6THD4
A	-8	HIS	-	expression tag	UNP B6THD4
A	-7	SER	-	expression tag	UNP B6THD4
A	-6	GLN	-	expression tag	UNP B6THD4
A	-5	ASP	-	expression tag	UNP B6THD4
A	-4	PRO	-	expression tag	UNP B6THD4
A	-3	ASN	-	expression tag	UNP B6THD4
A	-2	SER	-	expression tag	UNP B6THD4
A	-1	SER	-	expression tag	UNP B6THD4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP B6THD4
B	-17	MET	-	initiating methionine	UNP B6THD4
B	-16	GLY	-	expression tag	UNP B6THD4
B	-15	SER	-	expression tag	UNP B6THD4
B	-14	SER	-	expression tag	UNP B6THD4
B	-13	HIS	-	expression tag	UNP B6THD4
B	-12	HIS	-	expression tag	UNP B6THD4
B	-11	HIS	-	expression tag	UNP B6THD4
B	-10	HIS	-	expression tag	UNP B6THD4
B	-9	HIS	-	expression tag	UNP B6THD4
B	-8	HIS	-	expression tag	UNP B6THD4
B	-7	SER	-	expression tag	UNP B6THD4
B	-6	GLN	-	expression tag	UNP B6THD4
B	-5	ASP	-	expression tag	UNP B6THD4
B	-4	PRO	-	expression tag	UNP B6THD4
B	-3	ASN	-	expression tag	UNP B6THD4
B	-2	SER	-	expression tag	UNP B6THD4
B	-1	SER	-	expression tag	UNP B6THD4
B	0	SER	-	expression tag	UNP B6THD4
C	-17	MET	-	initiating methionine	UNP B6THD4
C	-16	GLY	-	expression tag	UNP B6THD4
C	-15	SER	-	expression tag	UNP B6THD4
C	-14	SER	-	expression tag	UNP B6THD4
C	-13	HIS	-	expression tag	UNP B6THD4
C	-12	HIS	-	expression tag	UNP B6THD4
C	-11	HIS	-	expression tag	UNP B6THD4
C	-10	HIS	-	expression tag	UNP B6THD4
C	-9	HIS	-	expression tag	UNP B6THD4
C	-8	HIS	-	expression tag	UNP B6THD4
C	-7	SER	-	expression tag	UNP B6THD4
C	-6	GLN	-	expression tag	UNP B6THD4
C	-5	ASP	-	expression tag	UNP B6THD4
C	-4	PRO	-	expression tag	UNP B6THD4
C	-3	ASN	-	expression tag	UNP B6THD4
C	-2	SER	-	expression tag	UNP B6THD4
C	-1	SER	-	expression tag	UNP B6THD4
C	0	SER	-	expression tag	UNP B6THD4
D	-17	MET	-	initiating methionine	UNP B6THD4
D	-16	GLY	-	expression tag	UNP B6THD4
D	-15	SER	-	expression tag	UNP B6THD4
D	-14	SER	-	expression tag	UNP B6THD4
D	-13	HIS	-	expression tag	UNP B6THD4

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-12	HIS	-	expression tag	UNP B6THD4
D	-11	HIS	-	expression tag	UNP B6THD4
D	-10	HIS	-	expression tag	UNP B6THD4
D	-9	HIS	-	expression tag	UNP B6THD4
D	-8	HIS	-	expression tag	UNP B6THD4
D	-7	SER	-	expression tag	UNP B6THD4
D	-6	GLN	-	expression tag	UNP B6THD4
D	-5	ASP	-	expression tag	UNP B6THD4
D	-4	PRO	-	expression tag	UNP B6THD4
D	-3	ASN	-	expression tag	UNP B6THD4
D	-2	SER	-	expression tag	UNP B6THD4
D	-1	SER	-	expression tag	UNP B6THD4
D	0	SER	-	expression tag	UNP B6THD4
E	-17	MET	-	initiating methionine	UNP B6THD4
E	-16	GLY	-	expression tag	UNP B6THD4
E	-15	SER	-	expression tag	UNP B6THD4
E	-14	SER	-	expression tag	UNP B6THD4
E	-13	HIS	-	expression tag	UNP B6THD4
E	-12	HIS	-	expression tag	UNP B6THD4
E	-11	HIS	-	expression tag	UNP B6THD4
E	-10	HIS	-	expression tag	UNP B6THD4
E	-9	HIS	-	expression tag	UNP B6THD4
E	-8	HIS	-	expression tag	UNP B6THD4
E	-7	SER	-	expression tag	UNP B6THD4
E	-6	GLN	-	expression tag	UNP B6THD4
E	-5	ASP	-	expression tag	UNP B6THD4
E	-4	PRO	-	expression tag	UNP B6THD4
E	-3	ASN	-	expression tag	UNP B6THD4
E	-2	SER	-	expression tag	UNP B6THD4
E	-1	SER	-	expression tag	UNP B6THD4
E	0	SER	-	expression tag	UNP B6THD4
F	-17	MET	-	initiating methionine	UNP B6THD4
F	-16	GLY	-	expression tag	UNP B6THD4
F	-15	SER	-	expression tag	UNP B6THD4
F	-14	SER	-	expression tag	UNP B6THD4
F	-13	HIS	-	expression tag	UNP B6THD4
F	-12	HIS	-	expression tag	UNP B6THD4
F	-11	HIS	-	expression tag	UNP B6THD4
F	-10	HIS	-	expression tag	UNP B6THD4
F	-9	HIS	-	expression tag	UNP B6THD4
F	-8	HIS	-	expression tag	UNP B6THD4
F	-7	SER	-	expression tag	UNP B6THD4

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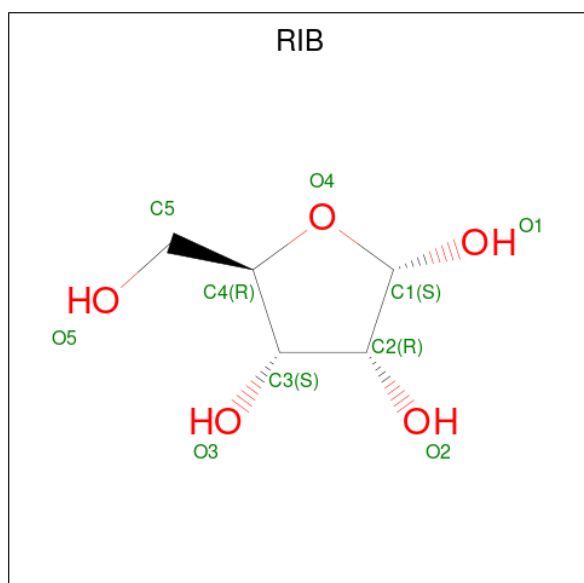
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Chain	Residue	Modelled	Actual	Comment	Reference
F	-6	GLN	-	expression tag	UNP B6THD4
F	-5	ASP	-	expression tag	UNP B6THD4
F	-4	PRO	-	expression tag	UNP B6THD4
F	-3	ASN	-	expression tag	UNP B6THD4
F	-2	SER	-	expression tag	UNP B6THD4
F	-1	SER	-	expression tag	UNP B6THD4
F	0	SER	-	expression tag	UNP B6THD4

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

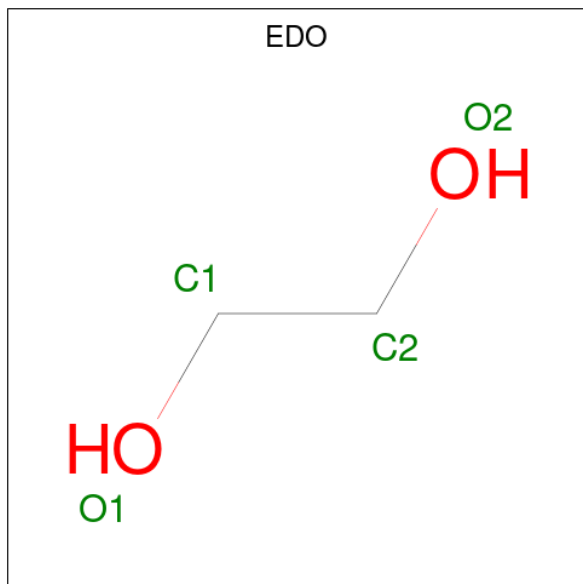
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Ca 1 1	0	0
2	B	1	Total Ca 1 1	0	0
2	C	1	Total Ca 1 1	0	0
2	D	1	Total Ca 1 1	0	0
2	E	1	Total Ca 1 1	0	0
2	F	1	Total Ca 1 1	0	0

- Molecule 3 is alpha-D-ribofuranose (three-letter code: RIB) (formula: C<sub>5</sub>H<sub>10</sub>O<sub>5</sub>) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



