



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 25, 2025 – 10:22 AM EDT

PDB ID : 9EFP / pdb\_00009efp  
Title : Crystal Structure of *Saccharomyces cerevisiae* Sec14p in complex with phosphatidylcholine  
Authors : Green, S.M.; Laganowsky, A.; Krieger, I.; Singh, P.K.; Sacchettini, J.; Igu-menova, T.I.; Bankaitis, V.A.  
Deposited on : 2024-11-20  
Resolution : 1.83 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.46

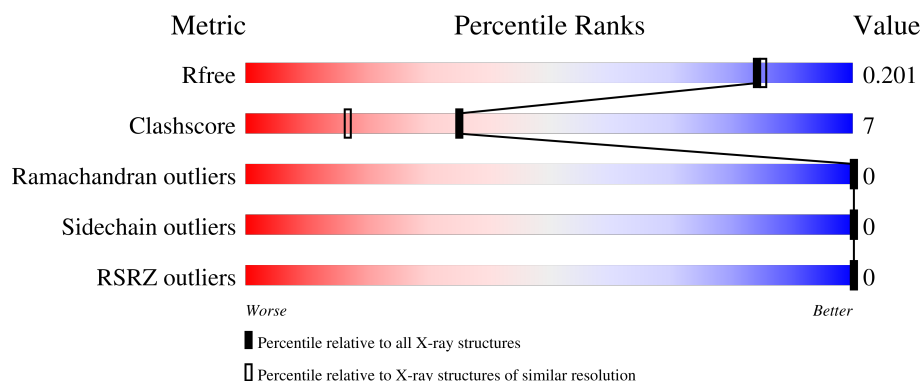
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.83 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))
$R_{free}$	164625	1150 (1.84-1.84)
Clashscore	180529	1248 (1.84-1.84)
Ramachandran outliers	177936	1240 (1.84-1.84)
Sidechain outliers	177891	1240 (1.84-1.84)
RSRZ outliers	164620	1149 (1.84-1.84)

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	304	
1	B	304	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	PEG	B	404	-	-	X	-

## 2 Entry composition [i](#)

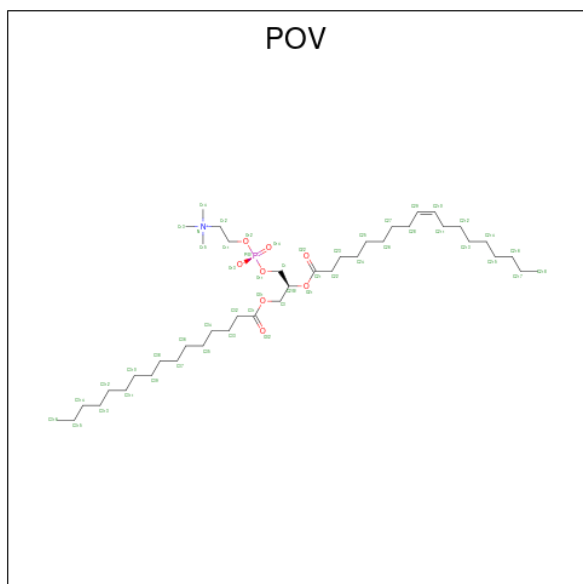
There are 6 unique types of molecules in this entry. The entry contains 5126 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SEC14 cytosolic factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	2	0
			2346	1508	382	445	11			
1	B	290	Total	C	N	O	S	0	0	0
			2346	1510	383	442	11			

- Molecule 2 is (2S)-3-(hexadecanoyloxy)-2-[(9Z)-octadec-9-enoyloxy]propyl 2-(trimethylammonio)ethyl phosphate (CCD ID: POV) (formula:  $C_{42}H_{82}NO_8P$ ) (labeled as "Ligand of Interest" by depositor).



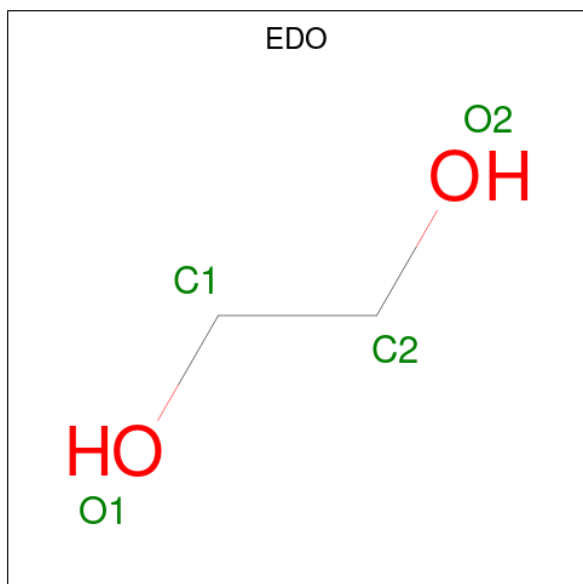
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			52	42	1	8	1		
2	B	1	Total	C	N	O	P	0	0
			52	42	1	8	1		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula:  $C_3H_8O_3$ ) (labeled as "Ligand of Interest" by depositor).



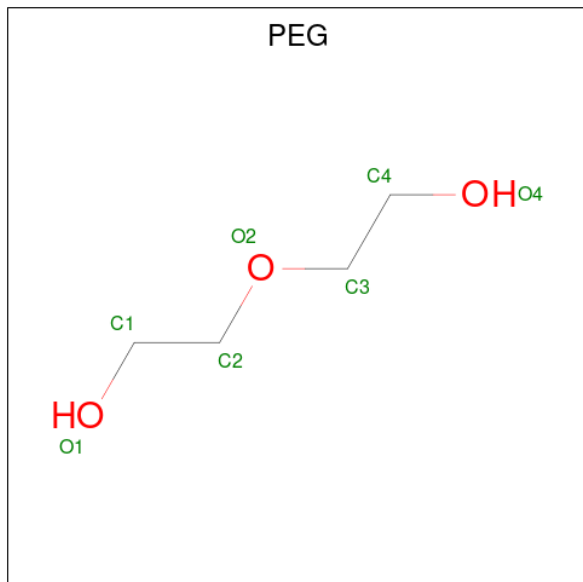
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula:  $C_2H_6O_2$ ) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula:  $C_4H_{10}O_3$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			7	4	3		
5	B	1	Total	C	O	0	0
			7	4	3		
5	B	1	Total	C	O	0	0
			7	4	3		

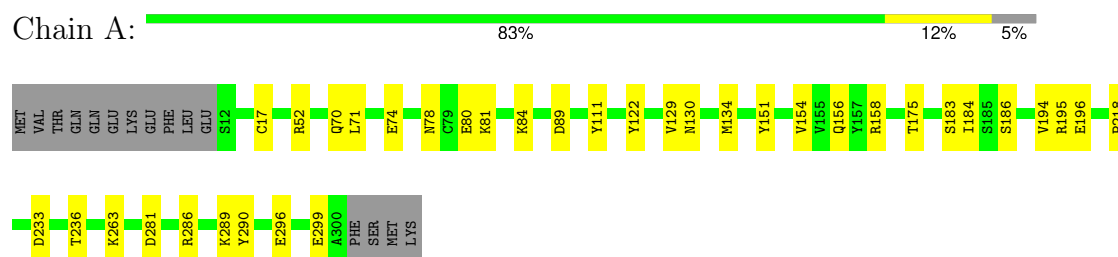
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	152	Total	O	0	0
			152	152		
6	B	141	Total	O	0	0
			141	141		

