



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jul 31, 2025 – 10:26 am BST

PDB ID : 9GTI / pdb\_00009gti  
Title : X-ray crystal structure of mouse NPTN N-terminal domain  
Authors : Vinayagam, D.; Raunser, S.; Sistel, O.; Shulte, U.; Constantin, C.E.;  
Prubaum, D.; Zolles, G.; Fakler, B.  
Deposited on : 2024-09-17  
Resolution : 2.03 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 Phenix2.0rc1  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 2.0rc1  
EDS : 3.0  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.006 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.45.1

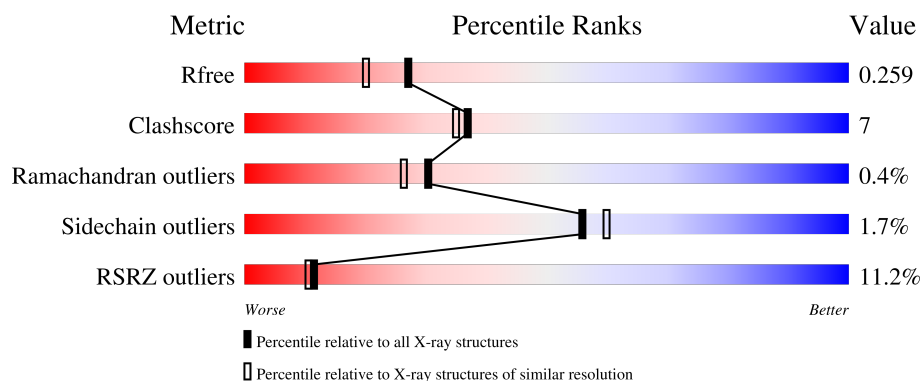
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.03 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	12358 (2.04-2.00)
Clashscore	180529	13897 (2.04-2.00)
Ramachandran outliers	177936	13770 (2.04-2.00)
Sidechain outliers	177891	13769 (2.04-2.00)
RSRZ outliers	164620	12358 (2.04-2.00)

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	161	
1	B	161	
1	C	161	
1	D	161	

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
2	PEG	A	205	-	-	X	-
2	PEG	D	202	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4059 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 Neuroplastin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	118	Total	C	N	O	S	0	0	0
			938	584	169	182	3			
1	B	120	Total	C	N	O	S	0	0	0
			958	595	175	185	3			
1	C	118	Total	C	N	O	S	0	0	0
			938	582	171	182	3			
1	D	117	Total	C	N	O	S	0	0	0
			926	576	167	180	3			

There are 168 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP P97300
A	1	GLU	-	expression tag	UNP P97300
A	2	THR	-	expression tag	UNP P97300
A	3	ASP	-	expression tag	UNP P97300
A	4	THR	-	expression tag	UNP P97300
A	5	LEU	-	expression tag	UNP P97300
A	6	LEU	-	expression tag	UNP P97300
A	7	LEU	-	expression tag	UNP P97300
A	8	TRP	-	expression tag	UNP P97300
A	9	VAL	-	expression tag	UNP P97300
A	10	LEU	-	expression tag	UNP P97300
A	11	LEU	-	expression tag	UNP P97300
A	12	LEU	-	expression tag	UNP P97300
A	13	TRP	-	expression tag	UNP P97300
A	14	VAL	-	expression tag	UNP P97300
A	15	PRO	-	expression tag	UNP P97300
A	16	GLY	-	expression tag	UNP P97300
A	17	SER	-	expression tag	UNP P97300
A	18	THR	-	expression tag	UNP P97300
A	19	GLY	-	expression tag	UNP P97300
A	20	ASP	-	expression tag	UNP P97300

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Chain	Residue	Modelled	Actual	Comment	Reference
A	21	ALA	-	expression tag	UNP P97300
A	22	ALA	-	expression tag	UNP P97300
A	23	GLN	-	expression tag	UNP P97300
A	24	PRO	-	expression tag	UNP P97300
A	25	ALA	-	expression tag	UNP P97300
A	26	ALA	-	expression tag	UNP P97300
A	27	ARG	-	expression tag	UNP P97300
A	28	ASP	-	expression tag	UNP P97300
A	148	GLU	-	expression tag	UNP P97300
A	149	ASP	-	expression tag	UNP P97300
A	150	LEU	-	expression tag	UNP P97300
A	151	TYR	-	expression tag	UNP P97300
A	152	PHE	-	expression tag	UNP P97300
A	153	GLN	-	expression tag	UNP P97300
A	154	SER	-	expression tag	UNP P97300
A	155	HIS	-	expression tag	UNP P97300
A	156	HIS	-	expression tag	UNP P97300
A	157	HIS	-	expression tag	UNP P97300
A	158	HIS	-	expression tag	UNP P97300
A	159	HIS	-	expression tag	UNP P97300
A	160	HIS	-	expression tag	UNP P97300
B	0	MET	-	initiating methionine	UNP P97300
B	1	GLU	-	expression tag	UNP P97300
B	2	THR	-	expression tag	UNP P97300
B	3	ASP	-	expression tag	UNP P97300
B	4	THR	-	expression tag	UNP P97300
B	5	LEU	-	expression tag	UNP P97300
B	6	LEU	-	expression tag	UNP P97300
B	7	LEU	-	expression tag	UNP P97300
B	8	TRP	-	expression tag	UNP P97300
B	9	VAL	-	expression tag	UNP P97300
B	10	LEU	-	expression tag	UNP P97300
B	11	LEU	-	expression tag	UNP P97300
B	12	LEU	-	expression tag	UNP P97300
B	13	TRP	-	expression tag	UNP P97300
B	14	VAL	-	expression tag	UNP P97300
B	15	PRO	-	expression tag	UNP P97300
B	16	GLY	-	expression tag	UNP P97300
B	17	SER	-	expression tag	UNP P97300
B	18	THR	-	expression tag	UNP P97300
B	19	GLY	-	expression tag	UNP P97300
B	20	ASP	-	expression tag	UNP P97300

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Chain	Residue	Modelled	Actual	Comment	Reference
B	21	ALA	-	expression tag	UNP P97300
B	22	ALA	-	expression tag	UNP P97300
B	23	GLN	-	expression tag	UNP P97300
B	24	PRO	-	expression tag	UNP P97300
B	25	ALA	-	expression tag	UNP P97300
B	26	ALA	-	expression tag	UNP P97300
B	27	ARG	-	expression tag	UNP P97300
B	28	ASP	-	expression tag	UNP P97300
B	148	GLU	-	expression tag	UNP P97300
B	149	ASP	-	expression tag	UNP P97300
B	150	LEU	-	expression tag	UNP P97300
B	151	TYR	-	expression tag	UNP P97300
B	152	PHE	-	expression tag	UNP P97300
B	153	GLN	-	expression tag	UNP P97300
B	154	SER	-	expression tag	UNP P97300
B	155	HIS	-	expression tag	UNP P97300
B	156	HIS	-	expression tag	UNP P97300
B	157	HIS	-	expression tag	UNP P97300
B	158	HIS	-	expression tag	UNP P97300
B	159	HIS	-	expression tag	UNP P97300
B	160	HIS	-	expression tag	UNP P97300
C	0	MET	-	initiating methionine	UNP P97300
C	1	GLU	-	expression tag	UNP P97300
C	2	THR	-	expression tag	UNP P97300
C	3	ASP	-	expression tag	UNP P97300
C	4	THR	-	expression tag	UNP P97300
C	5	LEU	-	expression tag	UNP P97300
C	6	LEU	-	expression tag	UNP P97300
C	7	LEU	-	expression tag	UNP P97300
C	8	TRP	-	expression tag	UNP P97300
C	9	VAL	-	expression tag	UNP P97300
C	10	LEU	-	expression tag	UNP P97300
C	11	LEU	-	expression tag	UNP P97300
C	12	LEU	-	expression tag	UNP P97300
C	13	TRP	-	expression tag	UNP P97300
C	14	VAL	-	expression tag	UNP P97300
C	15	PRO	-	expression tag	UNP P97300
C	16	GLY	-	expression tag	UNP P97300
C	17	SER	-	expression tag	UNP P97300
C	18	THR	-	expression tag	UNP P97300
C	19	GLY	-	expression tag	UNP P97300
C	20	ASP	-	expression tag	UNP P97300

