Full wwPDB X-ray Structure Validation Report (i)

Nov 5, 2023 – 10:48 pm GMT

PDB ID	:	6T5M
Title	:	Crystal structure of 2-methylisocitrate lyase (PrpB) from Pseudomonas aeruginosa in complex with Mg(II)-pyruvate.
Authors	:	Wijaya, A.J.; Brear, P.; Dolan, S.K.; Welch, M.
Deposited on	:	2019-10-16
Resolution	:	1.76 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 (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:

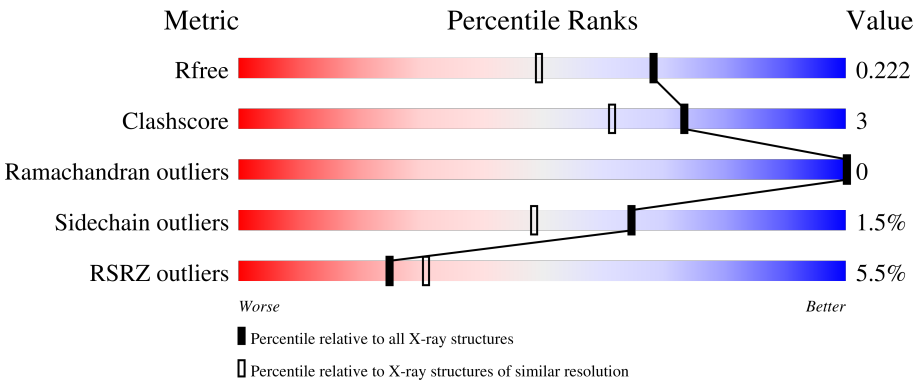
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 1.76 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 >=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	298	5% 88% 5% 7%
1	B	298	10% 83% 10% 7%
1	C	298	3% 84% 9% 7%
1	D	298	3% 84% 8% 7%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 8901 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.

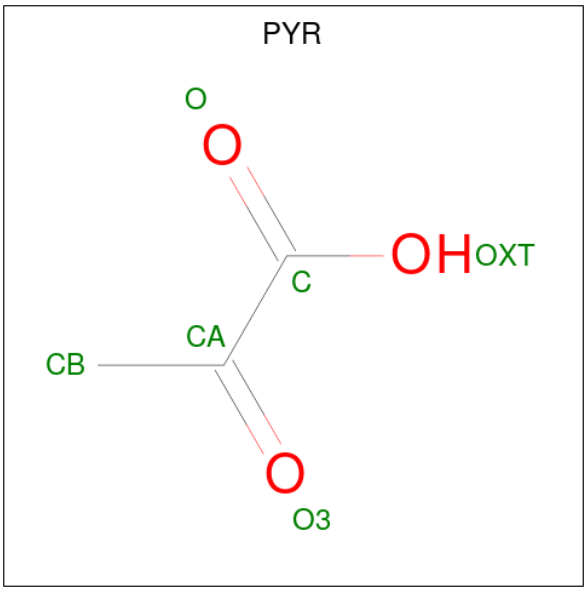
- Molecule 1 is a protein called 2-methylisocitrate lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	C	277	2120	1334	360	412	14	0	4	0
1	A	276	2085	1313	355	404	13	0	0	0
1	D	276	2085	1313	355	404	13	0	0	0
1	B	276	2093	1319	356	405	13	0	1	0

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 3 is PYRUVIC ACID (three-letter code: PYR) (formula: $C_3H_4O_3$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is water.

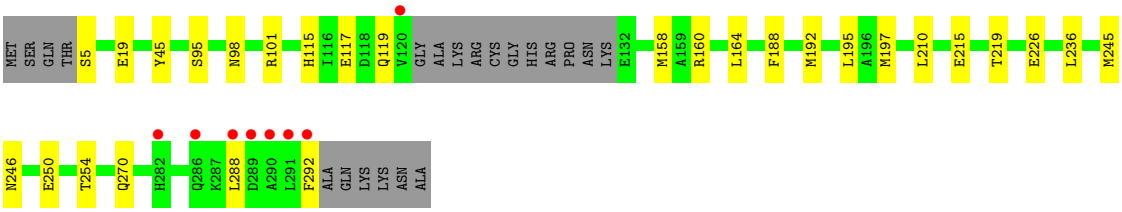
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	C	166	170	170	0	4
4	A	135	136	136	0	1
4	D	106	106	106	0	0
4	B	78	78	78	0	0

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.

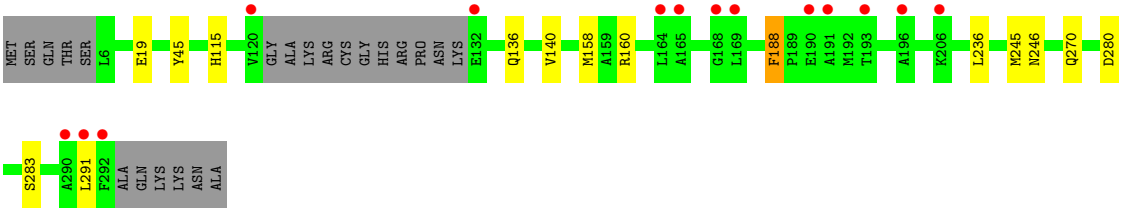
• Molecule 1: 2-methylisocitrate lyase

Chain C:



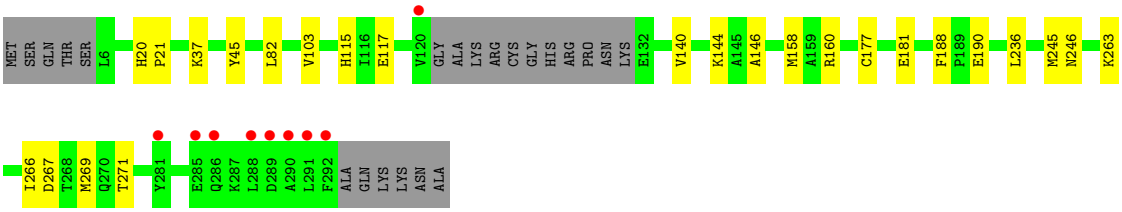
• Molecule 1: 2-methylisocitrate lyase

Chain A:



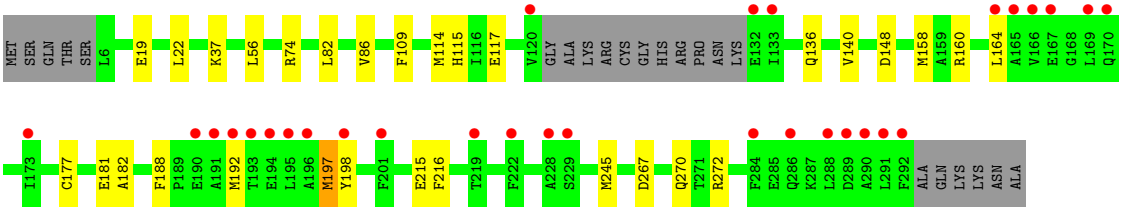
• Molecule 1: 2-methylisocitrate lyase

Chain D:



• Molecule 1: 2-methylisocitrate lyase

Chain B:


MET SER GLN THR SER L6 E19 L22 K37 L56 R74 L82 V86 F109 M114 H115 I116 E117 V120 GLY ALA LYS ARG CYS GLY HIS ARG PRO ASN LYS E132 I133 Q136 V140 D148 M158 A159 R160 L164 A165 V166 E167 G168 L169 Q170
I173 C177 E181 A182 F188 P189 E190 A191 M192 T193 E194 L195 A196 M197 Y198 F201 E215 F216 T219 F222 A228 S229 M245 D267 Q270 T271 R272 F284 E285 Q286 K287 L288 D289 A290 L291 F292 ALA GLN LYS LYS ASN ALA

