



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 27, 2024 – 11:44 pm BST

PDB ID : 6FL9  
Title : The active form of a pentameric ion channel (sTeLIC) gated by alkaline pH - Wild type 2.3 Angstrom resolution  
Authors : Hu, H.; Delarue, M.  
Deposited on : 2018-01-25  
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](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 types of validation reports are described at

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

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

MolProbity : 4.02b-467  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36.2  
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) : 2.36.2

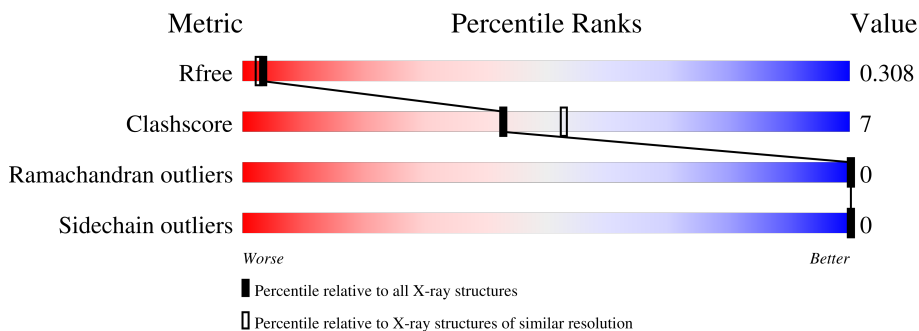
# 1 Overall quality at a glance

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)

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  $\leq 5\%$ .

Mol	Chain	Length	Quality of chain
1	A	320	85% 12% .
1	B	320	85% 12% .
1	C	320	84% 13% .
1	D	320	86% 11% .
1	E	320	85% 12% .

## 2 Entry composition [i](#)

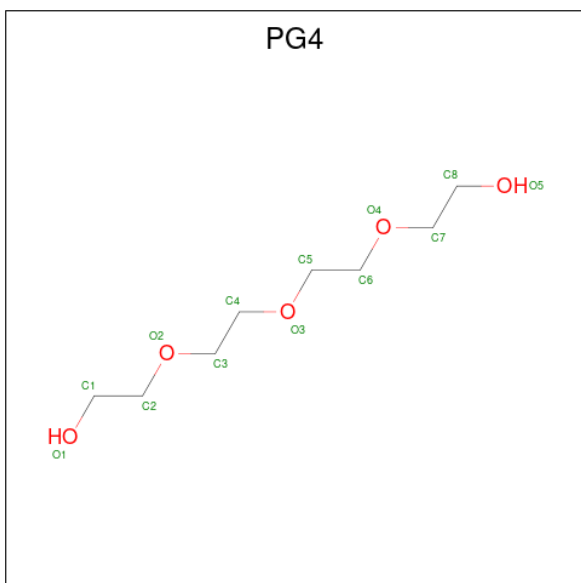
There are 4 unique types of molecules in this entry. The entry contains 13403 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Cys-loop ligand-gated ion channel.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	310	Total 2566	C 1690	N 423	O 448	S 5	0	0	0
1	B	310	Total 2566	C 1690	N 423	O 448	S 5	0	0	0
1	C	310	Total 2566	C 1690	N 423	O 448	S 5	0	0	0
1	D	310	Total 2566	C 1690	N 423	O 448	S 5	0	0	0
1	E	310	Total 2566	C 1690	N 423	O 448	S 5	0	0	0

- Molecule 2 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C<sub>8</sub>H<sub>18</sub>O<sub>5</sub>).



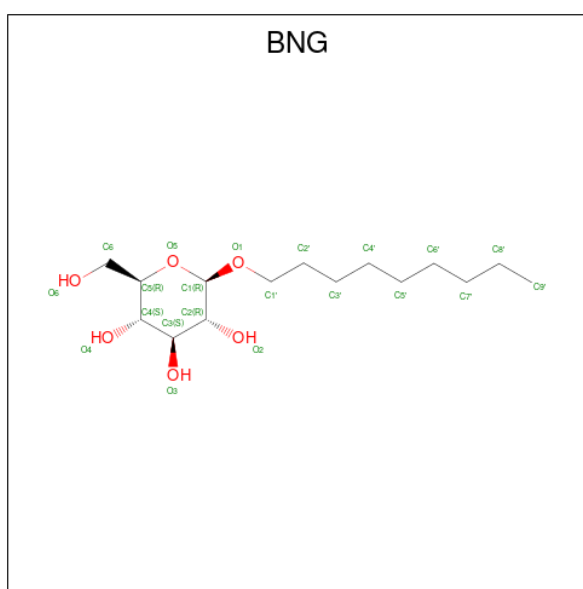
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	Total 13	C 8	O 5	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 13 8 5	0	0
2	C	1	Total C O 13 8 5	0	0
2	C	1	Total C O 13 8 5	0	0
2	E	1	Total C O 13 8 5	0	0

- Molecule 3 is nonyl beta-D-glucopyranoside (three-letter code: BNG) (formula: C<sub>15</sub>H<sub>30</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 21 15 6	0	0
3	B	1	Total C O 21 15 6	0	0
3	C	1	Total C O 21 15 6	0	0
3	D	1	Total C O 21 15 6	0	0
3	E	1	Total C O 21 15 6	0	0

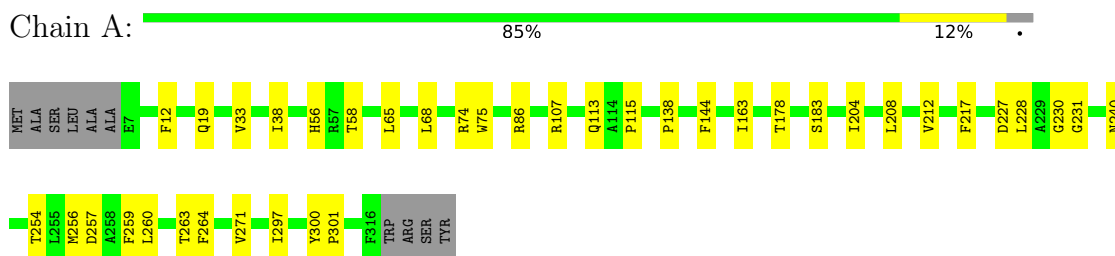
- Molecule 4 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
4	A	74	Total 74	O 74	0	0
4	B	99	Total 99	O 99	0	0
4	C	101	Total 101	O 101	0	0
4	D	88	Total 88	O 88	0	0
4	E	41	Total 41	O 41	0	0

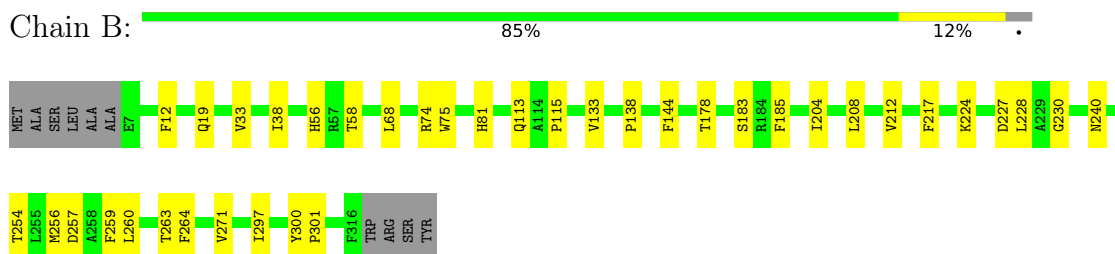
### 3 Residue-property plots [i](#)

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. 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. 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.