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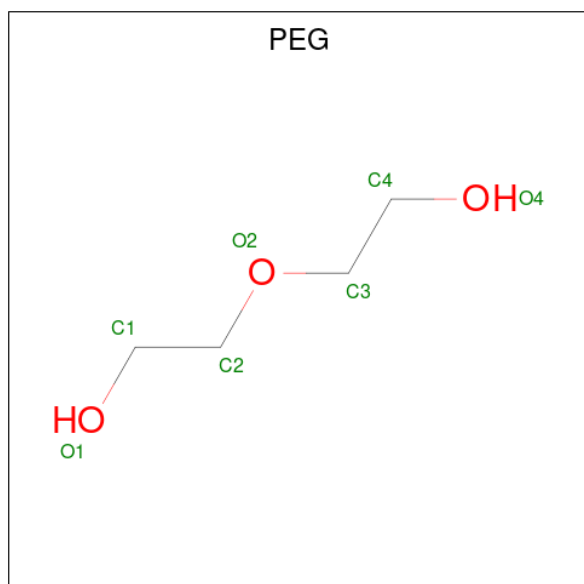
Chain	Residue	Modelled	Actual	Comment	Reference
C	21	ALA	-	expression tag	UNP P97300
C	22	ALA	-	expression tag	UNP P97300
C	23	GLN	-	expression tag	UNP P97300
C	24	PRO	-	expression tag	UNP P97300
C	25	ALA	-	expression tag	UNP P97300
C	26	ALA	-	expression tag	UNP P97300
C	27	ARG	-	expression tag	UNP P97300
C	28	ASP	-	expression tag	UNP P97300
C	148	GLU	-	expression tag	UNP P97300
C	149	ASP	-	expression tag	UNP P97300
C	150	LEU	-	expression tag	UNP P97300
C	151	TYR	-	expression tag	UNP P97300
C	152	PHE	-	expression tag	UNP P97300
C	153	GLN	-	expression tag	UNP P97300
C	154	SER	-	expression tag	UNP P97300
C	155	HIS	-	expression tag	UNP P97300
C	156	HIS	-	expression tag	UNP P97300
C	157	HIS	-	expression tag	UNP P97300
C	158	HIS	-	expression tag	UNP P97300
C	159	HIS	-	expression tag	UNP P97300
C	160	HIS	-	expression tag	UNP P97300
D	0	MET	-	initiating methionine	UNP P97300
D	1	GLU	-	expression tag	UNP P97300
D	2	THR	-	expression tag	UNP P97300
D	3	ASP	-	expression tag	UNP P97300
D	4	THR	-	expression tag	UNP P97300
D	5	LEU	-	expression tag	UNP P97300
D	6	LEU	-	expression tag	UNP P97300
D	7	LEU	-	expression tag	UNP P97300
D	8	TRP	-	expression tag	UNP P97300
D	9	VAL	-	expression tag	UNP P97300
D	10	LEU	-	expression tag	UNP P97300
D	11	LEU	-	expression tag	UNP P97300
D	12	LEU	-	expression tag	UNP P97300
D	13	TRP	-	expression tag	UNP P97300
D	14	VAL	-	expression tag	UNP P97300
D	15	PRO	-	expression tag	UNP P97300
D	16	GLY	-	expression tag	UNP P97300
D	17	SER	-	expression tag	UNP P97300
D	18	THR	-	expression tag	UNP P97300
D	19	GLY	-	expression tag	UNP P97300
D	20	ASP	-	expression tag	UNP P97300

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Chain	Residue	Modelled	Actual	Comment	Reference
D	21	ALA	-	expression tag	UNP P97300
D	22	ALA	-	expression tag	UNP P97300
D	23	GLN	-	expression tag	UNP P97300
D	24	PRO	-	expression tag	UNP P97300
D	25	ALA	-	expression tag	UNP P97300
D	26	ALA	-	expression tag	UNP P97300
D	27	ARG	-	expression tag	UNP P97300
D	28	ASP	-	expression tag	UNP P97300
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D	152	PHE	-	expression tag	UNP P97300
D	153	GLN	-	expression tag	UNP P97300
D	154	SER	-	expression tag	UNP P97300
D	155	HIS	-	expression tag	UNP P97300
D	156	HIS	-	expression tag	UNP P97300
D	157	HIS	-	expression tag	UNP P97300
D	158	HIS	-	expression tag	UNP P97300
D	159	HIS	-	expression tag	UNP P97300
D	160	HIS	-	expression tag	UNP P97300

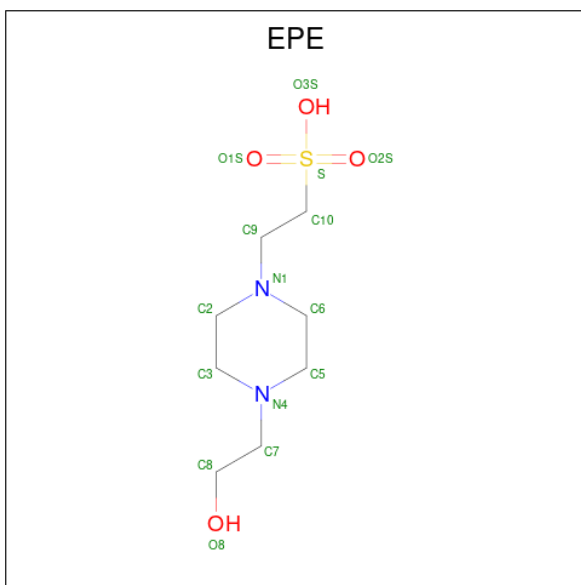
- Molecule 2 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula:  $C_4H_{10}O_3$ ) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 7 4 3	0	0
2	A	1	Total C O 7 4 3	0	0
2	A	1	Total C O 7 4 3	0	0
2	A	1	Total C O 7 4 3	0	0
2	A	1	Total C O 7 4 3	0	0
2	B	1	Total C O 7 4 3	0	0
2	B	1	Total C O 7 4 3	0	0
2	B	1	Total C O 7 4 3	0	0
2	B	1	Total C O 7 4 3	0	0
2	C	1	Total C O 7 4 3	0	0
2	C	1	Total C O 7 4 3	0	0
2	C	1	Total C O 7 4 3	0	0
2	C	1	Total C O 7 4 3	0	0
2	D	1	Total C O 7 4 3	0	0
2	D	1	Total C O 7 4 3	0	0

- Molecule 3 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: C<sub>8</sub>H<sub>18</sub>N<sub>2</sub>O<sub>4</sub>S) (labeled as "Ligand of Interest" by depositor).

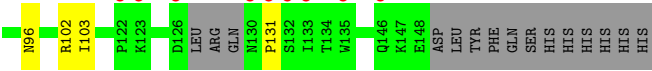
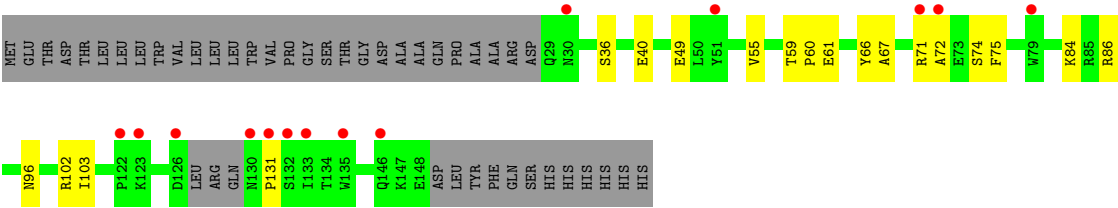


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	D	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	50	Total	O	0	0
			50	50		
4	B	58	Total	O	0	0
			58	58		
4	C	34	Total	O	0	0
			34	34		
4	D	37	Total	O	0	0
			37	37		





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.91Å 85.91Å 163.28Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.93 – 2.03 43.93 – 2.03	Depositor EDS
% Data completeness (in resolution range)	99.8 (43.93-2.03) 99.8 (43.93-2.03)	Depositor EDS
$R_{merge}$	0.20	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.32 (at 2.03Å)	Xtriage
Refinement program	PHENIX 1.18.2-3874	Depositor
R, $R_{free}$	0.215 , 0.248 0.227 , 0.259	Depositor DCC
$R_{free}$ test set	2286 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.1	Xtriage
Anisotropy	0.195	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 50.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4059	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.88% 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: EPE, PEG

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	0.35	0/956	0.49	0/1297
1	B	0.39	0/977	0.59	0/1326
1	C	0.34	0/955	0.50	0/1294
1	D	0.34	0/944	0.49	0/1281
All	All	0.35	0/3832	0.52	0/5198

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	C	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	128	ARG	Sidechain
1	C	27	ARG	Sidechain

## 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	938	0	912	14	0
1	B	958	0	934	17	0
1	C	938	0	910	8	0
1	D	926	0	895	14	0
2	A	35	0	50	8	0
2	B	28	0	40	8	0
2	C	28	0	40	3	0
2	D	14	0	20	5	0
3	D	15	0	18	0	0
4	A	50	0	0	1	0
4	B	58	0	0	0	0
4	C	34	0	0	1	0
4	D	37	0	0	2	0
All	All	4059	0	3819	54	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.

The worst 5 of 54 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:79:TRP:HE1	2:B:204:PEG:H12	1.25	1.01
1:A:131:PRO:HD2	2:A:205:PEG:H32	1.60	0.83
1:C:84:LYS:HG3	2:C:204:PEG:H42	1.62	0.81
1:A:75:PHE:O	2:A:204:PEG:H21	1.84	0.78
1:D:49:GLU:H	2:D:202:PEG:H42	1.57	0.70

There are no symmetry-related clashes.

