



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2022 – 06:10 PM JST

PDB ID : 7FIQ  
Title : The crystal structure of mannose-bound beta-1,2-mannobiose phosphorylase from *Thermoanaerobacter* sp.  
Authors : Dai, L.; Chang, Z.; Yang, J.; Liu, W.; Yang, Y.; Chen, C.-C.; Zhang, L.; Huang, J.; Sun, Y.; Guo, R.-T.  
Deposited on : 2021-08-01  
Resolution : 2.22 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](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.

---

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.27
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)	:	2.27

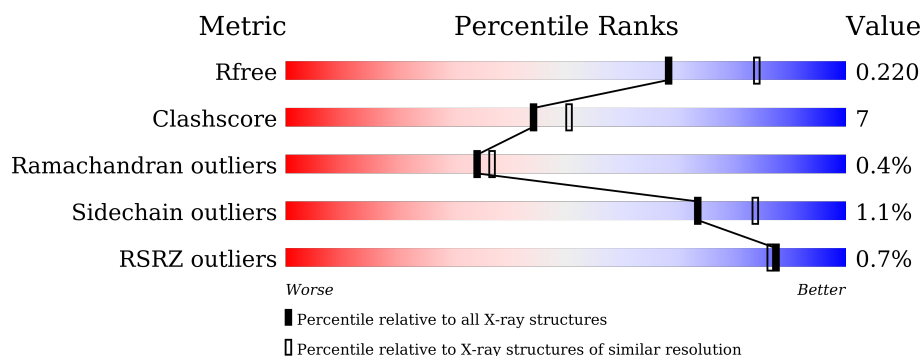
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.22 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	5912 (2.24-2.20)
Clashscore	141614	6646 (2.24-2.20)
Ramachandran outliers	138981	6543 (2.24-2.20)
Sidechain outliers	138945	6544 (2.24-2.20)
RSRZ outliers	127900	5797 (2.24-2.20)

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	313	 81% 15% . .
1	B	313	 82% 14% . .
1	C	313	 77% 19% .
1	D	313	 82% 14% .

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 11384 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 Beta-1,2-mannobiose phosphorylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	302	Total	C	N	O	S	0	0	0
			2485	1615	415	447	8			
1	B	302	Total	C	N	O	S	0	0	0
			2485	1615	415	447	8			
1	A	302	Total	C	N	O	S	0	0	0
			2485	1615	415	447	8			
1	D	302	Total	C	N	O	S	0	0	0
			2485	1615	415	447	8			

There are 44 discrepancies between the modelled and reference sequences:

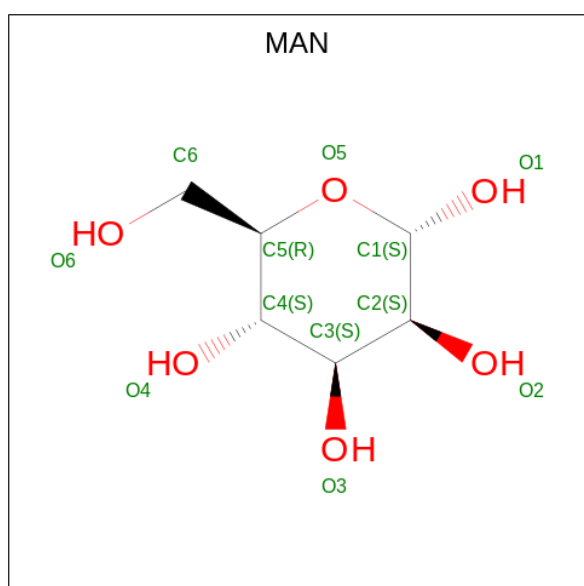
Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	GLY	-	expression tag	UNP B0K2C3
C	-9	ALA	-	expression tag	UNP B0K2C3
C	-8	GLY	-	expression tag	UNP B0K2C3
C	-7	ALA	-	expression tag	UNP B0K2C3
C	-6	GLY	-	expression tag	UNP B0K2C3
C	-5	ALA	-	expression tag	UNP B0K2C3
C	-4	GLY	-	expression tag	UNP B0K2C3
C	-3	ALA	-	expression tag	UNP B0K2C3
C	-2	GLY	-	expression tag	UNP B0K2C3
C	-1	ALA	-	expression tag	UNP B0K2C3
C	0	GLY	-	expression tag	UNP B0K2C3
B	-10	GLY	-	expression tag	UNP B0K2C3
B	-9	ALA	-	expression tag	UNP B0K2C3
B	-8	GLY	-	expression tag	UNP B0K2C3
B	-7	ALA	-	expression tag	UNP B0K2C3
B	-6	GLY	-	expression tag	UNP B0K2C3
B	-5	ALA	-	expression tag	UNP B0K2C3
B	-4	GLY	-	expression tag	UNP B0K2C3
B	-3	ALA	-	expression tag	UNP B0K2C3
B	-2	GLY	-	expression tag	UNP B0K2C3
B	-1	ALA	-	expression tag	UNP B0K2C3

*Continued on next page...*

Continued from previous page...

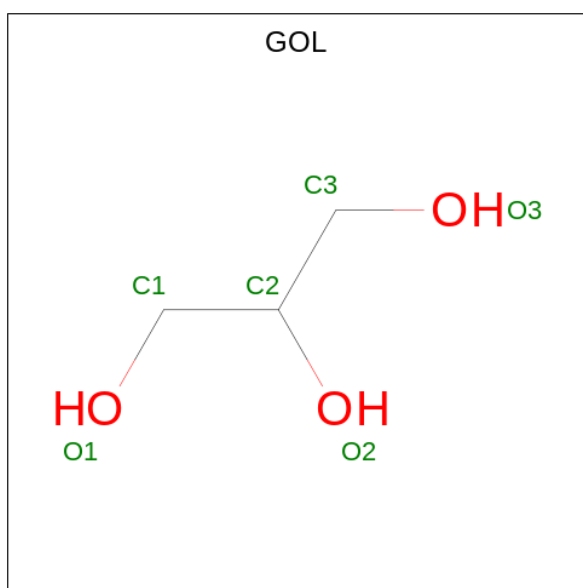
Chain	Residue	Modelled	Actual	Comment	Reference
B	0	GLY	-	expression tag	UNP B0K2C3
A	-10	GLY	-	expression tag	UNP B0K2C3
A	-9	ALA	-	expression tag	UNP B0K2C3
A	-8	GLY	-	expression tag	UNP B0K2C3
A	-7	ALA	-	expression tag	UNP B0K2C3
A	-6	GLY	-	expression tag	UNP B0K2C3
A	-5	ALA	-	expression tag	UNP B0K2C3
A	-4	GLY	-	expression tag	UNP B0K2C3
A	-3	ALA	-	expression tag	UNP B0K2C3
A	-2	GLY	-	expression tag	UNP B0K2C3
A	-1	ALA	-	expression tag	UNP B0K2C3
A	0	GLY	-	expression tag	UNP B0K2C3
D	-10	GLY	-	expression tag	UNP B0K2C3
D	-9	ALA	-	expression tag	UNP B0K2C3
D	-8	GLY	-	expression tag	UNP B0K2C3
D	-7	ALA	-	expression tag	UNP B0K2C3
D	-6	GLY	-	expression tag	UNP B0K2C3
D	-5	ALA	-	expression tag	UNP B0K2C3
D	-4	GLY	-	expression tag	UNP B0K2C3
D	-3	ALA	-	expression tag	UNP B0K2C3
D	-2	GLY	-	expression tag	UNP B0K2C3
D	-1	ALA	-	expression tag	UNP B0K2C3
D	0	GLY	-	expression tag	UNP B0K2C3

- Molecule 2 is alpha-D-mannopyranose (three-letter code: MAN) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>) (labeled as "Ligand of Interest" by depositor).



