



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 28, 2023 – 07:10 am BST

PDB ID : 8OIA
Title : Trichomonas vaginalis riboside hydrolase in complex with D-ribose
Authors : Patrone, M.; Stockman, B.J.; Degano, M.
Deposited on : 2023-03-22
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Mogul : 1.8.4, 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

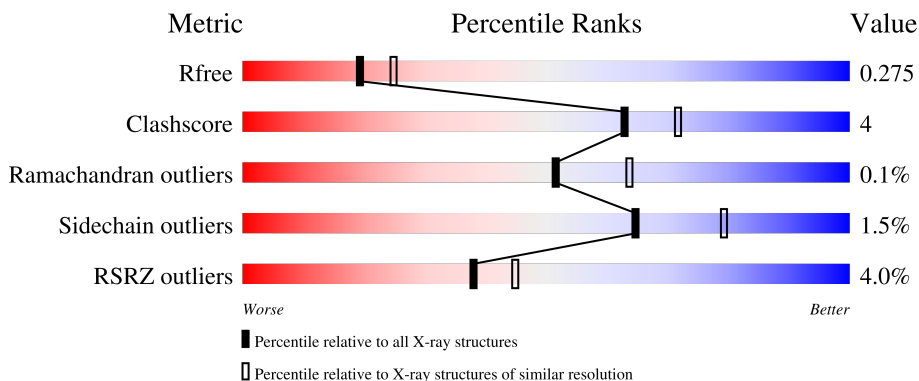
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


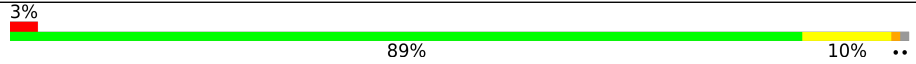
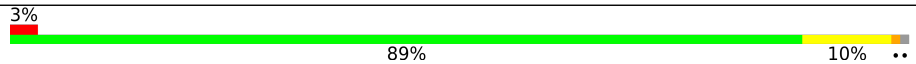

The reported resolution of this entry is 2.30 Å.

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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	347	 4% 85% 10% • 5%
1	B	347	 3% 89% 10% ••
1	C	347	 3% 89% 10% ••
1	D	347	 6% 88% 8% •

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 11006 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.

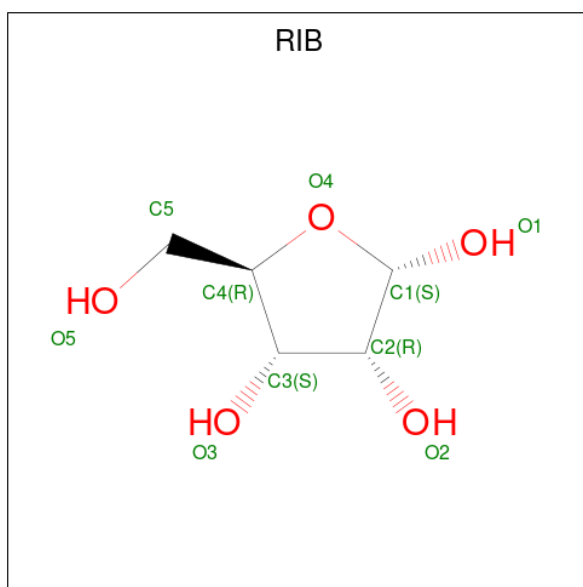
- Molecule 1 is a protein called Inosine-uridine preferring nucleoside hydrolase family protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	331	2654	1704	440	493	17	0	4	0
1	B	345	2761	1772	457	515	17	0	4	0
1	C	345	2735	1754	452	512	17	0	0	0
1	D	335	2676	1718	440	502	16	0	3	0

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

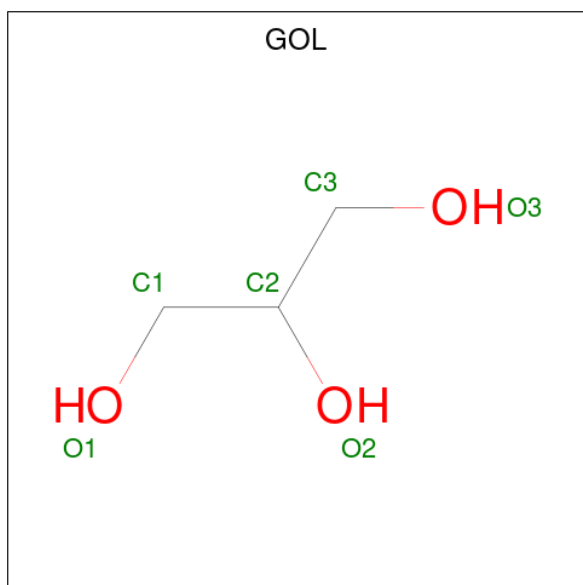
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Ca 1	0	0
2	B	1	Total 1	Ca 1	0	0
2	C	1	Total 1	Ca 1	0	0
2	D	1	Total 1	Ca 1	0	0

- Molecule 3 is alpha-D-ribofuranose (three-letter code: RIB) (formula: C₅H₁₀O₅) (labeled as "Ligand of Interest" by depositor).



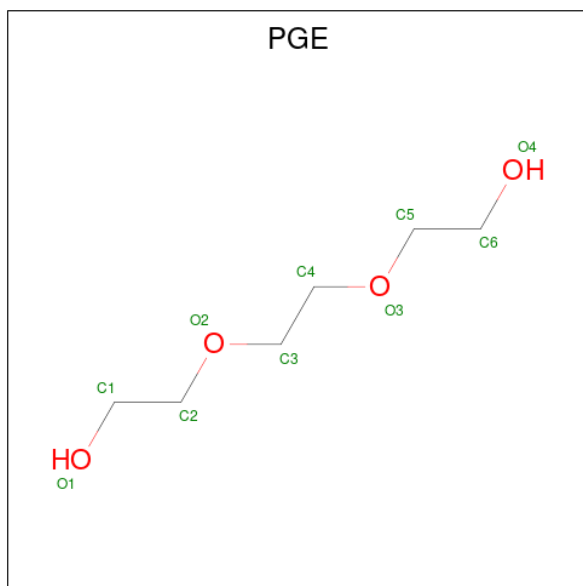
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 10 5 5	0	0
3	B	1	Total C O 10 5 5	0	0
3	C	1	Total C O 10 5 5	0	0
3	D	1	Total C O 10 5 5	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 5 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			10	6	4		

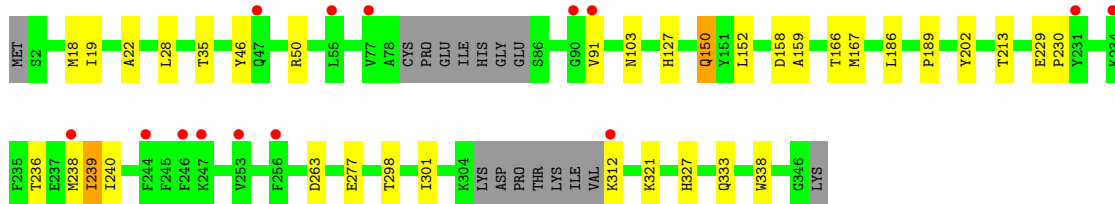
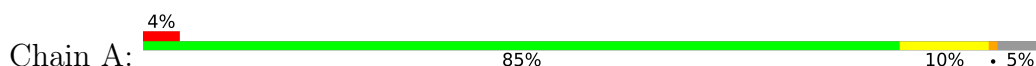
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	24	Total	O	0	0
			24	24		
6	B	28	Total	O	0	0
			28	28		
6	C	29	Total	O	0	0
			29	29		
6	D	27	Total	O	0	0
			27	27		

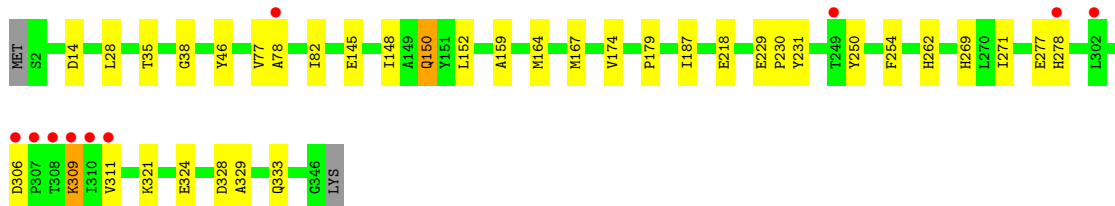
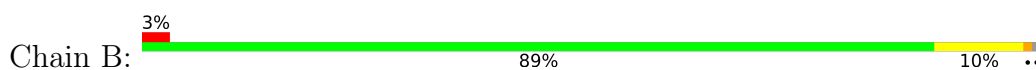
3 Residue-property plots [i](#)

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.