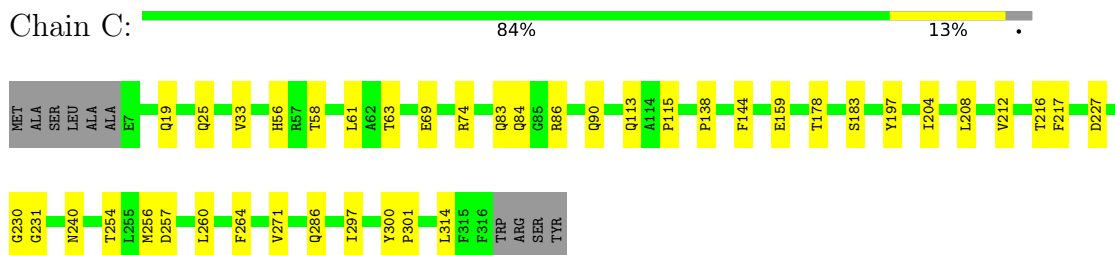
- Molecule 1: Cys-loop ligand-gated ion channel



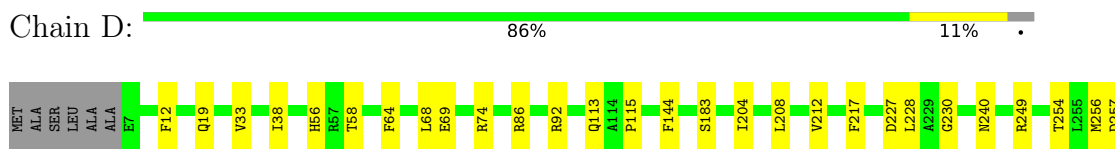
- Molecule 1: Cys-loop ligand-gated ion channel



- Molecule 1: Cys-loop ligand-gated ion channel



- Molecule 1: Cys-loop ligand-gated ion channel





- Molecule 1: Cys-loop ligand-gated ion channel

Chain E: 85% 12%



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	219.60Å 112.96Å 144.43Å 90.00° 112.05° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 19.97 – 2.30	Depositor EDS
% Data completeness (in resolution range)	89.7 (20.00-2.30) 89.7 (19.97-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.18 (at 2.30Å)	Xtrriage
Refinement program	BUSTER 2.10.3	Depositor
R, $R_{free}$	0.209 , 0.236 0.275 , 0.308	Depositor DCC
$R_{free}$ test set	6386 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.4	Xtrriage
Anisotropy	0.056	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 68.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	13403	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	80.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.52% 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: PG4, BNG

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	0.52	0/2639	0.68	0/3583
1	B	0.51	0/2639	0.67	0/3583
1	C	0.53	0/2639	0.69	0/3583
1	D	0.50	0/2639	0.68	0/3583
1	E	0.47	0/2639	0.67	0/3583
All	All	0.51	0/13195	0.68	0/17915

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	2566	0	2539	42	0
1	B	2566	0	2539	37	0
1	C	2566	0	2539	45	0
1	D	2566	0	2539	39	0
1	E	2566	0	2539	36	0
2	A	26	0	36	4	0
2	C	26	0	36	9	0
2	E	13	0	18	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	21	0	30	7	0
3	B	21	0	30	8	0
3	C	21	0	30	8	0
3	D	21	0	30	3	0
3	E	21	0	30	6	0
4	A	74	0	0	3	0
4	B	99	0	0	3	0
4	C	101	0	0	6	0
4	D	88	0	0	1	0
4	E	41	0	0	2	0
All	All	13403	0	12935	176	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 7.