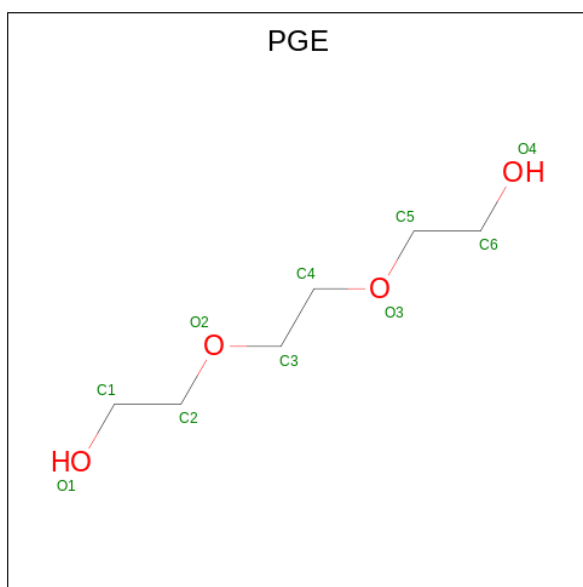
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			12	6	6		
2	B	1	Total	C	O	0	0
			12	6	6		
2	A	1	Total	C	O	0	0
			12	6	6		
2	D	1	Total	C	O	0	0
			12	6	6		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



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

- Molecule 4 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula:  $C_6H_{14}O_4$ ).

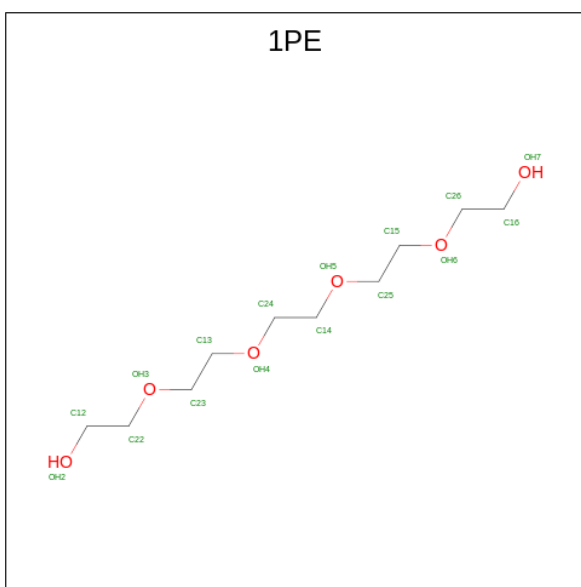


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			10	6	4		
4	D	1	Total	C	O	0	0
			10	6	4		

- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	4	Total	Zn	0	0
			4	4		
5	B	5	Total	Zn	0	0
			5	5		
5	A	5	Total	Zn	0	0
			5	5		
5	D	3	Total	Zn	0	0
			3	3		

- Molecule 6 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: C<sub>10</sub>H<sub>22</sub>O<sub>6</sub>).



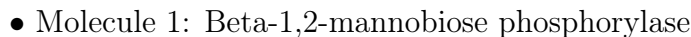
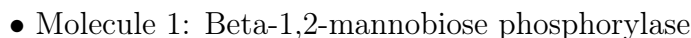
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			16	10	6		

- Molecule 7 is water.

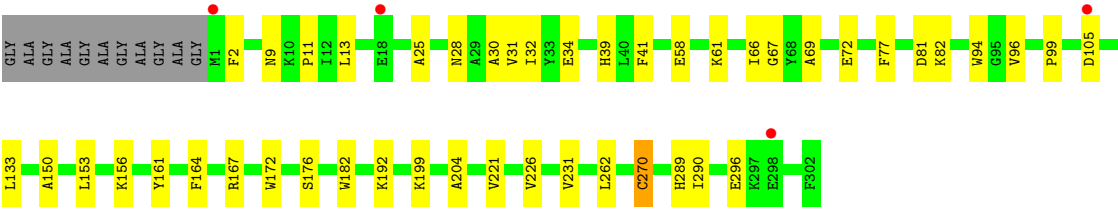
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	C	373	Total O 373 373	0	0
7	B	342	Total O 342 342	0	0
7	A	326	Total O 326 326	0	0
7	D	296	Total O 296 296	0	0



- Molecule 1: Beta-1,2-mannobiose phosphorylase







## 4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	203.19Å 203.19Å 77.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.85 – 2.22 36.03 – 2.22	Depositor EDS
% Data completeness (in resolution range)	97.6 (34.85-2.22) 84.4 (36.03-2.22)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.96 (at 2.22Å)	Xtriage
Refinement program	PHENIX 1.17.1	Depositor
R, $R_{free}$	0.173 , 0.220 0.173 , 0.220	Depositor DCC
$R_{free}$ test set	1928 reflections (2.57%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	18.1	Xtriage
Anisotropy	0.276	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 47.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.025 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	11384	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% 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: 1PE, MAN, GOL, ZN, PGE

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	0.45	1/2556 (0.0%)	0.62	1/3465 (0.0%)
1	B	0.45	0/2556	0.58	0/3465
1	C	0.46	0/2556	0.59	0/3465
1	D	0.43	1/2556 (0.0%)	0.56	0/3465
All	All	0.45	2/10224 (0.0%)	0.59	1/13860 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	270	CYS	CB-SG	-6.92	1.70	1.82
1	D	270	CYS	CB-SG	-5.72	1.72	1.81

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	9	ASN	C-N-CA	5.01	134.22	121.70

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	2485	0	2452	28	1

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2485	0	2452	29	0
1	C	2485	0	2452	43	1
1	D	2485	0	2452	30	0
2	A	12	0	12	0	0
2	B	12	0	12	0	0
2	C	12	0	12	0	0
2	D	12	0	12	0	0
3	C	6	0	8	3	0
4	C	10	0	14	1	0
4	D	10	0	14	0	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0
5	C	4	0	0	0	1
5	D	3	0	0	0	0
6	A	16	0	22	3	0
7	A	326	0	0	6	1
7	B	342	0	0	3	0
7	C	373	0	0	8	2
7	D	296	0	0	6	0
All	All	11384	0	9914	131	3

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (131) 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:C:296:GLU:OE1	1:C:298:GLU:HB2	1.62	0.99
1:C:296:GLU:OE1	1:C:298:GLU:CB	2.15	0.94
1:A:189:MET:SD	7:A:815:HOH:O	2.34	0.85
1:B:277:ASP:O	7:B:501:HOH:O	1.95	0.84
1:C:140:ARG:HB2	3:C:402:GOL:H2	1.62	0.79
1:D:9:ASN:OD1	7:D:501:HOH:O	2.07	0.73
1:C:17:LYS:NZ	7:C:501:HOH:O	2.06	0.70
1:A:1:MET:SD	7:A:823:HOH:O	2.50	0.70
1:D:72:GLU:OE2	7:D:502:HOH:O	2.13	0.66
1:D:32:ILE:HD11	1:D:133:LEU:HD11	1.78	0.66
1:C:296:GLU:OE1	1:C:298:GLU:HB3	1.95	0.66
1:C:81:ASP:OD2	7:C:503:HOH:O	2.14	0.65
1:C:296:GLU:CD	1:C:298:GLU:H	1.99	0.65
1:C:180:VAL:HG13	3:C:402:GOL:H12	1.78	0.64