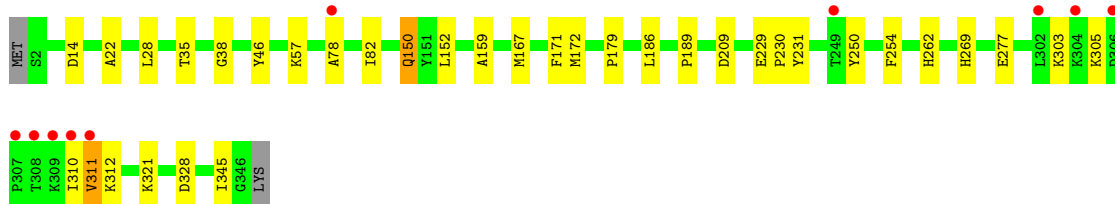
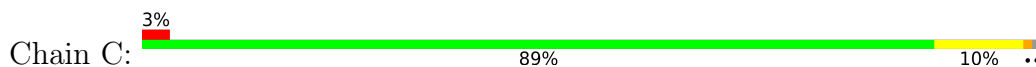
- Molecule 1: Inosine-uridine preferring nucleoside hydrolase family protein



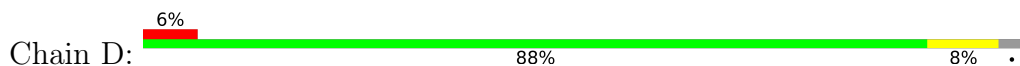
- Molecule 1: Inosine-uridine preferring nucleoside hydrolase family protein

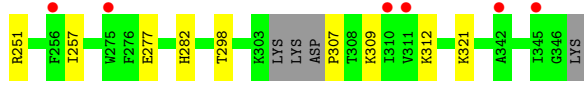
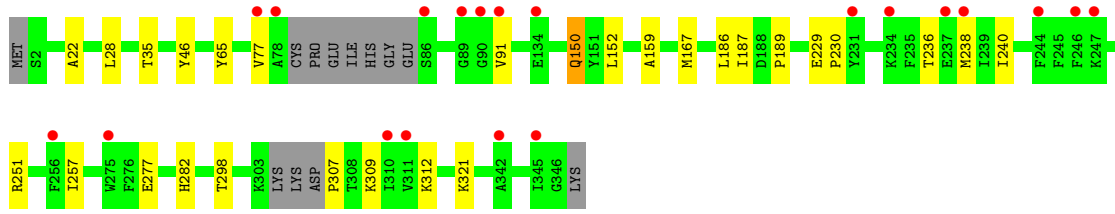


- Molecule 1: Inosine-uridine preferring nucleoside hydrolase family protein



- Molecule 1: Inosine-uridine preferring nucleoside hydrolase family protein





4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	117.00Å 135.62Å 94.08Å 90.00° 104.69° 90.00°	Depositor
Resolution (Å)	43.98 – 2.30 43.98 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (43.98-2.30) 99.9 (43.98-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.30	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 2.20Å)	Xtrriage
Refinement program	REFMAC 5.8.0405	Depositor
R, R_{free}	0.234 , 0.270 0.240 , 0.275	Depositor DCC
R_{free} test set	3666 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	22.8	Xtrriage
Anisotropy	0.602	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 39.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.000 for $-1/2^*h+1/2^*k-1,1/2^*h-1/2^*k-1,-1/2^*h-1/2^*k$ 0.000 for $-1/2^*h-1/2^*k-1,-1/2^*h-1/2^*k+1,-1/2^*h+1/2^*k$	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	11006	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 24.83 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.5586e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: RIB, CA, PGE, GOL

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	0.40	0/2727	0.69	0/3699
1	B	0.42	0/2838	0.69	0/3854
1	C	0.41	0/2802	0.69	0/3805
1	D	0.39	0/2745	0.67	0/3725
All	All	0.41	0/11112	0.68	0/15083