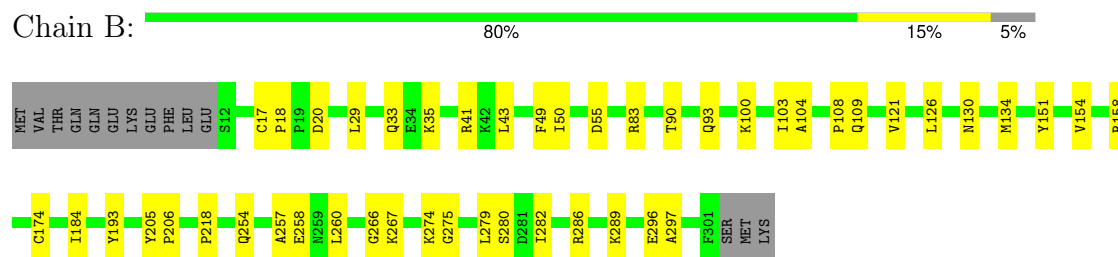
### 3 Residue-property plots

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: SEC14 cytosolic factor



- Molecule 1: SEC14 cytosolic factor



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.40Å 74.40Å 195.86Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.86 – 1.83 45.86 – 1.83	Depositor EDS
% Data completeness (in resolution range)	99.7 (45.86-1.83) 99.7 (45.86-1.83)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.83 (at 1.83Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487, PHENIX 1.20.1_4487	Depositor
R, $R_{free}$	0.172 , 0.208 0.175 , 0.201	Depositor DCC
$R_{free}$ test set	2814 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.9	Xtriage
Anisotropy	0.280	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 32.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.40$ , $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.299 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	5126	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, CME, EDO, POV, GOL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.42	0/2402	0.56	0/3247
1	B	0.39	0/2397	0.58	0/3239
All	All	0.40	0/4799	0.57	0/6486

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

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	2346	0	2299	30	0
1	B	2346	0	2298	36	0
2	A	52	0	82	9	0
2	B	52	0	82	5	0
3	A	6	0	8	2	0
3	B	6	0	8	0	0
4	A	4	0	6	0	0
5	B	21	0	30	6	0
6	A	152	0	0	4	2
6	B	141	0	0	9	2
All	All	5126	0	4813	70	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 (70) 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:B:184:ILE:HD11	1:B:218:PRO:HG2	1.54	0.90
1:B:109:GLN:HB2	5:B:404:PEG:H12	1.55	0.86
1:B:280:SER:HB2	1:B:282:ILE:HD12	1.75	0.68
1:A:129:VAL:HG11	2:A:401:POV:H210	1.75	0.68
1:A:156:GLN:O	6:A:501:HOH:O	2.13	0.67
1:B:205:TYR:O	6:B:503:HOH:O	2.14	0.66
1:A:194:VAL:HG22	2:A:401:POV:H26A	1.77	0.65
1:B:50:ILE:O	6:B:502:HOH:O	2.13	0.65
1:A:195:ARG:NH1	1:A:196[A]:GLU:OE2	2.28	0.64
1:B:100:LYS:HE2	5:B:404:PEG:H22	1.80	0.63
1:A:186:SER:HB3	2:A:401:POV:H313	1.81	0.63
1:B:151:TYR:OH	6:B:501:HOH:O	2.08	0.62
1:B:104:ALA:HB2	5:B:404:PEG:H21	1.80	0.62
1:B:130:ASN:O	1:B:134:MET:HG3	2.01	0.61
1:A:233[B]:ASP:OD1	1:A:236:THR:N	2.34	0.60
1:B:274:LYS:NZ	6:B:518:HOH:O	2.35	0.60
1:A:80:GLU:OE2	6:A:502:HOH:O	2.16	0.59
1:B:41:ARG:NH2	1:B:55:ASP:OD1	2.32	0.57
1:A:184:ILE:HD11	1:A:218:PRO:HG2	1.86	0.57
1:B:49:PHE:O	6:B:504:HOH:O	2.17	0.56
1:B:254:GLN:NE2	6:B:506:HOH:O	2.22	0.55
1:B:289:LYS:NZ	6:B:505:HOH:O	2.21	0.55
1:A:281:ASP:HB3	1:A:286:ARG:HH21	1.71	0.54
1:B:206:PRO:HA	5:B:403:PEG:H22	1.87	0.54
1:A:183:SER:O	1:A:186:SER:HB2	2.08	0.54
1:B:275:GLY:HA3	1:B:279:LEU:HD12	1.90	0.53
1:A:89:ASP:HB2	6:A:587:HOH:O	2.09	0.52
1:A:111:TYR:OH	2:A:401:POV:H13A	2.09	0.52
1:B:35:LYS:HE2	1:B:35:LYS:HA	1.92	0.52
1:B:90:THR:O	1:B:93:GLN:HG2	2.10	0.51
1:A:186:SER:CB	2:A:401:POV:H315	2.40	0.51
1:B:126:LEU:HD23	2:B:401:POV:H31E	1.91	0.50
1:A:286:ARG:HH11	1:A:296:GLU:HG2	1.76	0.50
1:B:43:LEU:HD23	6:B:509:HOH:O	2.11	0.50
1:B:154:VAL:HG22	1:B:158:ARG:HD2	1.93	0.49
1:A:286:ARG:NH2	1:A:299:GLU:HB2	2.27	0.49
1:B:83:ARG:NH1	6:B:514:HOH:O	2.32	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:286:ARG:NH1	1:B:297:ALA:O	2.33	0.49
1:A:281:ASP:OD2	1:A:286:ARG:NH2	2.47	0.48
1:A:158:ARG:HH11	1:A:158:ARG:HG3	1.79	0.48
1:B:257:ALA:HB1	1:B:267:LYS:HG2	1.96	0.47
1:B:280:SER:HB2	1:B:282:ILE:CD1	2.43	0.47
1:A:130:ASN:O	1:A:134:MET:HG3	2.14	0.47
1:B:193:TYR:CE2	2:B:401:POV:H212	2.50	0.47
1:A:263:LYS:HE2	3:A:402:GOL:H31	1.97	0.46
1:A:263:LYS:O	3:A:402:GOL:H32	2.15	0.46
1:A:84:LYS:HB2	1:A:84:LYS:HE2	1.57	0.45
1:A:175:THR:OG1	2:A:401:POV:O13	2.19	0.45
1:B:258:GLU:HA	1:B:267:LYS:HB2	1.99	0.45
1:A:289:LYS:HD3	1:A:290:TYR:CZ	2.51	0.44
1:A:122:TYR:CE2	2:A:401:POV:H13B	2.53	0.44
2:B:401:POV:H210	2:B:401:POV:H213	1.79	0.44
1:A:78:ASN:HA	1:A:81:LYS:HE3	2.00	0.43
1:B:29:LEU:HD22	1:B:33:GLN:HB3	2.00	0.43
1:A:151:TYR:O	1:A:154:VAL:HG12	2.18	0.43
1:A:70:GLN:O	1:A:74:GLU:HG3	2.19	0.43
1:B:286:ARG:HH11	1:B:296:GLU:HG2	1.84	0.42
2:A:401:POV:H21H	2:A:401:POV:H215	1.69	0.42
1:B:154:VAL:HG21	2:B:401:POV:H15B	2.01	0.42
2:A:401:POV:H15B	2:A:401:POV:H11A	1.73	0.42
1:B:108:PRO:HA	5:B:404:PEG:H42	2.00	0.42
1:B:18:PRO:HB2	1:B:20:ASP:OD1	2.20	0.42
1:B:103:ILE:HD13	1:B:103:ILE:HA	1.92	0.41
1:A:71:LEU:HD23	1:A:71:LEU:HA	1.85	0.41
1:A:184:ILE:HA	1:A:184:ILE:HD13	1.76	0.41
1:A:52:ARG:NH1	6:A:512:HOH:O	2.38	0.41
2:B:401:POV:H36	2:B:401:POV:H33A	1.85	0.41
1:B:104:ALA:HB1	5:B:404:PEG:O4	2.21	0.41
1:B:260:LEU:HB3	1:B:266:GLY:HA3	2.03	0.41
1:B:121:VAL:HA	1:B:174:CYS:O	2.21	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)
6:B:560:HOH:O	6:B:576:HOH:O[6_444]	2.02	0.18
6:A:576:HOH:O	6:B:518:HOH:O[6_554]	2.17	0.03
6:A:648:HOH:O	6:A:648:HOH:O[6_554]	2.19	0.01