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

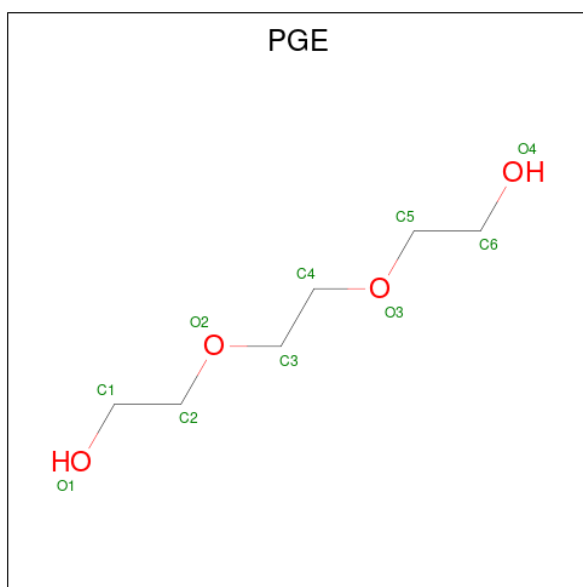
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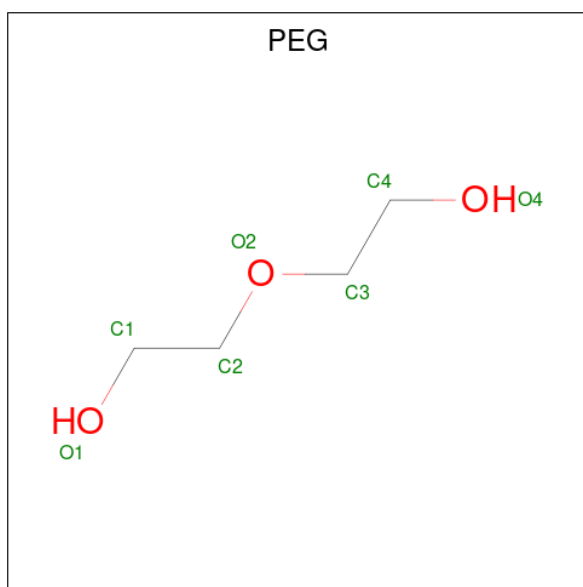
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	E	1	Total	C	O	0	0
			4	2	2		
4	E	1	Total	C	O	0	0
			4	2	2		
4	E	1	Total	C	O	0	0
			4	2	2		
4	E	1	Total	C	O	0	0
			4	2	2		
4	E	1	Total	C	O	0	0
			4	2	2		
4	F	1	Total	C	O	0	0
			4	2	2		
4	F	1	Total	C	O	0	0
			4	2	2		
4	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		
5	A	1	Total	C	O	0	0
			10	6	4		
5	B	1	Total	C	O	0	0
			10	6	4		
5	B	1	Total	C	O	0	0
			10	6	4		
5	B	1	Total	C	O	0	0
			10	6	4		
5	B	1	Total	C	O	0	0
			10	6	4		
5	C	1	Total	C	O	0	0
			10	6	4		
5	C	1	Total	C	O	0	0
			10	6	4		
5	E	1	Total	C	O	0	0
			10	6	4		
5	F	1	Total	C	O	0	0
			10	6	4		
5	F	1	Total	C	O	0	0
			10	6	4		
5	F	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	C	1	Total	C	O	0	0
			7	4	3		
6	C	1	Total	C	O	0	0
			7	4	3		
6	D	1	Total	C	O	0	0
			7	4	3		
6	D	1	Total	C	O	0	0
			7	4	3		
6	D	1	Total	C	O	0	0
			7	4	3		
6	D	1	Total	C	O	0	0
			7	4	3		
6	E	1	Total	C	O	0	0
			7	4	3		
6	E	1	Total	C	O	0	0
			7	4	3		
6	E	1	Total	C	O	0	0
			7	4	3		
6	E	1	Total	C	O	0	0
			7	4	3		
6	F	1	Total	C	O	0	0
			7	4	3		
6	F	1	Total	C	O	0	0
			7	4	3		


- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	297	Total 297	O 297	0	0
7	B	281	Total 281	O 281	0	0
7	C	339	Total 339	O 339	0	0
7	D	290	Total 290	O 290	0	0
7	E	267	Total 267	O 267	0	0
7	F	258	Total 258	O 258	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.

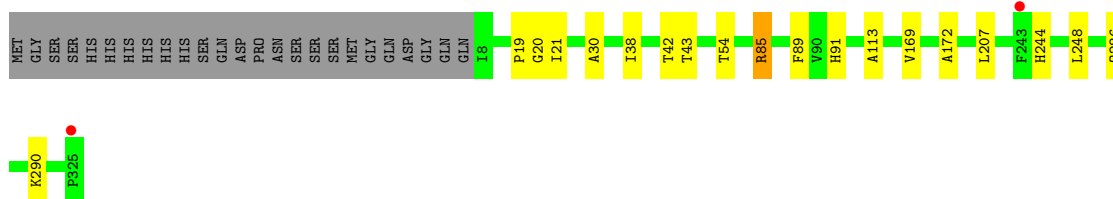
- Molecule 1: Pyrimidine-specific ribonucleoside hydrolase rihA

Chain A: 



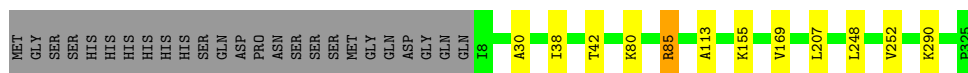
- Molecule 1: Pyrimidine-specific ribonucleoside hydrolase rihA

Chain B: 




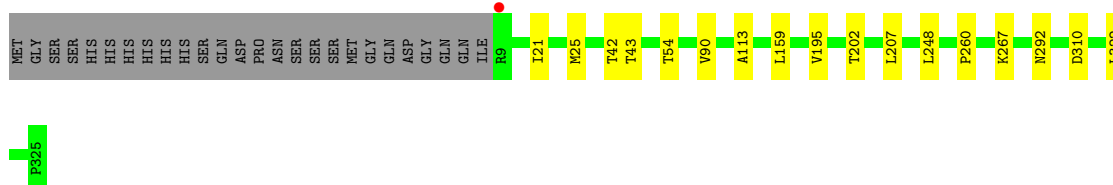
- Molecule 1: Pyrimidine-specific ribonucleoside hydrolase rihA

Chain C: 



- Molecule 1: Pyrimidine-specific ribonucleoside hydrolase rihA

Chain D: 



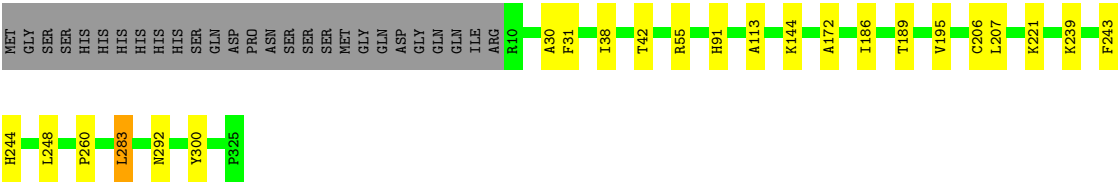
- Molecule 1: Pyrimidine-specific ribonucleoside hydrolase rihA

Chain E: 

85%

6%

8%



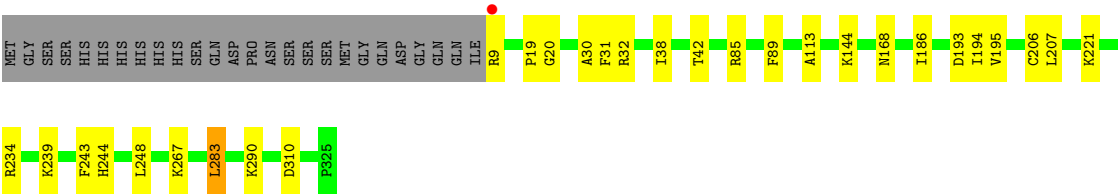
● Molecule 1: Pyrimidine-specific ribonucleoside hydrolase rihA

Chain F: 

84%

8%

8%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.81Å 86.71Å 131.91Å 90.00° 99.44° 90.00°	Depositor
Resolution (Å)	46.50 – 1.75 48.24 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.3 (46.50-1.75) 99.4 (48.24-1.75)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.81 (at 1.75Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, $R_{free}$	0.161 , 0.181 0.164 , 0.184	Depositor DCC
$R_{free}$ test set	10610 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	21.8	Xtriage
Anisotropy	0.272	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	(Not available) , (Not available)	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	16641	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.97% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: EDO, CA, PGE, PEG, RIB

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	0.49	0/2492	0.59	0/3383
1	B	0.49	0/2485	0.57	0/3374
1	C	0.51	0/2478	0.58	0/3364
1	D	0.49	0/2491	0.58	0/3382
1	E	0.50	0/2465	0.59	0/3348
1	F	0.50	0/2491	0.58	0/3381
All	All	0.50	0/14902	0.58	0/20232