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	2654	0	2629	29	0
1	B	2761	0	2743	32	1
1	C	2735	0	2712	19	1
1	D	2676	0	2653	19	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	10	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	10	0	0	0	0
3	C	10	0	0	0	0
3	D	10	0	0	0	0
4	A	6	0	8	0	0
4	B	6	0	8	0	0
4	C	6	0	8	1	0
5	C	10	0	14	1	0
6	A	24	0	0	0	0
6	B	28	0	0	1	0
6	C	29	0	0	0	0
6	D	27	0	0	1	0
All	All	11006	0	10775	94	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (94) 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:A:239:ILE:HD11	1:A:338:TRP:CE2	1.83	1.13
1:B:148:ILE:HD13	1:B:164:MET:SD	2.04	0.96
1:C:150:GLN:HE21	1:C:150:GLN:HA	1.40	0.87
1:B:77:VAL:HB	1:B:187[B]:ILE:HD13	1.57	0.86
1:B:150:GLN:HA	1:B:150:GLN:HE21	1.39	0.86
1:A:150:GLN:HE21	1:A:150:GLN:HA	1.40	0.85
1:D:150:GLN:HE21	1:D:150:GLN:HA	1.42	0.84
1:B:77:VAL:HB	1:B:187[B]:ILE:CD1	2.13	0.77
1:A:239:ILE:CD1	1:A:338:TRP:CE2	2.65	0.76
1:B:306:ASP:OD2	1:B:309:LYS:HG2	1.88	0.73
1:B:28:LEU:HD13	1:B:271:ILE:HG12	1.73	0.71
1:C:303:LYS:HG2	1:C:310:ILE:HD11	1.75	0.69
1:A:239:ILE:HD11	1:A:338:TRP:CZ2	2.28	0.68
1:D:251:ARG:HH22	1:D:257:ILE:HG23	1.59	0.67
1:A:239:ILE:HD11	1:A:338:TRP:NE1	2.10	0.65
1:B:269:HIS:NE2	1:B:278[B]:HIS:CE1	2.65	0.65
1:C:179:PRO:HG3	1:D:298:THR:O	1.99	0.62
1:B:77:VAL:N	1:B:187[B]:ILE:HD11	2.18	0.59
1:B:324:GLU:HA	1:B:324:GLU:OE1	2.02	0.59
1:A:298:THR:O	1:B:179:PRO:HG3	2.02	0.59
1:B:277:GLU:HG3	1:B:321:LYS:HG3	1.83	0.59
1:B:38:GLY:O	1:B:78:ALA:HA	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:150:GLN:HA	1:D:150:GLN:NE2	2.16	0.58
1:D:77:VAL:HG22	1:D:187:ILE:HD13	1.84	0.58
1:C:38:GLY:O	1:C:78:ALA:HA	2.04	0.57
1:B:77:VAL:H	1:B:187[B]:ILE:HD11	1.71	0.56
1:B:333[B]:GLN:NE2	1:B:333[B]:GLN:HA	2.21	0.55
1:B:329:ALA:O	1:B:333[B]:GLN:HG2	2.05	0.55
1:B:269:HIS:NE2	1:B:278[B]:HIS:HE1	2.05	0.55
1:D:277:GLU:HG3	1:D:321:LYS:HG3	1.88	0.55
1:C:150:GLN:HA	1:C:150:GLN:NE2	2.16	0.54
1:B:150:GLN:HA	1:B:150:GLN:NE2	2.16	0.54
1:B:277:GLU:HG3	1:B:321:LYS:CG	2.36	0.54
1:D:277:GLU:HG3	1:D:321:LYS:CG	2.38	0.54
1:D:307:PRO:HG2	1:D:309:LYS:HB2	1.89	0.54
1:C:171:PHE:HB3	1:C:172:MET:HE2	1.89	0.54
1:B:148:ILE:CD1	1:B:164:MET:SD	2.90	0.53
1:A:301:ILE:HD11	1:B:174:VAL:CG2	2.39	0.53
1:A:229:GLU:N	1:A:230:PRO:CD	2.73	0.52
1:A:239:ILE:CD1	1:A:338:TRP:NE1	2.73	0.52
1:C:303:LYS:CG	1:C:310:ILE:HD11	2.40	0.52
1:D:282:HIS:CG	1:D:312:LYS:HG3	2.45	0.52
1:A:239:ILE:HD12	1:A:338:TRP:CD1	2.45	0.51
1:D:229:GLU:N	1:D:230:PRO:CD	2.73	0.51
1:A:277:GLU:HG3	1:A:321:LYS:CG	2.40	0.51
1:A:150:GLN:HA	1:A:150:GLN:NE2	2.18	0.51
1:A:91:VAL:HG11	1:A:238:MET:CE	2.41	0.51
1:B:229:GLU:N	1:B:230:PRO:CD	2.74	0.51
1:C:229:GLU:N	1:C:230:PRO:CD	2.75	0.50
1:C:328:ASP:OD1	4:C:404:GOL:H12	2.11	0.50
1:B:152:LEU:HD22	1:B:159:ALA:HB2	1.94	0.50
1:A:239:ILE:CD1	1:A:338:TRP:CD2	2.95	0.49
1:D:236:THR:O	1:D:240:ILE:HG12	2.13	0.49
1:D:77:VAL:HG22	1:D:187:ILE:CD1	2.43	0.49
1:C:311:VAL:HG23	1:C:312:LYS:HG3	1.95	0.49
1:D:152:LEU:HD22	1:D:159:ALA:HB2	1.94	0.48
1:D:35:THR:HG21	1:D:46:TYR:HA	1.95	0.48
1:B:148:ILE:HD13	1:B:164:MET:CE	2.44	0.48
1:C:35:THR:HG21	1:C:46:TYR:HA	1.95	0.47
1:A:236:THR:O	1:A:240:ILE:HG12	2.14	0.47
1:C:152:LEU:HD22	1:C:159:ALA:HB2	1.95	0.47
1:B:35:THR:HG21	1:B:46:TYR:HA	1.95	0.47
1:A:152:LEU:HD22	1:A:159:ALA:HB2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:269:HIS:CE1	5:C:403:PGE:H12	2.50	0.47
1:A:35:THR:HG21	1:A:46:TYR:HA	1.95	0.47
1:A:91:VAL:HG11	1:A:238:MET:HE3	1.97	0.46
1:A:333:GLN:OE1	1:A:333:GLN:HA	2.16	0.46
1:D:91:VAL:HG11	1:D:238:MET:CE	2.46	0.45
1:D:65:TYR:HB3	6:D:527:HOH:O	2.16	0.45
1:A:213:THR:OG1	1:A:327:HIS:HE1	2.00	0.45
1:B:187[B]:ILE:HD12	1:B:187[B]:ILE:HA	1.80	0.45
1:B:150:GLN:HE22	1:D:150:GLN:HE22	1.65	0.44
1:A:127:HIS:HD2	1:A:158:ASP:OD2	1.99	0.44
1:D:22:ALA:HA	1:D:28:LEU:HD12	1.99	0.44
1:C:57:LYS:HE3	1:C:345:ILE:HG12	2.00	0.43
1:A:277:GLU:HG3	1:A:321:LYS:HG3	1.98	0.43
1:C:82:ILE:HD11	1:C:254:PHE:CE2	2.54	0.42
1:B:148:ILE:CD1	1:B:164:MET:CE	2.97	0.42
1:A:186:LEU:O	1:A:189:PRO:HD3	2.20	0.42
1:A:263:ASP:OD1	3:A:402:RIB:O3	2.37	0.42
1:A:22:ALA:HA	1:A:28:LEU:HD12	2.01	0.42
1:B:82:ILE:HD11	1:B:254:PHE:CE2	2.55	0.42
1:B:277:GLU:CG	1:B:321:LYS:HG3	2.49	0.42
1:D:186:LEU:O	1:D:189:PRO:HD3	2.20	0.41
1:A:18[B]:MET:HG3	1:A:19:ILE:N	2.33	0.41
1:B:14:ASP:OD2	1:B:262:HIS:HB2	2.21	0.41
1:B:145:GLU:HA	6:B:528:HOH:O	2.18	0.41
1:C:14:ASP:OD2	1:C:262:HIS:HB2	2.21	0.41
1:C:22:ALA:HA	1:C:28:LEU:HD12	2.02	0.41
1:A:50:ARG:CZ	1:A:103:ASN:HB2	2.51	0.41
1:C:277:GLU:HG3	1:C:321:LYS:HB2	2.01	0.41
1:C:186:LEU:O	1:C:189:PRO:HD3	2.21	0.40
1:A:166:THR:HG21	1:A:202:TYR:CZ	2.56	0.40
1:A:301:ILE:HD11	1:B:174:VAL:HG21	2.02	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231:TYR:OH	1:C:231:TYR:OH[1_455]	1.94	0.26

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	329/347 (95%)	315 (96%)	14 (4%)	0	100	100
1	B	347/347 (100%)	332 (96%)	15 (4%)	0	100	100
1	C	343/347 (99%)	328 (96%)	14 (4%)	1 (0%)	41	50
1	D	332/347 (96%)	317 (96%)	15 (4%)	0	100	100
All	All	1351/1388 (97%)	1292 (96%)	58 (4%)	1 (0%)	51	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	305	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	292/303 (96%)	288 (99%)	4 (1%)	67	81
1	B	305/303 (101%)	298 (98%)	7 (2%)	50	67
1	C	301/303 (99%)	296 (98%)	5 (2%)	60	76
1	D	295/303 (97%)	293 (99%)	2 (1%)	84	92
All	All	1193/1212 (98%)	1175 (98%)	18 (2%)	65	79

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	150	GLN
1	A	167	MET
1	A	239	ILE
1	A	312	LYS
1	B	150	GLN
1	B	167	MET
1	B	218	GLU
1	B	250	TYR
1	B	309	LYS
1	B	311	VAL
1	B	328	ASP
1	C	150	GLN
1	C	167	MET
1	C	209	ASP
1	C	250	TYR
1	C	311	VAL
1	D	150	GLN
1	D	167	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	40	GLN
1	A	127	HIS
1	A	150	GLN
1	A	327	HIS
1	B	150	GLN
1	C	39	ASN
1	C	150	GLN
1	D	31	GLN
1	D	150	GLN
1	D	262	HIS

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 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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
4	GOL	C	404	-	5,5,5	0.16	0	5,5,5	0.31	0
3	RIB	A	402	2	10,10,10	0.81	0	13,14,14	1.17	1 (7%)
5	PGE	C	403	-	9,9,9	0.30	0	8,8,8	0.21	0
4	GOL	A	403	-	5,5,5	0.19	0	5,5,5	0.39	0
3	RIB	B	402	2	10,10,10	1.40	1 (10%)	13,14,14	1.26	3 (23%)
3	RIB	D	402	2	10,10,10	0.75	0	13,14,14	1.18	1 (7%)
4	GOL	B	403	-	5,5,5	0.11	0	5,5,5	0.29	0
3	RIB	C	402	2	10,10,10	1.49	1 (10%)	13,14,14	1.56	3 (23%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	C	404	-	-	2/4/4/4	-
3	RIB	A	402	2	-	0/2/18/18	0/1/1/1
5	PGE	C	403	-	-	5/7/7/7	-
4	GOL	A	403	-	-	2/4/4/4	-
3	RIB	B	402	2	-	0/2/18/18	0/1/1/1
3	RIB	D	402	2	-	0/2/18/18	0/1/1/1
4	GOL	B	403	-	-	2/4/4/4	-
3	RIB	C	402	2	-	0/2/18/18	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	402	RIB	O4-C1	-4.16	1.38	1.43
3	B	402	RIB	O4-C1	-3.65	1.38	1.43

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	402	RIB	O1-C1-O4	-3.50	106.66	111.13
3	C	402	RIB	C5-C4-C3	-2.63	108.75	115.09
3	A	402	RIB	O1-C1-O4	-2.49	107.95	111.13
3	C	402	RIB	O2-C2-C1	2.45	118.57	111.82
3	B	402	RIB	O1-C1-O4	-2.32	108.16	111.13
3	D	402	RIB	O5-C5-C4	-2.31	103.36	111.29
3	B	402	RIB	C5-C4-C3	-2.27	109.62	115.09
3	B	402	RIB	O2-C2-C1	2.13	117.70	111.82