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:LYS:NZ	7:A:505:HOH:O	2.29	0.64
1:A:1:MET:SD	7:A:799:HOH:O	2.55	0.64
1:B:231:VAL:HG13	1:B:250:ILE:HD13	1.80	0.64
1:C:101:ILE:HD11	1:C:110:MET:HE3	1.83	0.59
1:C:296:GLU:CD	1:C:298:GLU:HB2	2.23	0.59
1:A:54:GLU:OE2	7:A:501:HOH:O	2.17	0.59
1:B:2:PHE:CZ	1:B:231:VAL:HG21	2.38	0.58
1:C:226:VAL:HG13	1:C:265:ASN:HB3	1.86	0.58
1:D:2:PHE:CZ	1:D:231:VAL:HG11	2.39	0.58
1:C:31:VAL:HG13	1:C:273:VAL:HG13	1.85	0.57
1:B:105:ASP:O	7:B:502:HOH:O	2.17	0.57
1:C:72:GLU:OE1	7:C:504:HOH:O	2.17	0.57
1:B:61:LYS:HA	1:B:94:TRP:CH2	2.41	0.56
1:A:31:VAL:HG13	1:A:273:VAL:HG13	1.87	0.56
1:A:11:PRO:HB3	1:A:289:HIS:HB3	1.86	0.56
1:D:167:ARG:HB2	1:D:172:TRP:NE1	2.21	0.56
1:C:1:MET:HE3	1:C:247:LYS:HB3	1.88	0.55
1:A:214:TRP:HB2	1:A:234:LEU:HB2	1.88	0.55
1:D:66:ILE:HD12	1:D:99:PRO:HG3	1.88	0.55
1:B:25:ALA:HB2	1:B:262:LEU:HD12	1.89	0.55
1:D:58:GLU:OE2	1:D:61:LYS:HD2	2.07	0.55
1:D:199:LYS:NZ	7:D:512:HOH:O	2.39	0.55
1:A:176:SER:HB2	1:A:182:TRP:CE3	2.41	0.54
1:A:61:LYS:HA	1:A:94:TRP:CH2	2.43	0.54
1:A:66:ILE:HG13	1:A:96:VAL:HG23	1.90	0.53
1:A:68:TYR:CD2	6:A:402:1PE:H152	2.43	0.53
1:D:99:PRO:O	1:D:270:CYS:HB3	2.08	0.52
1:B:217:ILE:HG12	1:B:231:VAL:HG12	1.91	0.52
1:D:58:GLU:OE2	7:D:503:HOH:O	2.19	0.52
1:D:61:LYS:HA	1:D:94:TRP:CH2	2.44	0.51
1:C:23:LYS:NZ	7:C:508:HOH:O	2.29	0.51
1:B:82:LYS:NZ	7:B:505:HOH:O	2.30	0.51
1:C:1:MET:HE2	7:C:748:HOH:O	2.11	0.50
1:B:27:PHE:CD2	1:B:287:ASP:HA	2.46	0.50
1:D:105:ASP:OD1	7:D:504:HOH:O	2.19	0.50
1:C:2:PHE:CZ	1:C:231:VAL:HG11	2.47	0.50
1:C:11:PRO:HB3	1:C:289:HIS:HB3	1.94	0.50
1:D:32:ILE:HD13	1:D:34:GLU:HG2	1.94	0.49
1:B:11:PRO:HB3	1:B:289:HIS:HB3	1.95	0.49
1:D:66:ILE:HG12	1:D:96:VAL:HG23	1.94	0.49
1:A:134:LYS:NZ	6:A:402:1PE:H162	2.28	0.49

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:176:SER:HB2	1:C:182:TRP:CE3	2.47	0.49
1:C:189:MET:HE2	1:C:201:ILE:HD12	1.94	0.48
1:C:36:ASN:ND2	7:C:502:HOH:O	2.13	0.48
1:B:61:LYS:HG3	1:B:94:TRP:CE2	2.48	0.48
1:C:91:GLN:NE2	1:C:118:ARG:HH12	2.11	0.48
1:C:134:LYS:HG2	4:C:403:PGE:H3	1.94	0.48
1:B:91:GLN:O	1:B:124:ARG:NH1	2.41	0.48
1:B:150:ALA:HA	1:B:164:PHE:O	2.14	0.48
1:C:278:MET:HE3	1:C:294:VAL:HB	1.96	0.48
1:B:214:TRP:HB2	1:B:234:LEU:HB2	1.97	0.47
1:D:25:ALA:HB2	1:D:262:LEU:HD12	1.97	0.47
1:D:150:ALA:HA	1:D:164:PHE:O	2.14	0.47
1:C:268:PHE:O	1:C:284:GLY:N	2.48	0.47
1:B:186:LYS:HE3	1:B:240:SER:HB2	1.95	0.47
1:C:1:MET:HE1	1:C:247:LYS:H	1.80	0.46
1:D:11:PRO:HB3	1:D:289:HIS:HB3	1.97	0.46
1:D:30:ALA:HB2	1:D:99:PRO:O	2.15	0.46
1:A:99:PRO:O	1:A:270:CYS:HB3	2.15	0.46
1:A:167:ARG:HB2	1:A:172:TRP:NE1	2.30	0.46
1:C:150:ALA:HA	1:C:164:PHE:O	2.14	0.46
1:C:44:ALA:O	1:C:64:SER:HA	2.16	0.46
1:C:101:ILE:HD11	1:C:110:MET:CE	2.44	0.45
1:B:101:ILE:HD11	1:B:110:MET:CE	2.45	0.45
1:D:192:LYS:NZ	7:D:519:HOH:O	2.48	0.45
1:D:32:ILE:HD11	1:D:133:LEU:CD1	2.44	0.45
1:B:21:TRP:CD2	1:B:83:PRO:HD3	2.51	0.45
1:D:31:VAL:HA	1:D:39:HIS:O	2.17	0.45
1:D:61:LYS:HG3	1:D:94:TRP:CE2	2.52	0.45
1:C:1:MET:CE	1:C:247:LYS:HB3	2.46	0.45
1:C:33:TYR:CZ	1:C:36:ASN:HA	2.52	0.45
1:C:66:ILE:HG12	1:C:96:VAL:HG23	1.99	0.44
1:D:156:LYS:HE3	1:D:161:TYR:CZ	2.53	0.44
1:A:25:ALA:HB2	1:A:262:LEU:HD12	1.99	0.43
3:C:402:GOL:H31	7:C:686:HOH:O	2.18	0.43
1:A:27:PHE:CD2	1:A:287:ASP:HA	2.53	0.43
1:A:21:TRP:CD2	1:A:83:PRO:HD3	2.54	0.43
1:A:127:MET:CE	1:A:140:ARG:HD2	2.49	0.43
1:A:268:PHE:O	1:A:284:GLY:N	2.50	0.43
1:C:21:TRP:CE2	1:C:83:PRO:HD3	2.54	0.43
1:C:94:TRP:CD1	1:C:117:GLY:HA3	2.54	0.43
1:B:7:LEU:HD23	1:B:7:LEU:HA	1.75	0.43