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2435	0	2448	9	0
1	B	2428	0	2447	16	0
1	C	2421	0	2439	10	0
1	D	2428	0	2450	11	0
1	E	2405	0	2420	20	0
1	F	2428	0	2450	27	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	10	0	0	0	0
3	B	10	0	0	0	0
3	C	10	0	0	0	0
3	D	10	0	0	0	0
3	E	10	0	0	0	0
3	F	10	0	0	0	0
4	A	12	0	18	0	0
4	B	4	0	6	0	0
4	C	12	0	18	3	0
4	D	20	0	30	3	0
4	E	20	0	30	5	0
4	F	12	0	18	7	0
5	A	20	0	28	1	0
5	B	40	0	56	8	0
5	C	20	0	28	3	0
5	E	10	0	14	6	0
5	F	30	0	42	4	0
6	A	7	0	10	0	0
6	B	7	0	10	0	0
6	C	14	0	20	3	0
6	D	28	0	40	3	0
6	E	28	0	40	6	0
6	F	14	0	20	5	0
7	A	297	0	0	0	0
7	B	281	0	0	1	0
7	C	339	0	0	4	0
7	D	290	0	0	2	0
7	E	267	0	0	2	0
7	F	258	0	0	0	0
All	All	16641	0	15082	97	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

The worst 5 of 97 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:267:LYS:HZ2	4:F:404:EDO:H22	1.18	1.07
1:F:267:LYS:NZ	4:F:404:EDO:H22	1.79	0.96
1:B:85:ARG:HH12	5:B:407:PGE:H2	1.41	0.83
6:C:409:PEG:H31	7:C:637:HOH:O	1.80	0.82
1:F:194:ILE:H	4:F:404:EDO:H12	1.47	0.80

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	318/343 (93%)	310 (98%)	7 (2%)	1 (0%)	41	22
1	B	317/343 (92%)	311 (98%)	6 (2%)	0	100	100
1	C	316/343 (92%)	309 (98%)	7 (2%)	0	100	100
1	D	318/343 (93%)	310 (98%)	8 (2%)	0	100	100
1	E	315/343 (92%)	309 (98%)	6 (2%)	0	100	100
1	F	318/343 (93%)	312 (98%)	6 (2%)	0	100	100
All	All	1902/2058 (92%)	1861 (98%)	40 (2%)	1 (0%)	51	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	154	LYS

### 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	262/283 (93%)	261 (100%)	1 (0%)	91	87
1	B	262/283 (93%)	260 (99%)	2 (1%)	81	72
1	C	261/283 (92%)	260 (100%)	1 (0%)	91	87
1	D	263/283 (93%)	262 (100%)	1 (0%)	91	87
1	E	260/283 (92%)	256 (98%)	4 (2%)	65	49
1	F	263/283 (93%)	258 (98%)	5 (2%)	57	37
All	All	1571/1698 (92%)	1557 (99%)	14 (1%)	78	67

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	244	HIS
1	E	283	LEU
1	F	283	LEU
1	F	234	ARG
1	F	244	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 58 ligands modelled in this entry, 6 are monoatomic - leaving 52 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
5	PGE	C	406	-	9,9,9	0.11	0	8,8,8	0.33	0
3	RIB	F	402	2	10,10,10	2.23	5 (50%)	13,14,14	1.58	2 (15%)
6	PEG	E	412	-	6,6,6	0.12	0	5,5,5	0.15	0
3	RIB	D	402	2	10,10,10	2.29	5 (50%)	13,14,14	1.53	3 (23%)
4	EDO	E	405	-	3,3,3	0.66	0	2,2,2	0.25	0
5	PGE	C	407	-	9,9,9	0.20	0	8,8,8	0.30	0
4	EDO	F	405	-	3,3,3	0.57	0	2,2,2	0.29	0
3	RIB	A	402	2	10,10,10	2.35	5 (50%)	13,14,14	1.55	3 (23%)
6	PEG	E	410	-	6,6,6	0.19	0	5,5,5	0.22	0
4	EDO	E	404	-	3,3,3	0.39	0	2,2,2	0.26	0
6	PEG	F	410	-	6,6,6	0.15	0	5,5,5	0.18	0
6	PEG	C	409	-	6,6,6	0.11	0	5,5,5	0.13	0
6	PEG	D	410	-	6,6,6	0.15	0	5,5,5	0.15	0
5	PGE	A	406	-	9,9,9	0.10	0	8,8,8	0.24	0
6	PEG	C	408	-	6,6,6	0.06	0	5,5,5	0.11	0
4	EDO	D	407	-	3,3,3	0.52	0	2,2,2	0.26	0
6	PEG	B	408	-	6,6,6	0.11	0	5,5,5	0.07	0
6	PEG	D	411	-	6,6,6	0.11	0	5,5,5	0.07	0
4	EDO	B	403	-	3,3,3	0.53	0	2,2,2	0.26	0
5	PGE	B	404	-	9,9,9	0.06	0	8,8,8	0.19	0
5	PGE	B	405	-	9,9,9	0.13	0	8,8,8	0.16	0
5	PGE	F	408	-	9,9,9	0.15	0	8,8,8	0.11	0
3	RIB	B	402	2	10,10,10	2.27	6 (60%)	13,14,14	1.64	2 (15%)
6	PEG	D	409	-	6,6,6	0.12	0	5,5,5	0.07	0
6	PEG	E	411	-	6,6,6	0.14	0	5,5,5	0.25	0
4	EDO	D	403	-	3,3,3	0.27	0	2,2,2	0.74	0
4	EDO	A	403	-	3,3,3	0.57	0	2,2,2	0.34	0
3	RIB	E	402	2	10,10,10	2.26	4 (40%)	13,14,14	1.48	2 (15%)
4	EDO	E	403	-	3,3,3	0.55	0	2,2,2	0.29	0
5	PGE	E	408	-	9,9,9	0.19	0	8,8,8	0.27	0
5	PGE	B	406	-	9,9,9	0.15	0	8,8,8	0.23	0
4	EDO	F	404	-	3,3,3	0.36	0	2,2,2	0.17	0
6	PEG	F	409	-	6,6,6	0.21	0	5,5,5	0.18	0
5	PGE	A	407	-	9,9,9	0.10	0	8,8,8	0.19	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	PGE	B	407	-	9,9,9	0.21	0	8,8,8	0.36	0
4	EDO	C	405	-	3,3,3	0.47	0	2,2,2	0.15	0
4	EDO	D	405	-	3,3,3	0.54	0	2,2,2	0.34	0
4	EDO	C	404	-	3,3,3	0.38	0	2,2,2	0.65	0
4	EDO	D	404	-	3,3,3	0.61	0	2,2,2	0.32	0
5	PGE	F	406	-	9,9,9	0.19	0	8,8,8	0.38	0
4	EDO	A	405	-	3,3,3	0.47	0	2,2,2	0.53	0
4	EDO	F	403	-	3,3,3	0.51	0	2,2,2	0.34	0
6	PEG	D	408	-	6,6,6	0.16	0	5,5,5	0.06	0
4	EDO	C	403	-	3,3,3	0.48	0	2,2,2	0.44	0
4	EDO	E	406	-	3,3,3	0.64	0	2,2,2	0.28	0
3	RIB	C	402	2	10,10,10	2.16	6 (60%)	13,14,14	1.62	2 (15%)
6	PEG	A	408	-	6,6,6	0.11	0	5,5,5	0.07	0
6	PEG	E	409	-	6,6,6	0.15	0	5,5,5	0.22	0
4	EDO	A	404	-	3,3,3	0.59	0	2,2,2	0.17	0
4	EDO	D	406	-	3,3,3	0.64	0	2,2,2	0.35	0
4	EDO	E	407	-	3,3,3	0.49	0	2,2,2	0.38	0
5	PGE	F	407	-	9,9,9	0.07	0	8,8,8	0.22	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PGE	C	406	-	-	4/7/7/7	-
3	RIB	F	402	2	-	0/2/18/18	0/1/1/1
6	PEG	E	412	-	-	2/4/4/4	-
3	RIB	D	402	2	-	0/2/18/18	0/1/1/1
4	EDO	E	405	-	-	1/1/1/1	-
5	PGE	C	407	-	-	6/7/7/7	-
4	EDO	F	405	-	-	0/1/1/1	-
3	RIB	A	402	2	-	0/2/18/18	0/1/1/1
6	PEG	E	410	-	-	1/4/4/4	-
4	EDO	E	404	-	-	1/1/1/1	-
6	PEG	F	410	-	-	2/4/4/4	-
6	PEG	C	409	-	-	2/4/4/4	-
6	PEG	D	410	-	-	2/4/4/4	-
5	PGE	A	406	-	-	0/7/7/7	-
6	PEG	C	408	-	-	3/4/4/4	-
4	EDO	D	407	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	PEG	B	408	-	-	1/4/4/4	-
6	PEG	D	411	-	-	3/4/4/4	-
4	EDO	B	403	-	-	1/1/1/1	-
5	PGE	B	404	-	-	3/7/7/7	-
5	PGE	B	405	-	-	0/7/7/7	-
5	PGE	F	408	-	-	6/7/7/7	-
3	RIB	B	402	2	-	0/2/18/18	0/1/1/1
6	PEG	D	409	-	-	2/4/4/4	-
6	PEG	E	411	-	-	4/4/4/4	-
4	EDO	D	403	-	-	0/1/1/1	-
4	EDO	A	403	-	-	0/1/1/1	-
3	RIB	E	402	2	-	0/2/18/18	0/1/1/1
4	EDO	E	403	-	-	1/1/1/1	-
5	PGE	E	408	-	-	3/7/7/7	-
5	PGE	B	406	-	-	5/7/7/7	-
4	EDO	F	404	-	-	0/1/1/1	-
6	PEG	F	409	-	-	1/4/4/4	-
5	PGE	A	407	-	-	1/7/7/7	-
5	PGE	B	407	-	-	4/7/7/7	-
4	EDO	C	405	-	-	0/1/1/1	-
4	EDO	D	405	-	-	0/1/1/1	-
4	EDO	C	404	-	-	0/1/1/1	-
4	EDO	D	404	-	-	1/1/1/1	-
5	PGE	F	406	-	-	3/7/7/7	-
4	EDO	A	405	-	-	0/1/1/1	-
4	EDO	F	403	-	-	0/1/1/1	-
6	PEG	D	408	-	-	0/4/4/4	-
4	EDO	C	403	-	-	1/1/1/1	-
4	EDO	E	406	-	-	0/1/1/1	-
3	RIB	C	402	2	-	0/2/18/18	0/1/1/1
6	PEG	A	408	-	-	2/4/4/4	-
6	PEG	E	409	-	-	1/4/4/4	-
4	EDO	A	404	-	-	0/1/1/1	-
4	EDO	D	406	-	-	0/1/1/1	-
4	EDO	E	407	-	-	0/1/1/1	-
5	PGE	F	407	-	-	3/7/7/7	-