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	288/304 (95%)	278 (96%)	10 (4%)	0	100	100
1	B	287/304 (94%)	275 (96%)	12 (4%)	0	100	100
All	All	575/608 (95%)	553 (96%)	22 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains

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	253/266 (95%)	253 (100%)	0	100	100
1	B	252/266 (95%)	252 (100%)	0	100	100
All	All	505/532 (95%)	505 (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 (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	B	130	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\# Z  > 2$
1	CME	A	17	1	8,9,10	0.76	0	6,9,11	3.03	2 (33%)
1	CME	B	17	1	8,9,10	0.66	0	6,9,11	4.95	2 (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
1	CME	A	17	1	-	1/5/8/10	-
1	CME	B	17	1	-	2/5/8/10	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	17	CME	CE-SD-SG	11.80	155.27	103.46
1	A	17	CME	CE-SD-SG	6.66	132.71	103.46
1	A	17	CME	CB-SG-SD	-2.72	96.83	103.86
1	B	17	CME	CB-SG-SD	2.23	109.62	103.86

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	17	CME	SD-CE-CZ-OH
1	A	17	CME	SD-CE-CZ-OH
1	B	17	CME	CZ-CE-SD-SG

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

8 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	$\# Z  > 2$
2	POV	A	401	-	51,51,51	0.55	0	57,59,59	0.71	0
5	PEG	B	403	-	6,6,6	0.08	0	5,5,5	0.15	0
5	PEG	B	404	-	6,6,6	0.26	0	5,5,5	0.33	0
3	GOL	B	405	-	5,5,5	0.99	0	5,5,5	1.12	0
4	EDO	A	403	-	3,3,3	0.42	0	2,2,2	0.34	0
2	POV	B	401	-	51,51,51	0.56	0	57,59,59	0.65	1 (1%)
3	GOL	A	402	-	5,5,5	1.06	0	5,5,5	1.02	0
5	PEG	B	402	-	6,6,6	0.22	0	5,5,5	0.12	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	POV	A	401	-	-	33/55/55/55	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PEG	B	403	-	-	1/4/4/4	-
5	PEG	B	404	-	-	1/4/4/4	-
3	GOL	B	405	-	-	0/4/4/4	-
4	EDO	A	403	-	-	0/1/1/1	-
2	POV	B	401	-	-	23/55/55/55	-
3	GOL	A	402	-	-	2/4/4/4	-
5	PEG	B	402	-	-	3/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	POV	O13-P-O14	2.19	122.63	112.44

There are no chirality outliers.

All (63) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	POV	C11-O12-P-O11
2	A	401	POV	C11-O12-P-O13
2	A	401	POV	C11-O12-P-O14
2	A	401	POV	C22-C21-O21-C2
2	B	401	POV	C22-C21-O21-C2
3	A	402	GOL	O1-C1-C2-O2
3	A	402	GOL	O1-C1-C2-C3
2	A	401	POV	O22-C21-O21-C2
2	B	401	POV	O22-C21-O21-C2
5	B	402	PEG	O1-C1-C2-O2
5	B	404	PEG	O2-C3-C4-O4
2	A	401	POV	C21-C22-C23-C24
2	B	401	POV	C23-C24-C25-C26
2	B	401	POV	C211-C212-C213-C214
2	A	401	POV	C211-C212-C213-C214
2	B	401	POV	C36-C37-C38-C39
2	A	401	POV	C34-C35-C36-C37
2	A	401	POV	C35-C36-C37-C38
2	B	401	POV	C37-C38-C39-C310
2	A	401	POV	C32-C33-C34-C35
2	A	401	POV	C36-C37-C38-C39
2	A	401	POV	C39-C310-C311-C312

*Continued on next page...*

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Mol	Chain	Res	Type	Atoms
2	A	401	POV	C312-C313-C314-C315
2	A	401	POV	C215-C216-C217-C218
2	A	401	POV	C23-C24-C25-C26
2	A	401	POV	C210-C211-C212-C213
2	A	401	POV	C214-C215-C216-C217
5	B	402	PEG	O2-C3-C4-O4
2	B	401	POV	C2-C3-O31-C31
2	A	401	POV	C313-C314-C315-C316
2	B	401	POV	C25-C26-C27-C28
2	B	401	POV	C1-C2-C3-O31
2	B	401	POV	C22-C23-C24-C25
2	B	401	POV	C213-C214-C215-C216
2	B	401	POV	C39-C310-C311-C312
2	A	401	POV	C25-C26-C27-C28
2	A	401	POV	C1-C2-C3-O31
2	A	401	POV	C37-C38-C39-C310
2	A	401	POV	C11-C12-N-C15
2	A	401	POV	O12-C11-C12-N
2	A	401	POV	C33-C34-C35-C36
2	B	401	POV	C35-C36-C37-C38
2	A	401	POV	O11-C1-C2-C3
2	A	401	POV	O11-C1-C2-O21
2	A	401	POV	C310-C311-C312-C313
2	A	401	POV	O21-C2-C3-O31
2	B	401	POV	O21-C2-C3-O31
2	B	401	POV	C32-C33-C34-C35
5	B	402	PEG	C1-C2-O2-C3
2	B	401	POV	C11-O12-P-O14
2	B	401	POV	C24-C25-C26-C27
2	B	401	POV	C32-C31-O31-C3
2	B	401	POV	O32-C31-O31-C3
2	B	401	POV	O11-C1-C2-O21
2	A	401	POV	C27-C28-C29-C210
2	A	401	POV	C11-C12-N-C14
2	B	401	POV	C26-C27-C28-C29
2	B	401	POV	C31-C32-C33-C34
2	A	401	POV	C11-C12-N-C13
5	B	403	PEG	C1-C2-O2-C3
2	A	401	POV	O31-C31-C32-C33
2	A	401	POV	O32-C31-C32-C33
2	B	401	POV	O11-C1-C2-C3

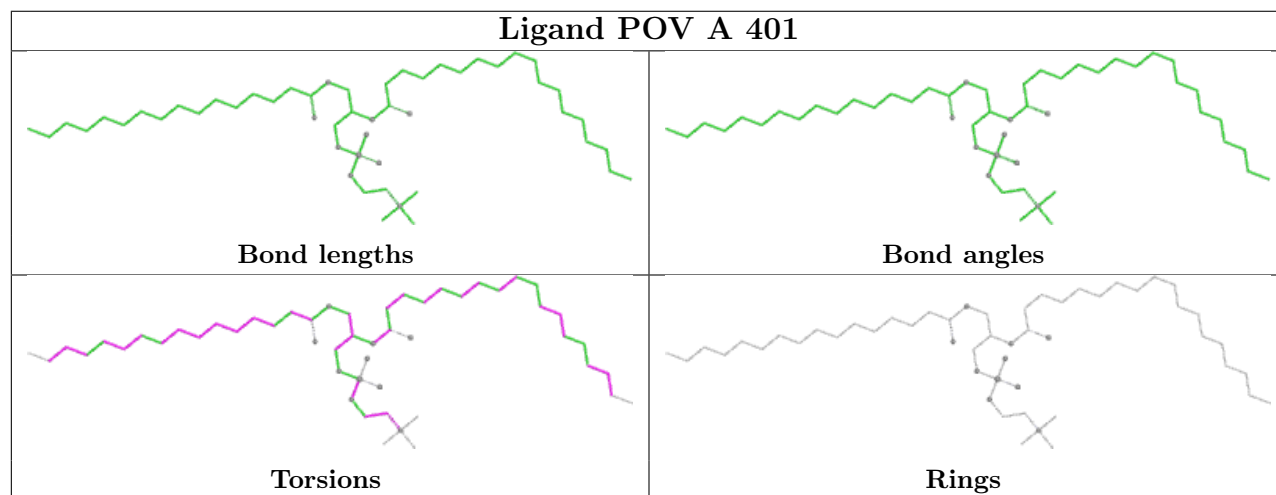
There are no ring outliers.