## 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	114/161 (71%)	113 (99%)	1 (1%)	0	100	100
1	B	118/161 (73%)	115 (98%)	3 (2%)	0	100	100
1	C	114/161 (71%)	111 (97%)	2 (2%)	1 (1%)	14	9
1	D	113/161 (70%)	109 (96%)	3 (3%)	1 (1%)	14	9
All	All	459/644 (71%)	448 (98%)	9 (2%)	2 (0%)	30	26

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	147	LYS
1	D	131	PRO

### 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	102/139 (73%)	102 (100%)	0	100	100
1	B	104/139 (75%)	100 (96%)	4 (4%)	28	26
1	C	100/139 (72%)	98 (98%)	2 (2%)	50	53
1	D	100/139 (72%)	99 (99%)	1 (1%)	73	77
All	All	406/556 (73%)	399 (98%)	7 (2%)	56	60

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	148	GLU
1	C	28	ASP
1	D	36	SER
1	C	148	GLU
1	B	147	LYS



Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	130	ASN
1	D	139	GLN
1	C	90	ASN
1	C	125	ASN
1	D	70	ASN

### 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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

16 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	PEG	B	203	-	6,6,6	0.18	0	5,5,5	0.37	0
2	PEG	B	202	-	6,6,6	0.24	0	5,5,5	0.54	0
2	PEG	A	205	-	6,6,6	0.28	0	5,5,5	0.73	0
2	PEG	C	202	-	6,6,6	0.22	0	5,5,5	0.31	0
2	PEG	A	201	-	6,6,6	0.22	0	5,5,5	0.32	0
2	PEG	C	203	-	6,6,6	0.21	0	5,5,5	0.39	0
3	EPE	D	203	-	15,15,15	0.80	1 (6%)	18,20,20	1.80	4 (22%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	PEG	B	201	-	6,6,6	0.24	0	5,5,5	0.59	0
2	PEG	A	203	-	6,6,6	0.19	0	5,5,5	0.54	0
2	PEG	A	202	-	6,6,6	0.19	0	5,5,5	0.38	0
2	PEG	C	201	-	6,6,6	0.25	0	5,5,5	0.60	0
2	PEG	C	204	-	6,6,6	0.37	0	5,5,5	0.43	0
2	PEG	D	201	-	6,6,6	0.20	0	5,5,5	0.31	0
2	PEG	D	202	-	6,6,6	0.43	0	5,5,5	0.86	0
2	PEG	B	204	-	6,6,6	0.27	0	5,5,5	0.37	0
2	PEG	A	204	-	6,6,6	0.26	0	5,5,5	0.73	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	PEG	B	203	-	-	3/4/4/4	-
2	PEG	B	202	-	-	3/4/4/4	-
2	PEG	A	205	-	-	3/4/4/4	-
2	PEG	C	202	-	-	1/4/4/4	-
2	PEG	A	201	-	-	1/4/4/4	-
2	PEG	C	203	-	-	0/4/4/4	-
3	EPE	D	203	-	-	5/9/19/19	0/1/1/1
2	PEG	B	201	-	-	1/4/4/4	-
2	PEG	A	203	-	-	1/4/4/4	-
2	PEG	A	202	-	-	2/4/4/4	-
2	PEG	C	201	-	-	4/4/4/4	-
2	PEG	C	204	-	-	2/4/4/4	-
2	PEG	D	201	-	-	1/4/4/4	-
2	PEG	D	202	-	-	3/4/4/4	-
2	PEG	B	204	-	-	3/4/4/4	-
2	PEG	A	204	-	-	1/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	203	EPE	C10-S	2.33	1.80	1.77

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	203	EPE	C2-C3-N4	4.30	119.47	110.64
3	D	203	EPE	C5-N4-C3	3.25	116.15	108.83
3	D	203	EPE	C3-C2-N1	2.77	116.33	110.64
3	D	203	EPE	O2S-S-C10	-2.48	103.93	106.92

There are no chirality outliers.

5 of 34 torsion outliers are listed below:

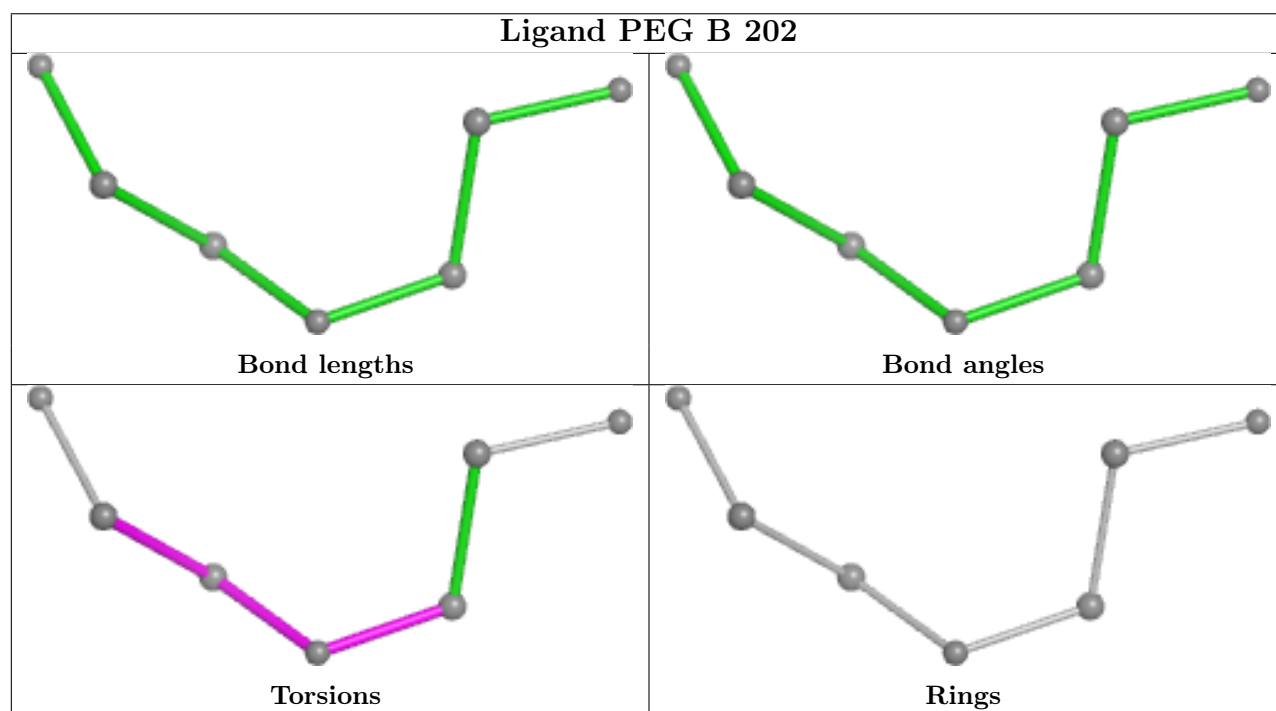
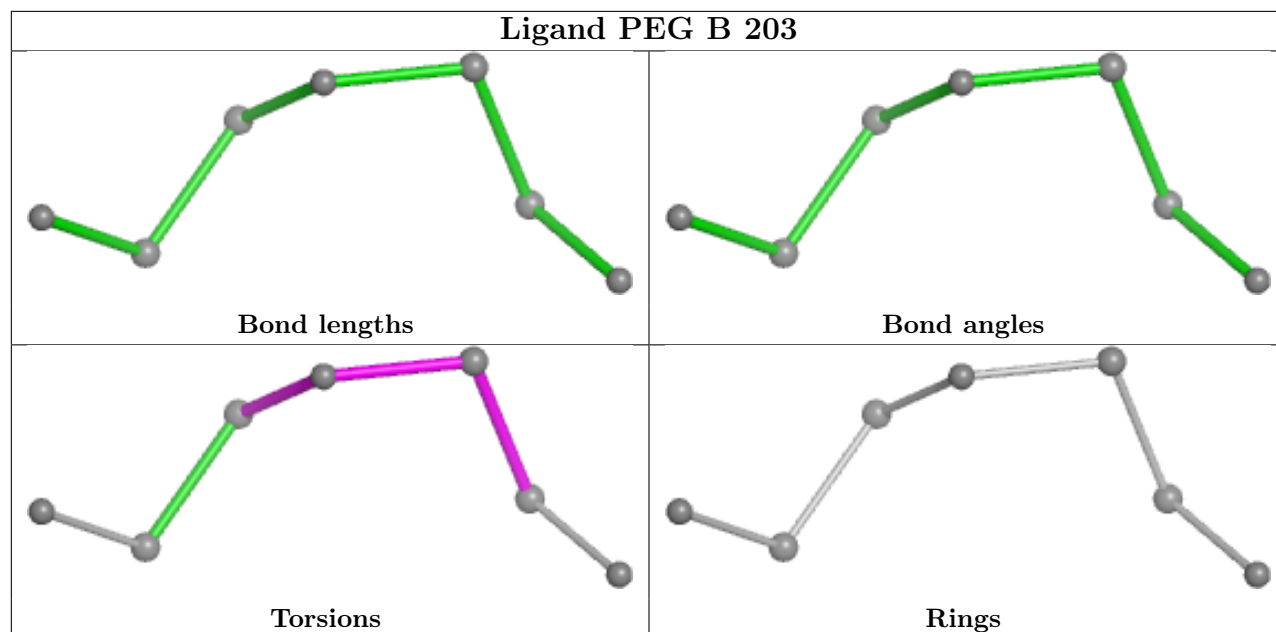
Mol	Chain	Res	Type	Atoms
2	A	202	PEG	O2-C3-C4-O4
2	A	205	PEG	O2-C3-C4-O4
2	B	203	PEG	O2-C3-C4-O4
2	C	204	PEG	O2-C3-C4-O4
2	A	203	PEG	O2-C3-C4-O4