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	402	RIB	O4-C1	4.20	1.48	1.43
3	E	402	RIB	O4-C1	3.94	1.48	1.43
3	A	402	RIB	O4-C1	3.86	1.47	1.43
3	A	402	RIB	O4-C4	-3.58	1.37	1.45
3	F	402	RIB	O4-C1	3.53	1.47	1.43

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	402	RIB	O4-C1-C2	-4.07	99.46	104.46
3	B	402	RIB	O4-C1-C2	-3.77	99.82	104.46
3	C	402	RIB	O4-C1-C2	-3.76	99.84	104.46
3	A	402	RIB	O4-C1-C2	-3.59	100.05	104.46
3	D	402	RIB	O4-C1-C2	-3.59	100.05	104.46

There are no chirality outliers.

5 of 71 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	407	PGE	O2-C3-C4-O3
6	C	408	PEG	O1-C1-C2-O2
5	B	404	PGE	O2-C3-C4-O3
6	E	411	PEG	O2-C3-C4-O4
6	E	410	PEG	O1-C1-C2-O2

There are no ring outliers.

23 monomers are involved in 57 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	406	PGE	1	0
6	E	412	PEG	3	0
5	C	407	PGE	2	0
6	E	410	PEG	2	0
4	E	404	EDO	4	0
6	F	410	PEG	1	0
6	C	409	PEG	2	0
6	D	410	PEG	2	0
6	C	408	PEG	1	0
5	F	408	PGE	1	0
6	E	411	PEG	1	0
4	D	403	EDO	3	0
5	E	408	PGE	6	0

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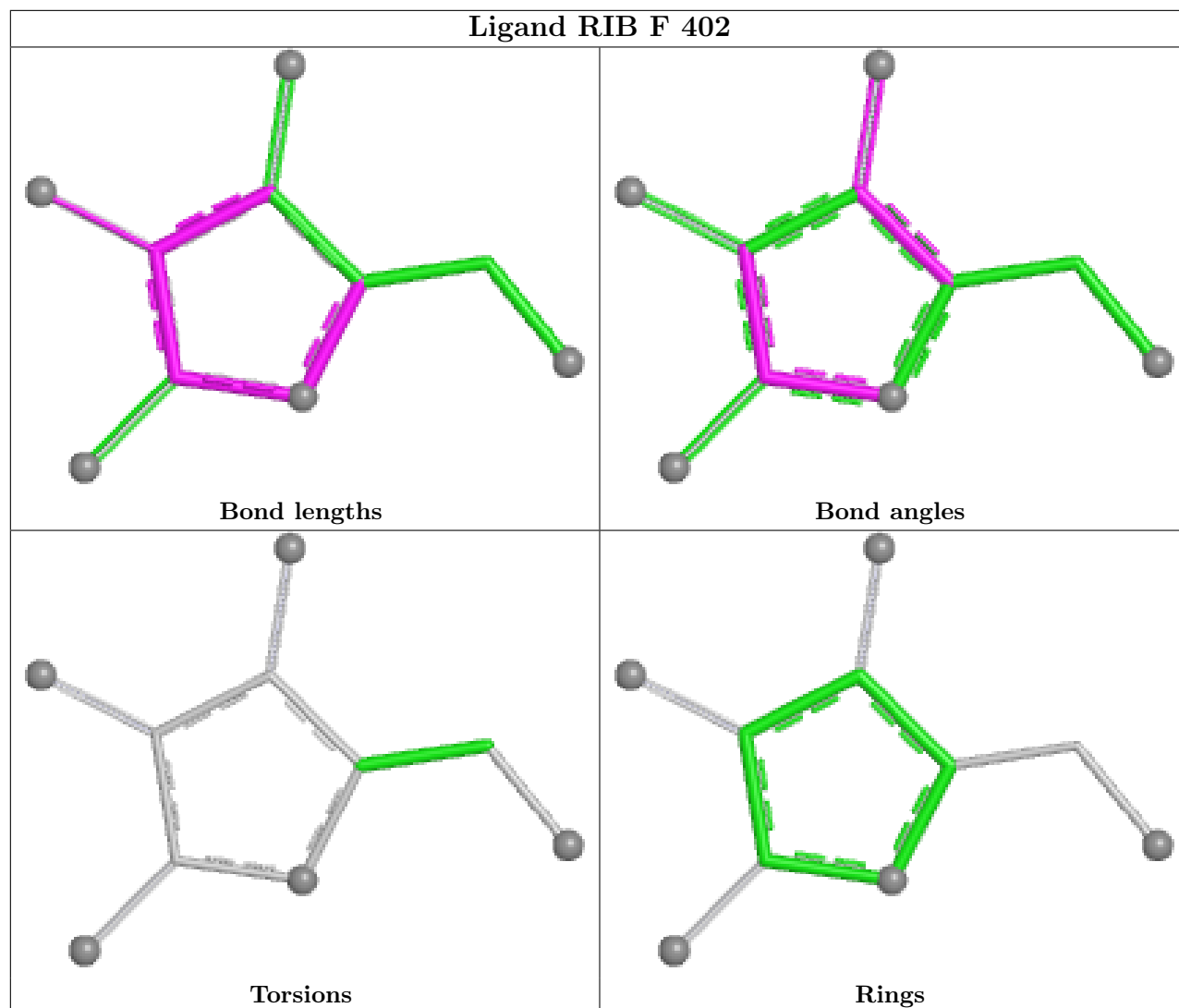
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	406	PGE	1	0
4	F	404	EDO	7	0
6	F	409	PEG	4	0
5	A	407	PGE	1	0
5	B	407	PGE	7	0
4	C	405	EDO	3	0
5	F	406	PGE	2	0
6	D	408	PEG	1	0
4	E	406	EDO	1	0
5	F	407	PGE	1	0

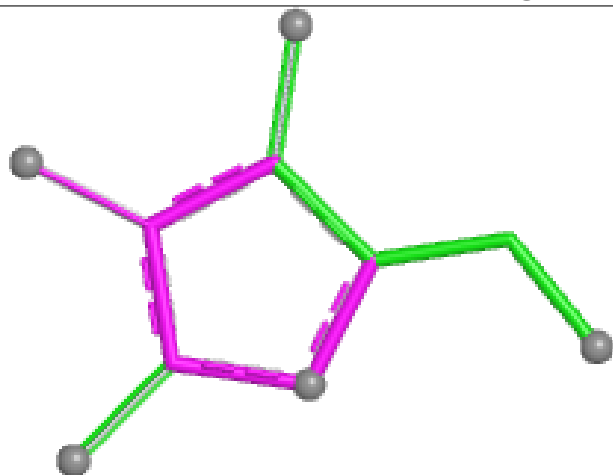
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