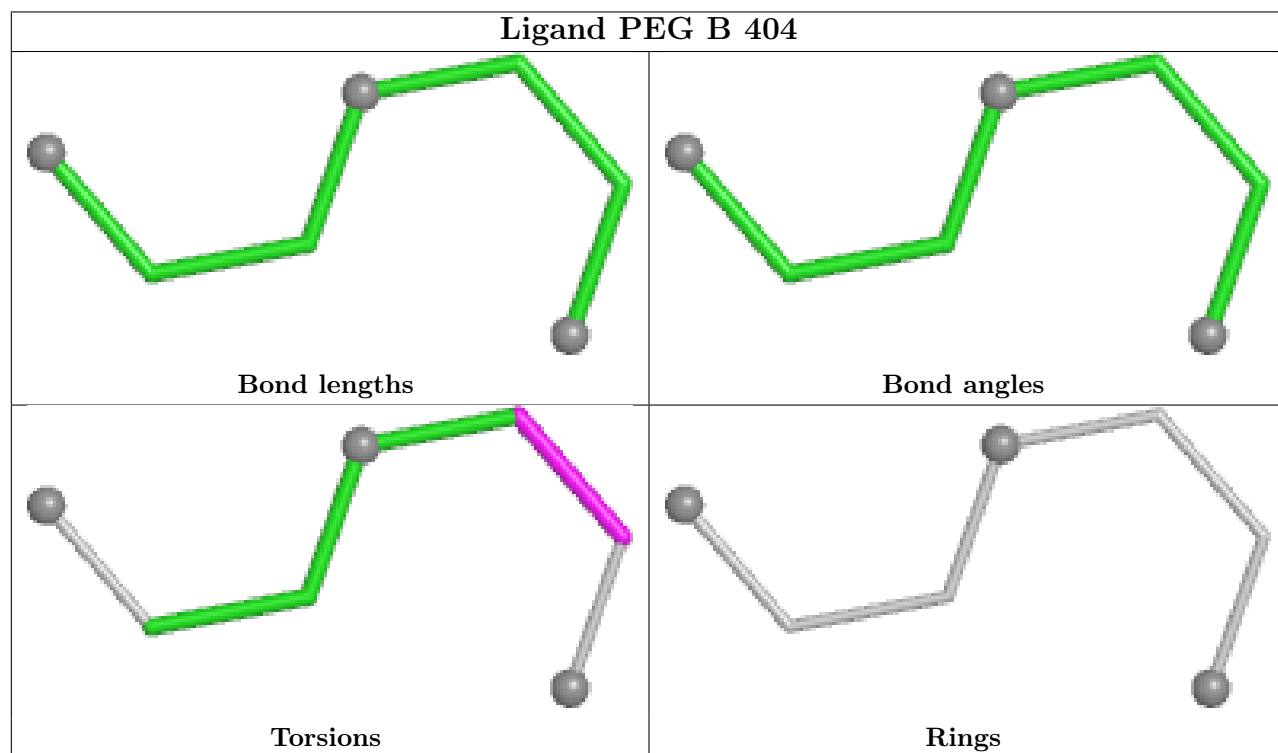
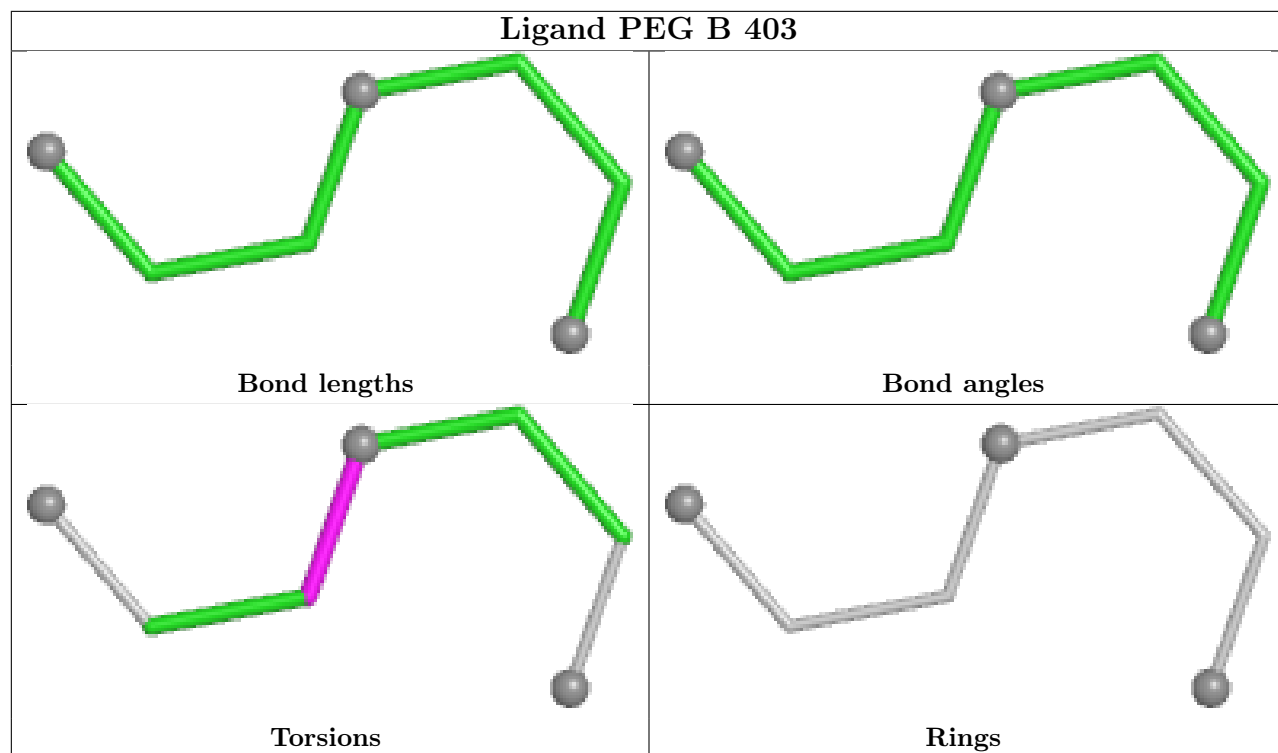