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	153.49Å 59.40Å 147.65Å 90.00° 120.30° 90.00°	Depositor
Resolution (Å)	42.49 – 1.76 63.74 – 1.76	Depositor EDS
% Data completeness (in resolution range)	85.4 (42.49-1.76) 85.4 (63.74-1.76)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.42 (at 1.76Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.211 , 0.227 0.208 , 0.222	Depositor DCC
R_{free} test set	4808 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å^2)	27.2	Xtriage
Anisotropy	0.340	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 37.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.006 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	8901	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.32% 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 (i)

5.1 Standard geometry (i)

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

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	0.79	1/2114 (0.0%)	0.83	0/2864
1	B	0.74	0/2122	0.81	0/2875
1	C	0.84	3/2149 (0.1%)	0.87	0/2913
1	D	0.78	0/2114	0.82	0/2864
All	All	0.79	4/8499 (0.0%)	0.83	0/11516

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	19	GLU	CD-OE2	-6.47	1.18	1.25
1	C	19	GLU	CD-OE1	-6.06	1.19	1.25
1	A	19	GLU	CD-OE2	-5.41	1.19	1.25
1	C	215	GLU	CD-OE2	-5.09	1.20	1.25

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	2085	0	2084	10	0
1	B	2093	0	2094	21	0
1	C	2120	0	2118	20	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2085	0	2084	15	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	6	0	0	0	0
3	B	6	0	0	0	0
3	C	6	0	0	0	0
3	D	6	0	0	0	0
4	A	136	0	0	0	0
4	B	78	0	0	0	0
4	C	170	0	0	4	0
4	D	106	0	0	0	0
All	All	8901	0	8380	56	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.

All (56) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:246:ASN:HD22	1:B:270:GLN:HE22	1.29	0.76
1:C:5:SER:HA	4:C:449:HOH:O	1.86	0.76
1:C:101:ARG:NH2	1:B:56[B]:LEU:HD11	2.06	0.70
1:C:270:GLN:HE22	1:A:246:ASN:HD22	1.39	0.69
1:C:246:ASN:HD22	1:A:270:GLN:HE22	1.41	0.67
1:B:192:MET:HB2	1:B:198:TYR:CZ	2.31	0.66
1:B:192:MET:HB2	1:B:198:TYR:CE2	2.34	0.62
1:D:246:ASN:HD22	1:B:270:GLN:NE2	1.98	0.60
1:B:37:LYS:HA	1:B:82:LEU:HD11	1.90	0.54
1:C:250:GLU:O	1:C:254[B]:THR:HG23	2.08	0.54
1:C:119:GLN:NE2	4:C:402[A]:HOH:O	2.38	0.53
1:C:5:SER:CA	4:C:449:HOH:O	2.51	0.52
1:C:164:LEU:HD22	1:C:192:MET:HE3	1.92	0.52
1:D:103:VAL:HG21	1:D:146:ALA:HA	1.92	0.51
1:C:164:LEU:HD22	1:C:192:MET:CE	2.40	0.50
1:C:195:LEU:HD11	1:C:226:GLU:HB3	1.94	0.49
1:C:246:ASN:HD22	1:A:270:GLN:NE2	2.07	0.49
1:B:117:GLU:HA	1:B:160:ARG:O	2.12	0.49
1:D:177:CYS:O	1:D:181:GLU:HG2	2.13	0.48
1:A:45:TYR:CG	1:A:236:LEU:HD11	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:177:CYS:O	1:B:181:GLU:HG2	2.14	0.47
1:D:140:VAL:HG12	1:D:144:LYS:HD2	1.96	0.47
1:D:263:LYS:O	1:D:266:ILE:HG12	2.14	0.47
1:A:136:GLN:O	1:A:140:VAL:HG23	2.14	0.47
1:C:115:HIS:HA	1:C:158:MET:O	2.16	0.46
1:C:270:GLN:NE2	1:A:246:ASN:HD22	2.08	0.46
1:B:164:LEU:HD22	1:B:192:MET:HE2	1.98	0.45
1:C:119:GLN:NE2	4:C:402[B]:HOH:O	2.50	0.44
1:B:115:HIS:HA	1:B:158:MET:O	2.18	0.44
1:C:45:TYR:CG	1:C:236:LEU:HD11	2.53	0.44
1:B:192:MET:SD	1:B:197:MET:HG2	2.57	0.44
1:A:115:HIS:HA	1:A:158:MET:O	2.17	0.44
1:D:269:MET:SD	1:B:215:GLU:HG2	2.58	0.44
1:B:19:GLU:HG2	1:B:22:LEU:HA	1.99	0.44
1:C:288:LEU:HD22	1:C:292:PHE:HE2	1.82	0.43
1:B:74:ARG:HB2	1:B:109:PHE:HD2	1.82	0.43
1:D:117:GLU:HA	1:D:160:ARG:O	2.18	0.43
1:C:210:LEU:C	1:C:210:LEU:HD23	2.38	0.43
1:D:45:TYR:CG	1:D:236:LEU:HD11	2.54	0.43
1:B:140:VAL:HG13	1:B:182:ALA:HB2	2.01	0.43
1:B:197:MET:HE2	1:B:197:MET:HB2	1.90	0.43
1:D:37:LYS:HA	1:D:82:LEU:HD11	2.00	0.43
1:D:115:HIS:HA	1:D:158:MET:O	2.19	0.43
1:D:160:ARG:NH2	1:D:190:GLU:CD	2.71	0.43
1:B:86:VAL:O	1:B:114:MET:HA	2.19	0.42
1:B:136:GLN:O	1:B:140:VAL:HG23	2.19	0.42
1:C:219[A]:THR:HG23	1:C:219[A]:THR:O	2.20	0.42
1:B:74:ARG:HB2	1:B:109:PHE:CD2	2.54	0.42
1:D:20:HIS:HA	1:D:21:PRO:HA	1.81	0.41
1:A:291:LEU:HD21	1:B:148:ASP:CB	2.51	0.41
1:C:117:GLU:HA	1:C:160:ARG:O	2.21	0.41
1:D:271:THR:HA	1:B:216:PHE:CE2	2.55	0.41
1:C:95:SER:HB3	1:C:98:ASN:OD1	2.21	0.40
1:A:280:ASP:OD1	1:A:283:SER:HB3	2.21	0.40
1:A:160:ARG:HA	1:A:188:PHE:O	2.22	0.40
1:D:263:LYS:HB2	1:D:263:LYS:HE3	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	272/298 (91%)	265 (97%)	7 (3%)	0	100	100
1	B	273/298 (92%)	265 (97%)	8 (3%)	0	100	100
1	C	277/298 (93%)	270 (98%)	7 (2%)	0	100	100
1	D	272/298 (91%)	265 (97%)	7 (3%)	0	100	100
All	All	1094/1192 (92%)	1065 (97%)	29 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	218/235 (93%)	216 (99%)	2 (1%)	78	67
1	B	219/235 (93%)	214 (98%)	5 (2%)	50	28
1	C	223/235 (95%)	220 (99%)	3 (1%)	69	54
1	D	218/235 (93%)	215 (99%)	3 (1%)	67	52
All	All	878/940 (93%)	865 (98%)	13 (2%)	65	49