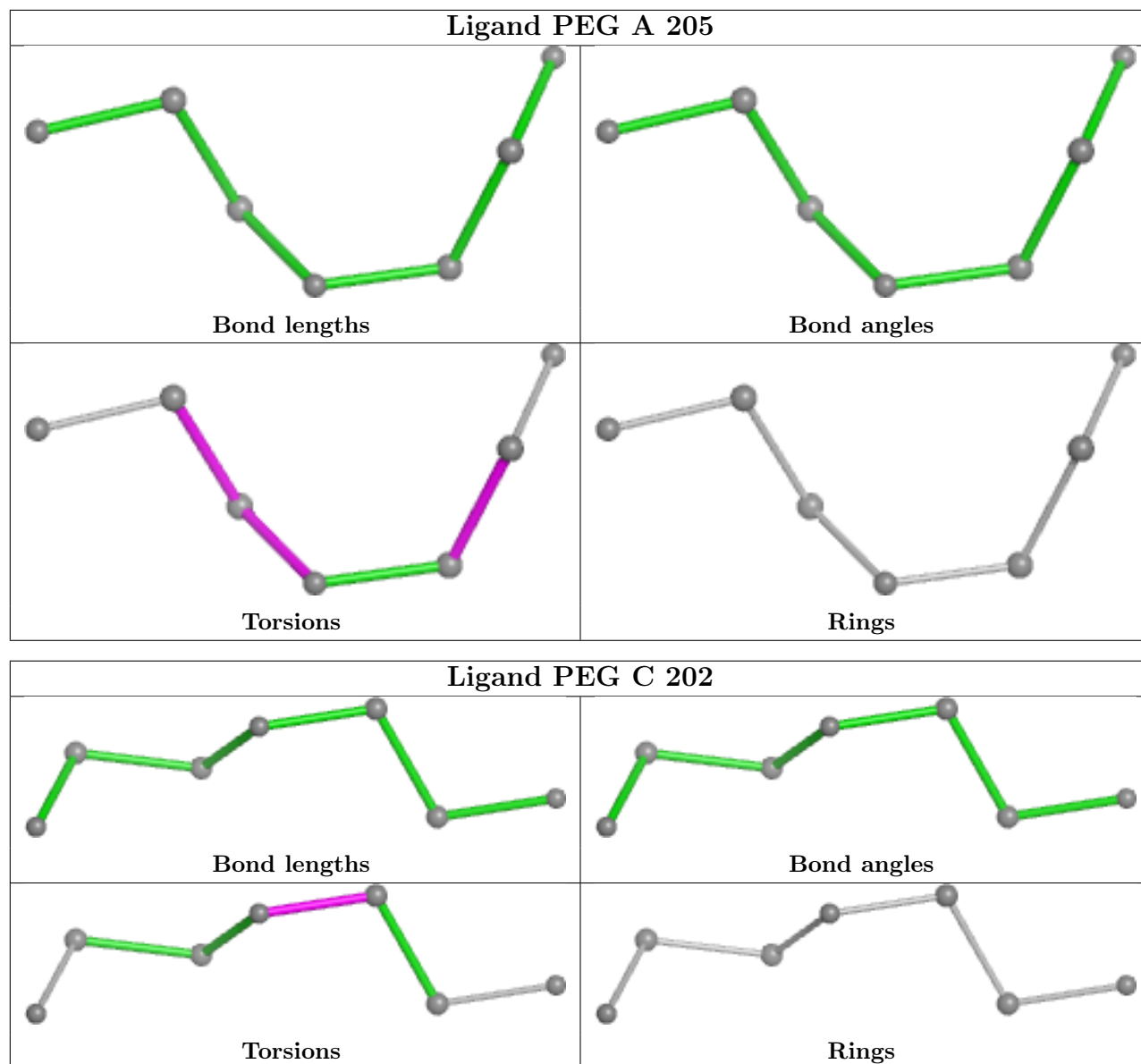
There are no ring outliers.

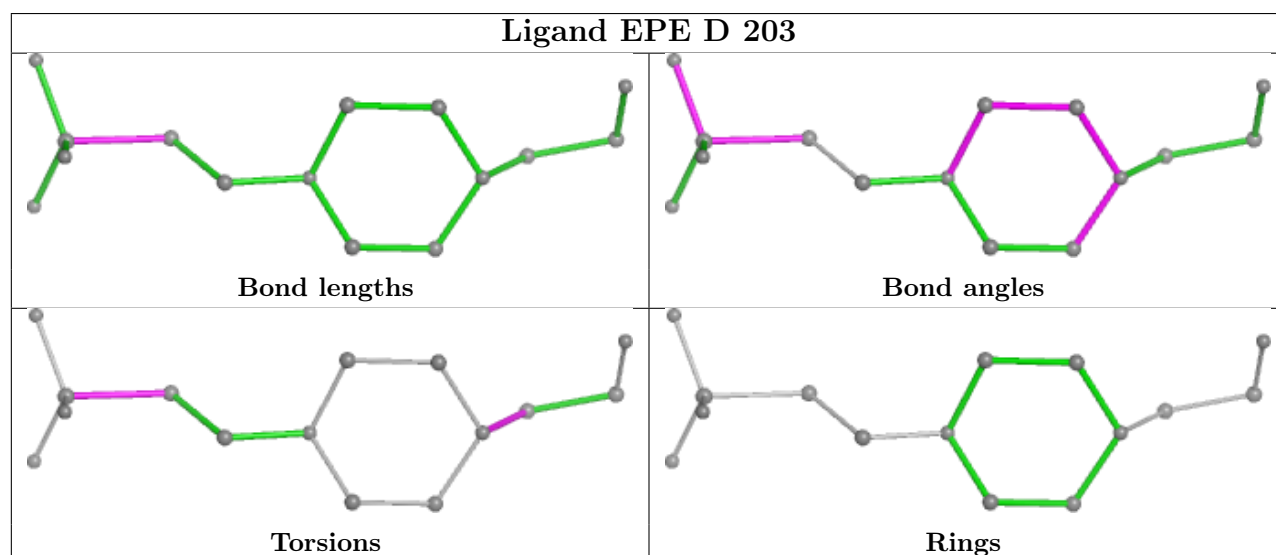
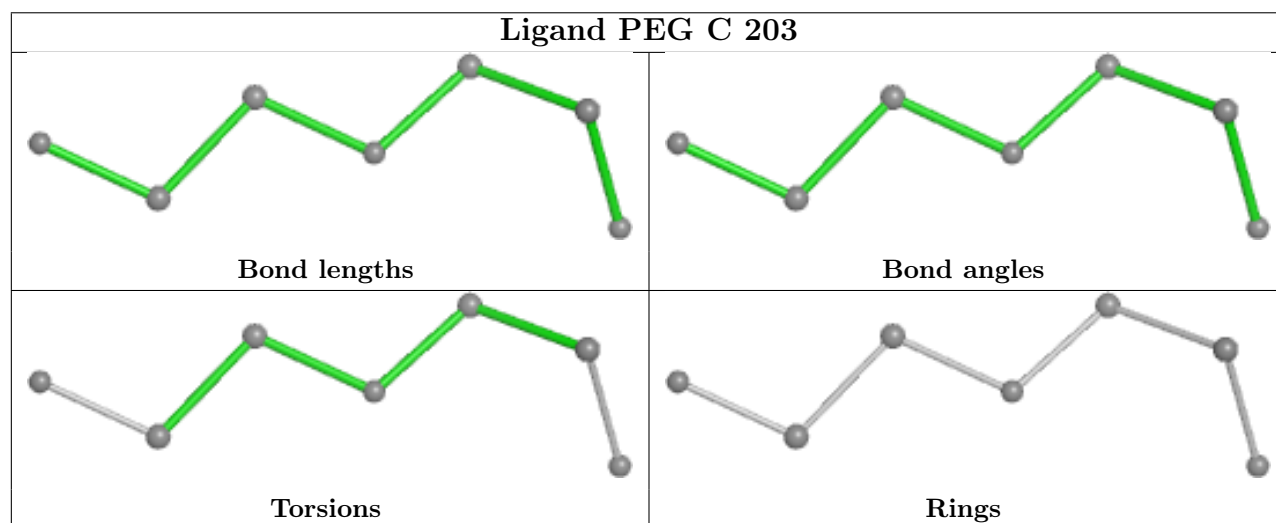
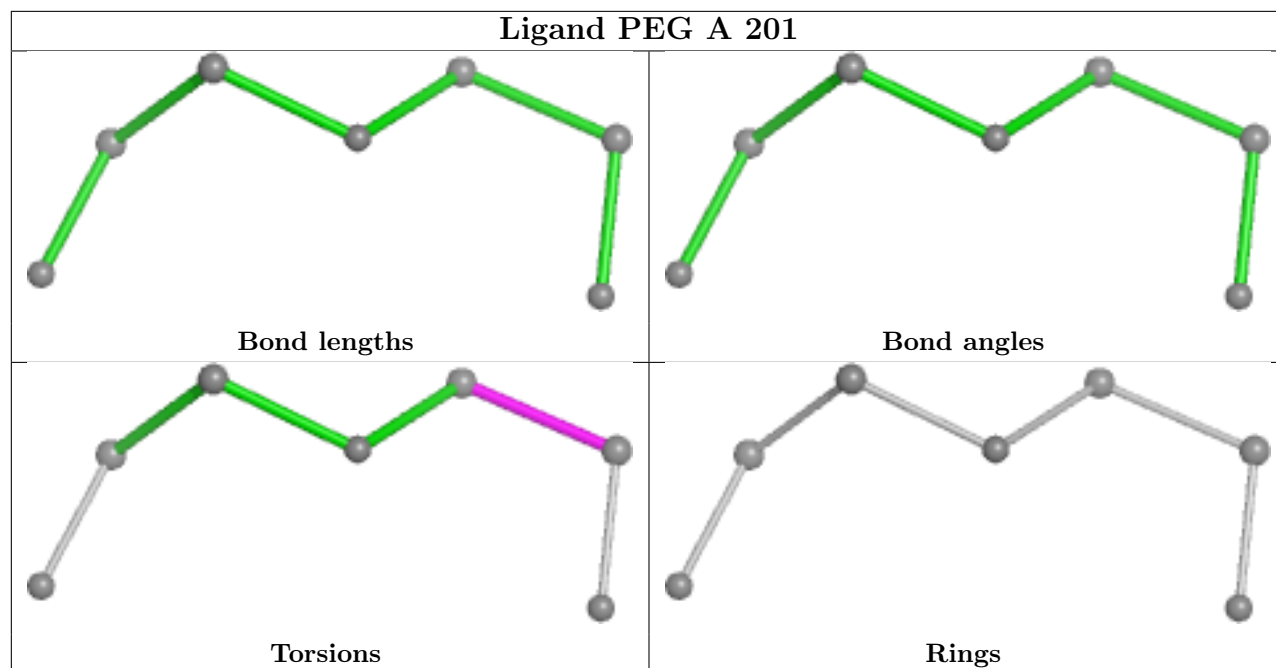
11 monomers are involved in 23 short contacts:

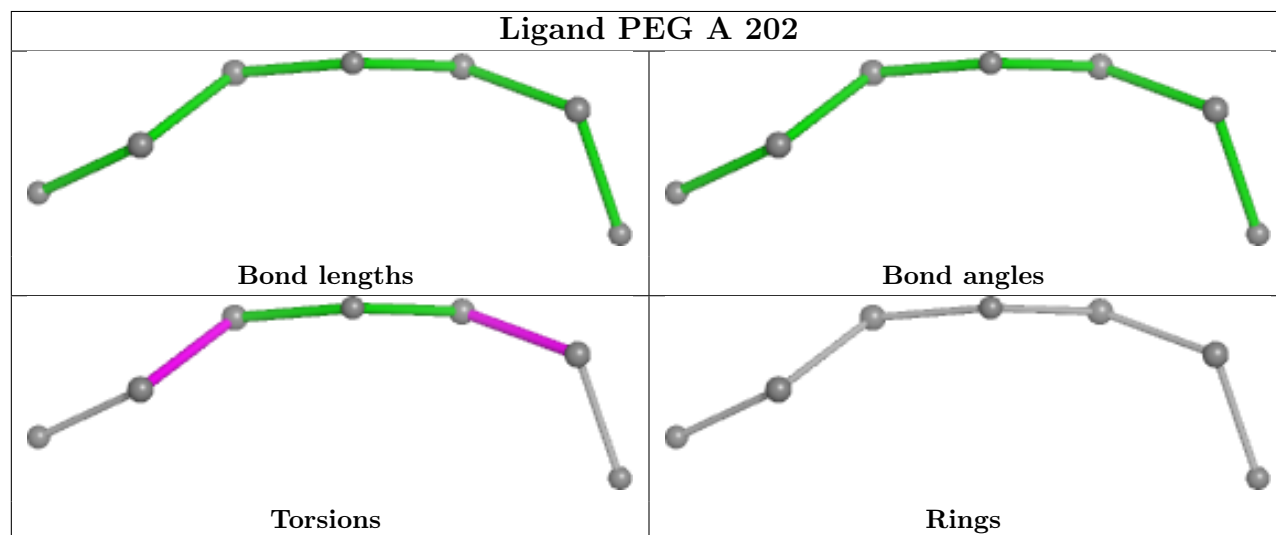
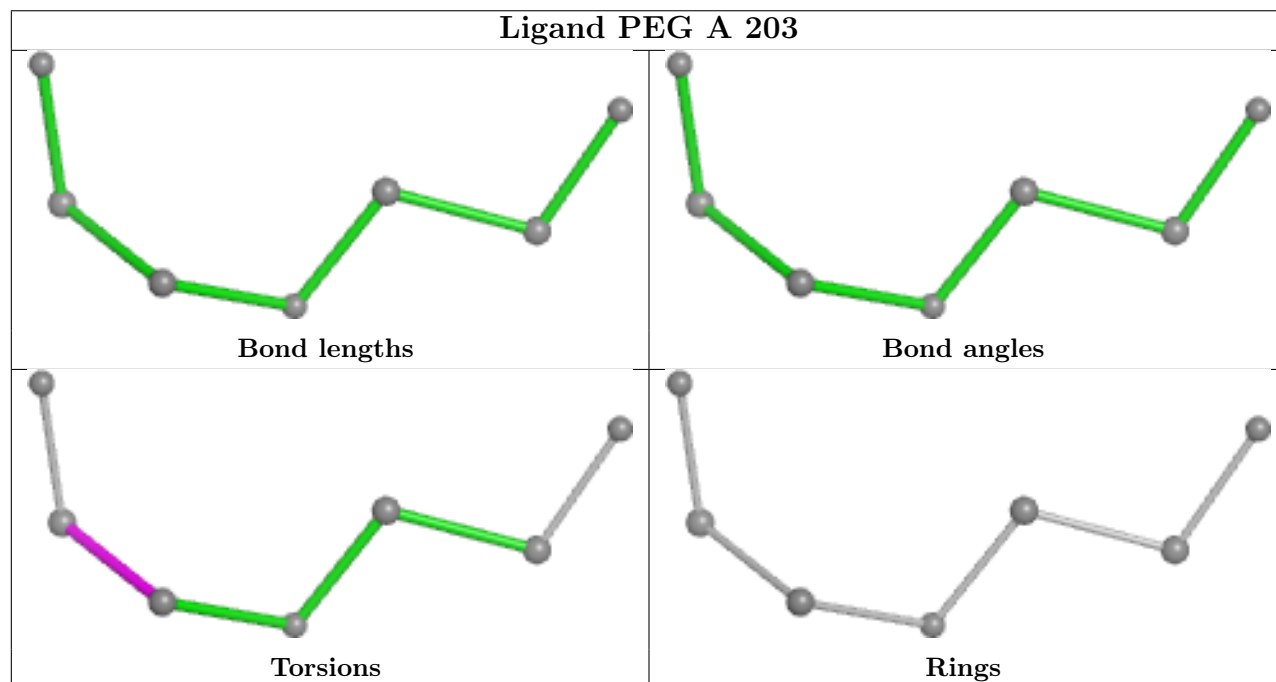
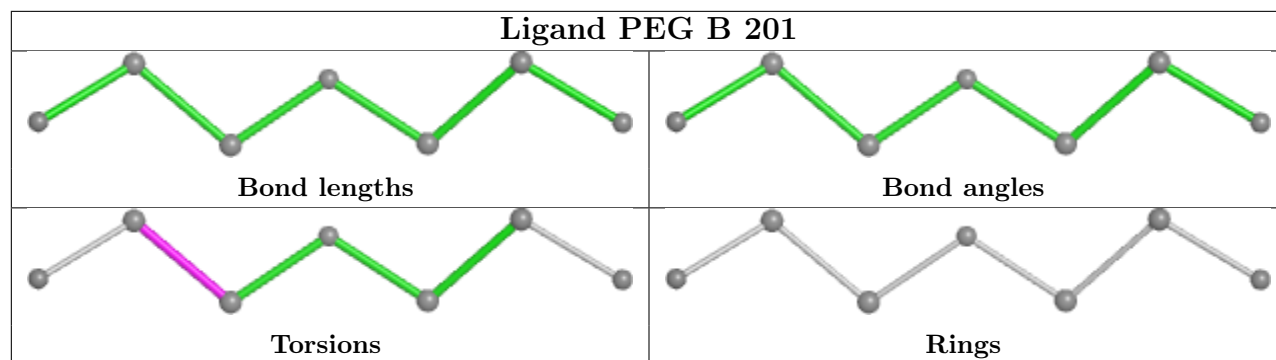
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	203	PEG	1	0
2	B	202	PEG	3	0
2	A	205	PEG	5	0
2	B	201	PEG	1	0
2	A	202	PEG	1	0
2	C	201	PEG	1	0
2	C	204	PEG	2	0
2	D	201	PEG	1	0
2	D	202	PEG	4	0
2	B	204	PEG	3	0
2	A	204	PEG	2	0