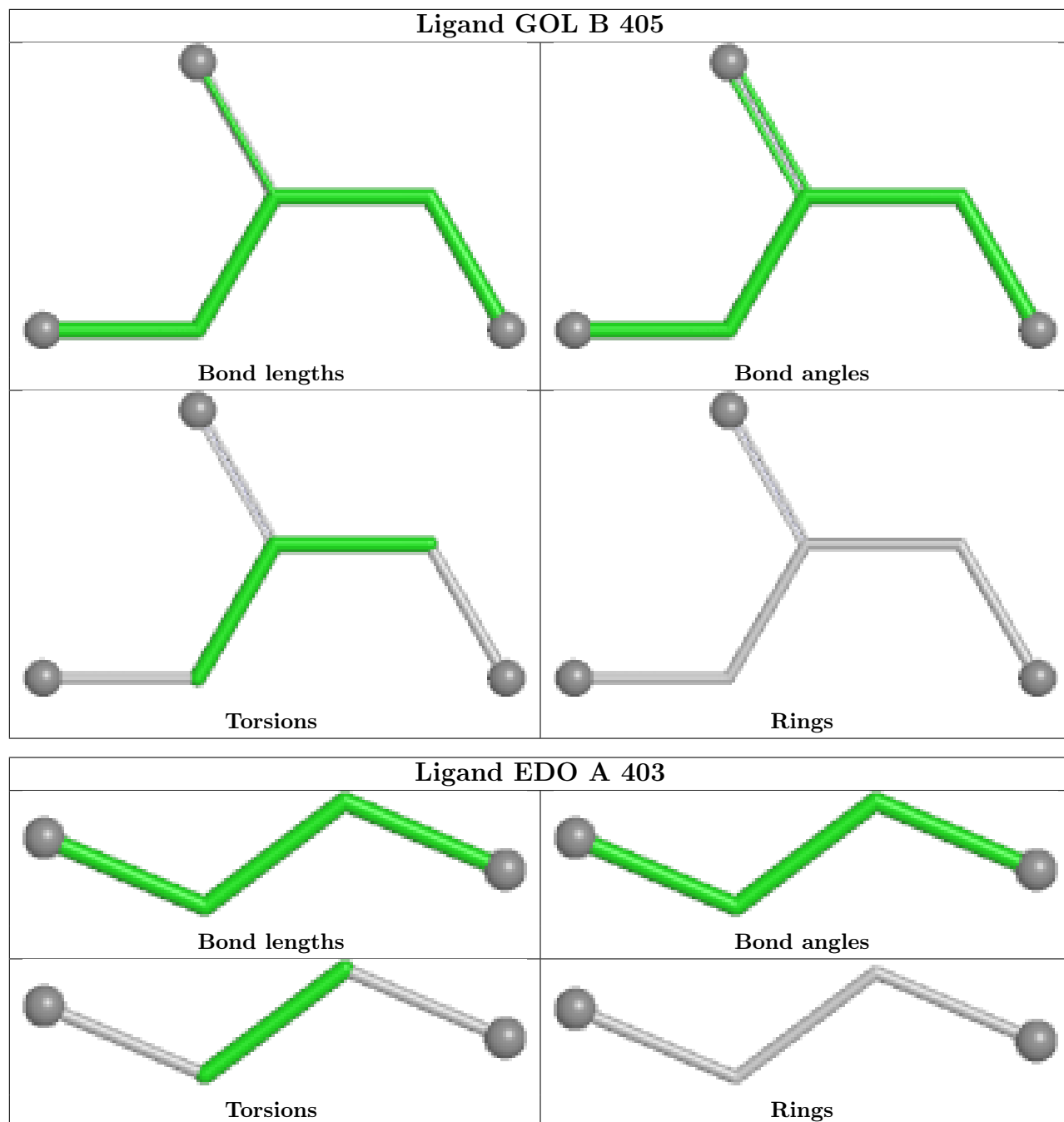
5 monomers are involved in 22 short contacts:

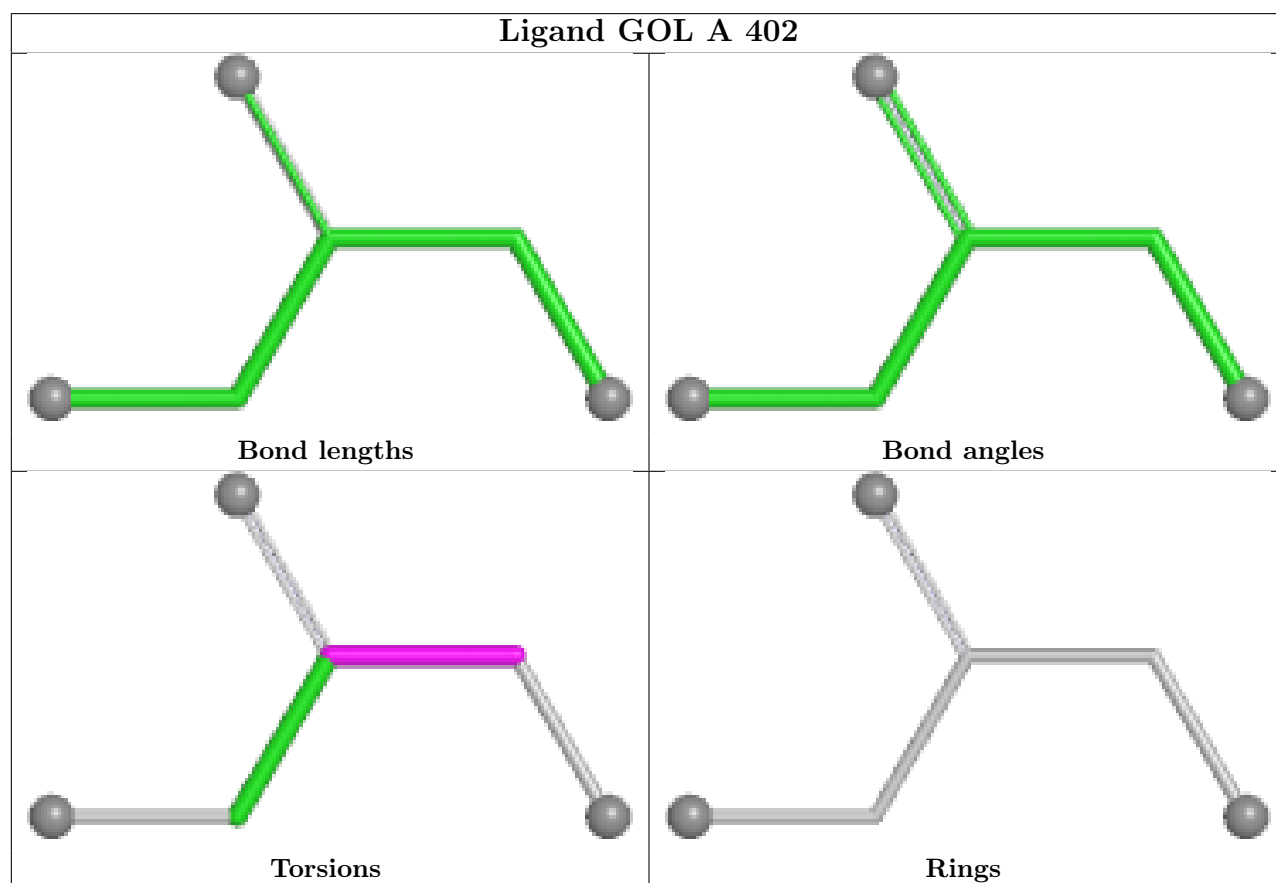
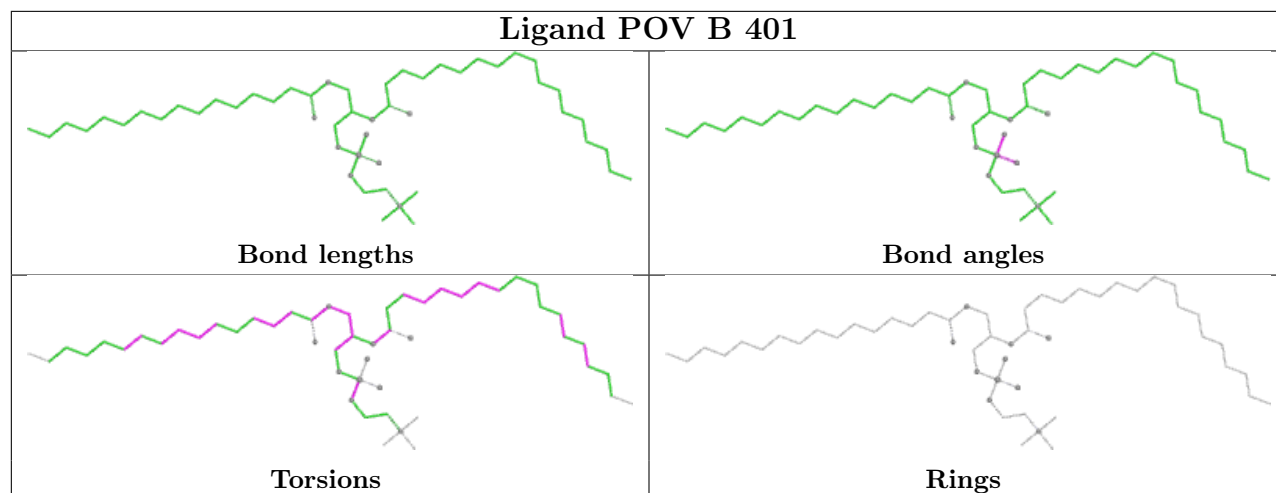
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	POV	9	0
5	B	403	PEG	1	0
5	B	404	PEG	5	0
2	B	401	POV	5	0
3	A	402	GOL	2	0