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	404	GOL	C1-C2-C3-O3
5	C	403	PGE	O1-C1-C2-O2
5	C	403	PGE	O2-C3-C4-O3
4	A	403	GOL	O1-C1-C2-C3
4	B	403	GOL	C1-C2-C3-O3
4	B	403	GOL	O2-C2-C3-O3
4	C	404	GOL	O2-C2-C3-O3
5	C	403	PGE	C1-C2-O2-C3
4	A	403	GOL	O1-C1-C2-O2
5	C	403	PGE	C3-C4-O3-C5
5	C	403	PGE	C6-C5-O3-C4

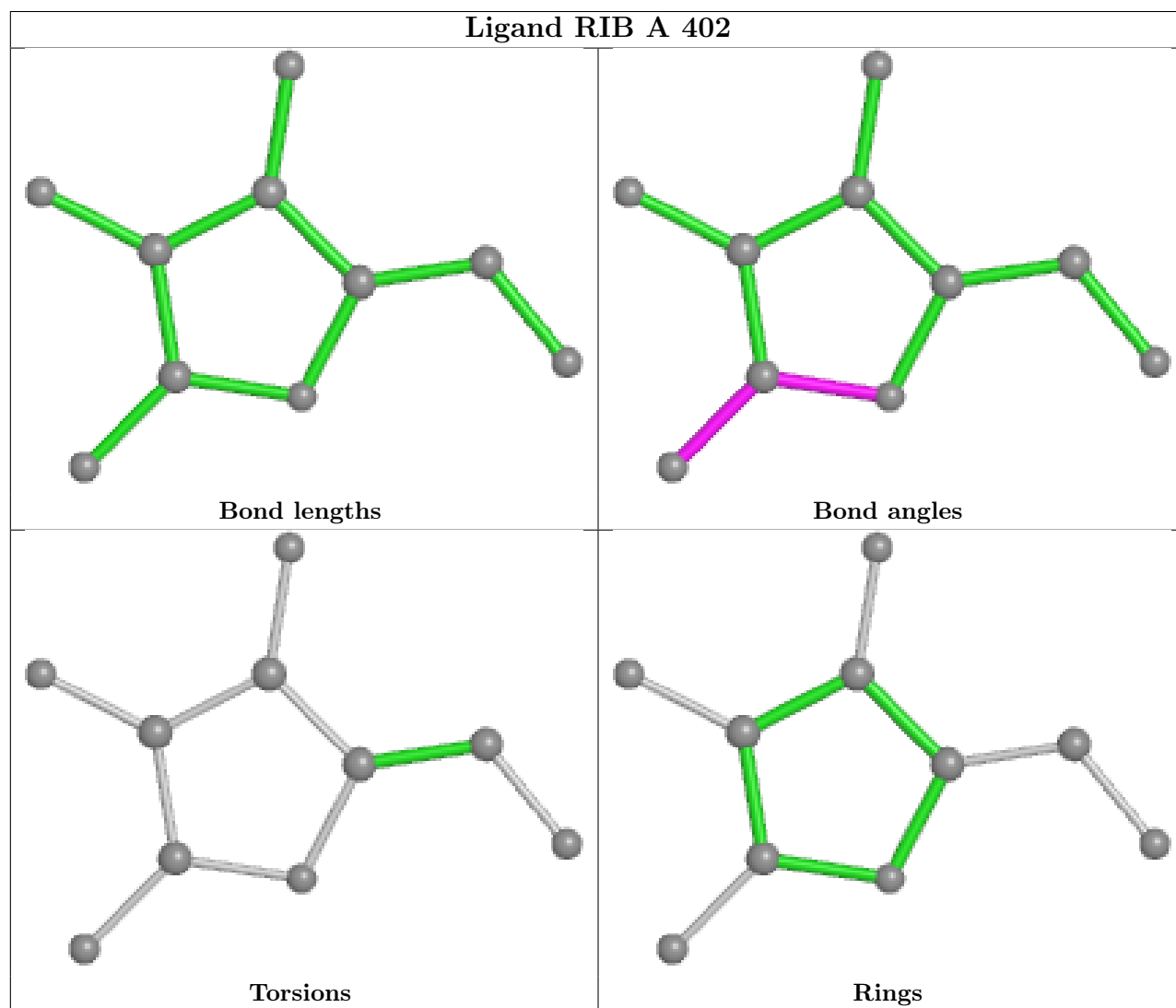
There are no ring outliers.

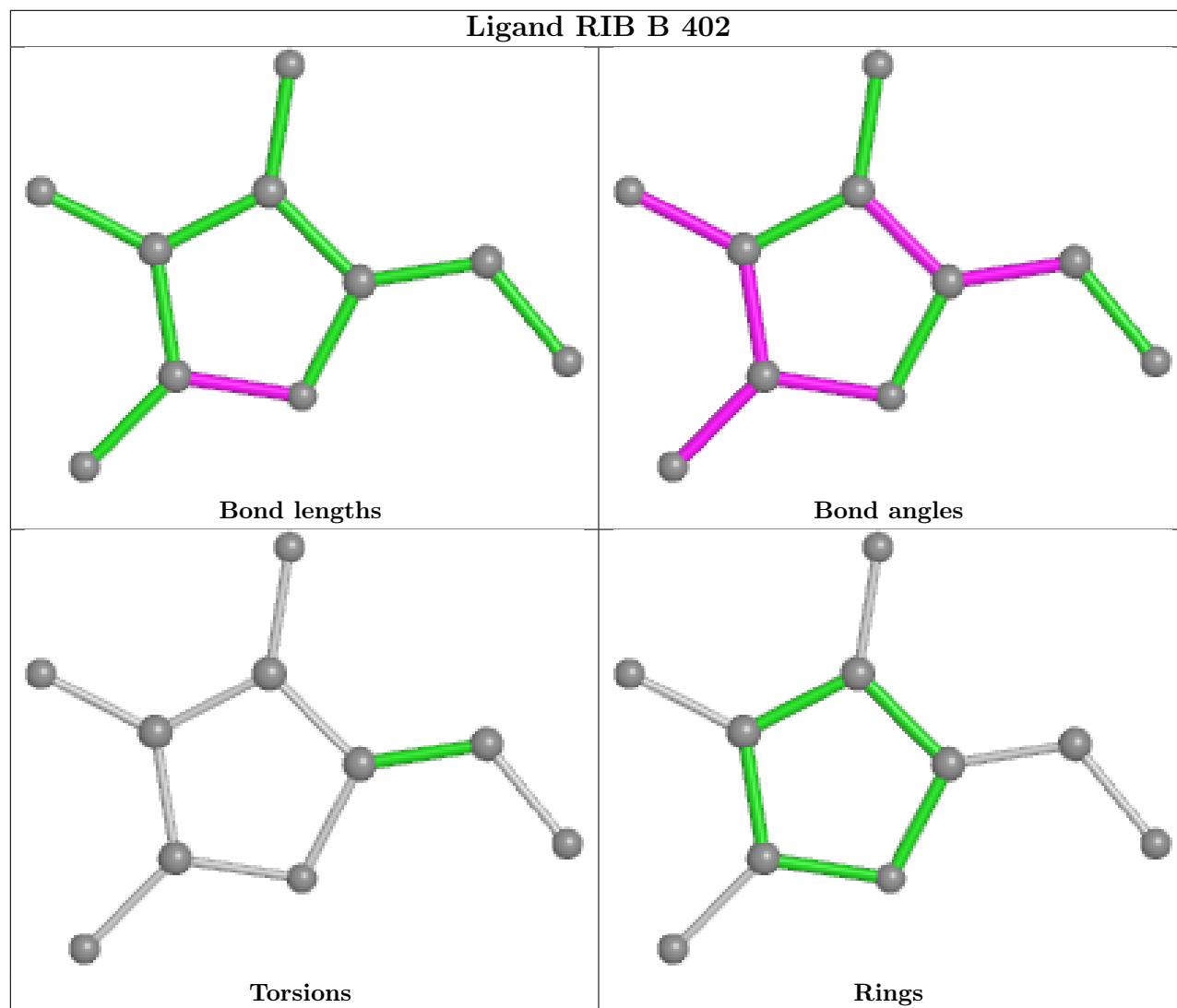
3 monomers are involved in 3 short contacts:

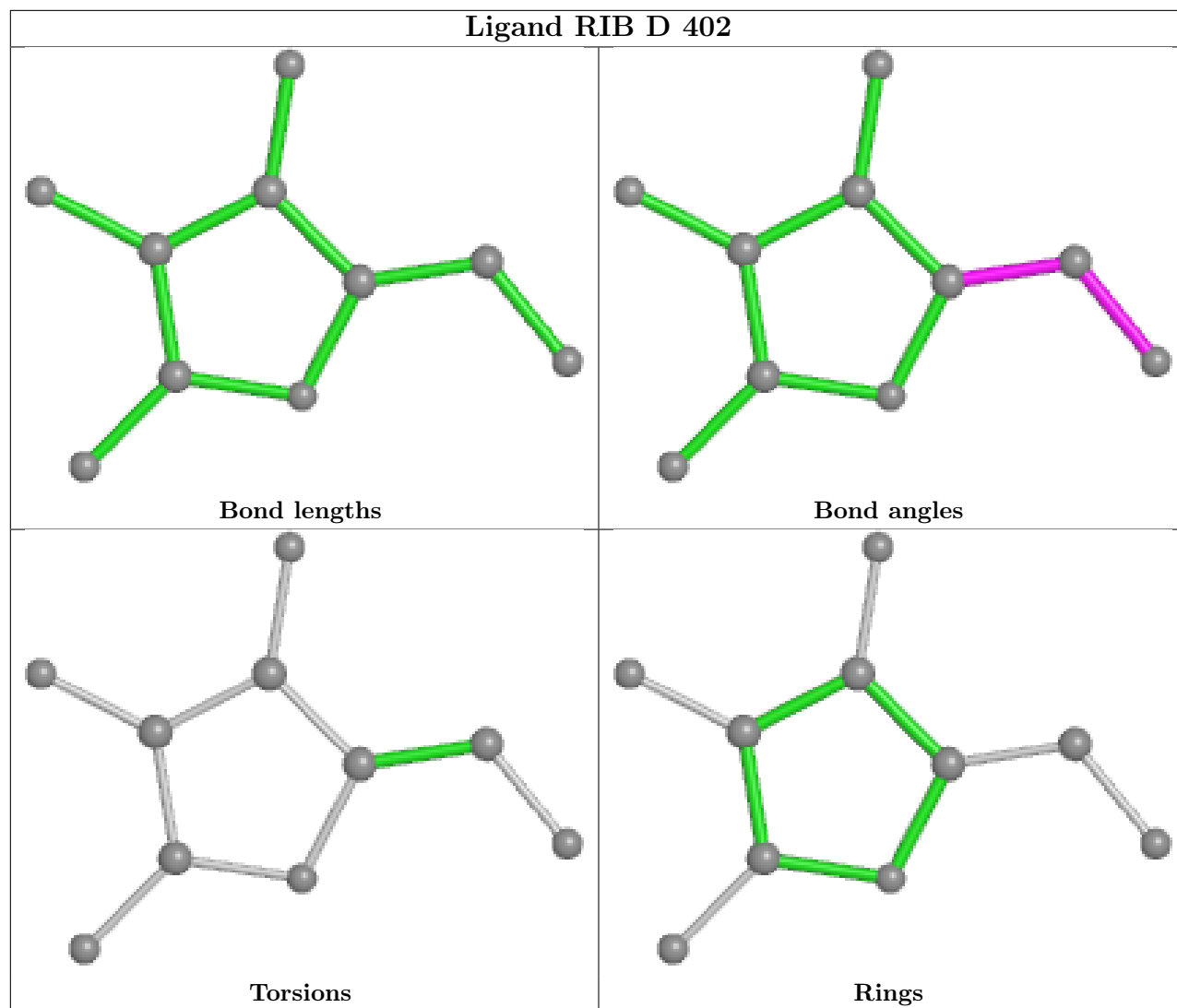
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	404	GOL	1	0
3	A	402	RIB	1	0
5	C	403	PGE	1	0

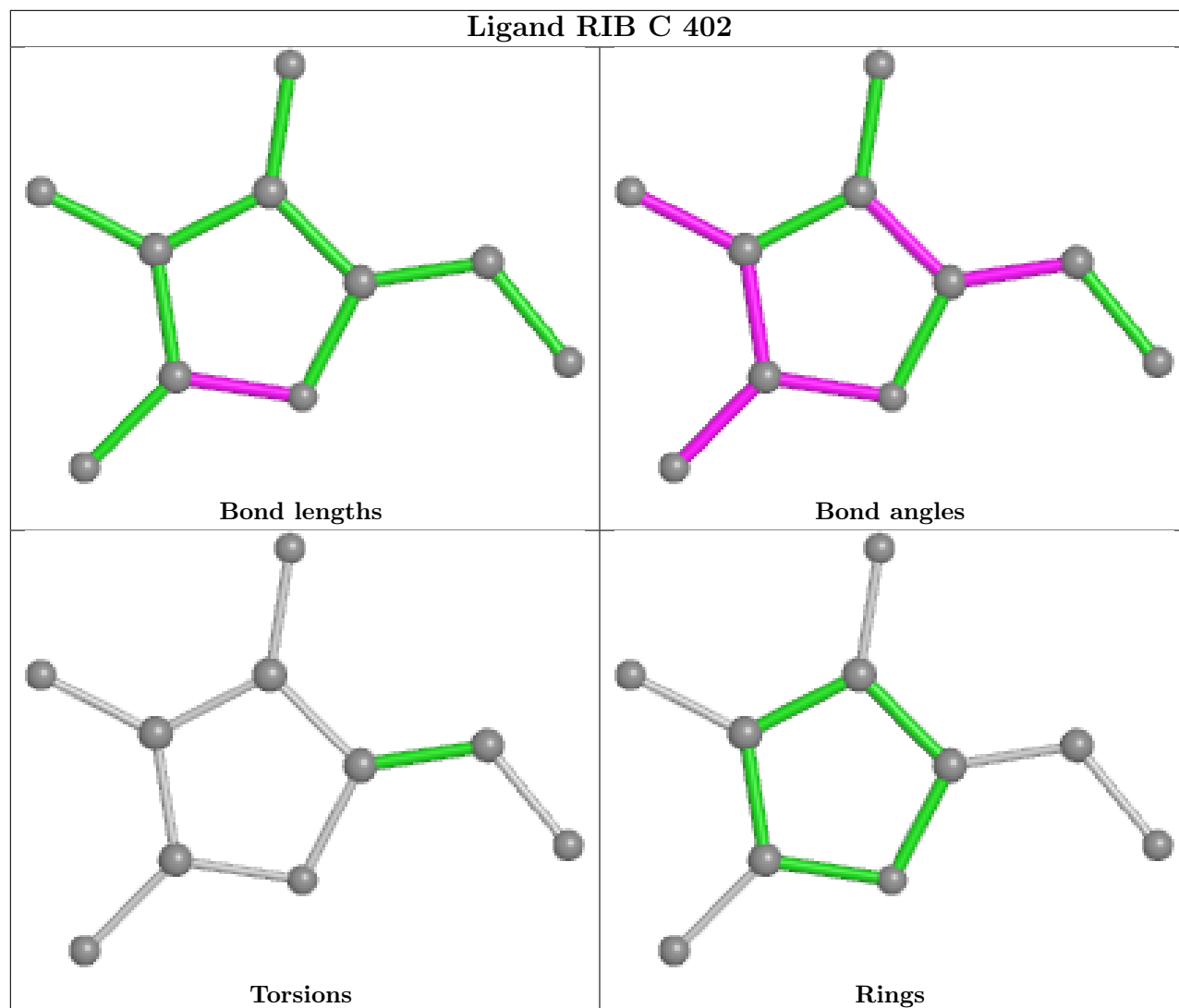
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

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.