All (176) 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)
3:C:403:BNG:H2'2	1:D:227:ASP:HA	1.61	0.83
1:C:61:LEU:HD21	2:C:401:PG4:C3	2.11	0.80
1:C:144:PHE:HB2	1:C:183:SER:OG	1.81	0.80
1:E:144:PHE:HB2	1:E:183:SER:OG	1.82	0.79
1:A:144:PHE:HB2	1:A:183:SER:OG	1.82	0.79
1:D:144:PHE:HB2	1:D:183:SER:OG	1.84	0.77
1:B:144:PHE:HB2	1:B:183:SER:OG	1.85	0.76
1:A:230:GLY:HA3	3:E:402:BNG:H1'2	1.67	0.73
1:B:144:PHE:HB2	1:B:183:SER:HG	1.53	0.73
1:C:61:LEU:HD21	2:C:401:PG4:H31	1.70	0.72
1:A:228:LEU:HD11	3:A:403:BNG:H6'1	1.72	0.72
1:C:144:PHE:HB2	1:C:183:SER:HG	1.55	0.72
3:C:403:BNG:H1'2	1:D:230:GLY:HA3	1.71	0.72
1:A:144:PHE:HB2	1:A:183:SER:HG	1.55	0.71
1:E:231:GLY:HA3	3:E:402:BNG:H61	1.71	0.71
1:D:144:PHE:HB2	1:D:183:SER:HG	1.54	0.71
2:C:402:PG4:C2	2:C:402:PG4:H82	2.21	0.70
1:E:144:PHE:HB2	1:E:183:SER:HG	1.52	0.70
1:A:58:THR:HG21	1:B:74:ARG:HG3	1.73	0.70
1:B:212:VAL:HG13	1:C:271:VAL:HG11	1.74	0.69
1:C:58:THR:HG21	1:D:74:ARG:HG3	1.74	0.68
1:C:300:TYR:HB3	1:C:301:PRO:HD3	1.75	0.67
1:A:271:VAL:HG11	1:E:212:VAL:HG13	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:212:VAL:HG13	1:E:271:VAL:HG11	1.77	0.66
1:A:212:VAL:HG13	1:B:271:VAL:HG11	1.78	0.66
1:A:231:GLY:HA3	3:A:403:BNG:O4	1.95	0.65
1:B:300:TYR:HB3	1:B:301:PRO:HD3	1.77	0.65
2:C:402:PG4:H82	2:C:402:PG4:H21	1.79	0.64
1:D:300:TYR:HB3	1:D:301:PRO:HD3	1.79	0.63
3:B:401:BNG:H1'2	1:C:230:GLY:HA3	1.79	0.63
1:A:300:TYR:HB3	1:A:301:PRO:HD3	1.79	0.63
1:B:58:THR:HG21	1:C:74:ARG:HG3	1.81	0.62
1:E:300:TYR:HB3	1:E:301:PRO:HD3	1.80	0.62
1:B:228:LEU:HD11	3:B:401:BNG:H6'1	1.81	0.62
1:A:74:ARG:HG3	1:E:58:THR:HG21	1.81	0.62
1:A:227:ASP:HA	3:E:402:BNG:H2'2	1.80	0.61
2:E:401:PG4:H41	2:E:401:PG4:H81	1.81	0.61
2:A:402:PG4:H11	4:A:544:HOH:O	2.02	0.60
1:C:208:LEU:CD2	1:D:264:PHE:HB3	2.31	0.60
2:C:401:PG4:O4	2:C:401:PG4:H42	2.01	0.60
1:B:212:VAL:CG1	1:C:271:VAL:HG11	2.32	0.60
1:D:204:ILE:O	1:D:208:LEU:HD13	2.03	0.59
2:C:402:PG4:H82	2:C:402:PG4:O2	2.03	0.59
1:C:204:ILE:O	1:C:208:LEU:HD13	2.01	0.59
3:C:403:BNG:H3'2	1:D:230:GLY:HA3	1.83	0.58
1:C:61:LEU:HD21	2:C:401:PG4:H32	1.84	0.58
1:B:204:ILE:O	1:B:208:LEU:HD13	2.04	0.57
1:C:208:LEU:HD21	1:D:264:PHE:HB3	1.86	0.57
1:A:208:LEU:CD2	1:B:264:PHE:HB3	2.35	0.57
1:A:204:ILE:O	1:A:208:LEU:HD13	2.04	0.57
1:E:204:ILE:O	1:E:208:LEU:HD13	2.04	0.57
1:A:75:TRP:HB2	4:A:553:HOH:O	2.05	0.56
1:C:159:GLU:HB3	4:C:533:HOH:O	2.06	0.56
3:A:403:BNG:H1'2	1:B:230:GLY:HA3	1.88	0.55
1:E:144:PHE:CB	1:E:183:SER:HG	2.20	0.55
1:B:144:PHE:CB	1:B:183:SER:HG	2.19	0.55
1:A:65:LEU:HD11	2:A:401:PG4:H32	1.88	0.54
1:A:271:VAL:HG11	1:E:212:VAL:CG1	2.37	0.54
1:A:163:ILE:HD12	4:A:526:HOH:O	2.06	0.54
1:A:231:GLY:HA3	3:A:403:BNG:H61	1.89	0.54
1:C:212:VAL:HG13	1:D:271:VAL:HG11	1.90	0.54
3:B:401:BNG:H9'1	4:B:548:HOH:O	2.07	0.54
1:D:208:LEU:CD2	1:E:264:PHE:HB3	2.37	0.54
1:A:230:GLY:HA3	3:E:402:BNG:H3'2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:401:BNG:H1'2	1:E:230:GLY:HA3	1.91	0.53
1:A:56:HIS:HB2	1:B:74:ARG:HH22	1.73	0.53
1:D:144:PHE:CB	1:D:183:SER:HG	2.22	0.53
2:E:401:PG4:H41	2:E:401:PG4:C8	2.39	0.53
1:C:231:GLY:HA3	3:C:403:BNG:O4	2.09	0.52
1:A:212:VAL:CG1	1:B:271:VAL:HG11	2.38	0.52
1:A:264:PHE:HB3	1:E:208:LEU:CD2	2.40	0.52
1:A:240:ASN:HA	1:A:260:LEU:HD13	1.90	0.52
1:B:208:LEU:CD2	1:C:264:PHE:HB3	2.40	0.52
1:A:208:LEU:HD21	1:B:264:PHE:HB3	1.91	0.52
1:D:58:THR:HG21	1:E:74:ARG:HG3	1.91	0.51
1:E:240:ASN:HA	1:E:260:LEU:HD13	1.91	0.51
1:B:240:ASN:HA	1:B:260:LEU:HD13	1.92	0.51
1:D:212:VAL:CG1	1:E:271:VAL:HG11	2.39	0.51
1:A:74:ARG:HH22	1:E:56:HIS:HB2	1.75	0.51
1:E:279:ARG:HA	4:E:502:HOH:O	2.09	0.51
1:C:240:ASN:HA	1:C:260:LEU:HD13	1.93	0.50
1:C:144:PHE:CB	1:C:183:SER:HG	2.22	0.50
3:B:401:BNG:H5	4:B:548:HOH:O	2.11	0.50
1:D:240:ASN:HA	1:D:260:LEU:HD13	1.92	0.49
1:A:144:PHE:CB	1:A:183:SER:HG	2.23	0.49
1:B:208:LEU:HD21	1:C:264:PHE:HB3	1.95	0.48
1:B:212:VAL:HG12	3:B:401:BNG:H5'1	1.94	0.48
1:B:113:GLN:OE1	1:B:115:PRO:HG3	2.14	0.48
3:D:401:BNG:H5	4:D:520:HOH:O	2.13	0.48
3:C:403:BNG:C1'	1:D:230:GLY:HA3	2.42	0.48
3:B:401:BNG:H2'2	1:C:227:ASP:HA	1.95	0.48
1:B:56:HIS:HB2	1:C:74:ARG:HH22	1.79	0.47
3:B:401:BNG:H9'3	3:B:401:BNG:H1'1	1.96	0.47
1:C:69:GLU:HA	4:C:510:HOH:O	2.13	0.47
3:A:403:BNG:H2'2	1:B:227:ASP:HA	1.96	0.47
1:B:212:VAL:HG13	1:C:271:VAL:CG1	2.42	0.46
3:A:403:BNG:H3'2	1:B:230:GLY:HA3	1.97	0.46
1:D:208:LEU:HD21	1:E:264:PHE:HB3	1.95	0.46
1:C:197:TYR:CD2	1:D:249:ARG:HG2	2.50	0.46
1:C:212:VAL:CG1	1:D:271:VAL:HG11	2.45	0.46
1:C:56:HIS:HB2	1:D:74:ARG:HH22	1.80	0.46
1:E:216:THR:HG21	3:E:402:BNG:H4'2	1.98	0.46
1:A:271:VAL:CG1	1:E:212:VAL:HG13	2.46	0.45
1:C:25:GLN:HG2	4:C:513:HOH:O	2.15	0.45
1:A:264:PHE:HB3	1:E:208:LEU:HD21	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:19:GLN:HB2	1:C:33:VAL:HB	1.99	0.45
1:A:212:VAL:HG13	1:B:271:VAL:CG1	2.45	0.45
1:A:113:GLN:OE1	1:A:115:PRO:HG3	2.17	0.44
1:B:12:PHE:O	1:B:38:ILE:HA	2.17	0.44
1:E:254:THR:HB	1:E:257:ASP:H	1.83	0.44
1:D:228:LEU:HD11	3:D:401:BNG:H8'1	1.99	0.44
1:E:113:GLN:OE1	1:E:115:PRO:HG3	2.18	0.44
1:C:86:ARG:HE	1:C:86:ARG:HB3	1.61	0.44
2:A:401:PG4:H52	2:A:401:PG4:H72	1.79	0.44
1:E:19:GLN:HB2	1:E:33:VAL:HB	2.00	0.44
1:E:144:PHE:CB	1:E:183:SER:OG	2.62	0.44
1:A:254:THR:HB	1:A:257:ASP:H	1.83	0.43
1:C:217:PHE:HB3	1:C:297:ILE:HG12	2.00	0.43
3:C:403:BNG:H1'2	1:D:230:GLY:CA	2.44	0.43
1:A:19:GLN:HB2	1:A:33:VAL:HB	2.00	0.43
1:A:138:PRO:HD3	1:A:178:THR:HG21	2.01	0.43
3:B:401:BNG:H3'2	1:C:230:GLY:HA3	2.00	0.43
1:D:113:GLN:OE1	1:D:115:PRO:HG3	2.17	0.43
1:D:212:VAL:HG13	1:E:271:VAL:CG1	2.46	0.43
1:E:217:PHE:HB3	1:E:297:ILE:HG12	2.00	0.43
1:D:259:PHE:O	1:D:263:THR:HG23	2.19	0.43
1:C:138:PRO:HD3	1:C:178:THR:HG21	2.01	0.43
1:C:254:THR:HB	1:C:257:ASP:H	1.84	0.43
2:A:402:PG4:H21	2:A:402:PG4:O5	2.19	0.43
1:C:113:GLN:OE1	1:C:115:PRO:HG3	2.18	0.43
1:D:86:ARG:HE	1:D:86:ARG:HB3	1.60	0.43
1:D:144:PHE:CB	1:D:183:SER:OG	2.63	0.43
1:D:19:GLN:HB2	1:D:33:VAL:HB	2.00	0.42
1:A:86:ARG:HE	1:A:86:ARG:HB3	1.61	0.42
1:B:138:PRO:HD3	1:B:178:THR:HG21	2.00	0.42
1:C:254:THR:HG22	1:C:256:MET:H	1.84	0.42
1:D:217:PHE:HB3	1:D:297:ILE:HG12	2.01	0.42
2:C:402:PG4:H62	2:C:402:PG4:H42	1.97	0.42
1:E:254:THR:HG22	1:E:256:MET:H	1.85	0.42
2:C:402:PG4:H41	1:D:92:ARG:NH2	2.35	0.42
1:B:224:LYS:HD2	4:B:575:HOH:O	2.20	0.42
1:C:286:GLN:HB2	4:C:542:HOH:O	2.20	0.42
1:D:56:HIS:HB2	1:E:74:ARG:HH22	1.85	0.42
1:A:107:ARG:NH2	1:B:81:HIS:HA	2.35	0.41
1:A:259:PHE:O	1:A:263:THR:HG23	2.21	0.41
1:B:68:LEU:HD13	1:B:75:TRP:HB3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:231:GLY:HA3	3:C:403:BNG:H61	2.01	0.41
1:D:64:PHE:CE2	1:D:68:LEU:HD11	2.55	0.41
1:E:12:PHE:O	1:E:38:ILE:HA	2.20	0.41
1:E:138:PRO:HD3	1:E:178:THR:HG21	2.02	0.41
1:A:12:PHE:O	1:A:38:ILE:HA	2.21	0.41
1:A:217:PHE:HB3	1:A:297:ILE:HG12	2.02	0.41
1:B:254:THR:HB	1:B:257:ASP:H	1.85	0.41
1:E:68:LEU:HD13	1:E:75:TRP:HB3	2.02	0.41
1:E:83:GLN:HG2	1:E:84:GLN:N	2.36	0.41
1:C:90:GLN:HG3	4:C:502:HOH:O	2.20	0.41
1:C:216:THR:HG21	3:C:403:BNG:H4'2	2.01	0.41
1:D:254:THR:HG22	1:D:256:MET:H	1.85	0.41
3:E:402:BNG:H9'3	3:E:402:BNG:H1'1	2.03	0.41
1:B:19:GLN:HB2	1:B:33:VAL:HB	2.02	0.41
1:C:58:THR:CG2	1:D:74:ARG:NH1	2.83	0.41
1:B:133:VAL:HB	1:B:185:PHE:HB3	2.03	0.41
1:B:217:PHE:HB3	1:B:297:ILE:HG12	2.02	0.41
1:B:259:PHE:O	1:B:263:THR:HG23	2.21	0.41
1:D:12:PHE:O	1:D:38:ILE:HA	2.20	0.41
1:A:68:LEU:HD13	1:A:75:TRP:HB3	2.03	0.41
1:C:63:THR:HG21	1:D:69:GLU:OE1	2.21	0.41
1:D:254:THR:HB	1:D:257:ASP:H	1.86	0.41
1:A:144:PHE:CB	1:A:183:SER:OG	2.61	0.40
1:B:254:THR:HG22	1:B:256:MET:H	1.86	0.40
1:E:133:VAL:HB	1:E:185:PHE:HB3	2.03	0.40
1:A:231:GLY:CA	3:A:403:BNG:H61	2.52	0.40
1:C:83:GLN:HG2	1:C:84:GLN:N	2.37	0.40
1:C:314:LEU:HD22	4:C:556:HOH:O	2.20	0.40
1:A:254:THR:HG22	1:A:256:MET:H	1.85	0.40
1:E:224:LYS:HG3	4:E:533:HOH:O	2.21	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	308/320 (96%)	303 (98%)	5 (2%)	0	100	100
1	B	308/320 (96%)	303 (98%)	5 (2%)	0	100	100
1	C	308/320 (96%)	302 (98%)	6 (2%)	0	100	100
1	D	308/320 (96%)	302 (98%)	6 (2%)	0	100	100
1	E	308/320 (96%)	303 (98%)	5 (2%)	0	100	100
All	All	1540/1600 (96%)	1513 (98%)	27 (2%)	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	275/282 (98%)	275 (100%)	0	100	100
1	B	275/282 (98%)	275 (100%)	0	100	100
1	C	275/282 (98%)	275 (100%)	0	100	100
1	D	275/282 (98%)	275 (100%)	0	100	100
1	E	275/282 (98%)	275 (100%)	0	100	100
All	All	1375/1410 (98%)	1375 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	GLN
1	B	49	HIS
1	B	83	GLN
1	C	83	GLN
1	D	49	HIS