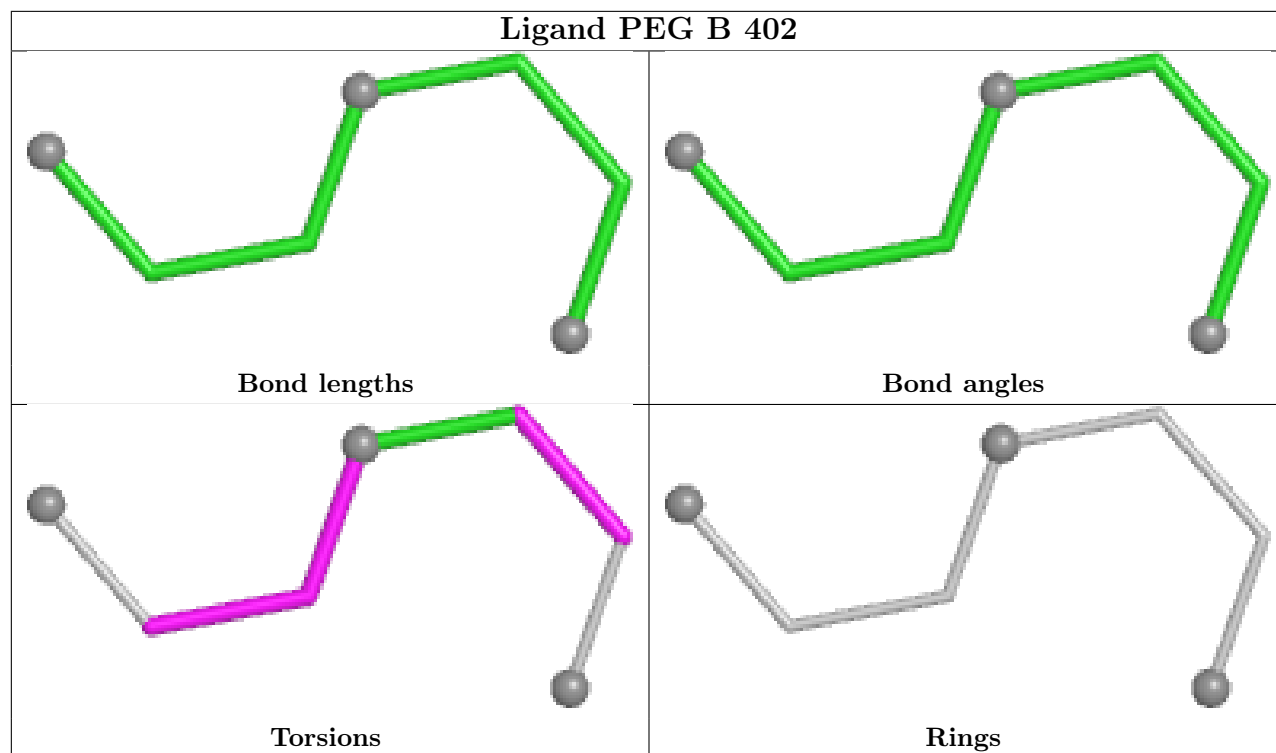
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	288/304 (94%)	-1.36	0 100 100	16, 27, 37, 60	2 (0%)
1	B	289/304 (95%)	-1.36	0 100 100	20, 28, 37, 52	0
All	All	577/608 (94%)	-1.36	0 100 100	16, 28, 37, 60	2 (0%)

There are no RSRZ outliers to report.

### 6.2 Non-standard residues in protein, DNA, RNA chains [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, 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
1	CME	A	17	10/11	0.99	0.03	20,25,31,34	0
1	CME	B	17	10/11	0.99	0.04	28,32,44,53	0

### 6.3 Carbohydrates [i](#)

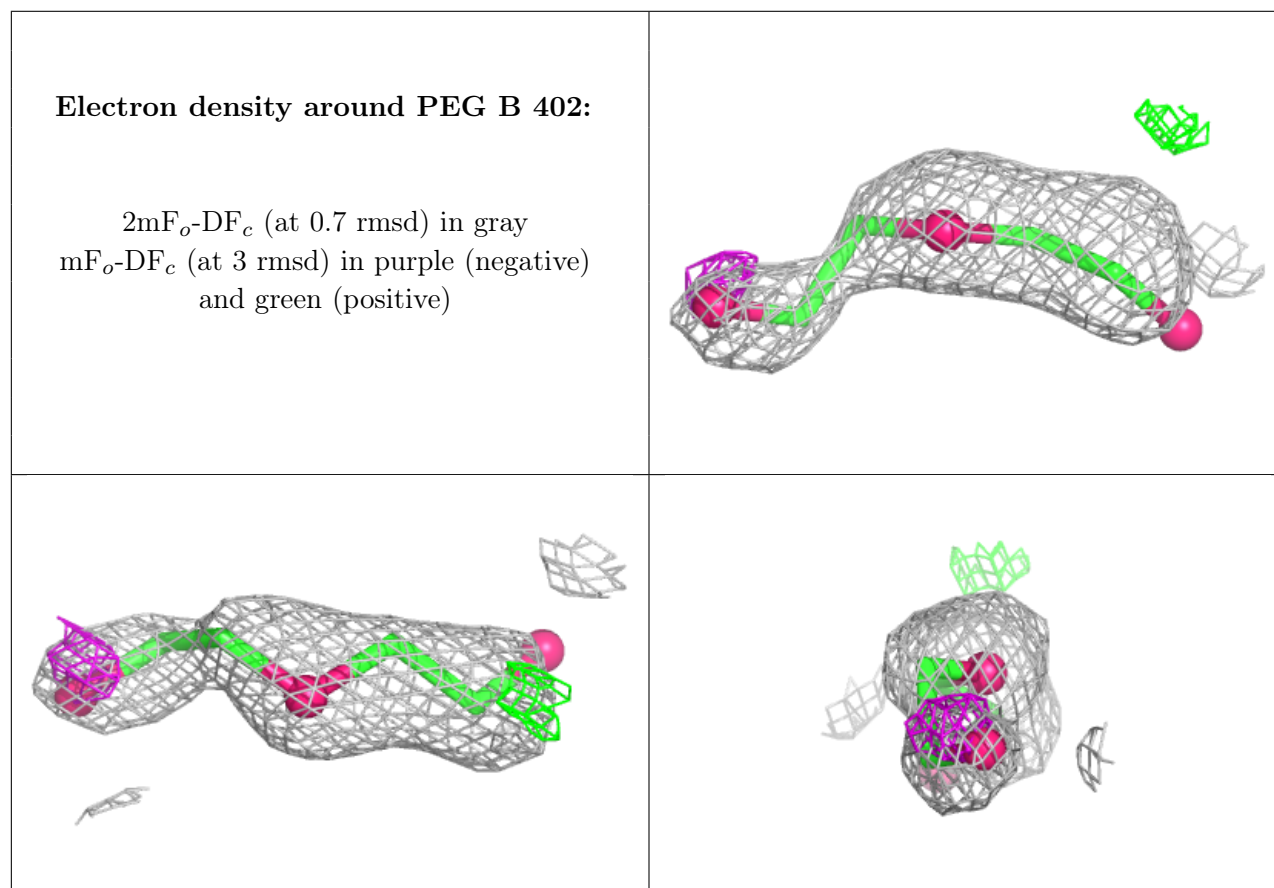
There are no oligosaccharides in this entry.