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:PHE:CZ	1:A:231:VAL:HG11	2.54	0.43
1:C:4:LEU:HD23	1:C:295:ILE:HB	2.01	0.42
1:B:107:LYS:HD3	1:B:129:TRP:CZ3	2.54	0.42
1:D:69:ALA:HB1	1:D:77:PHE:HB3	2.00	0.42
1:D:153:LEU:HD12	1:D:153:LEU:HA	1.91	0.42
1:A:25:ALA:HB3	1:A:45:SER:HB3	2.02	0.42
1:C:99:PRO:O	1:C:270:CYS:HB3	2.19	0.42
1:B:107:LYS:HD3	1:B:129:TRP:CE3	2.54	0.42
1:A:48:LYS:NZ	7:A:506:HOH:O	2.42	0.42
1:C:61:LYS:HA	1:C:94:TRP:CH2	2.55	0.42
1:A:150:ALA:HA	1:A:164:PHE:O	2.19	0.42
1:A:21:TRP:CE2	1:A:83:PRO:HD3	2.55	0.42
1:C:278:MET:HE3	1:C:278:MET:HB3	1.91	0.41
1:C:179:LEU:HD23	1:C:179:LEU:HA	1.89	0.41
1:C:27:PHE:CD2	1:C:287:ASP:HA	2.55	0.41
1:D:41:PHE:HA	1:D:67:GLY:O	2.21	0.41
1:B:173:ILE:HG22	1:B:186:LYS:HB3	2.01	0.41
1:A:232:ALA:HB2	1:A:245:ARG:HG3	2.01	0.41
1:B:279:TYR:O	1:B:294:VAL:HA	2.21	0.41
1:D:176:SER:HB2	1:D:182:TRP:CE3	2.56	0.41
1:B:6:ARG:NH2	1:B:252:GLU:O	2.54	0.41
1:B:99:PRO:O	1:B:270:CYS:HB3	2.20	0.41
1:C:29:ALA:HA	1:C:42:TYR:HA	2.03	0.41
1:B:168:MET:HA	1:B:169:PRO:HA	1.90	0.41
1:A:6:ARG:NH2	1:A:252:GLU:O	2.54	0.41
1:D:81:ASP:O	1:D:82:LYS:HD2	2.21	0.41
1:B:176:SER:HB2	1:B:182:TRP:CE3	2.55	0.40
1:A:68:TYR:CZ	6:A:402:1PE:H262	2.56	0.40
1:B:129:TRP:CH2	1:B:137:LYS:HE3	2.56	0.40
1:C:30:ALA:HB2	1:C:99:PRO:O	2.22	0.40
1:C:155:GLU:HG3	7:C:576:HOH:O	2.21	0.40
1:B:163:LEU:HB2	1:B:182:TRP:CZ3	2.57	0.40
1:B:216:LEU:HD13	1:B:234:LEU:HD11	2.02	0.40
1:D:13:LEU:HB2	1:D:290:ILE:HB	2.04	0.40
1:D:221:VAL:HA	1:D:226:VAL:O	2.22	0.40

All (3) 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:C:54:GLU:OE2	5:C:407:ZN:ZN[6_554]	1.68	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:GLU:OE1	7:C:501:HOH:O[6_554]	2.09	0.11
7:C:602:HOH:O	7:A:510:HOH:O[6_555]	2.14	0.06

## 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	300/313 (96%)	287 (96%)	13 (4%)	0	100	100
1	B	300/313 (96%)	288 (96%)	11 (4%)	1 (0%)	41	45
1	C	300/313 (96%)	289 (96%)	9 (3%)	2 (1%)	22	21
1	D	300/313 (96%)	284 (95%)	14 (5%)	2 (1%)	22	21
All	All	1200/1252 (96%)	1148 (96%)	47 (4%)	5 (0%)	34	37

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	28	ASN
1	C	204	ALA
1	B	28	ASN
1	D	204	ALA
1	C	28	ASN

### 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	267/267 (100%)	261 (98%)	6 (2%)	52	64
1	B	267/267 (100%)	265 (99%)	2 (1%)	84	91
1	C	267/267 (100%)	264 (99%)	3 (1%)	73	84
1	D	267/267 (100%)	266 (100%)	1 (0%)	91	95
All	All	1068/1068 (100%)	1056 (99%)	12 (1%)	73	84

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	27	PHE
1	C	124	ARG
1	C	137	LYS
1	B	124	ARG
1	B	277	ASP
1	A	1	MET
1	A	55	LYS
1	A	124	ARG
1	A	131	ASP
1	A	160	LYS
1	A	216	LEU
1	D	296	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	91	GLN
1	D	276	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry

Of 25 ligands modelled in this entry, 17 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	$\# Z  > 2$
2	MAN	B	401	-	12,12,12	0.59	0	17,17,17	0.81	0
4	PGE	D	402	-	9,9,9	0.45	0	8,8,8	0.69	0
6	1PE	A	402	-	15,15,15	0.48	0	14,14,14	0.50	0
2	MAN	C	401	-	12,12,12	0.53	0	17,17,17	1.93	2 (11%)
4	PGE	C	403	-	9,9,9	0.41	0	8,8,8	0.56	0
3	GOL	C	402	-	5,5,5	0.60	0	5,5,5	0.50	0
2	MAN	D	401	-	12,12,12	0.51	0	17,17,17	0.63	0
2	MAN	A	401	-	12,12,12	0.56	0	17,17,17	0.65	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
2	MAN	B	401	-	-	0/2/22/22	0/1/1/1
4	PGE	D	402	-	-	6/7/7/7	-
6	1PE	A	402	-	-	6/13/13/13	-
2	MAN	C	401	-	-	0/2/22/22	0/1/1/1
4	PGE	C	403	-	-	2/7/7/7	-
3	GOL	C	402	-	-	4/4/4/4	-
2	MAN	D	401	-	-	0/2/22/22	0/1/1/1
2	MAN	A	401	-	-	0/2/22/22	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	MAN	O5-C1-C2	4.94	119.10	110.28
2	C	401	MAN	C1-O5-C5	4.80	122.71	113.66

There are no chirality outliers.

All (18) torsion outliers are listed below:

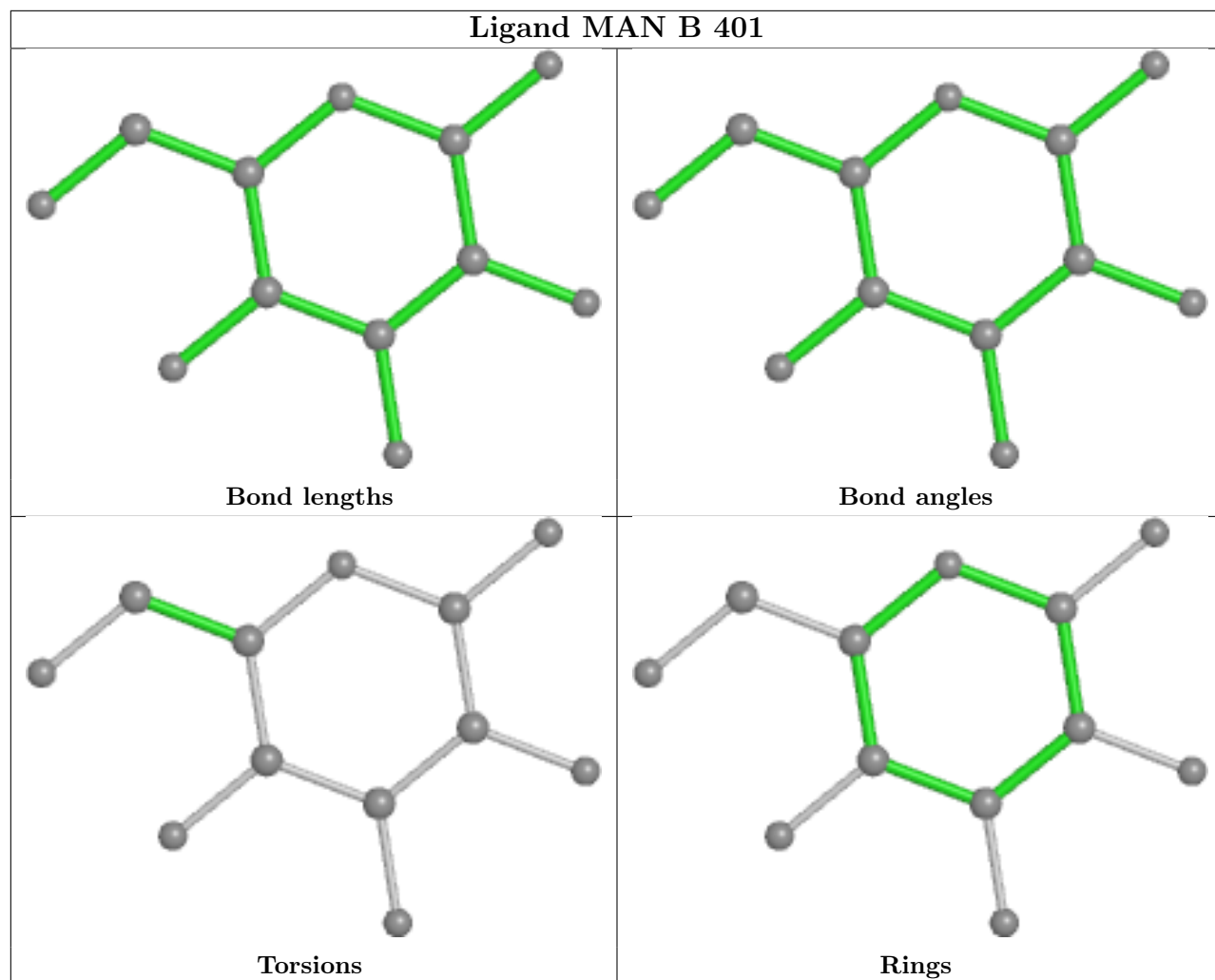
Mol	Chain	Res	Type	Atoms
3	C	402	GOL	O1-C1-C2-C3
3	C	402	GOL	C1-C2-C3-O3
6	A	402	1PE	OH5-C14-C24-OH4
3	C	402	GOL	O1-C1-C2-O2
6	A	402	1PE	OH2-C12-C22-OH3
6	A	402	1PE	OH7-C16-C26-OH6
6	A	402	1PE	OH6-C15-C25-OH5
4	C	403	PGE	C4-C3-O2-C2
6	A	402	1PE	C24-C14-OH5-C25
3	C	402	GOL	O2-C2-C3-O3
4	D	402	PGE	C1-C2-O2-C3
4	D	402	PGE	C3-C4-O3-C5
6	A	402	1PE	C16-C26-OH6-C15
4	C	403	PGE	C3-C4-O3-C5
4	D	402	PGE	O1-C1-C2-O2
4	D	402	PGE	O3-C5-C6-O4
4	D	402	PGE	C6-C5-O3-C4
4	D	402	PGE	O2-C3-C4-O3

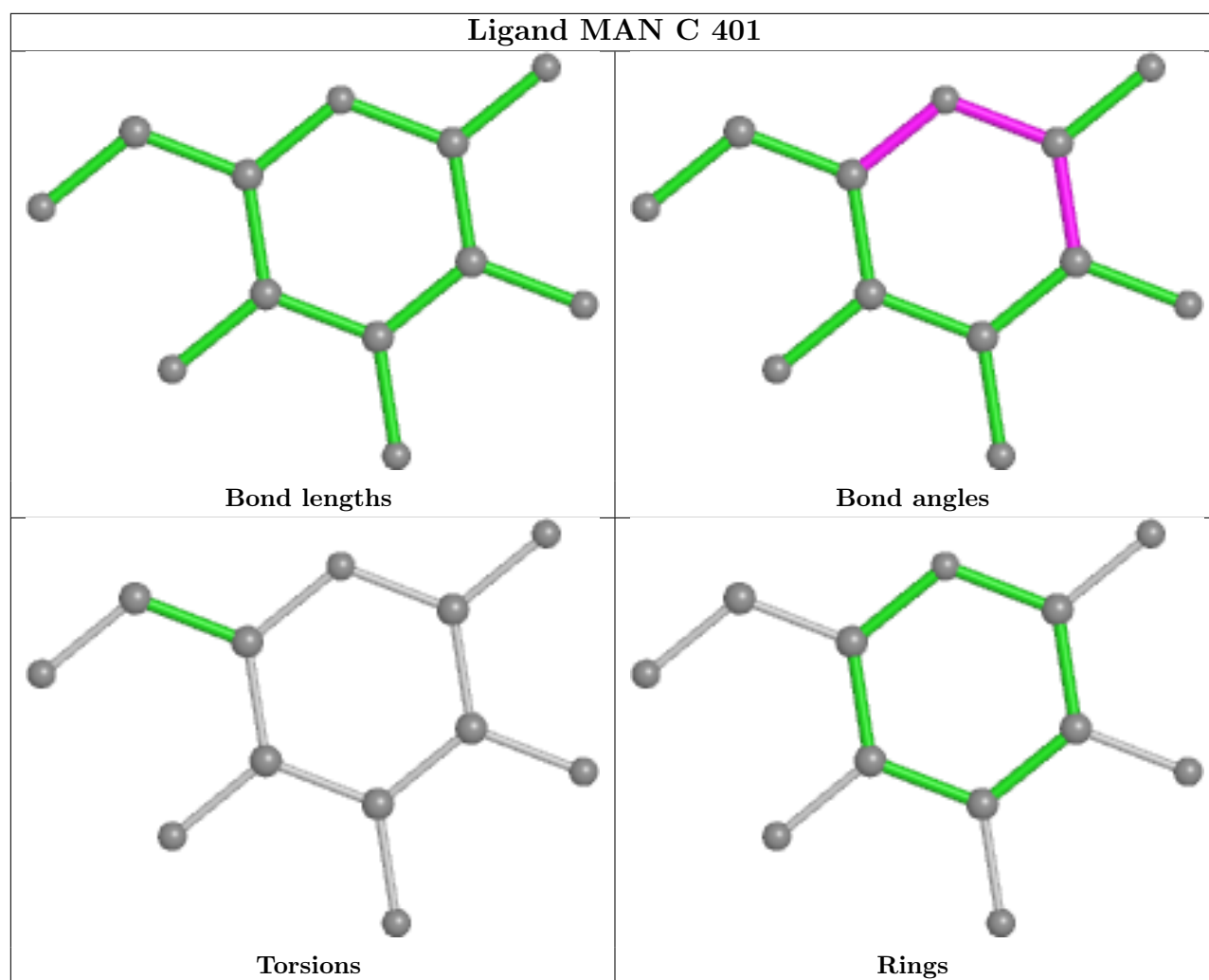
There are no ring outliers.