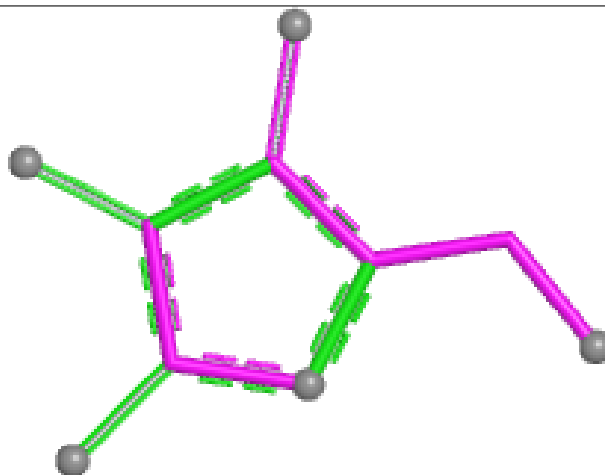
## Ligand RIB F 402



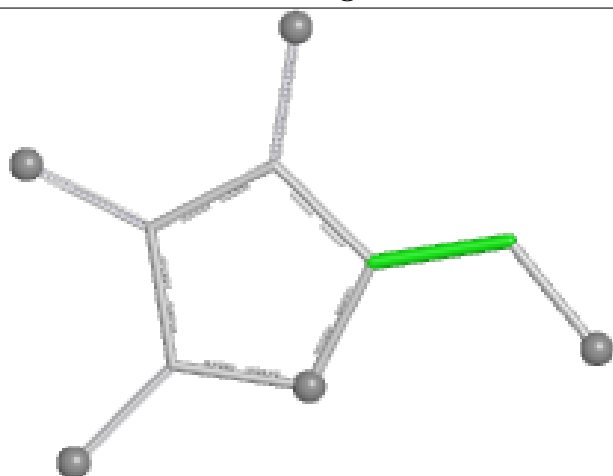
## Ligand RIB D 402



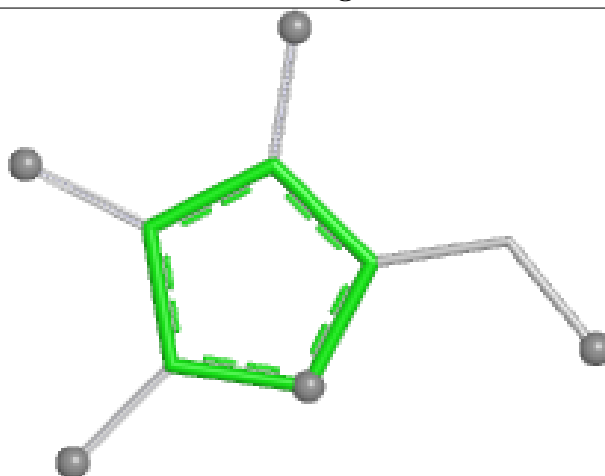
Bond lengths



Bond angles

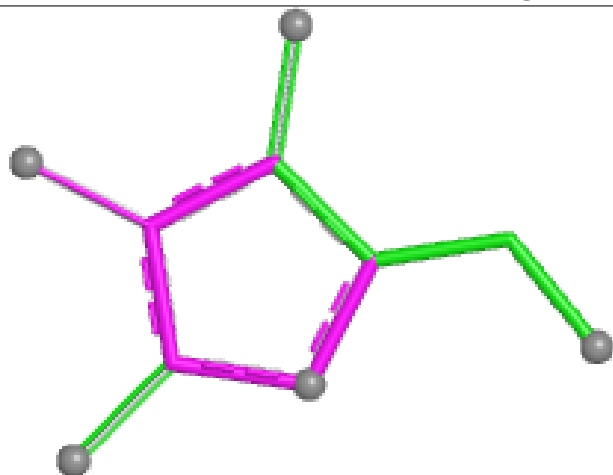


Torsions

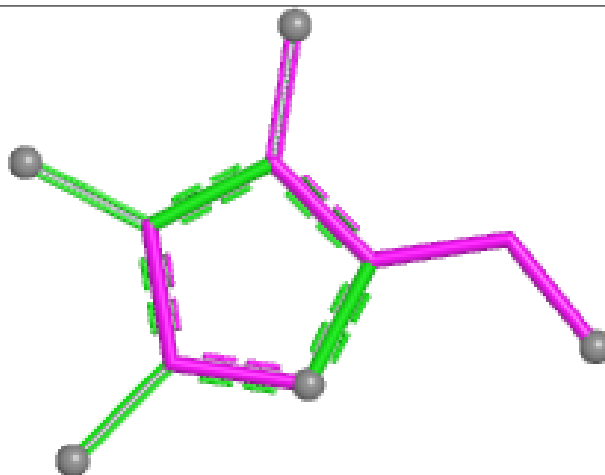


Rings

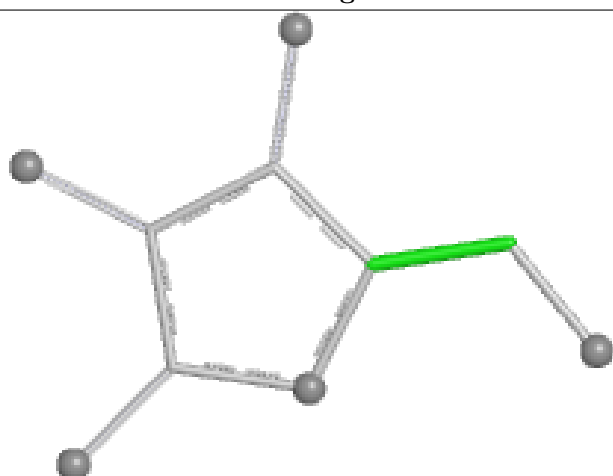
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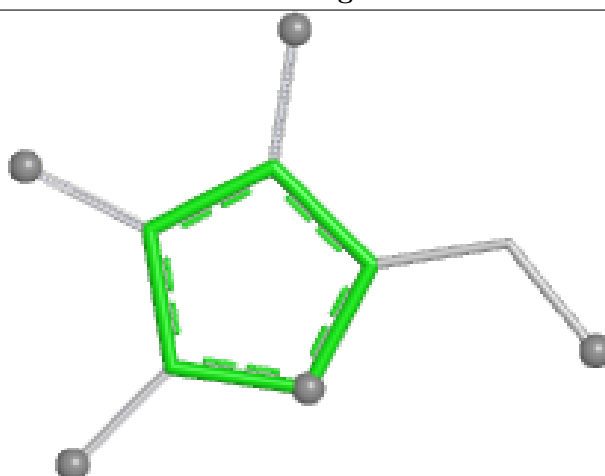
Bond lengths



Bond angles

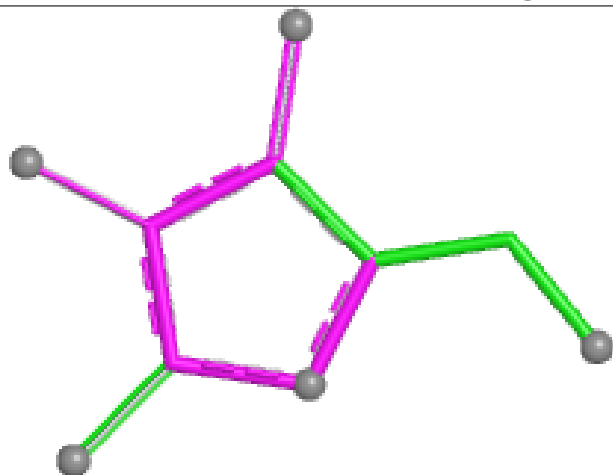


Torsions

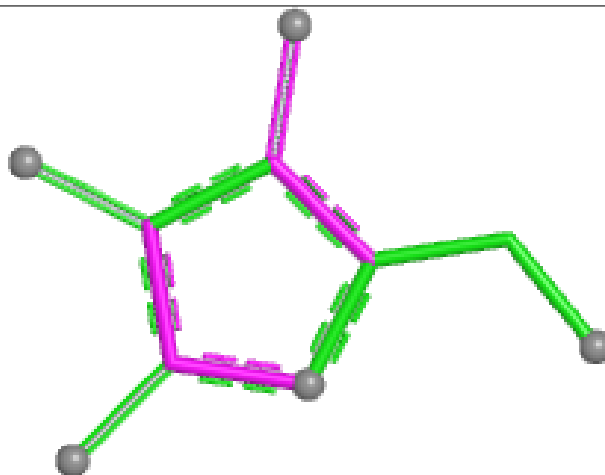


Rings

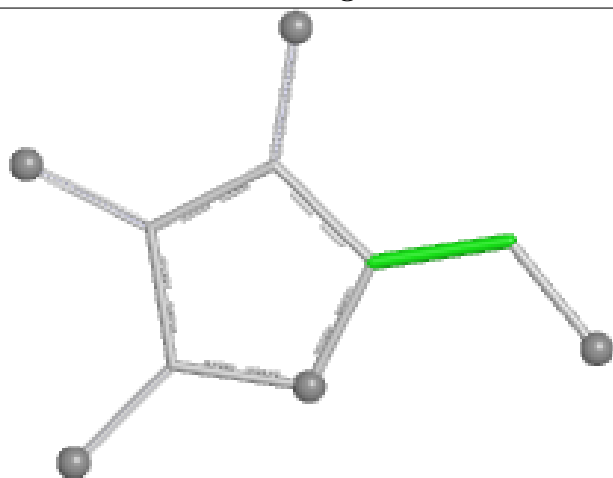
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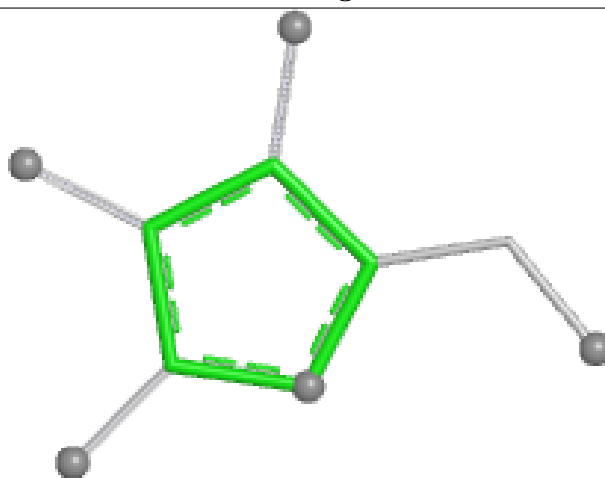
Bond lengths