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Mol	Chain	Res	Type
1	E	49	HIS
1	E	83	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](#)

10 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	BNG	C	403	-	21,21,21	0.29	0	26,26,26	0.86	1 (3%)
2	PG4	A	401	-	12,12,12	0.29	0	11,11,11	0.51	0
3	BNG	A	403	-	21,21,21	0.30	0	26,26,26	0.61	0
2	PG4	A	402	-	12,12,12	0.20	0	11,11,11	0.28	0
3	BNG	D	401	-	21,21,21	0.28	0	26,26,26	0.63	0
2	PG4	C	401	-	12,12,12	0.28	0	11,11,11	0.27	0
3	BNG	B	401	-	21,21,21	0.30	0	26,26,26	0.42	0
2	PG4	E	401	-	12,12,12	0.15	0	11,11,11	0.16	0
2	PG4	C	402	-	12,12,12	0.25	0	11,11,11	0.29	0
3	BNG	E	402	-	21,21,21	0.20	0	26,26,26	0.47	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
3	BNG	C	403	-	-	8/12/32/32	0/1/1/1
2	PG4	A	401	-	-	7/10/10/10	-
3	BNG	A	403	-	-	7/12/32/32	0/1/1/1
2	PG4	A	402	-	-	6/10/10/10	-
3	BNG	D	401	-	-	8/12/32/32	0/1/1/1
2	PG4	C	401	-	-	5/10/10/10	-
3	BNG	B	401	-	-	8/12/32/32	0/1/1/1
2	PG4	E	401	-	-	5/10/10/10	-
2	PG4	C	402	-	-	7/10/10/10	-
3	BNG	E	402	-	-	6/12/32/32	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	403	BNG	O5-C1-O1	2.55	116.00	109.97

There are no chirality outliers.

All (67) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	403	BNG	C2-C1-O1-C1'
3	A	403	BNG	O5-C1-O1-C1'
3	B	401	BNG	C2-C1-O1-C1'
3	B	401	BNG	O5-C1-O1-C1'
3	C	403	BNG	C2-C1-O1-C1'
3	D	401	BNG	C2-C1-O1-C1'
3	D	401	BNG	O5-C1-O1-C1'
2	C	402	PG4	C6-C5-O3-C4
3	B	401	BNG	O5-C5-C6-O6
2	C	401	PG4	O2-C3-C4-O3
2	C	402	PG4	O2-C3-C4-O3
2	A	402	PG4	O1-C1-C2-O2
3	C	403	BNG	O5-C1-O1-C1'
2	C	402	PG4	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	A	403	BNG	C3'-C4'-C5'-C6'
3	C	403	BNG	C3'-C4'-C5'-C6'
3	D	401	BNG	C3'-C4'-C5'-C6'
3	D	401	BNG	C4'-C5'-C6'-C7'
3	E	402	BNG	C3'-C4'-C5'-C6'
3	E	402	BNG	C2'-C3'-C4'-C5'
3	A	403	BNG	C4'-C5'-C6'-C7'
3	C	403	BNG	C2'-C3'-C4'-C5'
2	A	401	PG4	O1-C1-C2-O2
3	B	401	BNG	C3'-C4'-C5'-C6'
3	B	401	BNG	C2'-C1'-O1-C1
3	D	401	BNG	C2'-C3'-C4'-C5'
2	A	401	PG4	C1-C2-O2-C3
2	E	401	PG4	O2-C3-C4-O3
3	D	401	BNG	O5-C5-C6-O6
3	B	401	BNG	C4'-C5'-C6'-C7'
3	B	401	BNG	C2'-C3'-C4'-C5'
2	E	401	PG4	O4-C7-C8-O5
3	A	403	BNG	C2'-C3'-C4'-C5'
3	B	401	BNG	C4-C5-C6-O6
2	A	401	PG4	C5-C6-O4-C7
2	C	401	PG4	O4-C7-C8-O5
3	E	402	BNG	C4'-C5'-C6'-C7'
2	A	402	PG4	C3-C4-O3-C5
3	D	401	BNG	C5'-C6'-C7'-C8'
3	A	403	BNG	C2'-C1'-O1-C1
3	C	403	BNG	C2'-C1'-O1-C1
3	D	401	BNG	C2'-C1'-O1-C1
3	E	402	BNG	C2'-C1'-O1-C1
3	C	403	BNG	C4'-C5'-C6'-C7'
2	A	402	PG4	O3-C5-C6-O4
2	A	401	PG4	C4-C3-O2-C2
3	E	402	BNG	C6'-C7'-C8'-C9'
2	C	401	PG4	C6-C5-O3-C4
2	E	401	PG4	C4-C3-O2-C2
3	C	403	BNG	C6'-C7'-C8'-C9'
2	C	401	PG4	C3-C4-O3-C5
2	C	402	PG4	C3-C4-O3-C5
2	A	401	PG4	O2-C3-C4-O3
2	A	402	PG4	C6-C5-O3-C4
2	A	401	PG4	C3-C4-O3-C5
2	C	402	PG4	C1-C2-O2-C3