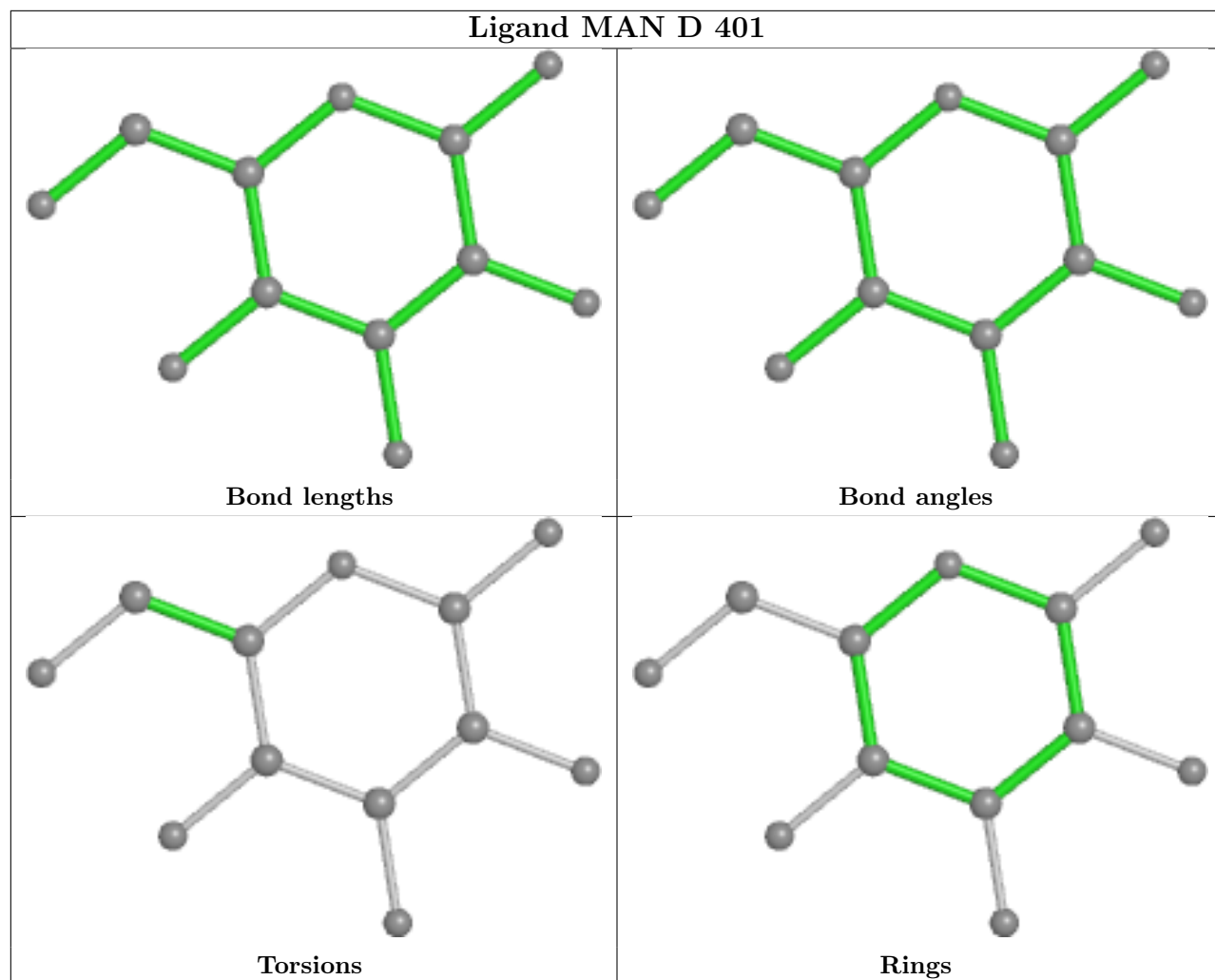
3 monomers are involved in 7 short contacts:

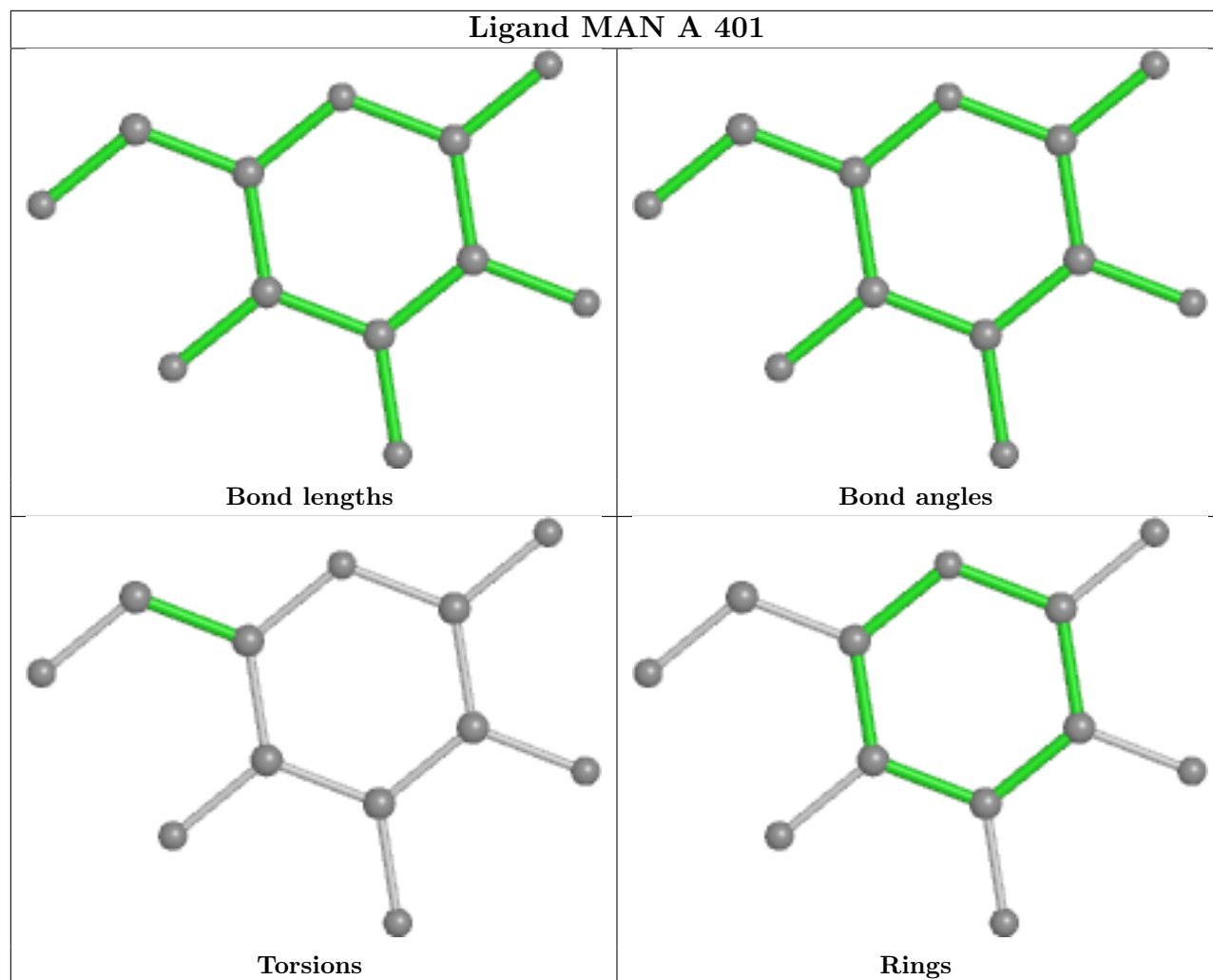
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	402	1PE	3	0
4	C	403	PGE	1	0
3	C	402	GOL	3	0