Bond angles

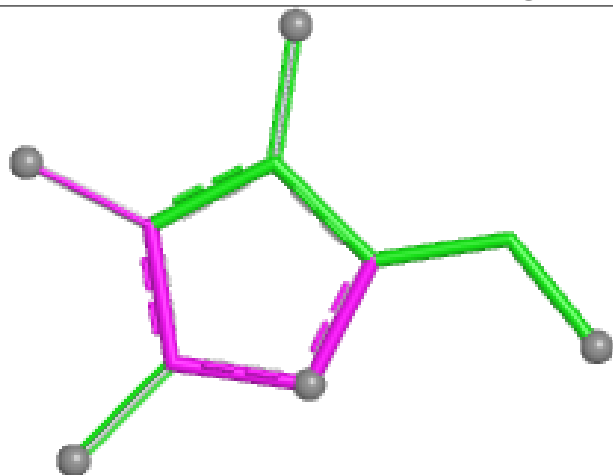


Torsions

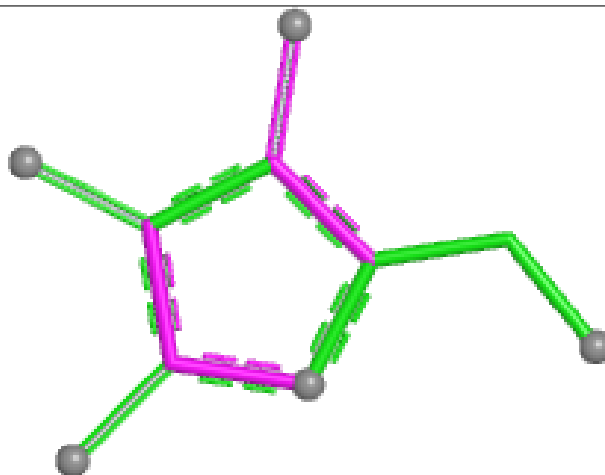


Rings

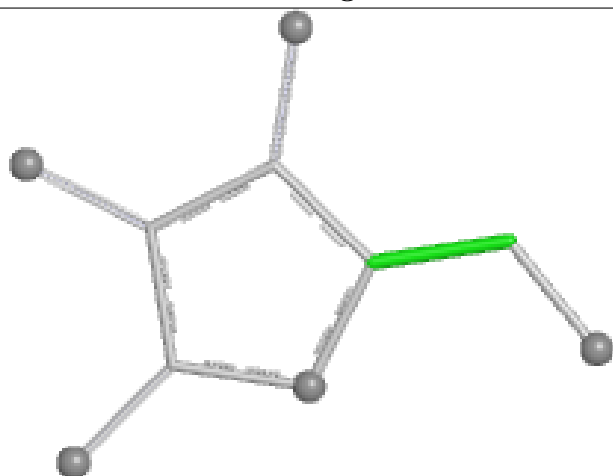
## Ligand RIB E 402



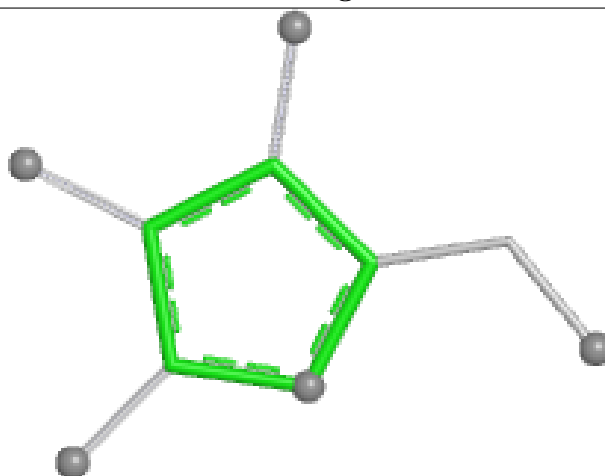
Bond lengths



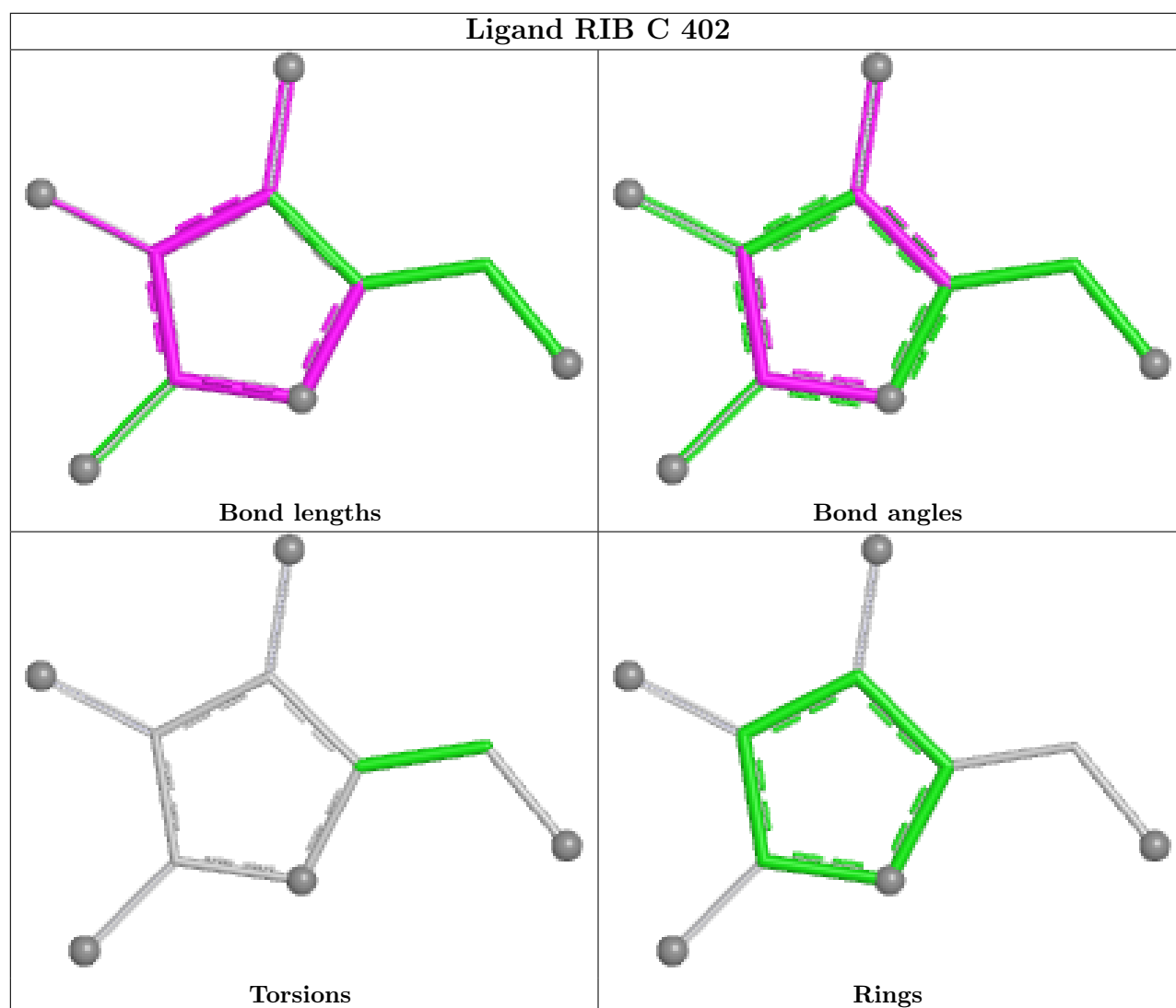
Bond angles



Torsions



Rings



## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	319/343 (93%)	-0.41	1 (0%) 94 95	16, 23, 40, 56	0
1	B	318/343 (92%)	-0.18	2 (0%) 89 92	16, 24, 42, 58	0
1	C	318/343 (92%)	-0.46	0 100 100	15, 21, 38, 55	0
1	D	317/343 (92%)	-0.49	1 (0%) 94 95	16, 23, 38, 73	0
1	E	316/343 (92%)	-0.30	0 100 100	16, 25, 44, 56	0
1	F	317/343 (92%)	-0.24	1 (0%) 94 95	17, 26, 47, 80	0
All	All	1905/2058 (92%)	-0.35	5 (0%) 94 95	15, 23, 42, 80	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	9	ARG	4.7
1	B	243	PHE	4.0
1	D	9	ARG	3.7
1	A	243	PHE	3.1
1	B	325	PRO	2.5