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	188	PHE
1	C	197	MET

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Mol	Chain	Res	Type
1	C	245	MET
1	A	188	PHE
1	A	245	MET
1	D	188	PHE
1	D	245	MET
1	D	267	ASP
1	B	188	PHE
1	B	197	MET
1	B	245	MET
1	B	267	ASP
1	B	272	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	251	ASN
1	C	264	ASN
1	C	270	GLN
1	C	282	HIS
1	A	264	ASN
1	A	270	GLN
1	D	264	ASN
1	D	270	GLN
1	B	264	ASN
1	B	270	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	$\#\|Z\| > 2$
3	PYR	C	302	2	5,5,5	2.65	1 (20%)	3,6,6	2.25	1 (33%)
3	PYR	D	302	2	5,5,5	2.62	1 (20%)	3,6,6	1.34	0
3	PYR	A	302	2	5,5,5	3.02	2 (40%)	3,6,6	1.12	0
3	PYR	B	302	2	5,5,5	2.43	1 (20%)	3,6,6	1.81	1 (33%)

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
3	PYR	C	302	2	-	0/4/4/4	-
3	PYR	D	302	2	-	0/4/4/4	-
3	PYR	A	302	2	-	0/4/4/4	-
3	PYR	B	302	2	-	0/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	PYR	CA-C	-6.22	1.32	1.54
3	D	302	PYR	CA-C	-5.50	1.34	1.54
3	C	302	PYR	CA-C	-5.22	1.35	1.54
3	B	302	PYR	CA-C	-4.99	1.36	1.54
3	A	302	PYR	OXT-C	-2.59	1.23	1.30

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	302	PYR	OXT-C-CA	3.66	123.98	113.97
3	B	302	PYR	OXT-C-CA	2.40	120.53	113.97

There are no chirality outliers.

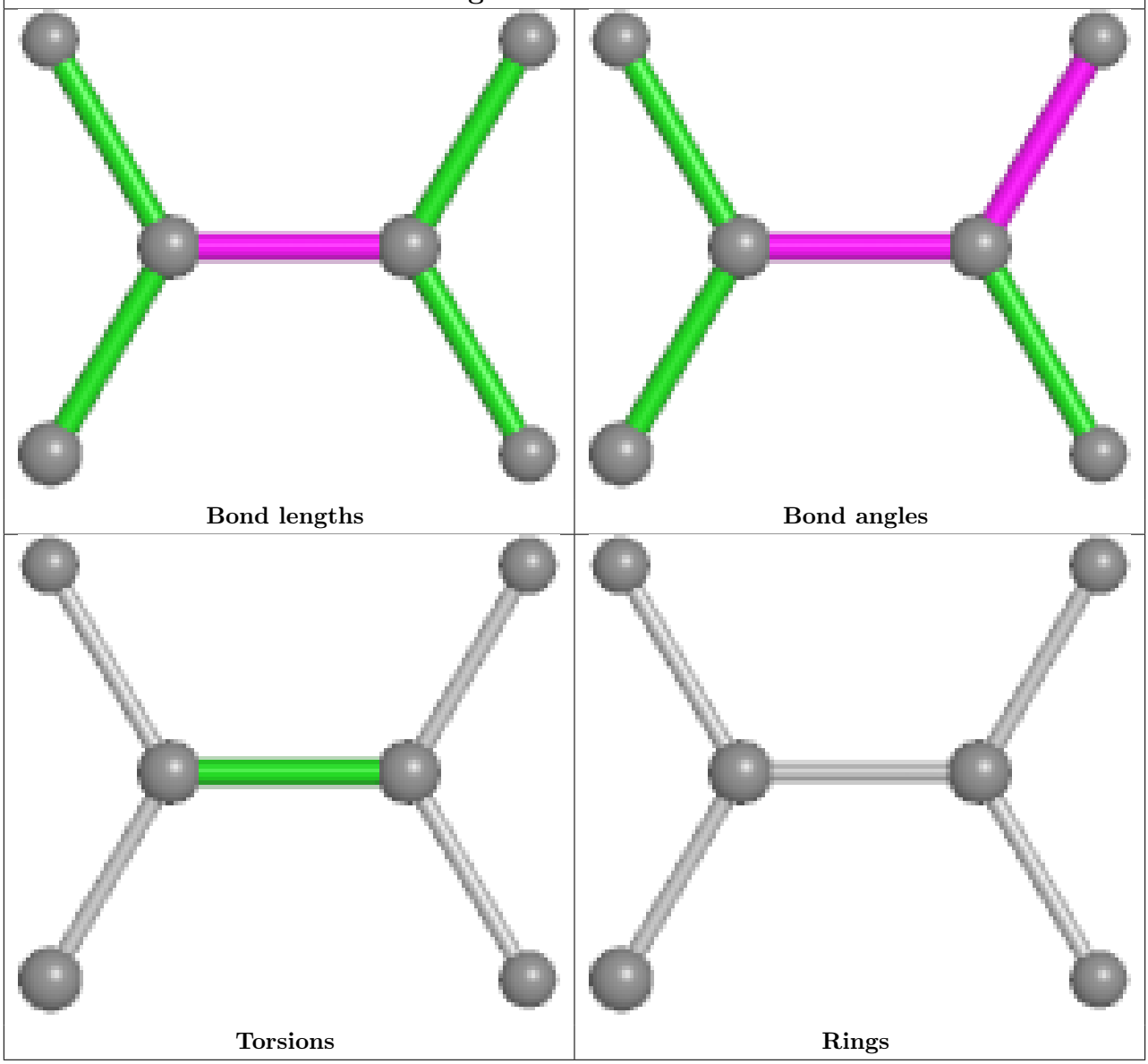
There are no torsion outliers.

There are no ring outliers.