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









## 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	302/313 (96%)	-0.36	1 (0%) 94 93	14, 21, 36, 47	0
1	B	302/313 (96%)	-0.45	3 (0%) 82 81	14, 22, 36, 58	0
1	C	302/313 (96%)	-0.49	0 100 100	13, 18, 33, 48	0
1	D	302/313 (96%)	-0.21	4 (1%) 77 75	16, 26, 45, 61	0
All	All	1208/1252 (96%)	-0.38	8 (0%) 87 86	13, 22, 39, 61	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	89	ILE	3.6
1	B	1	MET	2.6
1	D	1	MET	2.3
1	D	105	ASP	2.2
1	A	90	PRO	2.2
1	D	18	GLU	2.1
1	B	90	PRO	2.1
1	D	298	GLU	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 ⓘ

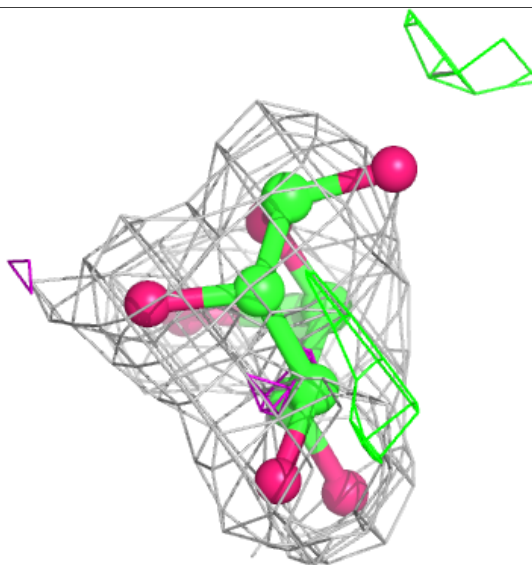
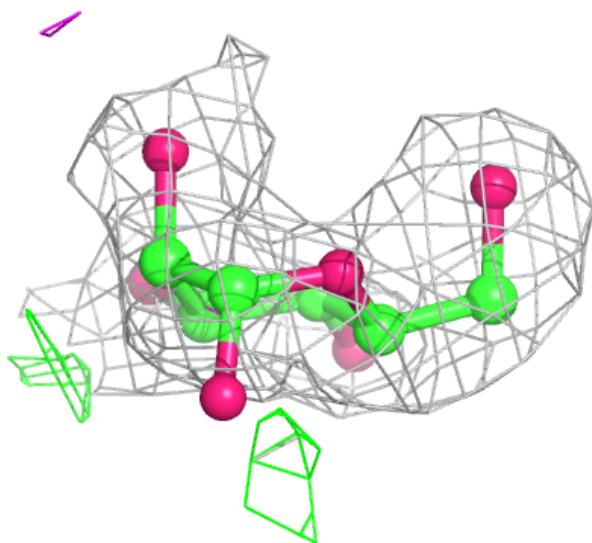
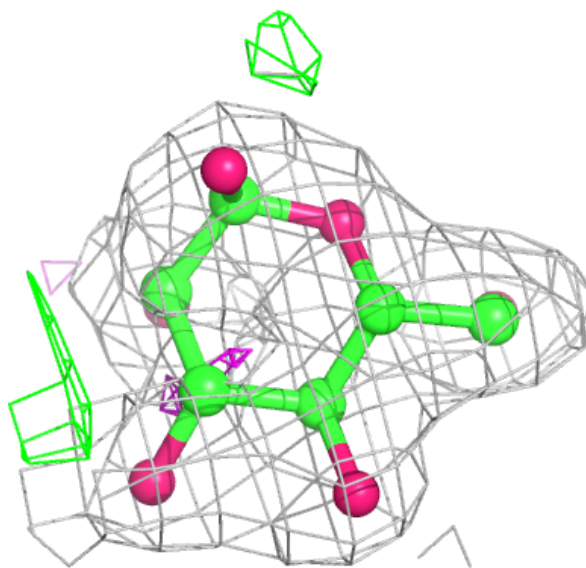
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
2	MAN	B	401	12/12	0.85	0.18	26,34,41,50	0
3	GOL	C	402	6/6	0.86	0.24	24,26,36,36	0
2	MAN	C	401	12/12	0.91	0.12	20,25,36,43	0
4	PGE	D	402	10/10	0.91	0.20	33,43,46,46	0
5	ZN	A	405	1/1	0.91	0.08	67,67,67,67	0
6	1PE	A	402	16/16	0.91	0.13	27,33,40,43	0
5	ZN	A	407	1/1	0.92	0.06	72,72,72,72	0
5	ZN	D	403	1/1	0.92	0.08	66,66,66,66	0
4	PGE	C	403	10/10	0.92	0.16	21,30,39,42	0
2	MAN	A	401	12/12	0.93	0.11	22,28,33,42	0
5	ZN	C	406	1/1	0.94	0.11	71,71,71,71	0
2	MAN	D	401	12/12	0.95	0.12	19,24,28,32	0
5	ZN	A	406	1/1	0.97	0.04	55,55,55,55	0
5	ZN	B	406	1/1	0.98	0.07	23,23,23,23	0
5	ZN	D	404	1/1	0.98	0.06	25,25,25,25	0
5	ZN	D	405	1/1	0.98	0.06	43,43,43,43	0
5	ZN	B	402	1/1	0.98	0.02	30,30,30,30	0
5	ZN	C	404	1/1	0.99	0.06	20,20,20,20	0
5	ZN	B	403	1/1	0.99	0.06	38,38,38,38	0
5	ZN	B	404	1/1	0.99	0.02	30,30,30,30	0
5	ZN	B	405	1/1	0.99	0.03	27,27,27,27	0
5	ZN	C	407	1/1	0.99	0.03	30,30,30,30	0
5	ZN	A	403	1/1	0.99	0.07	23,23,23,23	0
5	ZN	A	404	1/1	0.99	0.02	21,21,21,21	0
5	ZN	C	405	1/1	1.00	0.04	25,25,25,25	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.