### 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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
5	PGE	E	408	10/10	0.65	0.24	42,48,57,58	0
4	EDO	E	406	4/4	0.67	0.20	41,42,43,45	0
6	PEG	E	412	7/7	0.67	0.25	31,37,58,62	0
4	EDO	F	404	4/4	0.71	0.33	41,44,44,47	0
4	EDO	A	404	4/4	0.76	0.13	50,55,58,61	0
5	PGE	F	406	10/10	0.77	0.19	49,51,60,62	0
6	PEG	D	409	7/7	0.77	0.15	70,70,71,72	0
5	PGE	B	407	10/10	0.77	0.22	37,42,46,46	0
5	PGE	C	407	10/10	0.78	0.18	41,50,62,63	0
5	PGE	F	408	10/10	0.79	0.16	60,62,66,67	0
6	PEG	A	408	7/7	0.79	0.16	55,57,60,61	0
4	EDO	D	404	4/4	0.79	0.13	62,62,62,62	0
5	PGE	B	406	10/10	0.79	0.24	49,54,62,63	0
6	PEG	F	409	7/7	0.79	0.23	34,37,49,50	0
4	EDO	F	405	4/4	0.80	0.12	53,53,54,54	0
6	PEG	D	410	7/7	0.81	0.21	54,54,57,59	0
6	PEG	D	411	7/7	0.81	0.12	60,61,62,63	0
6	PEG	E	409	7/7	0.82	0.15	44,45,48,51	0
6	PEG	E	411	7/7	0.82	0.19	46,48,51,53	0
4	EDO	A	405	4/4	0.82	0.25	50,52,54,54	0
4	EDO	D	403	4/4	0.82	0.29	44,46,48,48	0
4	EDO	E	405	4/4	0.83	0.22	45,45,48,53	0
4	EDO	A	403	4/4	0.83	0.10	62,62,63,63	0
4	EDO	C	405	4/4	0.83	0.33	51,52,55,58	0
6	PEG	F	410	7/7	0.83	0.17	53,54,56,57	0
4	EDO	D	405	4/4	0.84	0.18	55,56,57,60	0
6	PEG	E	410	7/7	0.84	0.12	51,52,58,59	0
4	EDO	D	407	4/4	0.86	0.08	61,61,61,62	0
4	EDO	E	404	4/4	0.86	0.28	47,48,48,51	0
4	EDO	C	404	4/4	0.87	0.14	41,42,42,44	0
5	PGE	F	407	10/10	0.88	0.20	46,50,54,55	0
5	PGE	B	404	10/10	0.88	0.17	47,52,61,64	0
4	EDO	E	407	4/4	0.88	0.23	58,58,59,60	0
4	EDO	F	403	4/4	0.88	0.13	48,48,49,51	0
6	PEG	D	408	7/7	0.89	0.13	44,45,48,51	0
5	PGE	C	406	10/10	0.89	0.16	42,48,58,62	0
3	RIB	D	402	10/10	0.90	0.09	18,23,27,27	0

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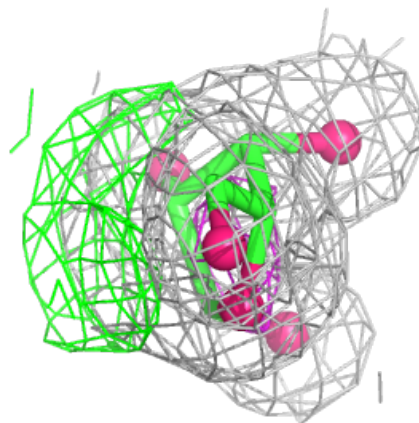
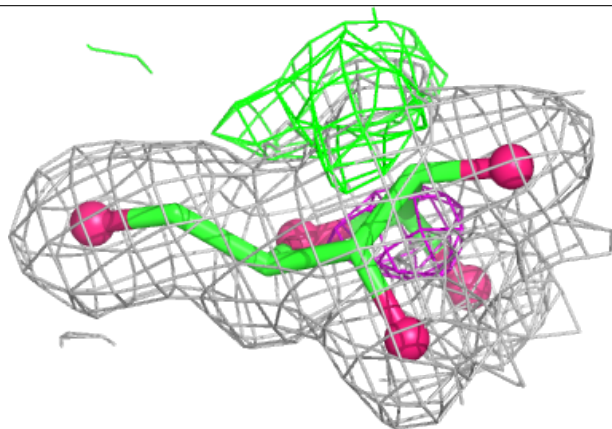
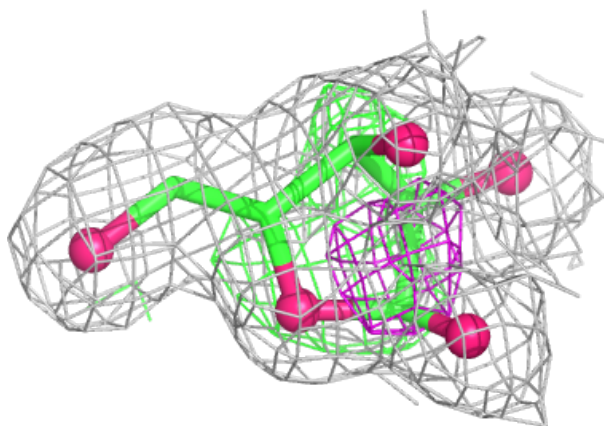
*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	RIB	B	402	10/10	0.90	0.10	20,25,27,28	0
3	RIB	C	402	10/10	0.90	0.13	17,23,26,29	0
6	PEG	B	408	7/7	0.90	0.17	53,55,60,63	0
6	PEG	C	409	7/7	0.90	0.18	51,53,54,54	0
3	RIB	F	402	10/10	0.91	0.09	19,23,27,29	0
4	EDO	E	403	4/4	0.91	0.10	48,49,50,52	0
5	PGE	B	405	10/10	0.91	0.10	44,47,53,54	0
5	PGE	A	406	10/10	0.92	0.15	35,37,44,48	0
5	PGE	A	407	10/10	0.92	0.14	48,51,55,57	0
6	PEG	C	408	7/7	0.92	0.13	44,45,51,55	0
4	EDO	B	403	4/4	0.92	0.07	50,52,52,54	0
4	EDO	C	403	4/4	0.92	0.23	48,48,49,49	0
3	RIB	A	402	10/10	0.92	0.07	18,23,28,28	0
3	RIB	E	402	10/10	0.92	0.09	20,22,26,28	0
4	EDO	D	406	4/4	0.95	0.10	29,37,43,47	0
2	CA	E	401	1/1	1.00	0.06	17,17,17,17	0
2	CA	F	401	1/1	1.00	0.06	17,17,17,17	0
2	CA	A	401	1/1	1.00	0.06	17,17,17,17	0
2	CA	B	401	1/1	1.00	0.06	18,18,18,18	0
2	CA	C	401	1/1	1.00	0.07	17,17,17,17	0
2	CA	D	401	1/1	1.00	0.06	17,17,17,17	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.

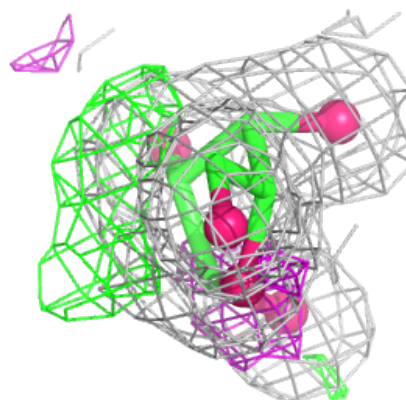
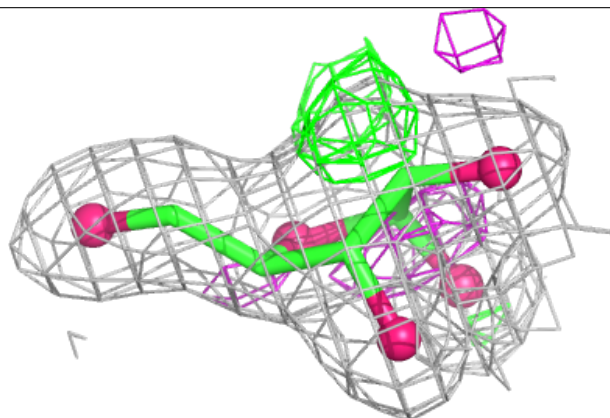
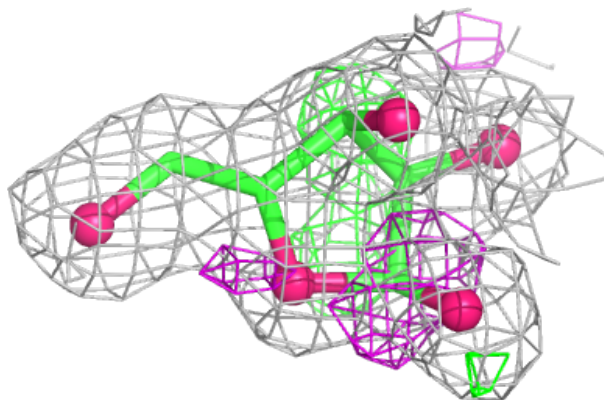
**Electron density around RIB D 402:**