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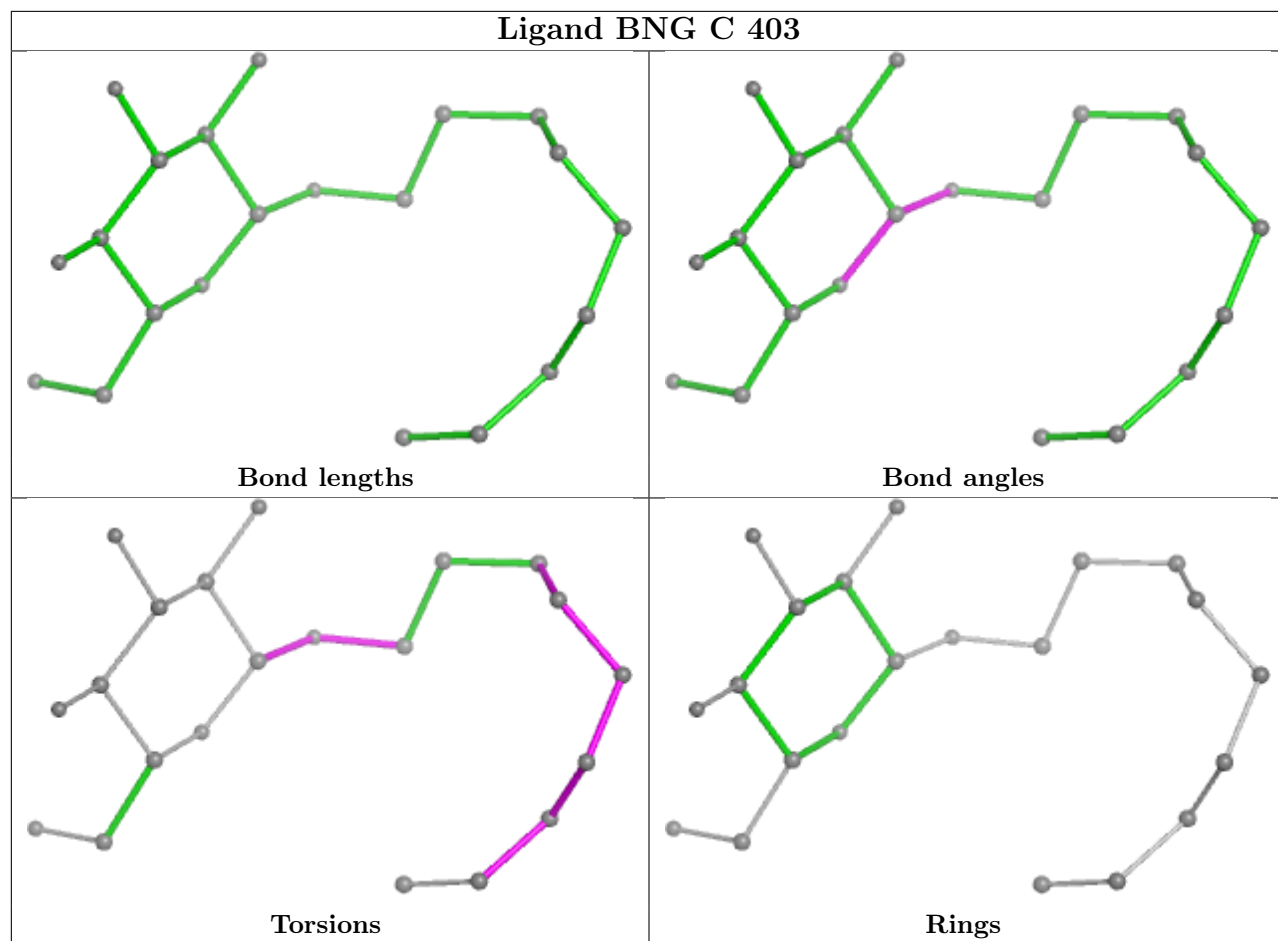
Mol	Chain	Res	Type	Atoms
2	A	402	PG4	O4-C7-C8-O5
2	E	401	PG4	C8-C7-O4-C6
3	A	403	BNG	C6'-C7'-C8'-C9'
2	C	402	PG4	O3-C5-C6-O4
3	C	403	BNG	C5'-C6'-C7'-C8'
2	C	401	PG4	C4-C3-O2-C2
2	C	402	PG4	C4-C3-O2-C2
2	A	402	PG4	O2-C3-C4-O3
2	A	401	PG4	C6-C5-O3-C4
3	E	402	BNG	C5'-C6'-C7'-C8'
2	E	401	PG4	C5-C6-O4-C7

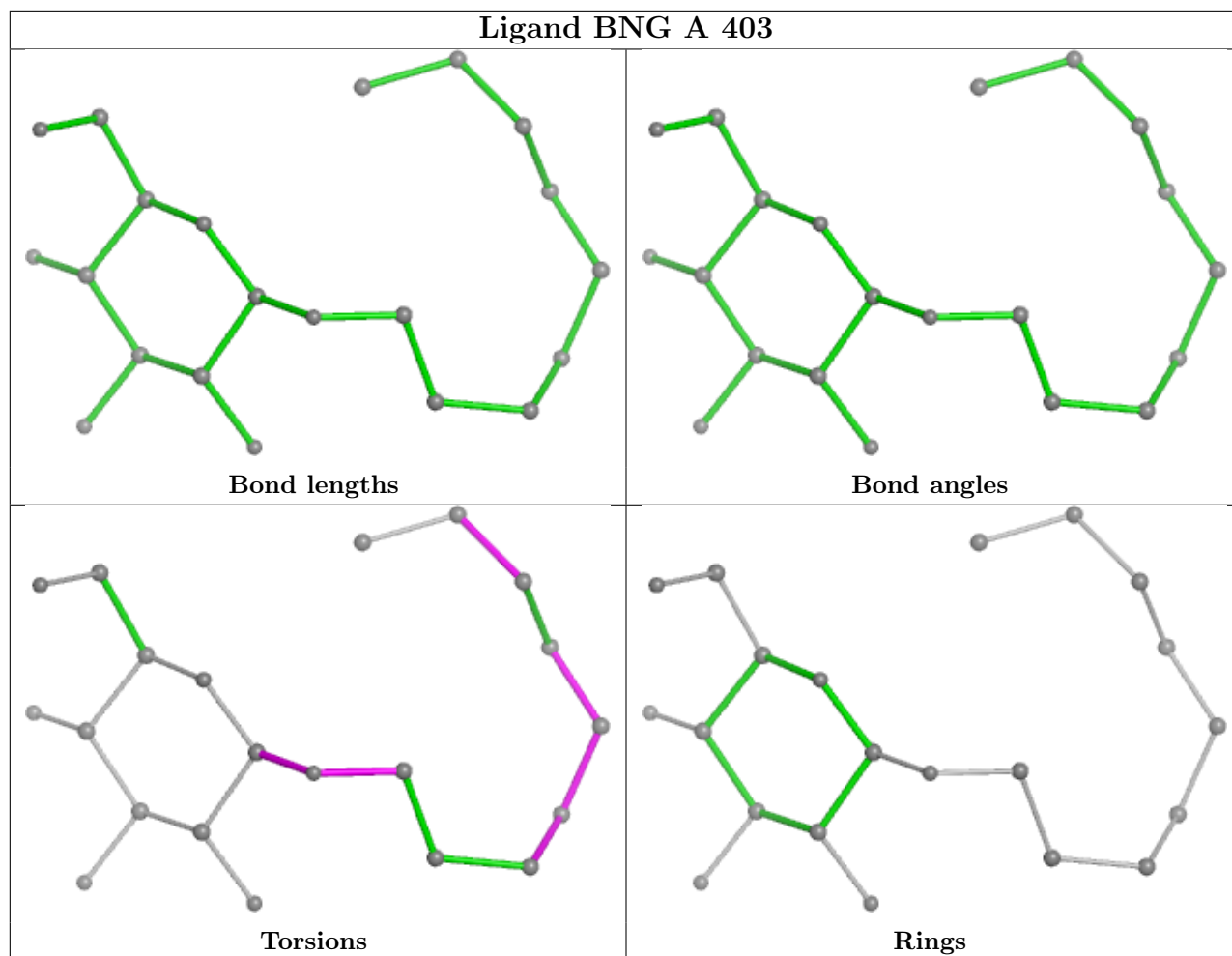
There are no ring outliers.