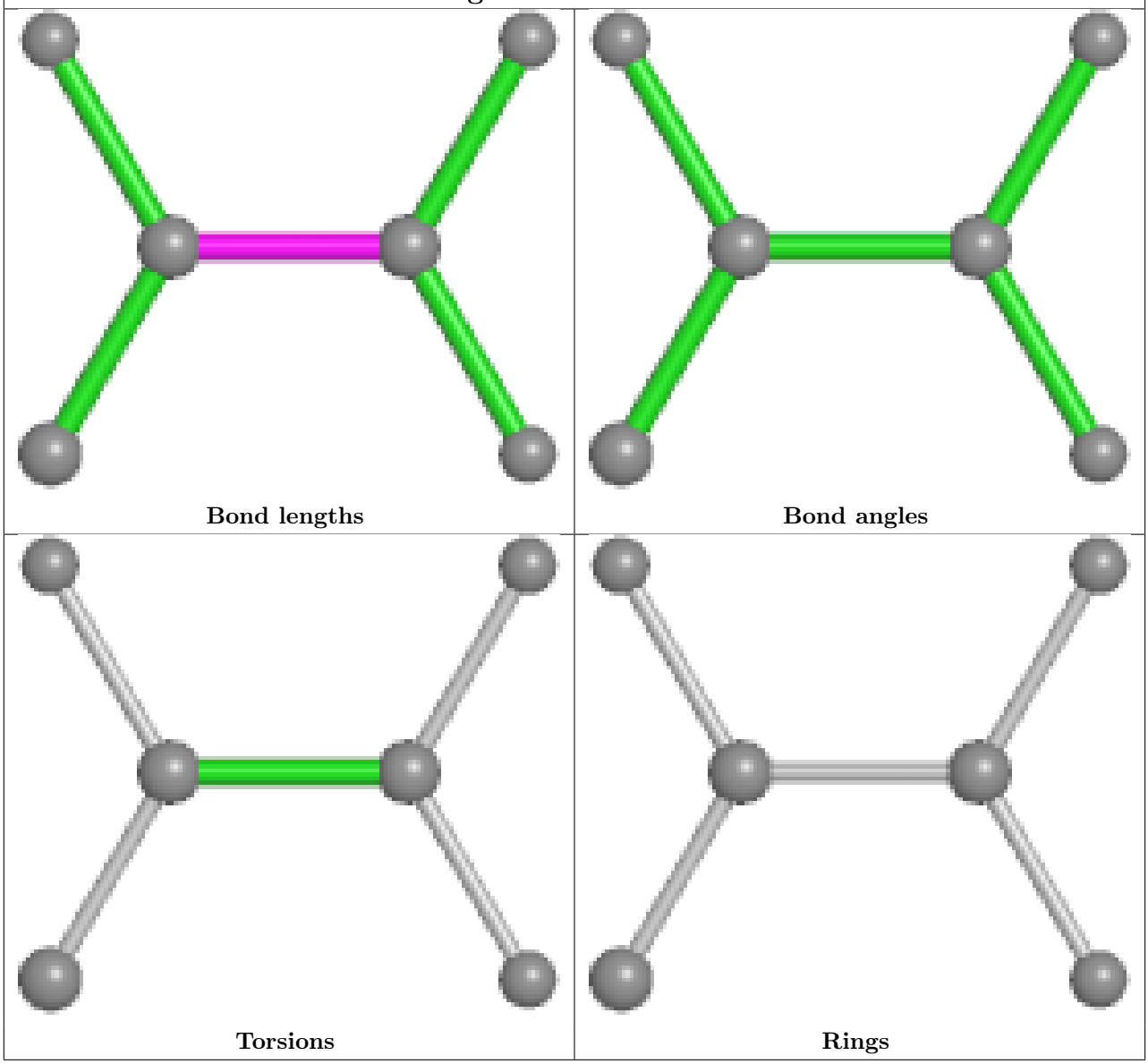
No monomer is involved in short contacts.

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.