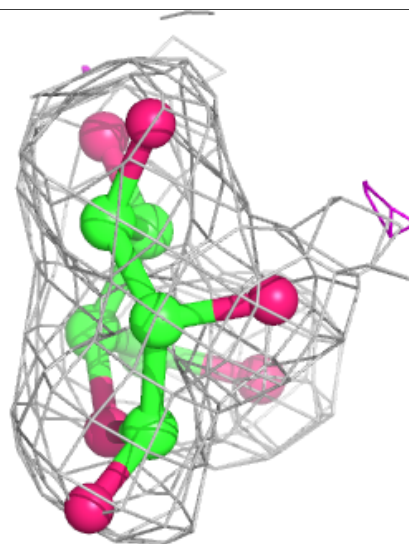
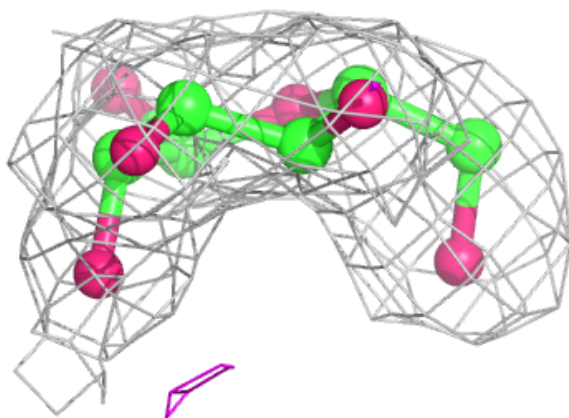
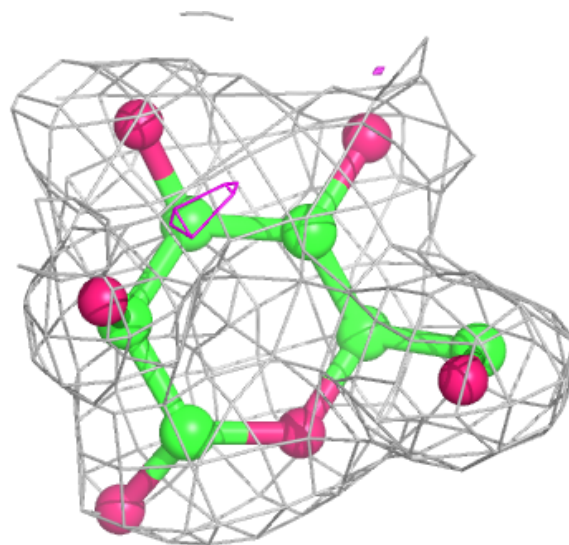
**Electron density around MAN B 401:**

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



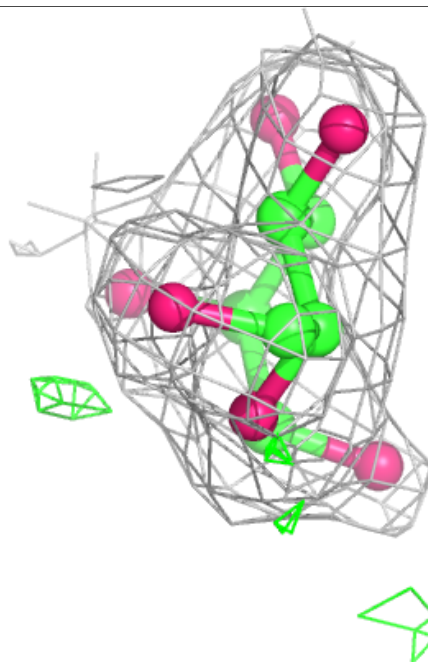
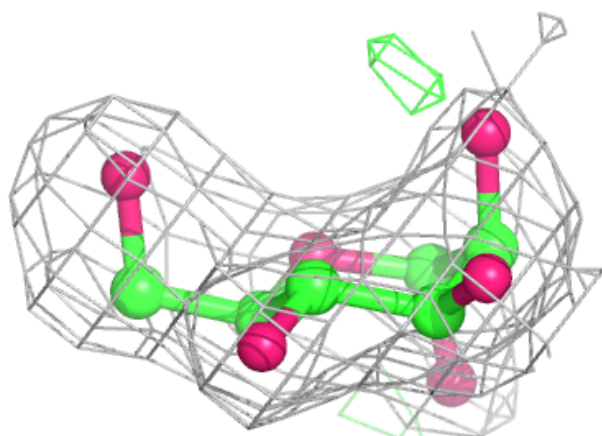
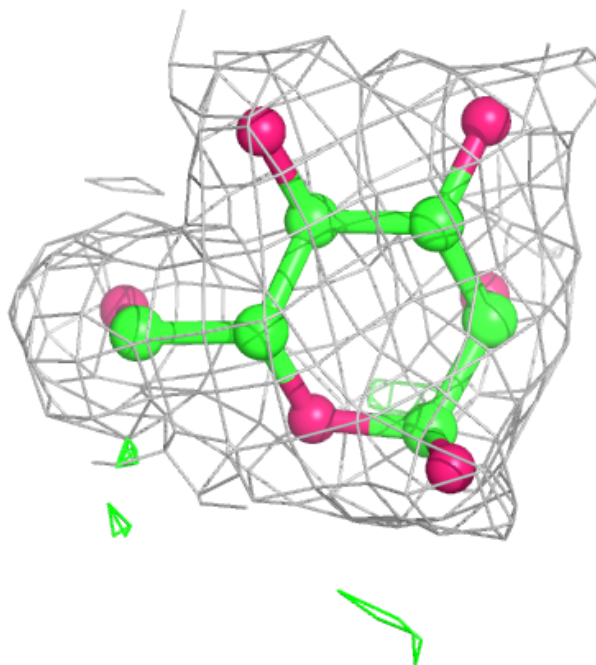
**Electron density around MAN C 401:**

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



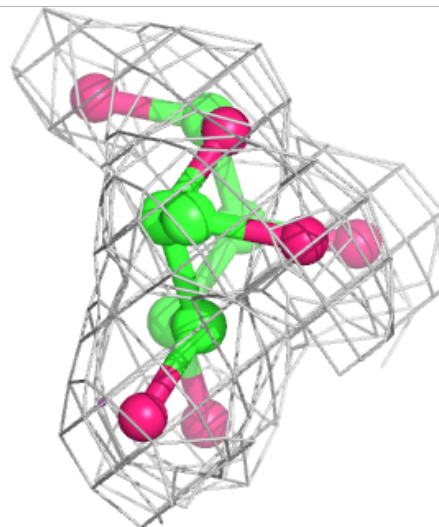
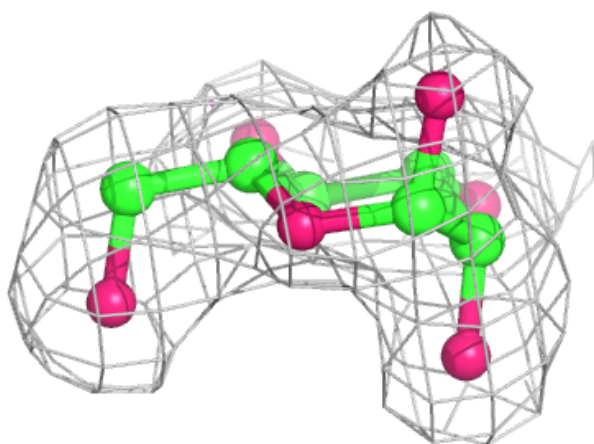
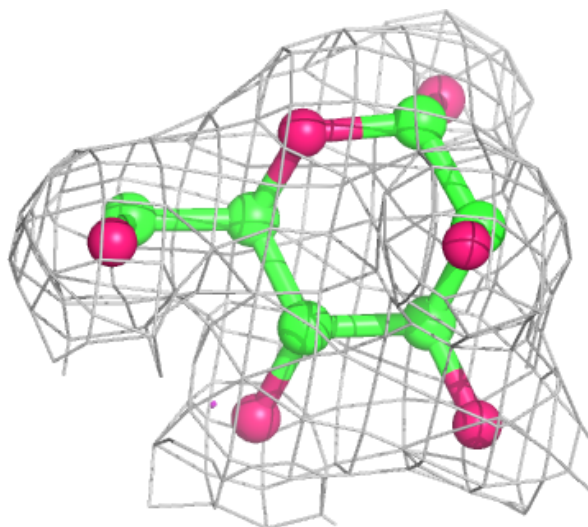
**Electron density around MAN A 401:**

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



**Electron density around MAN D 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.