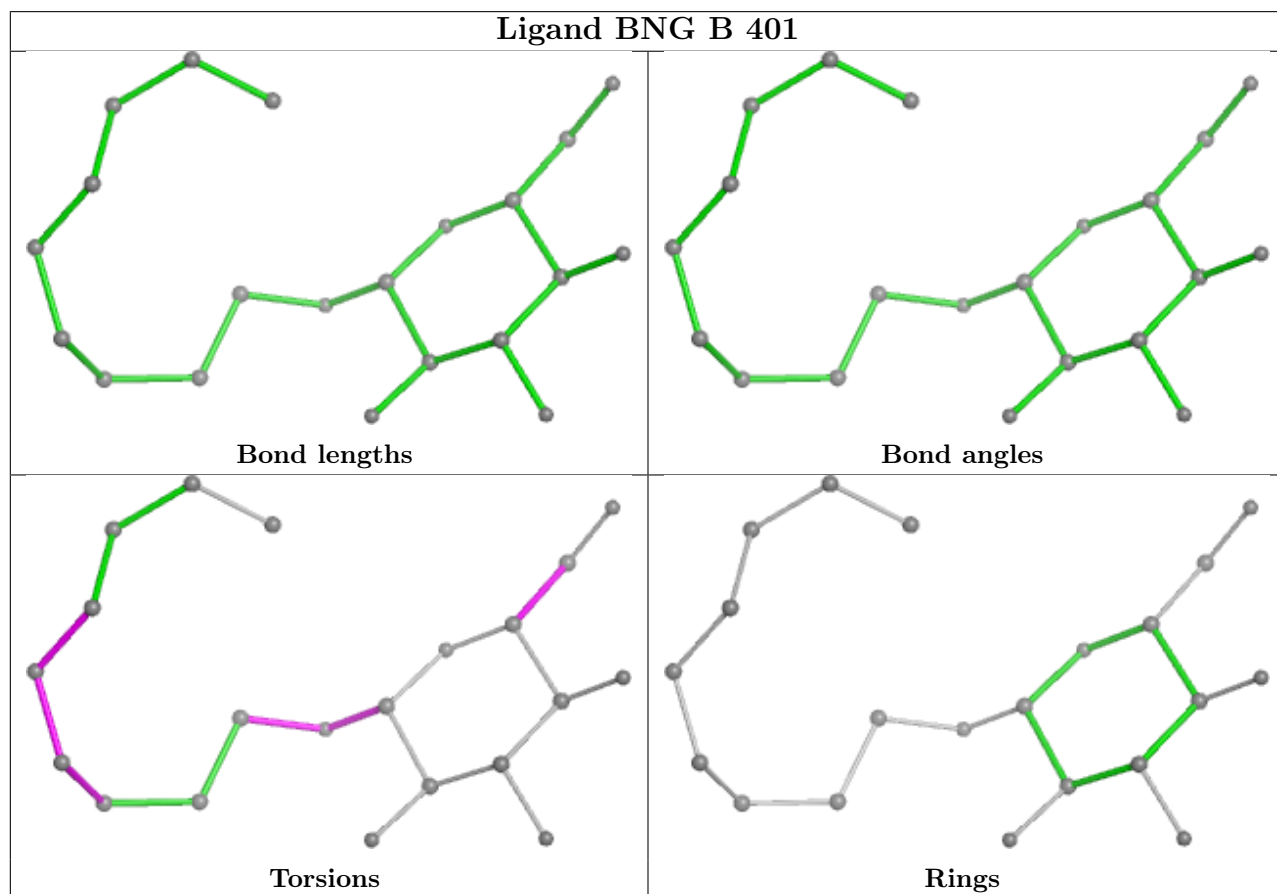
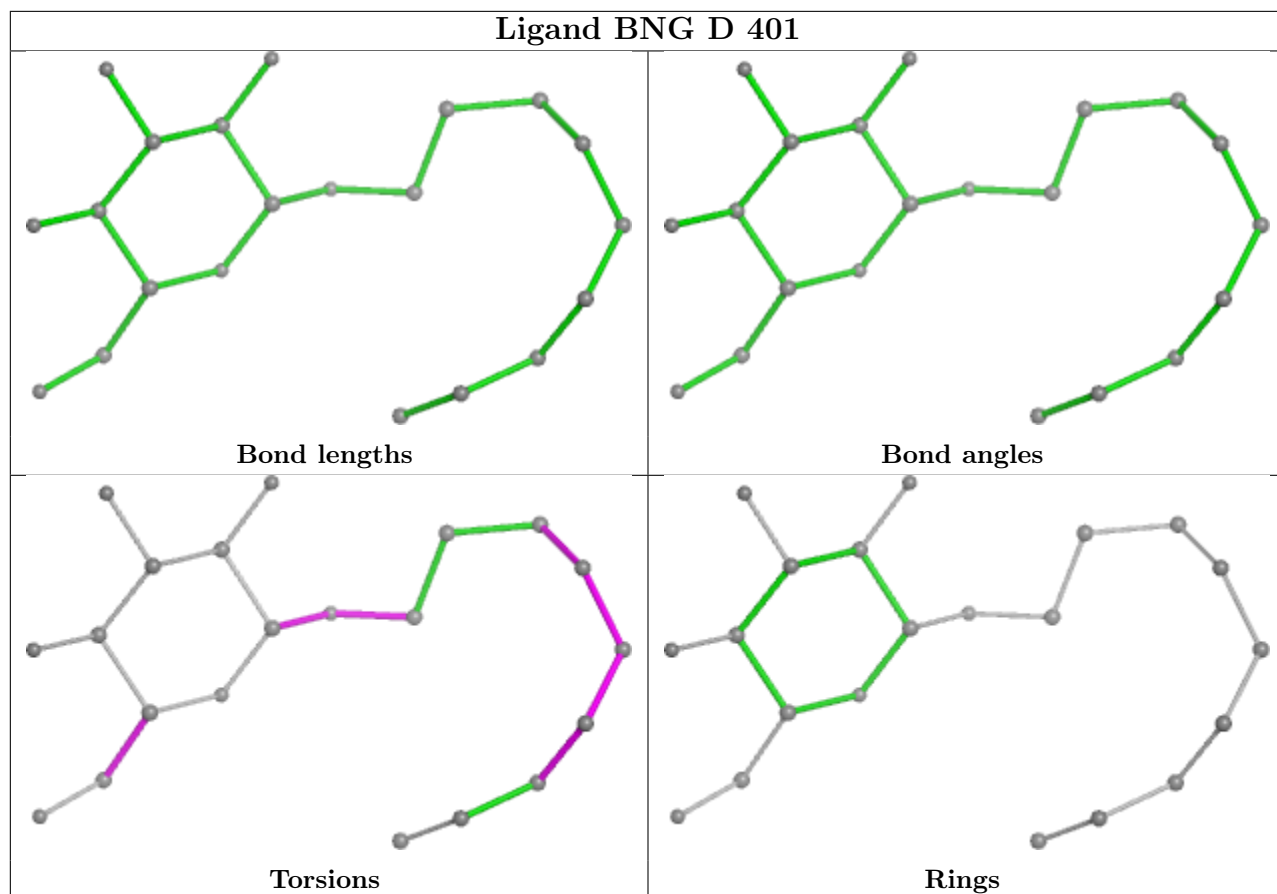
10 monomers are involved in 47 short contacts:

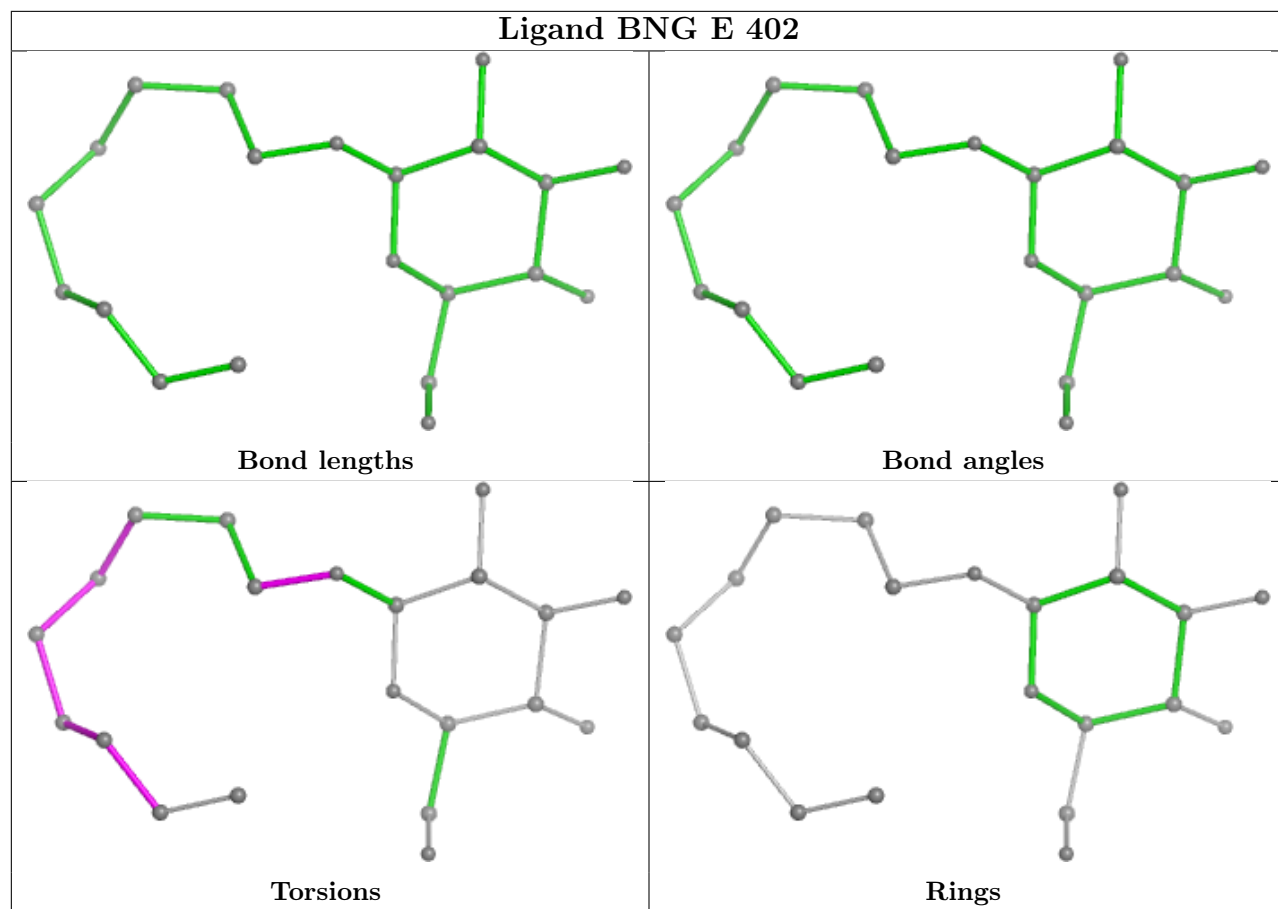
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	403	BNG	8	0
2	A	401	PG4	2	0
3	A	403	BNG	7	0
2	A	402	PG4	2	0
3	D	401	BNG	3	0
2	C	401	PG4	4	0
3	B	401	BNG	8	0
2	E	401	PG4	2	0
2	C	402	PG4	5	0
3	E	402	BNG	6	0

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

### 6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.4 Ligands

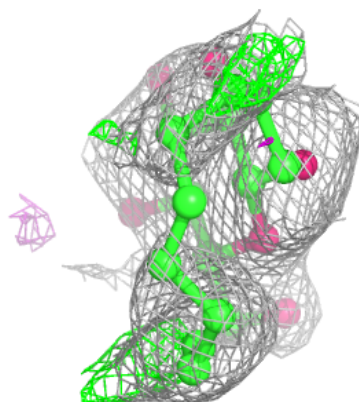
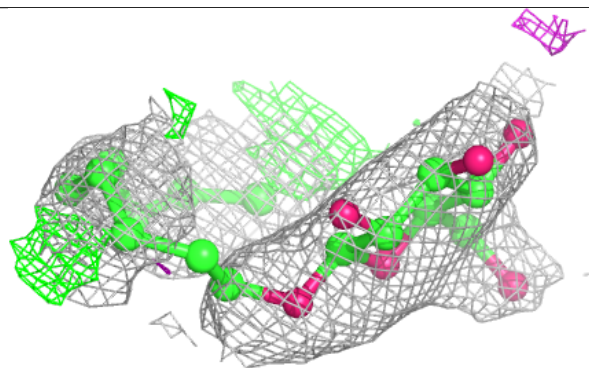
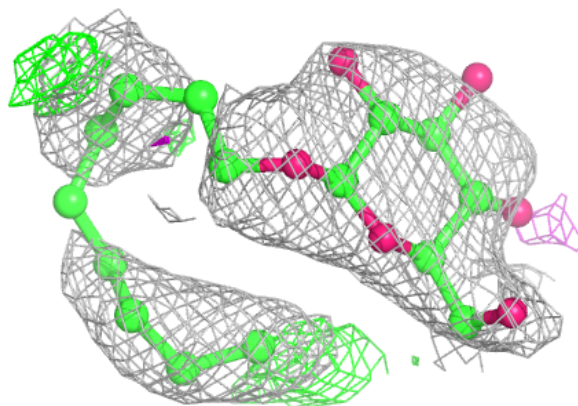
Unable to reproduce the depositors R factor - this section is therefore empty.

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.

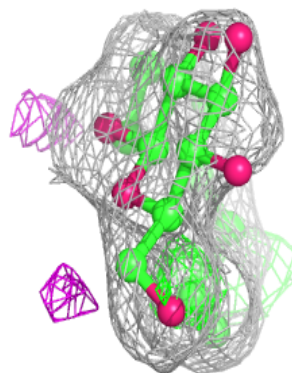
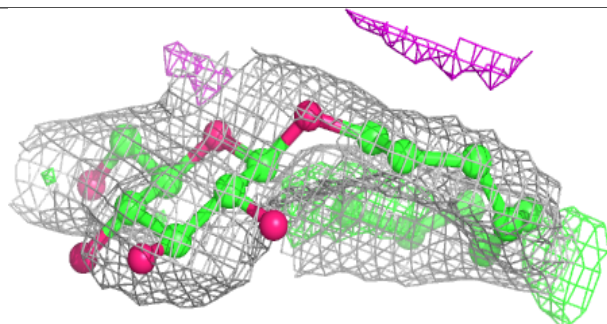
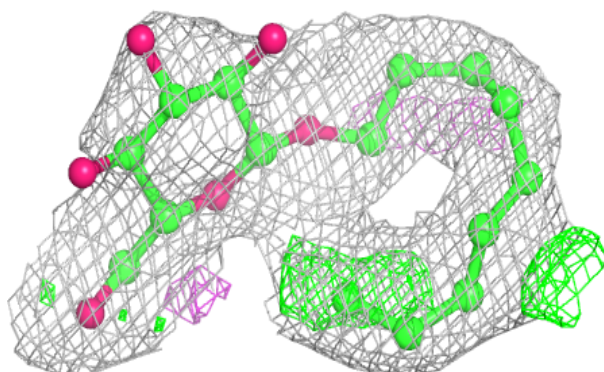