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



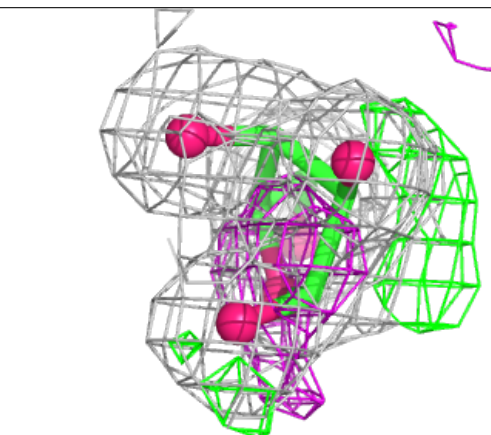
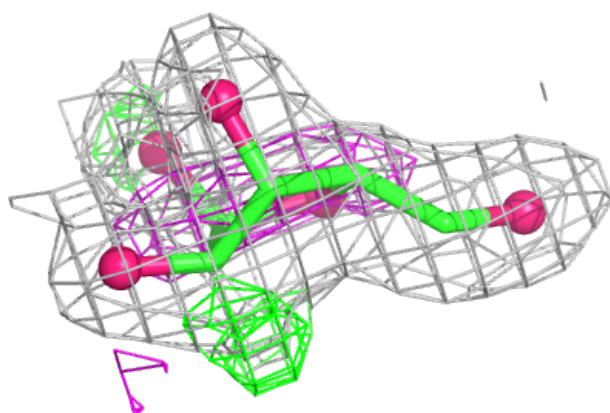
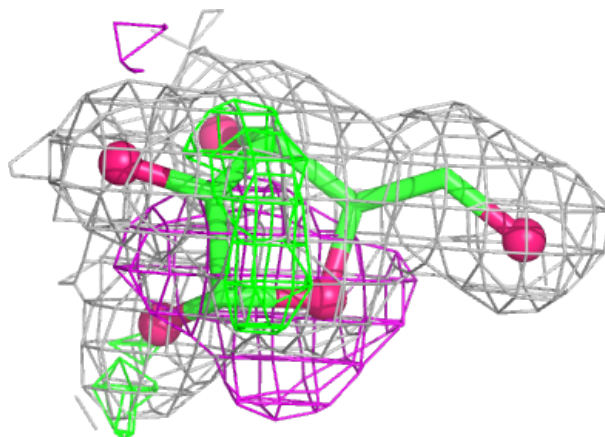
**Electron density around RIB B 402:**

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

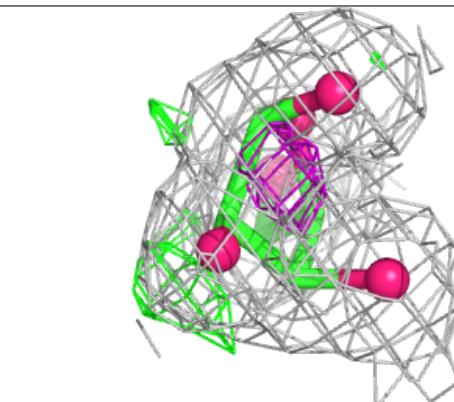
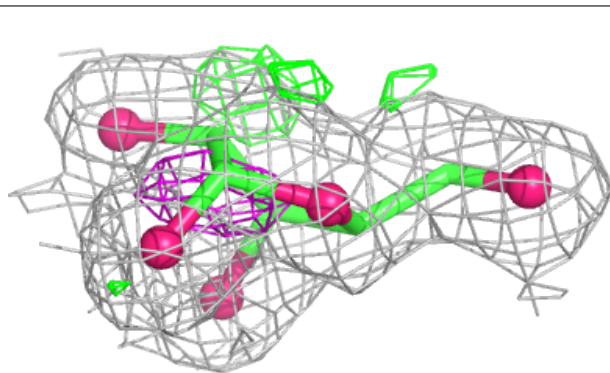
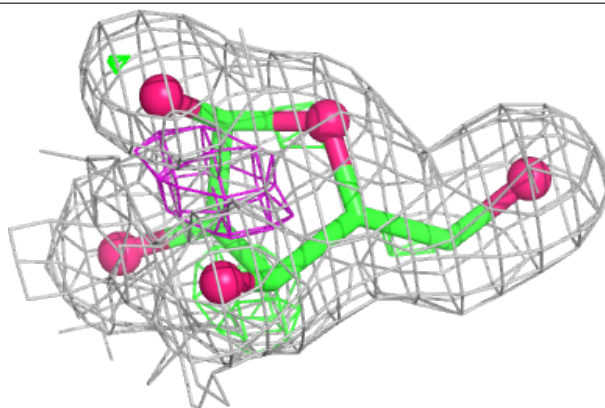


**Electron density around RIB C 402:**

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

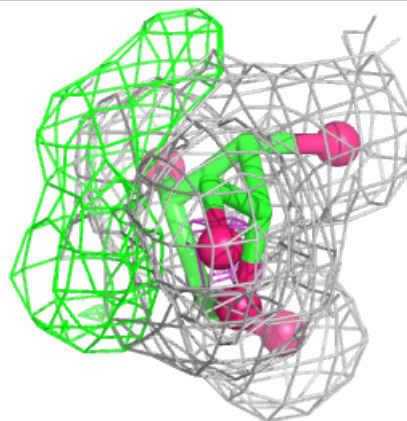
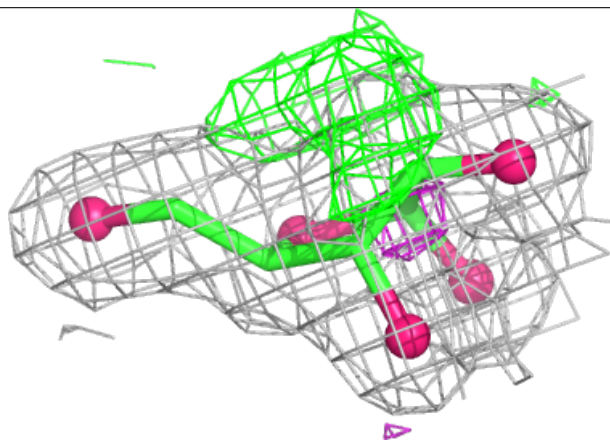
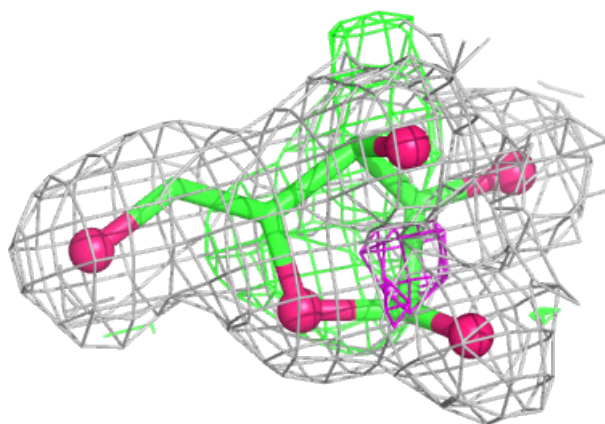
**Electron density around RIB F 402:**

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



**Electron density around RIB A 402:**

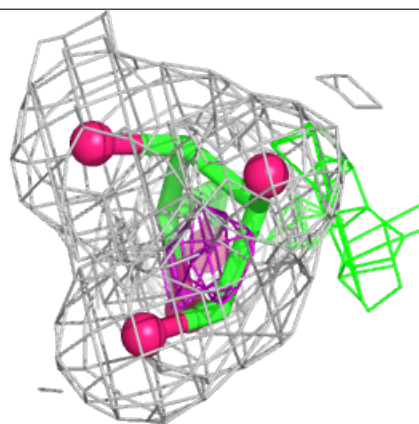
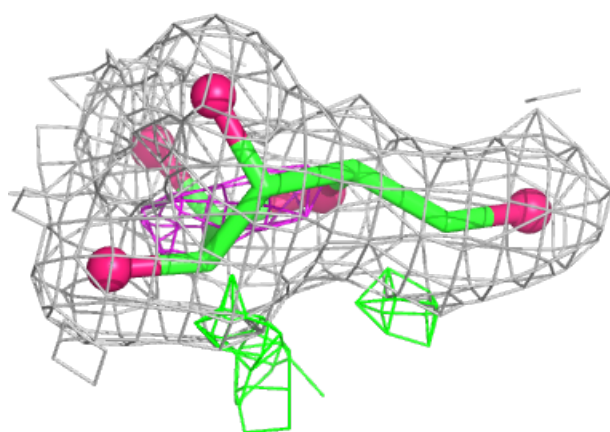
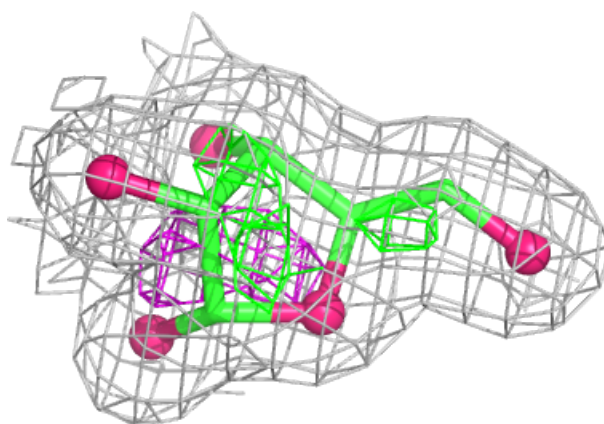
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around RIB E 402:**

$2mF_o-DF_c$  (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.