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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	331/347 (95%)	0.39	14 (4%) 36 43	18, 30, 55, 72	0
1	B	345/347 (99%)	0.35	10 (2%) 51 58	18, 28, 45, 79	0
1	C	345/347 (99%)	0.42	10 (2%) 51 58	18, 29, 46, 81	0
1	D	335/347 (96%)	0.48	20 (5%) 21 28	20, 31, 59, 71	0
All	All	1356/1388 (97%)	0.41	54 (3%) 38 45	18, 29, 55, 81	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	246	PHE	6.5
1	C	310	ILE	6.2
1	B	306	ASP	5.7
1	C	306	ASP	5.6
1	B	309	LYS	5.1
1	B	311	VAL	5.0
1	D	246	PHE	4.9
1	D	247	LYS	4.7
1	A	247	LYS	4.6
1	A	90	GLY	4.3
1	D	77	VAL	4.0
1	D	231	TYR	3.9
1	C	311	VAL	3.8
1	C	309	LYS	3.7
1	C	308	THR	3.6
1	B	78	ALA	3.6
1	B	307	PRO	3.5
1	D	90	GLY	3.4
1	D	89	GLY	3.2
1	C	307	PRO	3.2
1	D	91	VAL	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	91	VAL	3.0
1	A	47	GLN	3.0
1	D	234	LYS	2.9
1	B	310	ILE	2.9
1	D	311	VAL	2.9
1	A	238	MET	2.8
1	B	308	THR	2.8
1	C	249	THR	2.7
1	D	86	SER	2.6
1	D	238	MET	2.6
1	A	55	LEU	2.5
1	A	231	TYR	2.5
1	B	302	LEU	2.4
1	D	78	ALA	2.4
1	A	244	PHE	2.4
1	D	256	PHE	2.4
1	D	275	TRP	2.3
1	C	302	LEU	2.3
1	B	278[A]	HIS	2.2
1	C	304	LYS	2.2
1	A	253	VAL	2.2
1	A	312	LYS	2.2
1	D	310	ILE	2.2
1	A	77	VAL	2.1
1	D	244	PHE	2.1
1	D	134	GLU	2.1
1	B	249	THR	2.1
1	D	342	ALA	2.1
1	D	345	ILE	2.0
1	C	78	ALA	2.0
1	D	237	GLU	2.0
1	A	256	PHE	2.0
1	A	234	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands

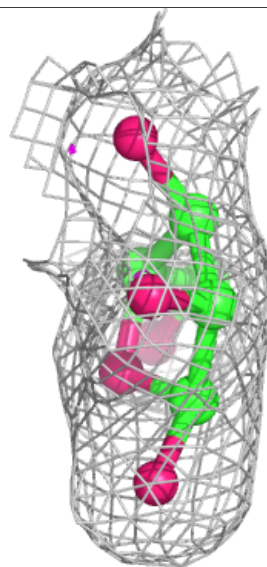
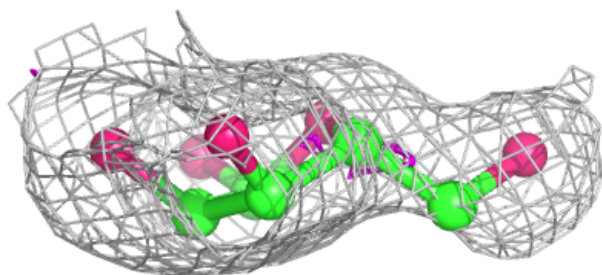
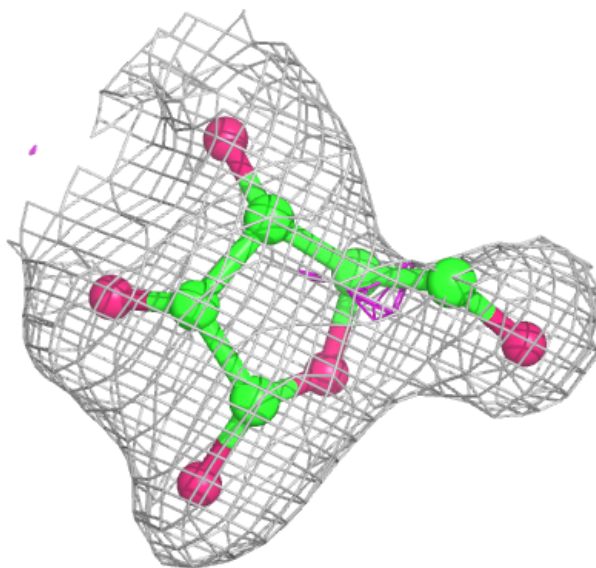
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	B	403	6/6	0.69	0.17	47,48,54,56	0
4	GOL	C	404	6/6	0.80	0.19	37,43,46,52	0
4	GOL	A	403	6/6	0.87	0.17	29,34,36,36	0
5	PGE	C	403	10/10	0.87	0.20	30,41,46,47	0
3	RIB	A	402	10/10	0.91	0.14	20,28,32,36	0
3	RIB	C	402	10/10	0.94	0.12	19,21,22,24	0
3	RIB	D	402	10/10	0.94	0.10	23,27,29,29	0
3	RIB	B	402	10/10	0.95	0.16	17,22,27,30	0
2	CA	A	401	1/1	0.98	0.05	23,23,23,23	0
2	CA	C	401	1/1	0.98	0.06	22,22,22,22	0
2	CA	B	401	1/1	0.99	0.03	21,21,21,21	0
2	CA	D	401	1/1	0.99	0.04	23,23,23,23	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.

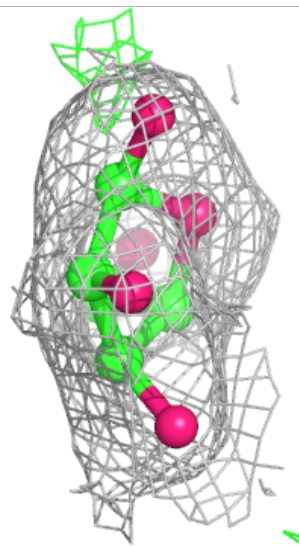
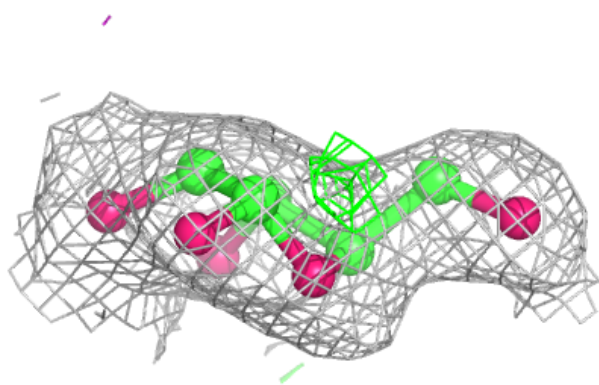
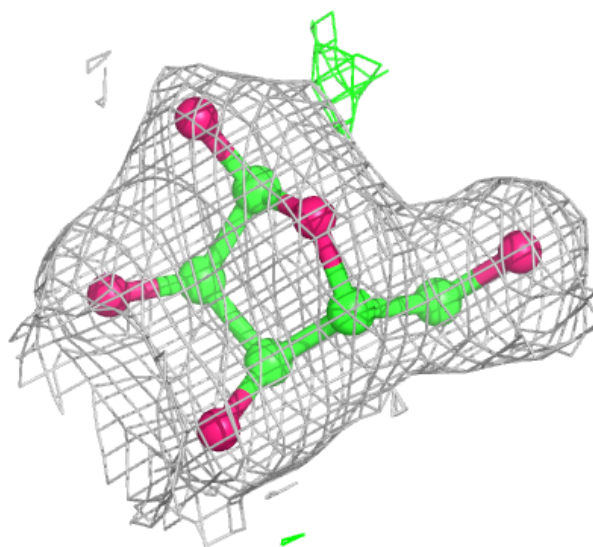
Electron density around RIB A 402:

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



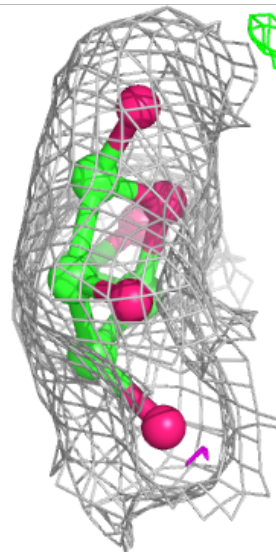
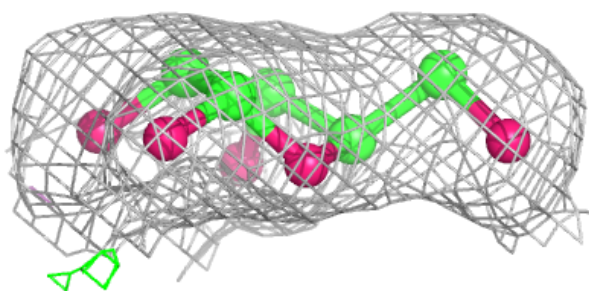
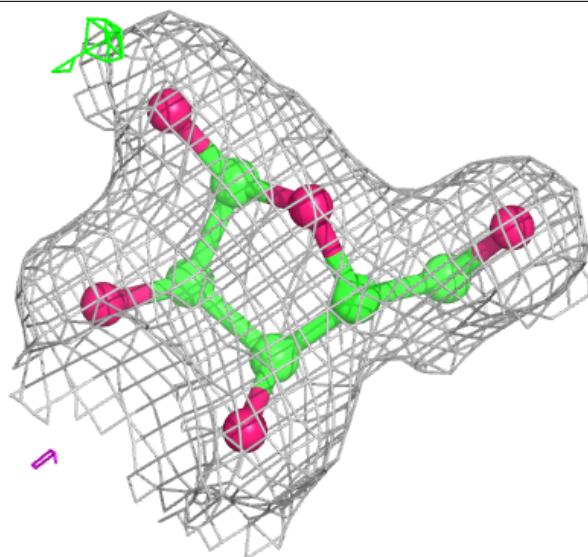
Electron density around RIB C 402:

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



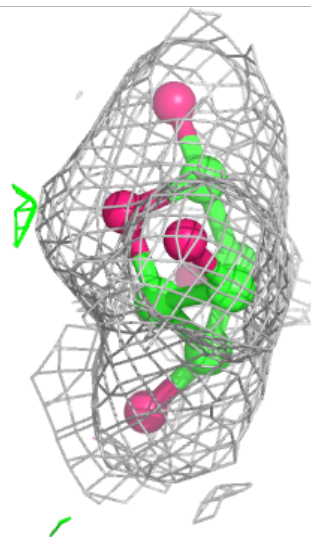
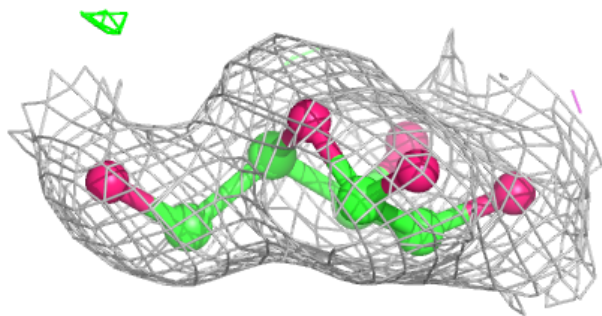
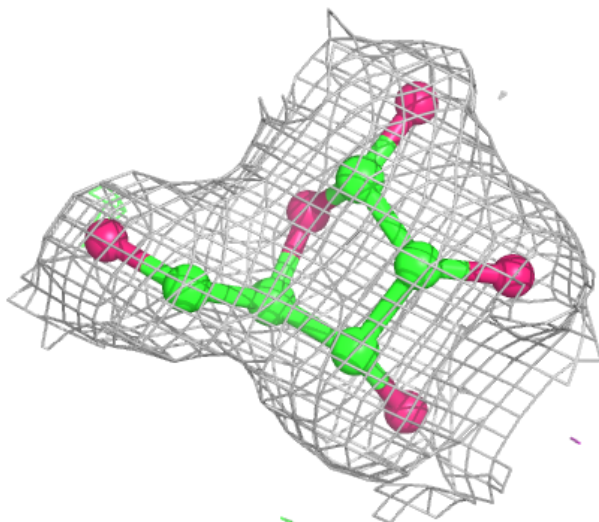
Electron density around RIB D 402:

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and green (positive)



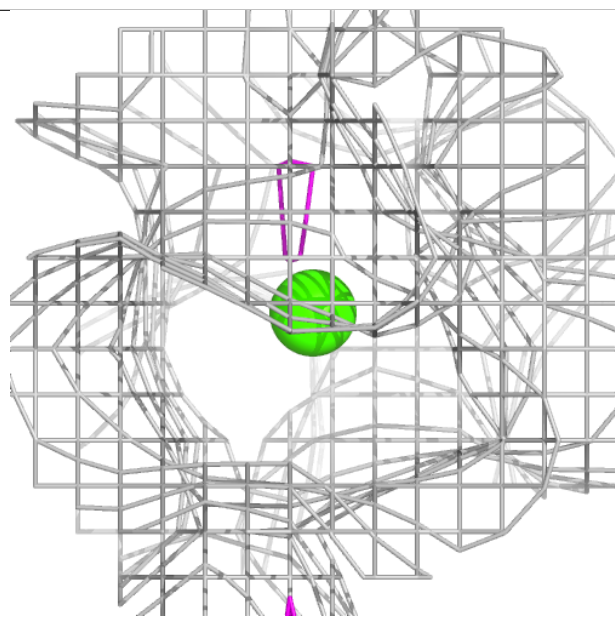
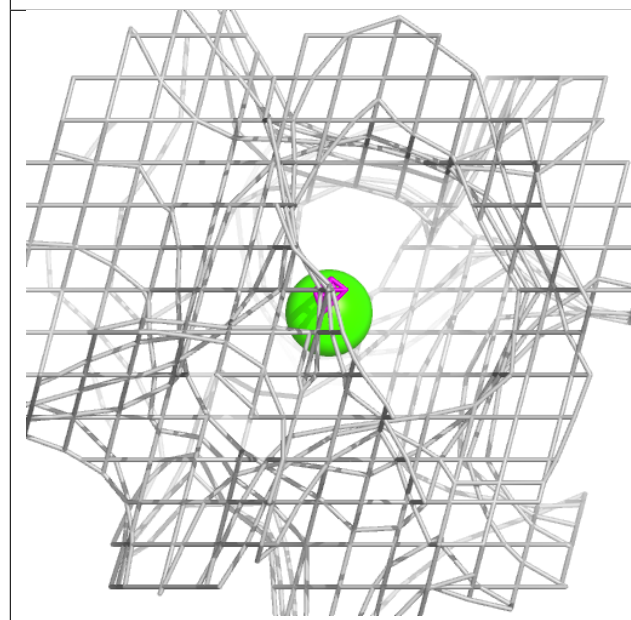
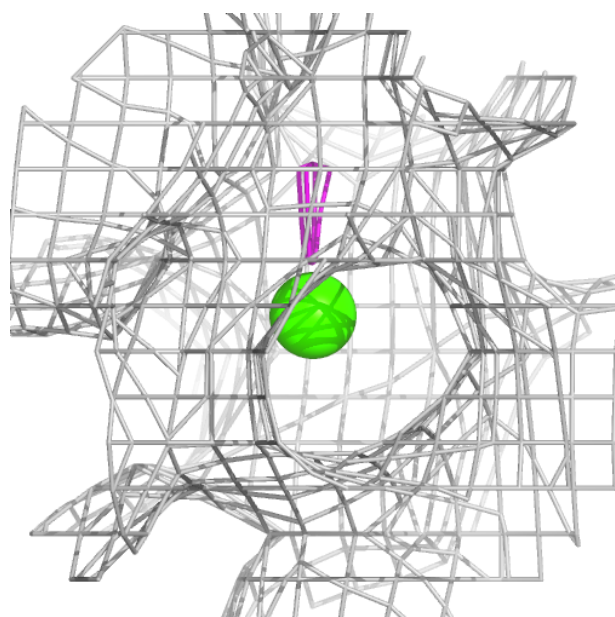
Electron density around RIB B 402:

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and green (positive)