### 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, 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.

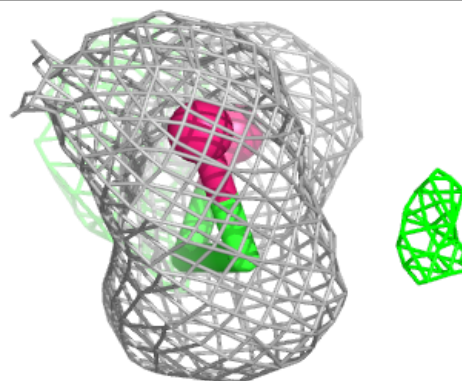
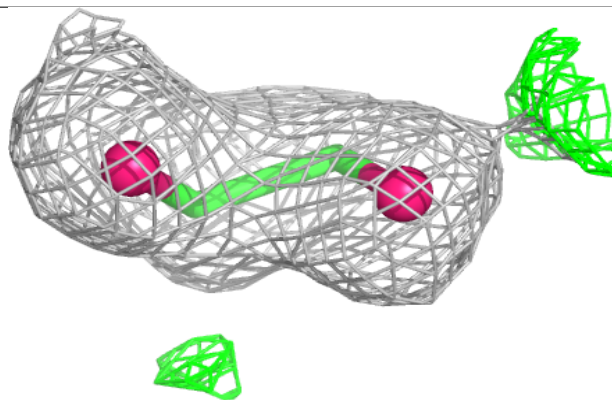
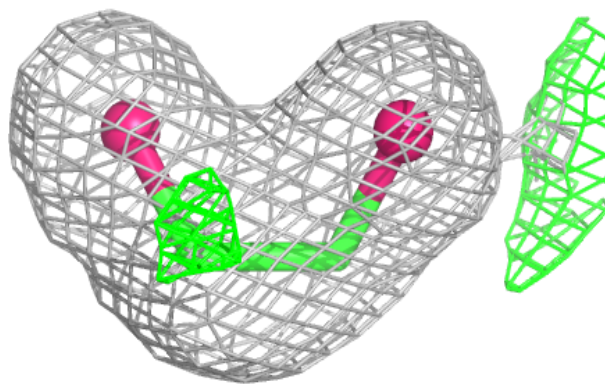
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	PEG	B	402	7/7	0.97	0.06	33,37,44,45	0
4	EDO	A	403	4/4	0.98	0.04	29,30,31,34	0
3	GOL	A	402	6/6	0.98	0.04	34,36,39,45	0
5	PEG	B	403	7/7	0.98	0.05	30,32,35,37	0
2	POV	B	401	52/52	0.99	0.04	26,32,36,41	0
2	POV	A	401	52/52	0.99	0.05	24,31,38,42	0
3	GOL	B	405	6/6	0.99	0.04	31,33,35,41	0
5	PEG	B	404	7/7	0.99	0.04	25,30,33,38	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 EDO A 403:**

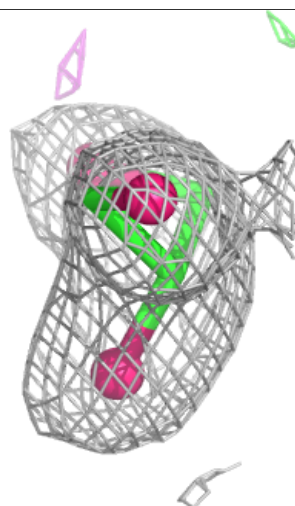
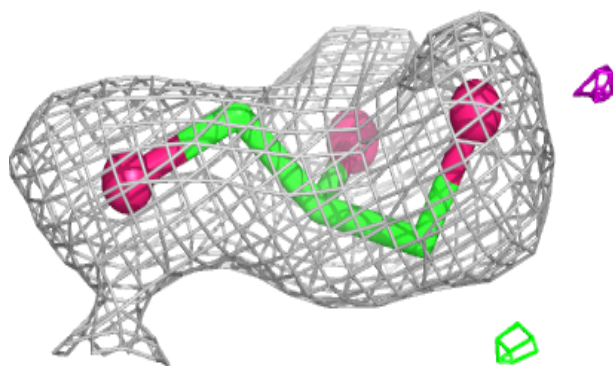
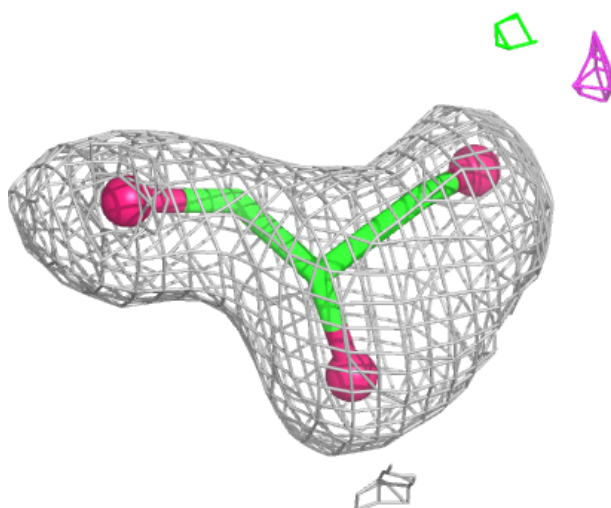
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





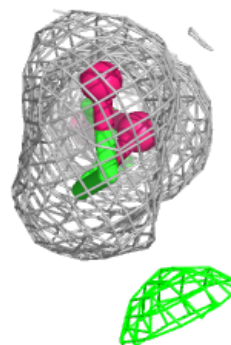
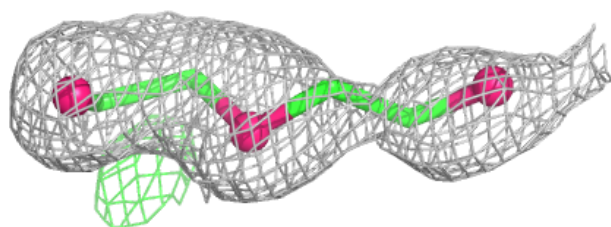
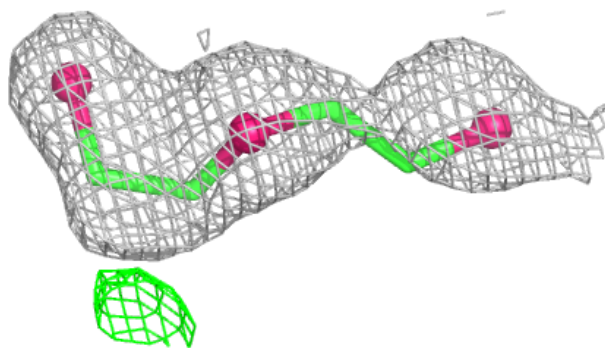
**Electron density around GOL A 402:**

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

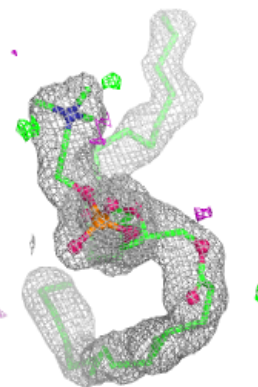
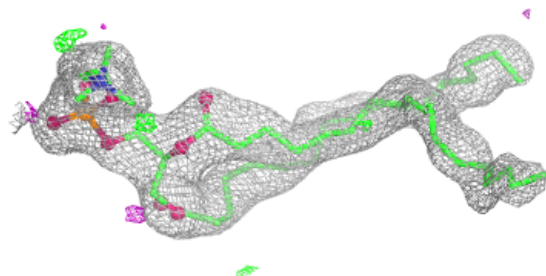
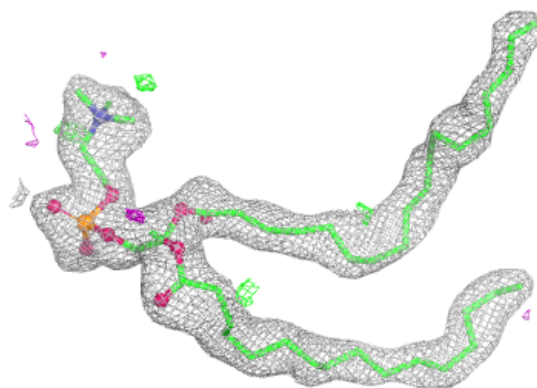