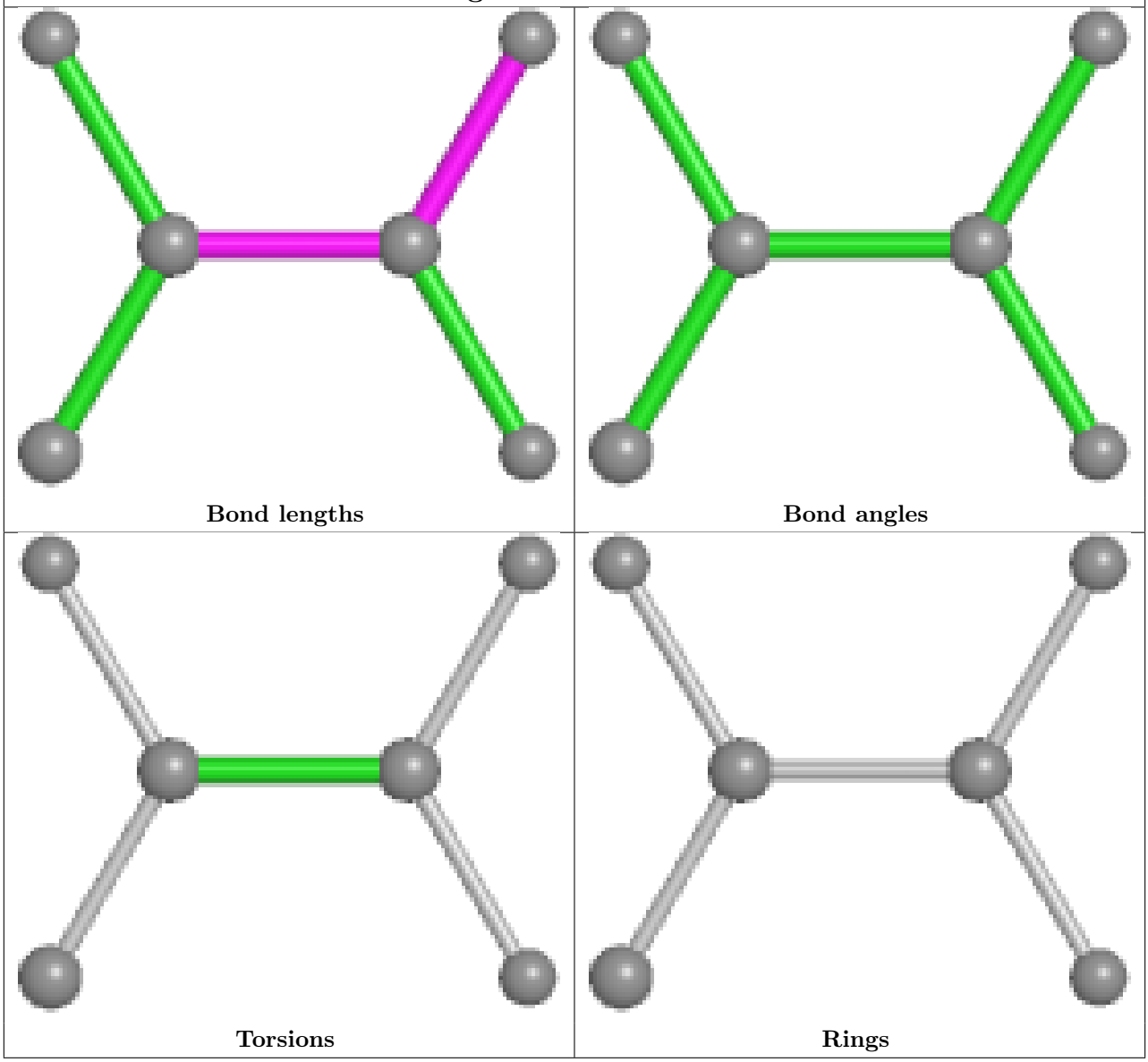
Ligand PYR C 302



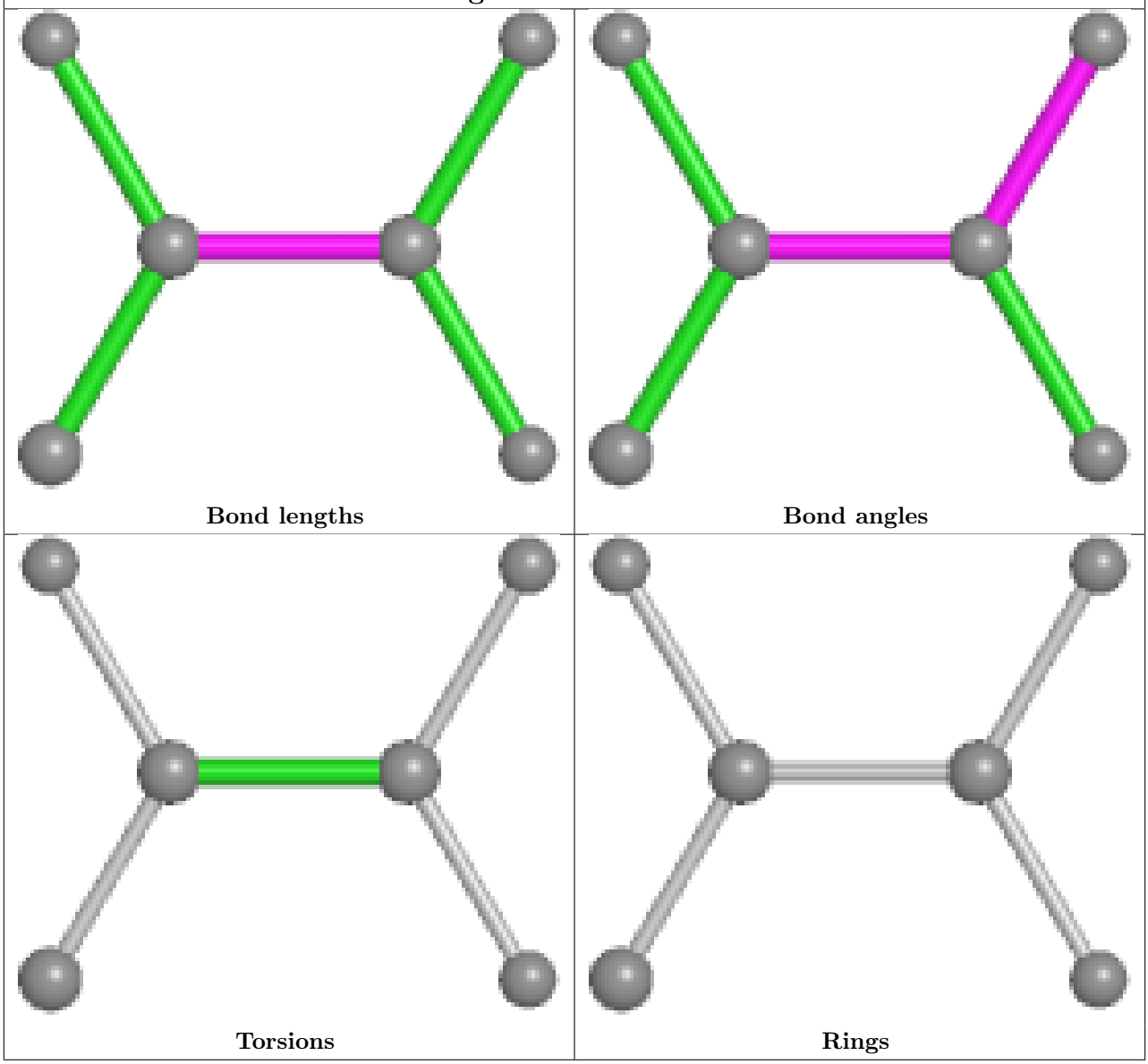
Ligand PYR D 302



Ligand PYR A 302



Ligand PYR B 302



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	276/298 (92%)	0.31	14 (5%) 28 34	16, 27, 56, 68	0
1	B	276/298 (92%)	0.57	30 (10%) 5 7	22, 35, 63, 97	0
1	C	277/298 (92%)	0.12	8 (2%) 51 57	16, 24, 42, 72	0
1	D	276/298 (92%)	0.16	9 (3%) 46 53	21, 31, 48, 70	0
All	All	1105/1192 (92%)	0.29	61 (5%) 25 31	16, 29, 56, 97	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	292	PHE	13.3
1	B	191	ALA	9.8
1	C	292	PHE	9.7
1	B	290	ALA	9.6
1	D	292	PHE	8.9
1	B	291	LEU	8.5
1	B	288	LEU	8.5
1	A	191	ALA	8.3
1	A	292	PHE	6.8
1	B	289	ASP	6.2
1	A	291	LEU	6.2
1	A	120	VAL	5.7
1	D	290	ALA	5.5
1	A	164	LEU	5.3
1	D	289	ASP	4.8
1	C	291	LEU	4.6
1	D	120	VAL	4.6
1	B	166	VAL	4.4
1	B	120	VAL	4.2
1	C	288	LEU	4.1
1	D	286	GLN	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	165	ALA	3.9
1	B	190	GLU	3.9
1	B	196	ALA	3.9
1	A	290	ALA	3.8
1	B	164	LEU	3.7
1	D	291	LEU	3.5
1	A	190	GLU	3.5
1	B	229	SER	3.5
1	C	289	ASP	3.5
1	B	201	PHE	3.4
1	C	120	VAL	3.4
1	C	286	GLN	3.3
1	B	173	ILE	3.2
1	B	169	LEU	3.1
1	B	193	THR	3.0
1	C	290	ALA	2.9
1	A	193	THR	2.9
1	B	170	GLN	2.9
1	D	285	GLU	2.8
1	A	168	GLY	2.8
1	B	133	ILE	2.7
1	C	282	HIS	2.6
1	B	195	LEU	2.6
1	A	165	ALA	2.6
1	B	198	TYR	2.5
1	B	194	GLU	2.5
1	A	196	ALA	2.4
1	B	284	PHE	2.4
1	B	219	THR	2.3
1	B	167	GLU	2.3
1	A	132	GLU	2.3
1	B	286	GLN	2.3
1	D	288	LEU	2.3
1	B	222	PHE	2.2
1	B	192	MET	2.2
1	D	281	TYR	2.2
1	B	132	GLU	2.2
1	A	169	LEU	2.2
1	B	228	ALA	2.1
1	A	206	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 (i)

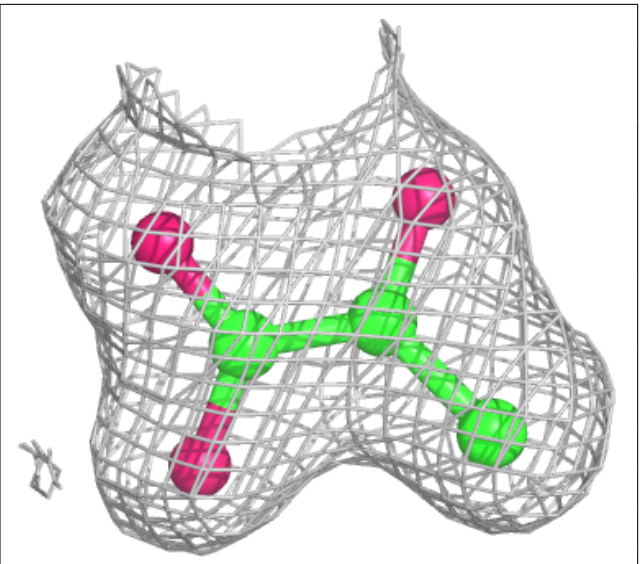
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

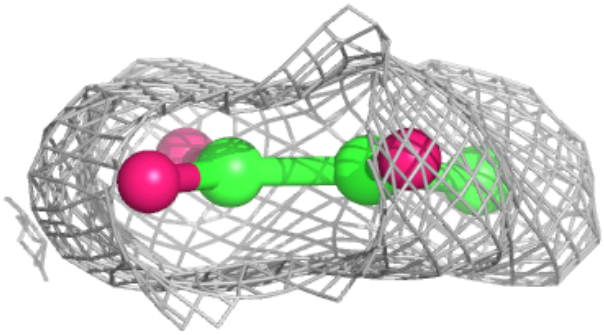
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	PYR	D	302	6/6	0.87	0.12	31,32,37,39	0
3	PYR	A	302	6/6	0.89	0.13	29,31,34,34	0
2	MG	B	301	1/1	0.91	0.07	44,44,44,44	0
3	PYR	B	302	6/6	0.92	0.16	32,36,37,41	0
3	PYR	C	302	6/6	0.95	0.09	21,23,26,29	0
2	MG	D	301	1/1	0.98	0.06	33,33,33,33	0
2	MG	A	301	1/1	0.98	0.05	33,33,33,33	0
2	MG	C	301	1/1	0.99	0.03	26,26,26,26	0