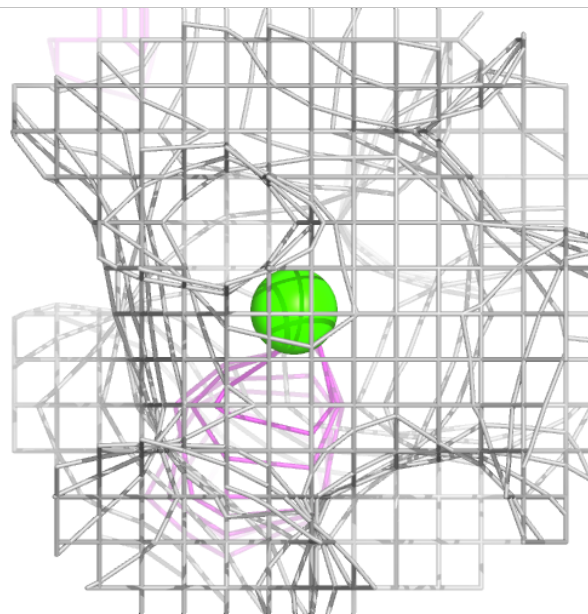
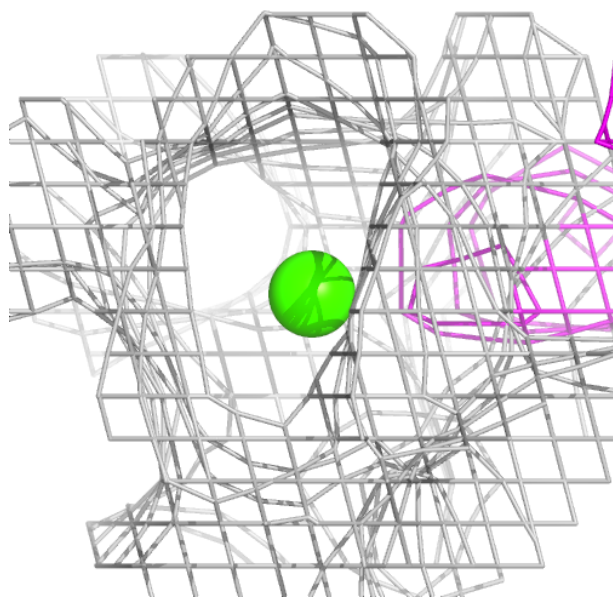
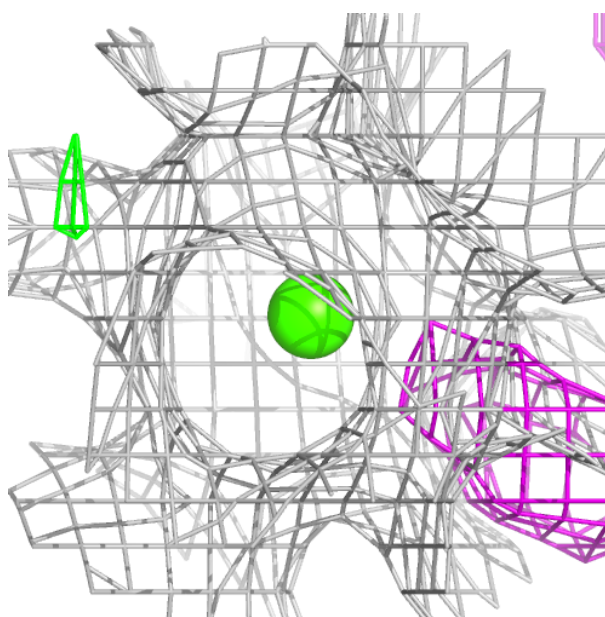
Electron density around CA A 401:

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



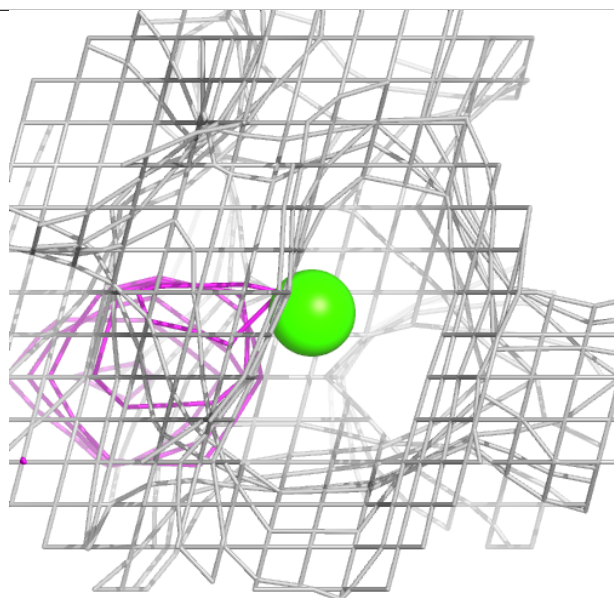
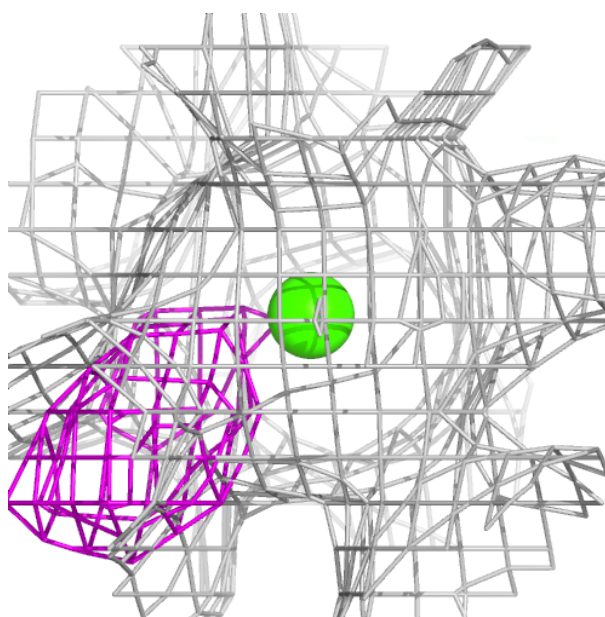
Electron density around CA C 401:

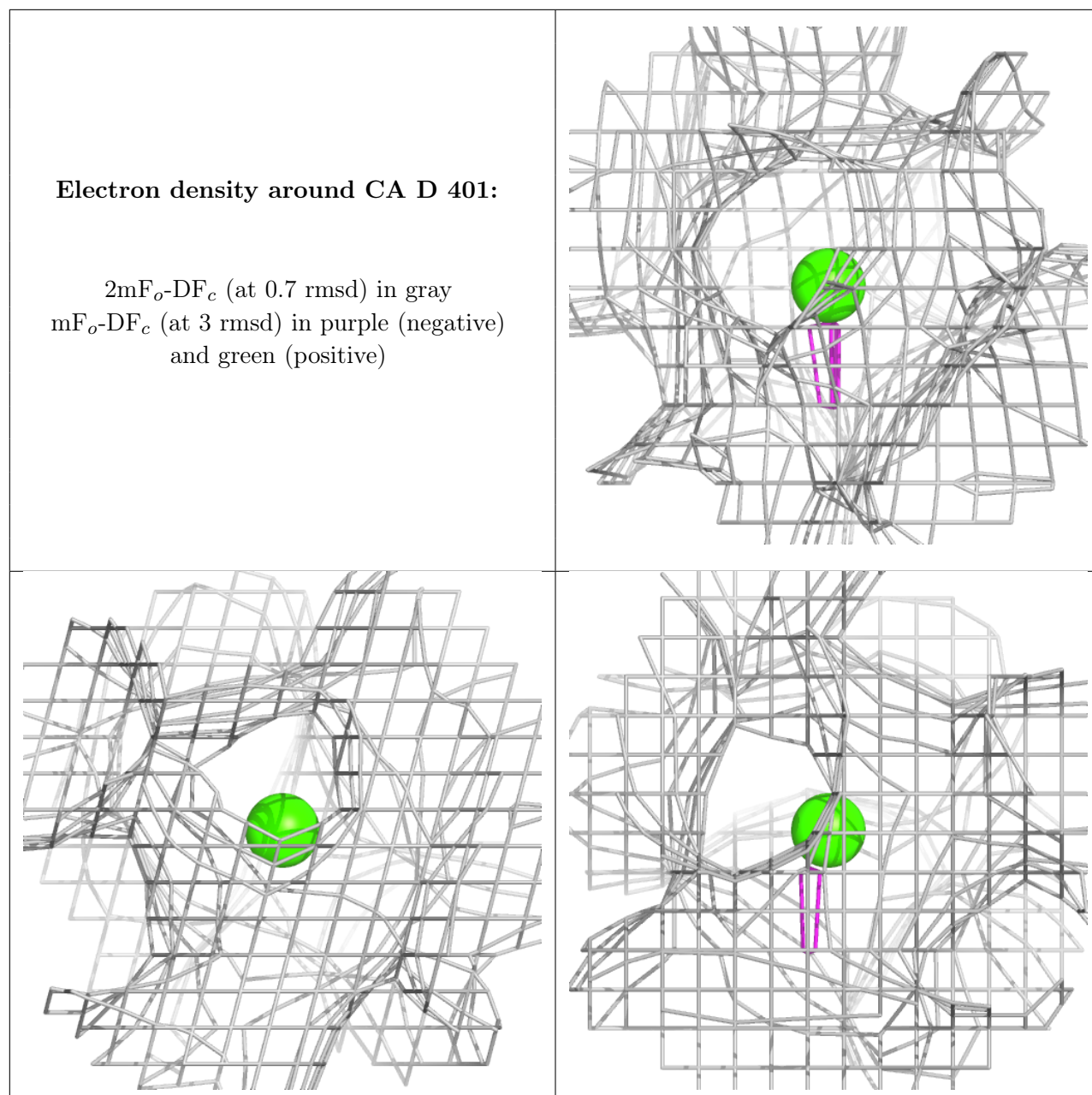
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around CA B 401:

$2mF_o-DF_c$ (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.