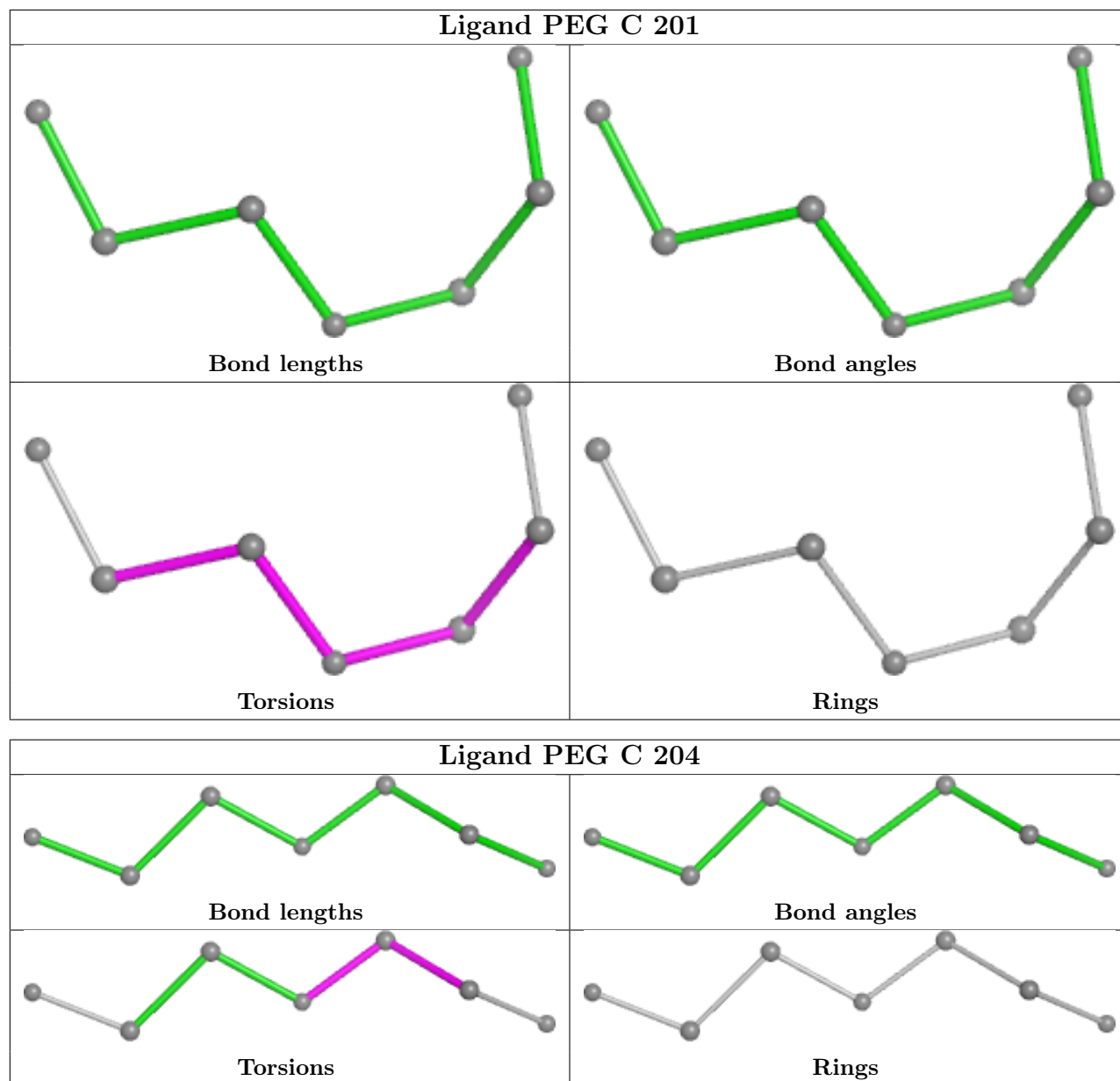
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



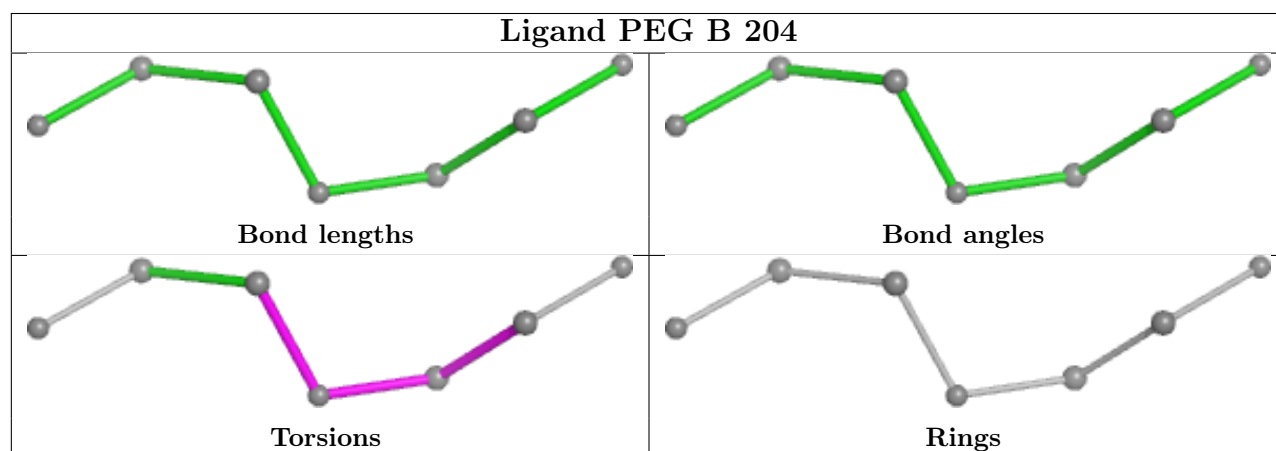
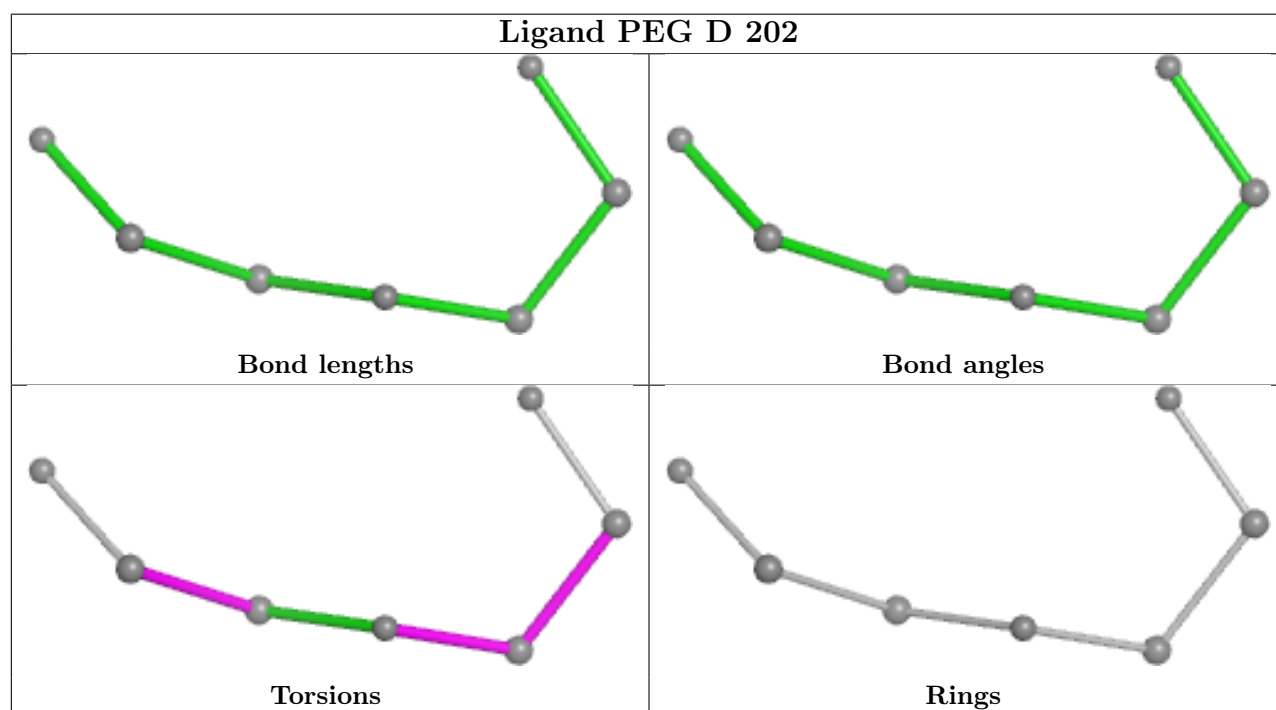
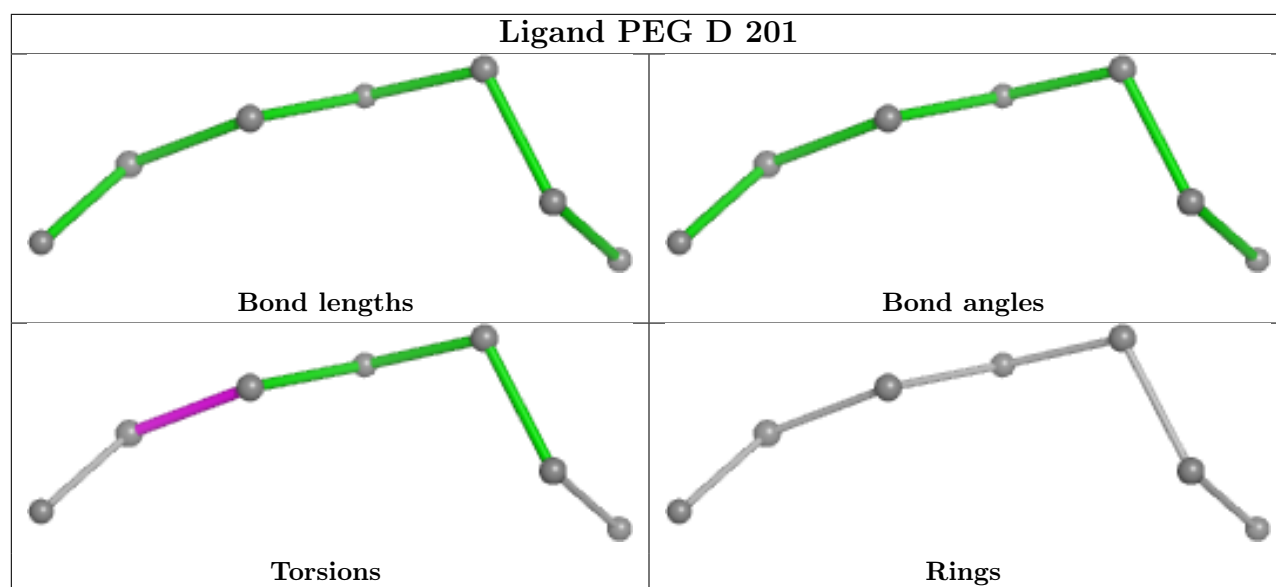