**Electron density around PEG B 403:**

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

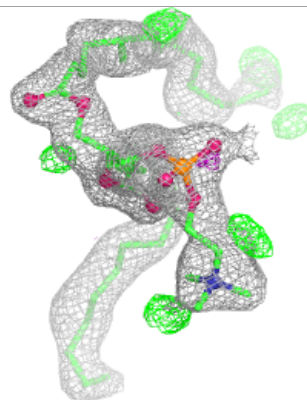
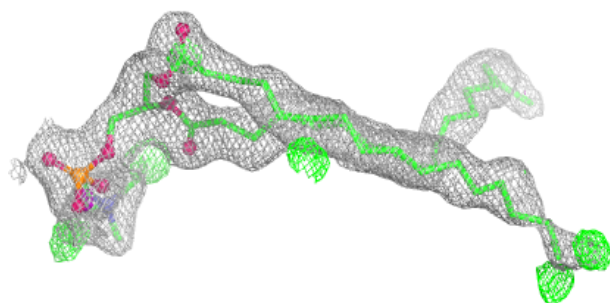
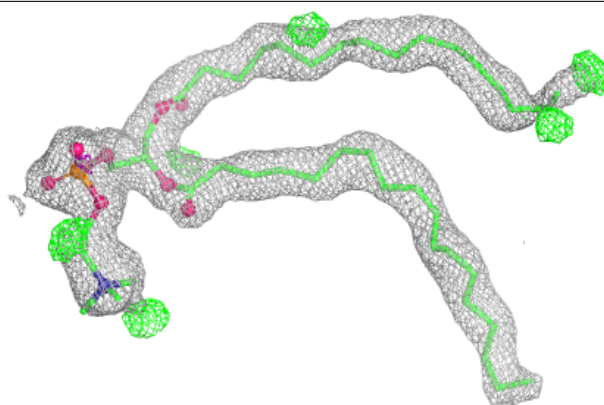
**Electron density around POV B 401:**

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



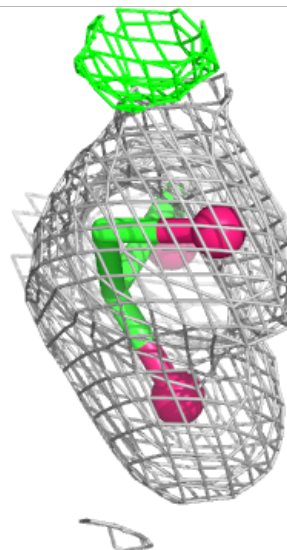
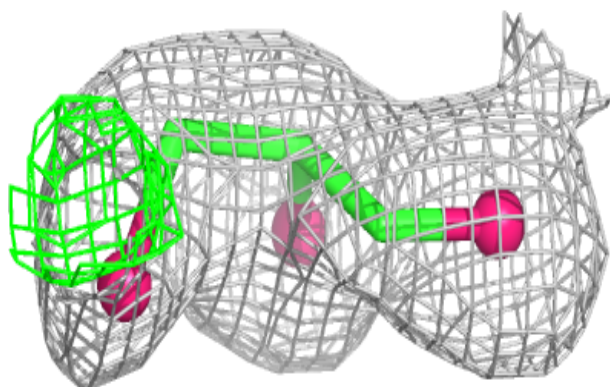
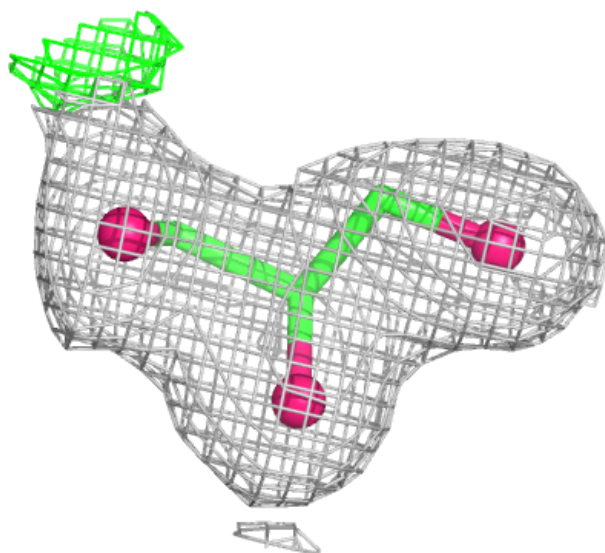
**Electron density around POV A 401:**

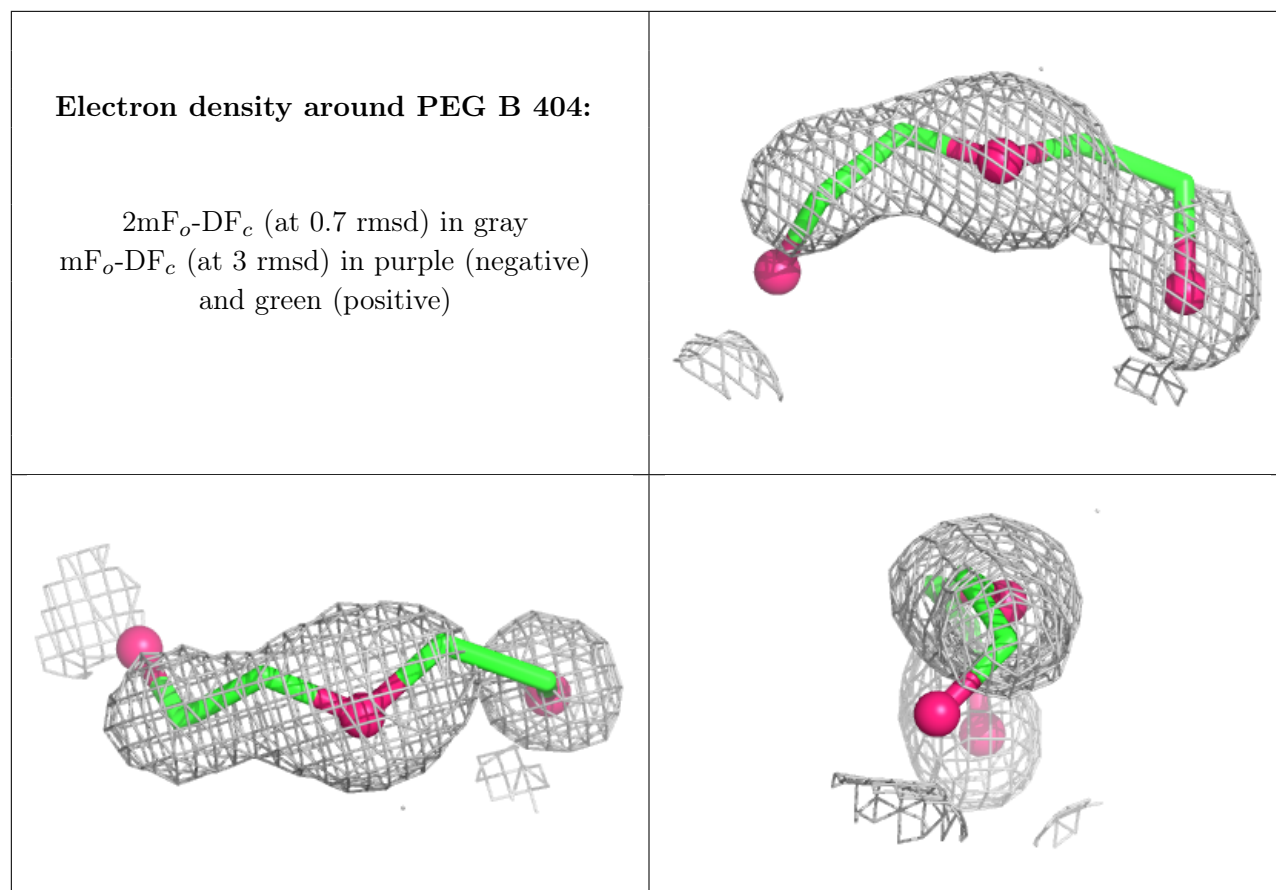
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around GOL B 405:**

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





## 6.5 Other polymers ⓘ

There are no such residues in this entry.