**Electron density around BNG A 403:**

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

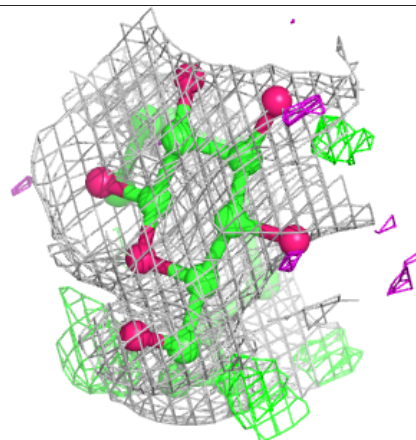
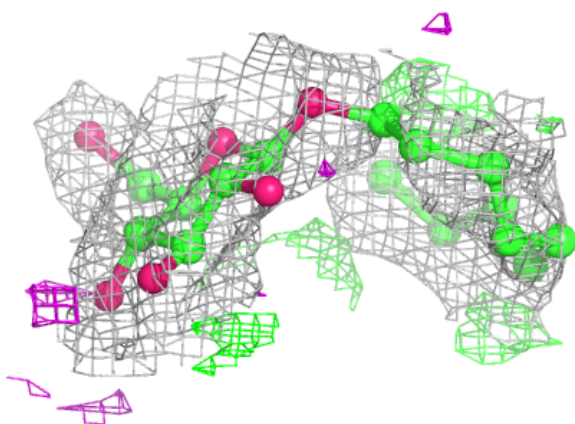
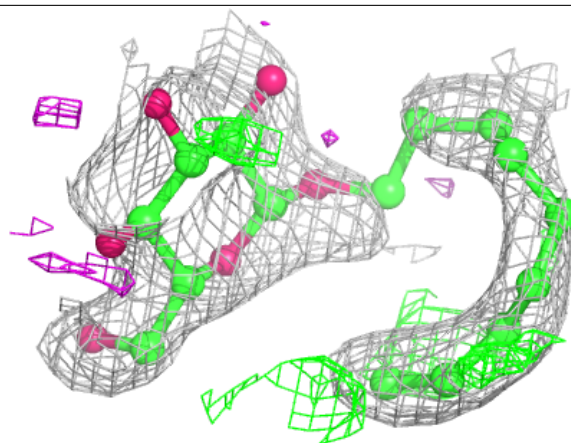
**Electron density around BNG B 401:**

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

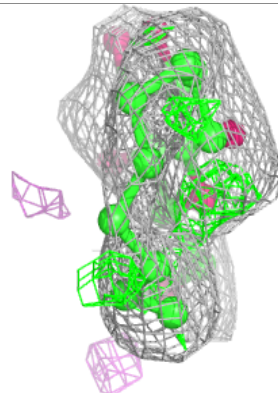
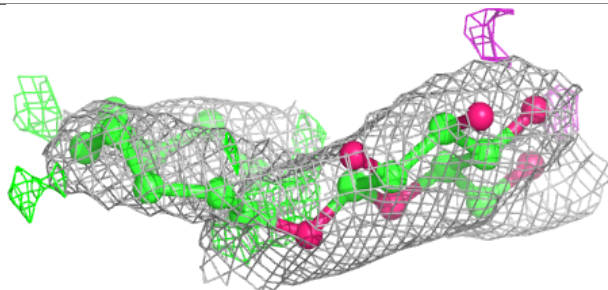
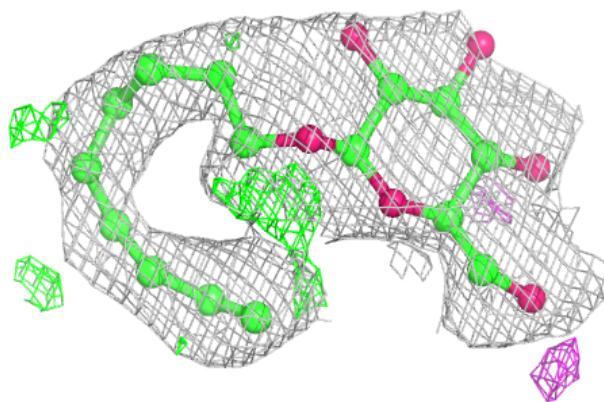


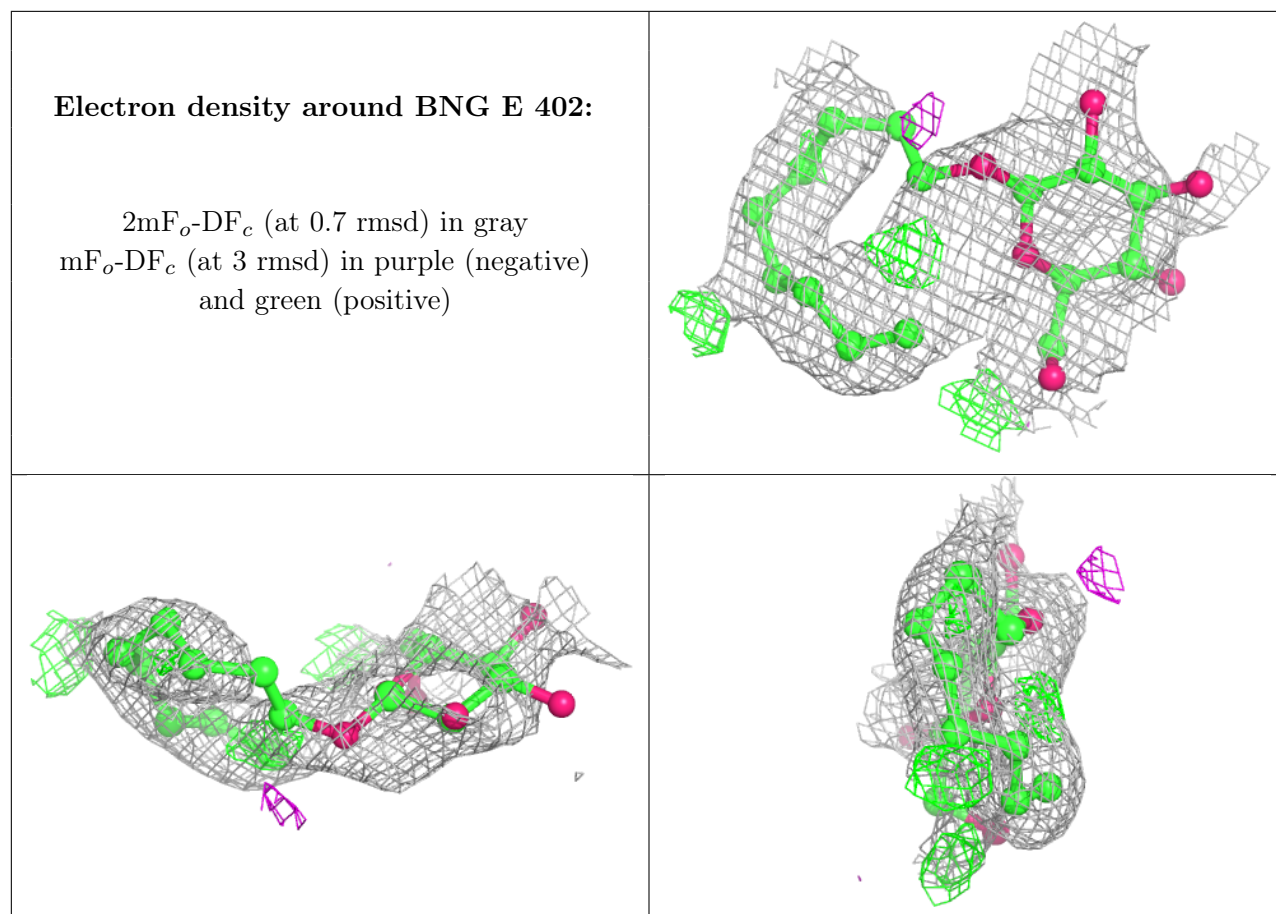
**Electron density around BNG C 403:**

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

**Electron density around BNG 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](#)

Unable to reproduce the depositor's R factor - this section is therefore empty.