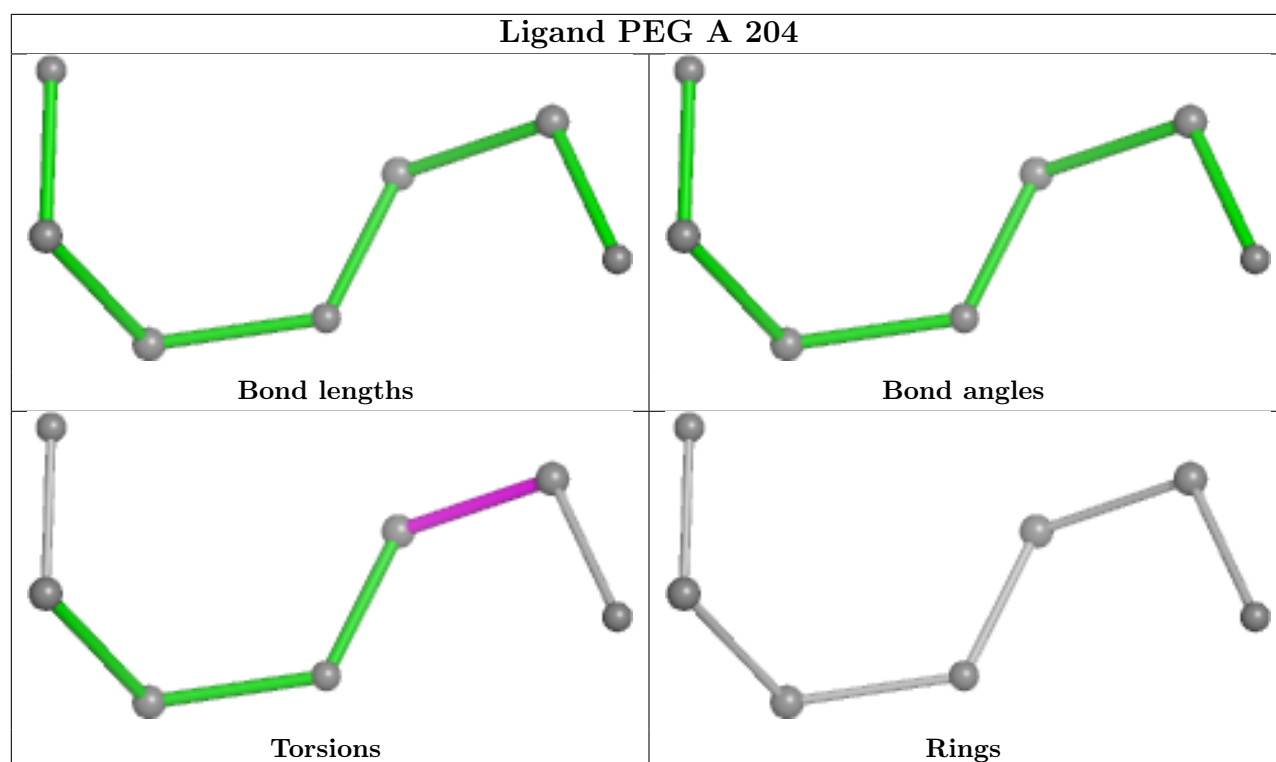












## 5.7 Other polymers [i](#)

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	118/161 (73%)	0.56	12 (10%) 13 13	27, 35, 56, 77	0
1	B	120/161 (74%)	0.67	12 (10%) 14 13	27, 36, 55, 70	0
1	C	118/161 (73%)	0.98	15 (12%) 9 8	30, 43, 60, 86	0
1	D	117/161 (72%)	0.89	14 (11%) 10 9	32, 41, 66, 75	0
All	All	473/644 (73%)	0.78	53 (11%) 11 10	27, 39, 59, 86	0

The worst 5 of 53 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	127	LEU	6.1
1	C	28	ASP	5.5
1	B	128	ARG	5.3
1	D	130	ASN	5.3
1	D	126	ASP	5.0

### 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 oligosaccharides in this entry.

### 6.4 Ligands [i](#)

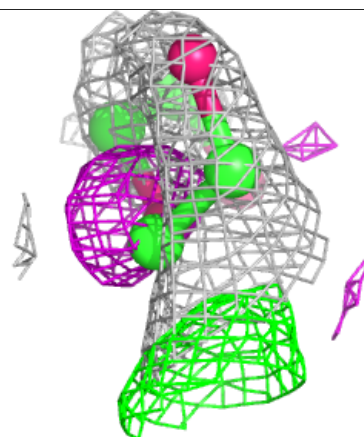
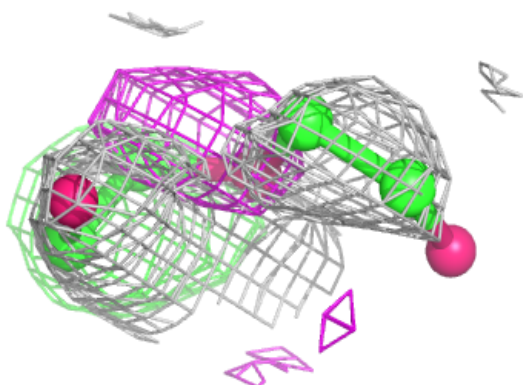
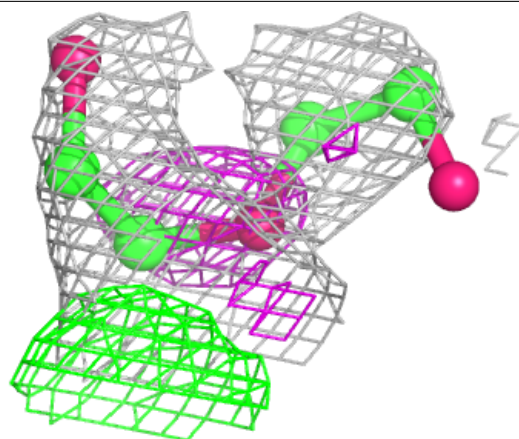
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 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
2	PEG	A	204	7/7	0.55	0.22	54,55,65,71	0
2	PEG	B	203	7/7	0.71	0.18	47,52,59,63	0
2	PEG	B	201	7/7	0.72	0.22	50,54,59,60	0
2	PEG	D	202	7/7	0.72	0.21	43,55,58,64	0
2	PEG	B	204	7/7	0.74	0.22	48,53,57,59	0
2	PEG	C	204	7/7	0.75	0.25	43,45,63,63	0
2	PEG	A	201	7/7	0.75	0.19	55,62,70,72	0
2	PEG	A	205	7/7	0.81	0.19	43,52,57,58	0
2	PEG	C	201	7/7	0.82	0.15	34,42,46,51	0
2	PEG	A	202	7/7	0.82	0.17	48,52,58,62	0
2	PEG	B	202	7/7	0.82	0.18	45,49,65,66	0
2	PEG	C	203	7/7	0.85	0.15	43,49,57,58	0
2	PEG	A	203	7/7	0.85	0.14	47,48,56,58	0
2	PEG	C	202	7/7	0.85	0.14	48,53,65,65	0
3	EPE	D	203	15/15	0.86	0.18	53,60,83,85	0
2	PEG	D	201	7/7	0.87	0.13	53,54,63,69	0

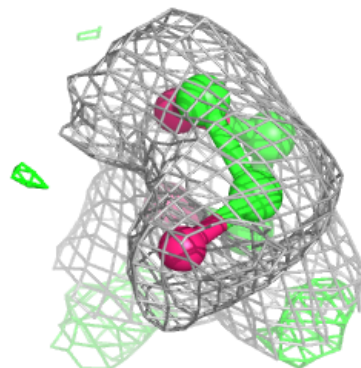
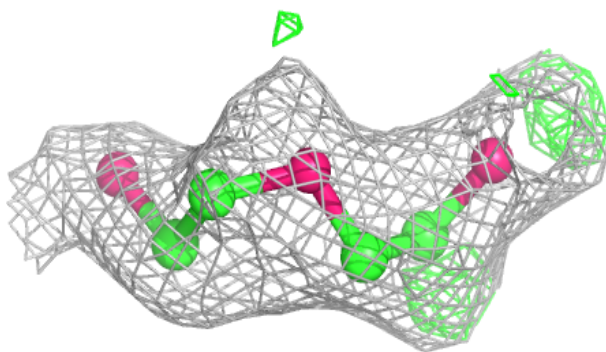
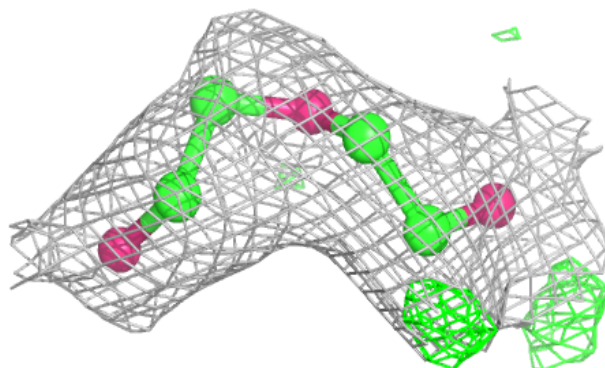
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around PEG A 204:**