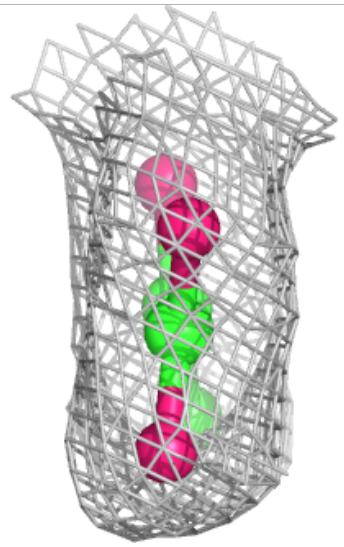
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around PYR D 302:

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

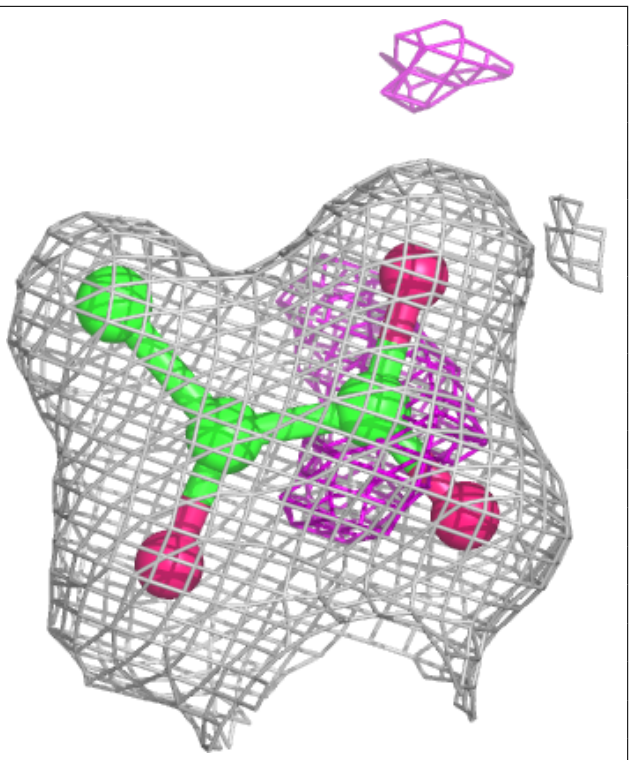


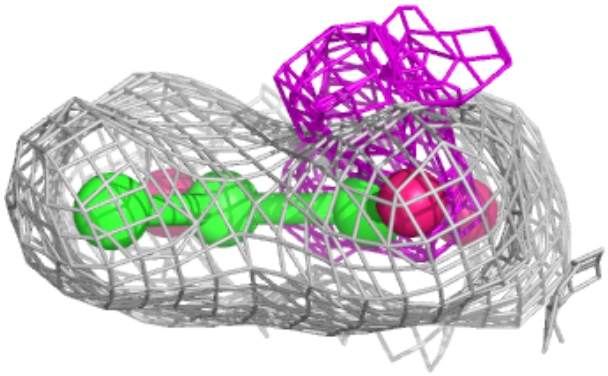


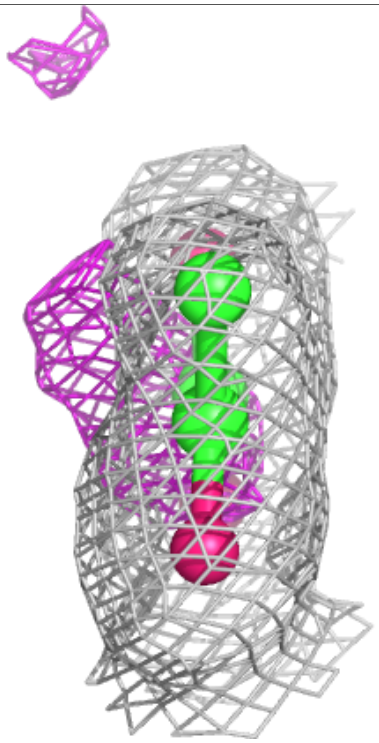


Electron density around PYR A 302:

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

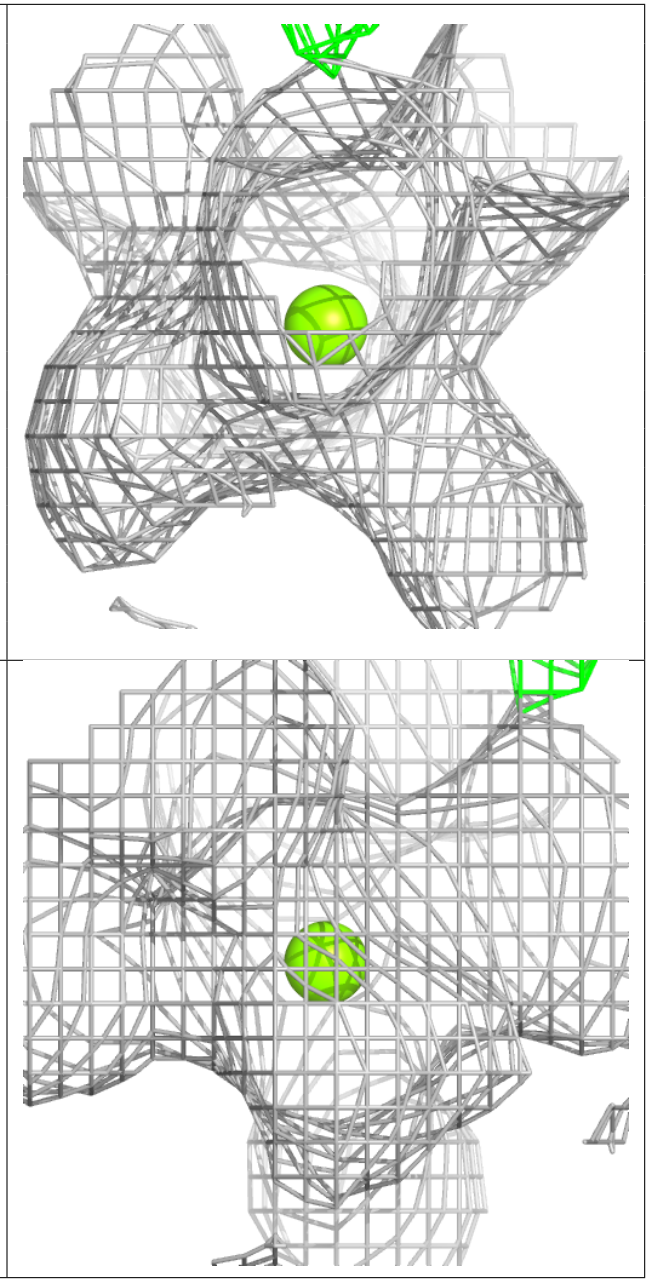






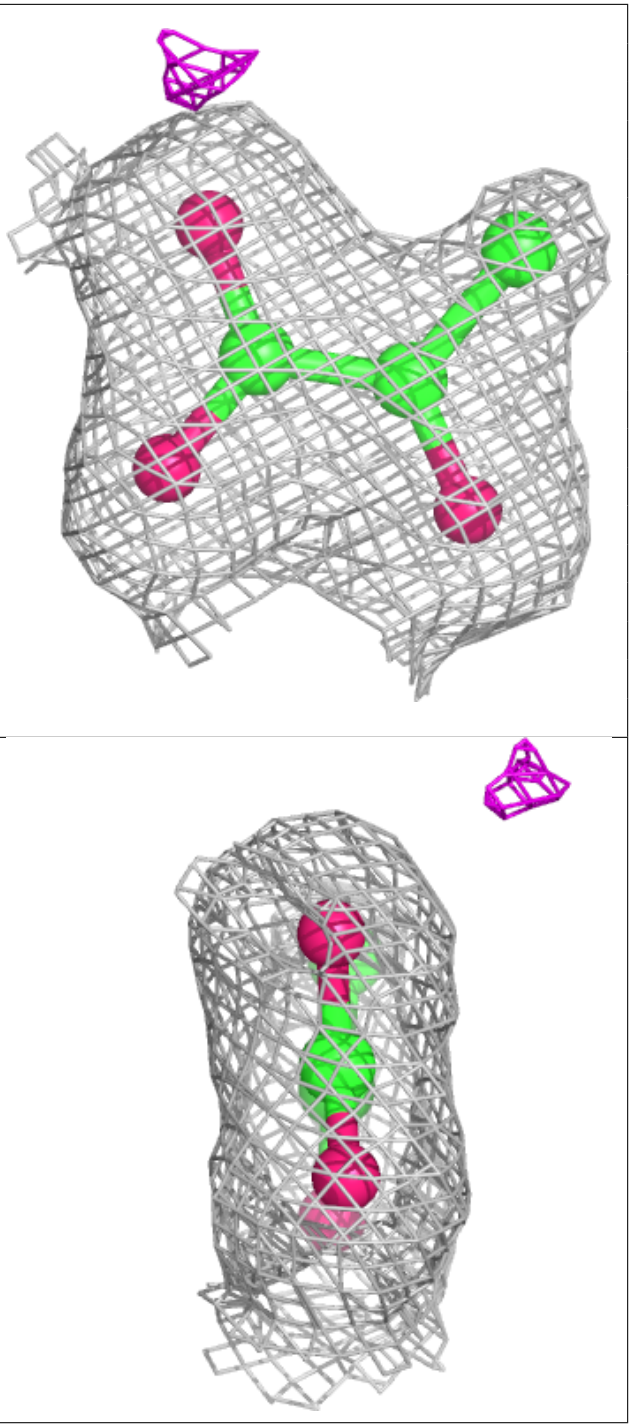
Electron density around MG B 301:

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



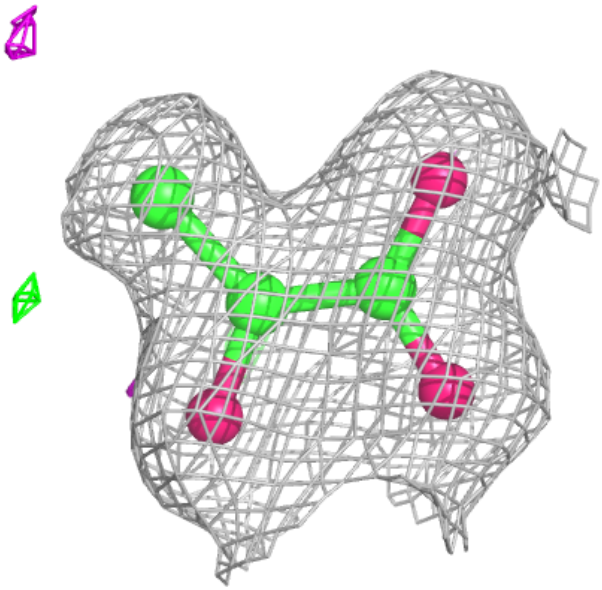
Electron density around PYR B 302:

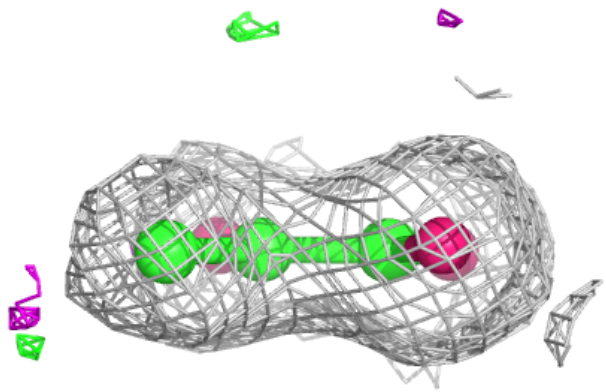
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

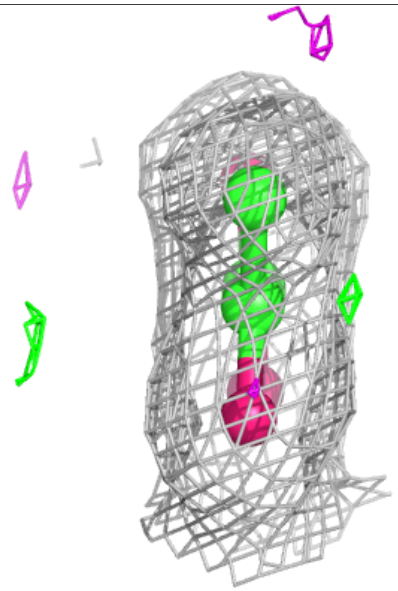


Electron density around PYR C 302:

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

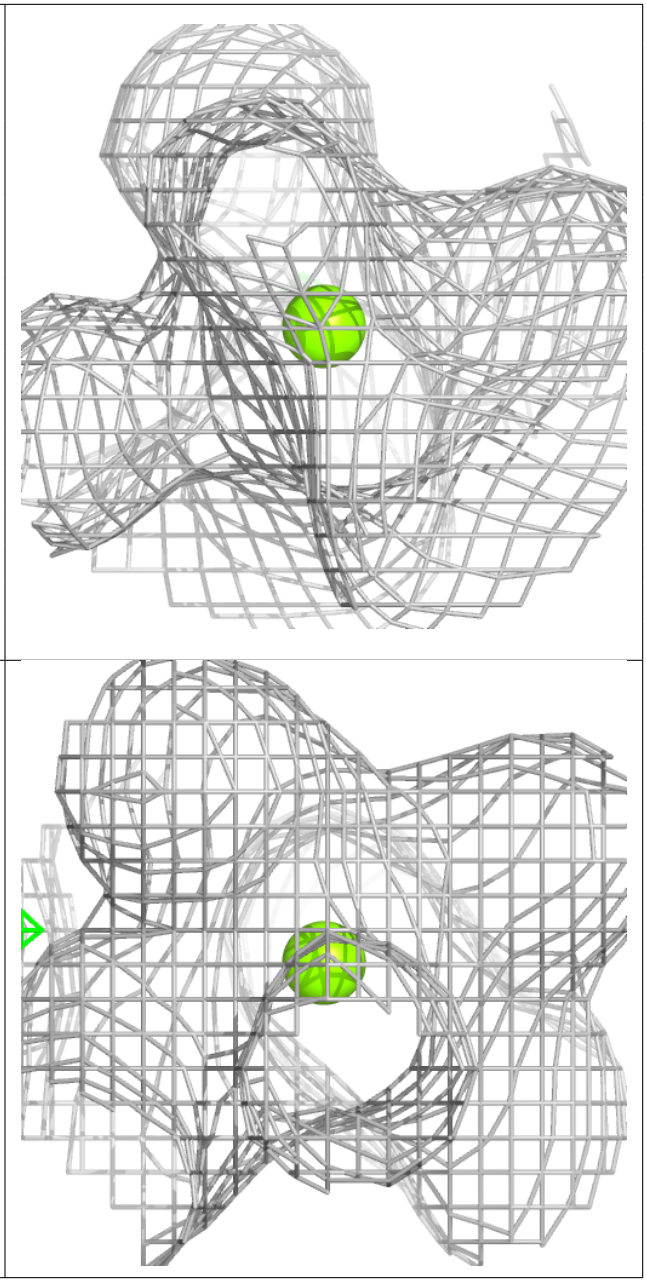






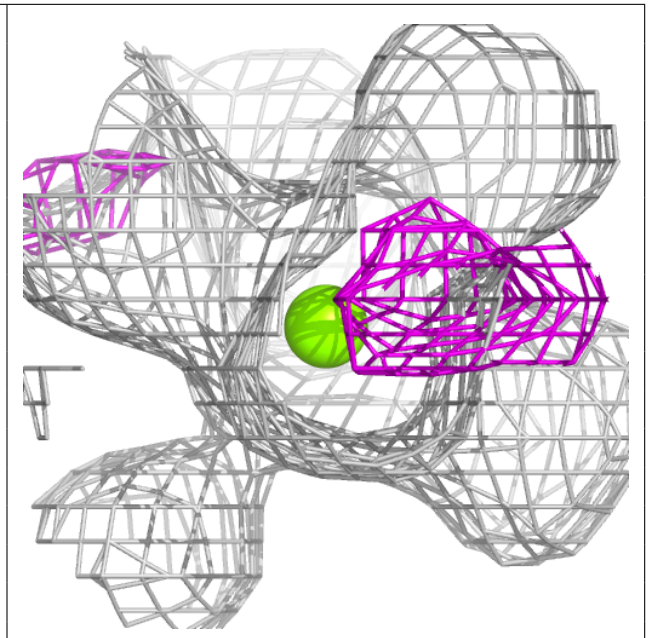
Electron density around MG D 301:

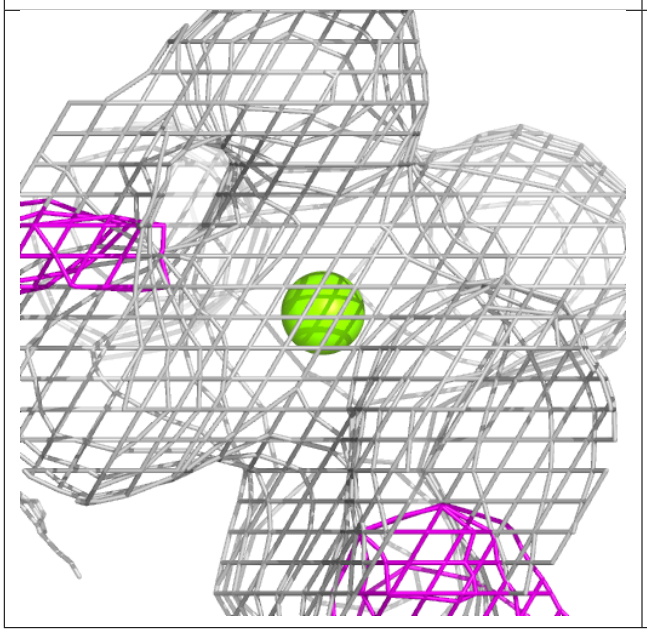
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

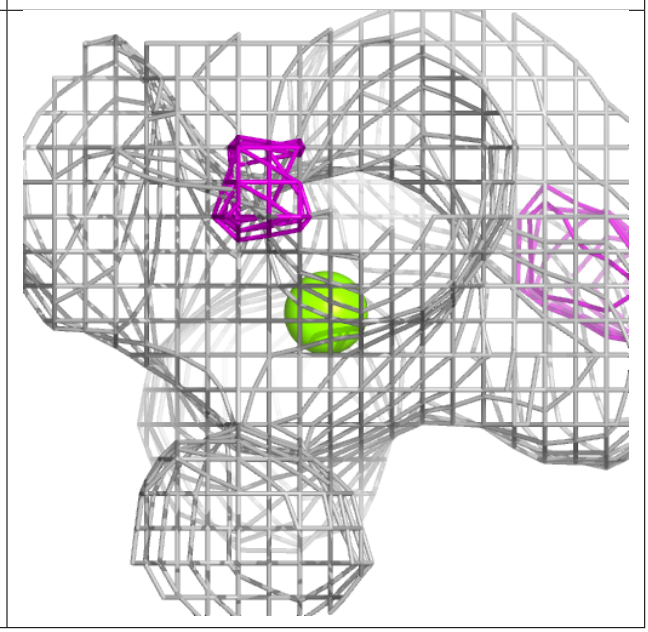


Electron density around MG A 301:

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

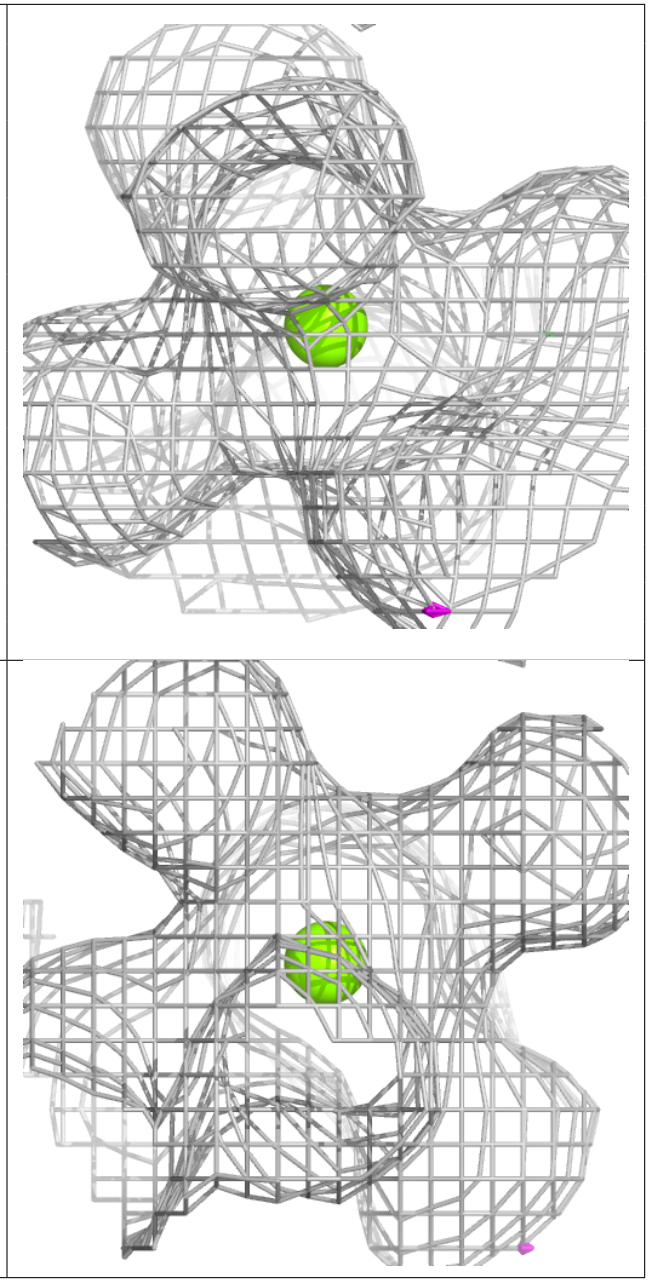






Electron density around MG C 301:

$2mF_o$-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.