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

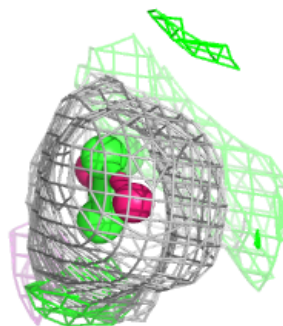
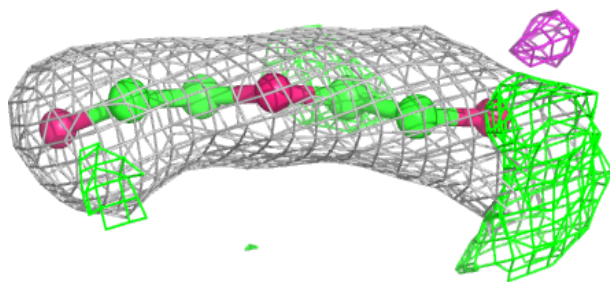
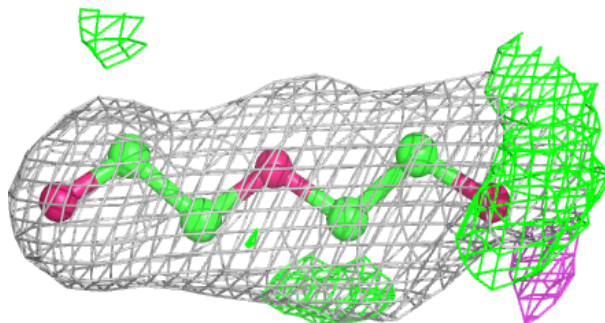
**Electron density around PEG B 203:**

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



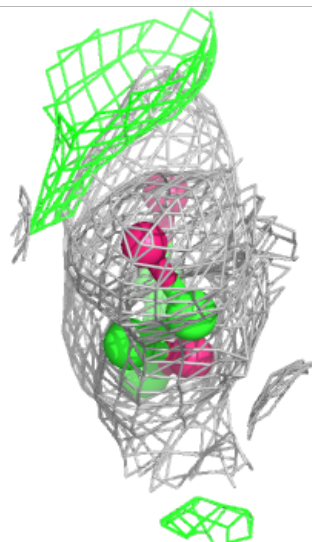
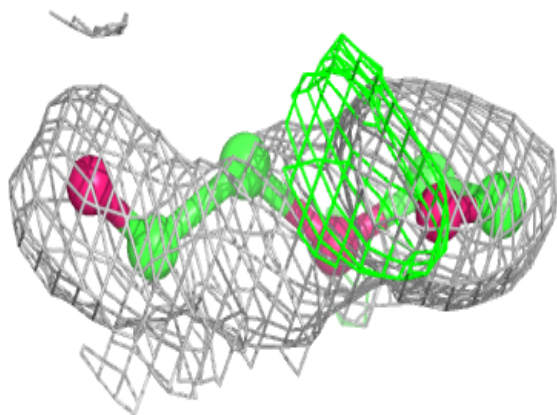
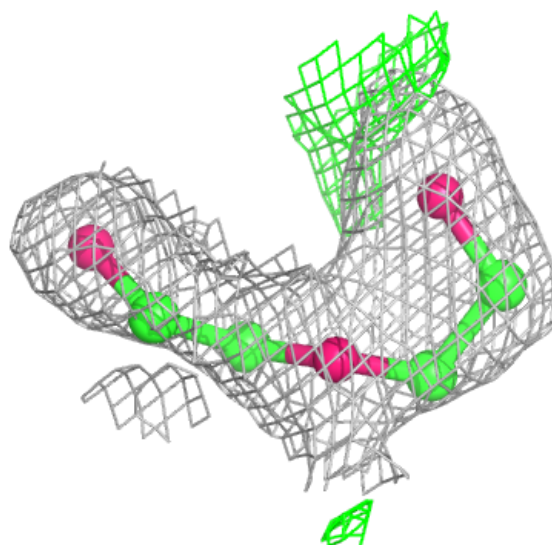
**Electron density around PEG B 201:**

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



**Electron density around PEG D 202:**

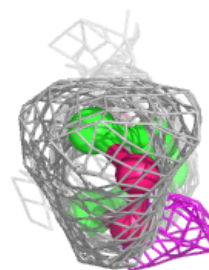
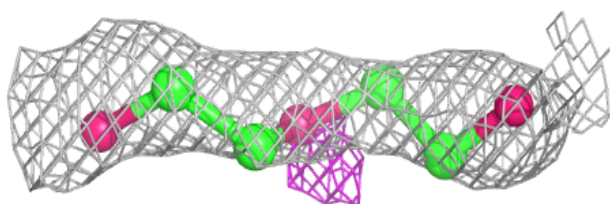
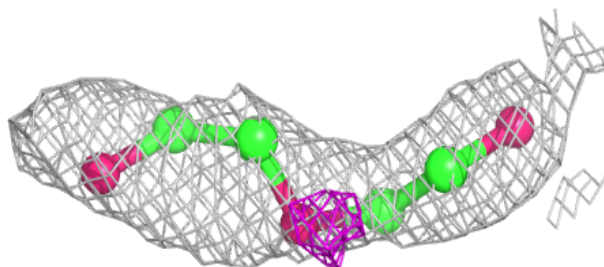
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



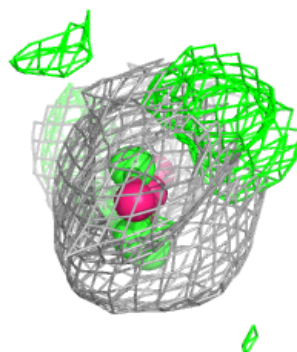
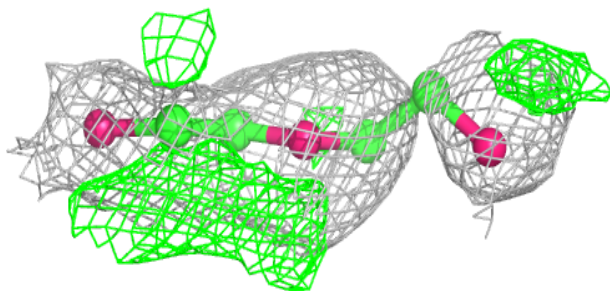
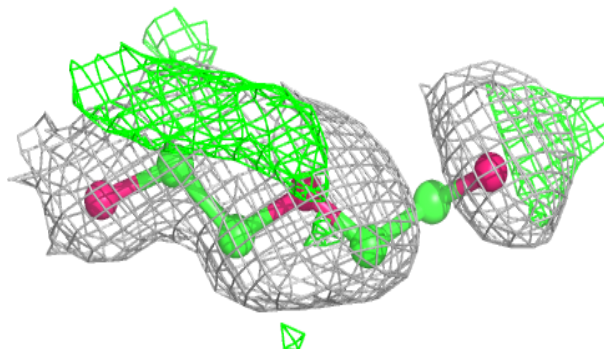


**Electron density around PEG B 204:**

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

**Electron density around PEG C 204:**

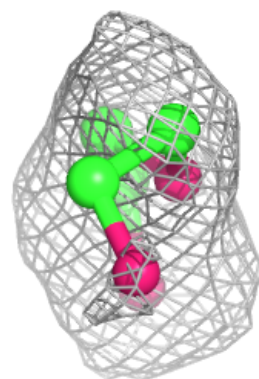
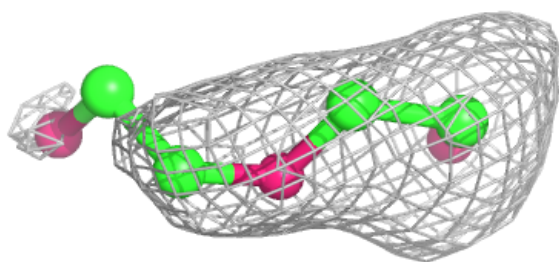
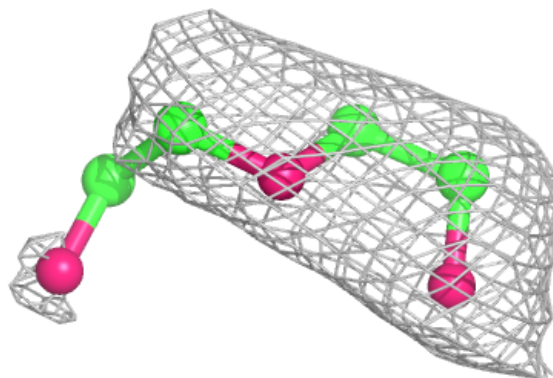
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



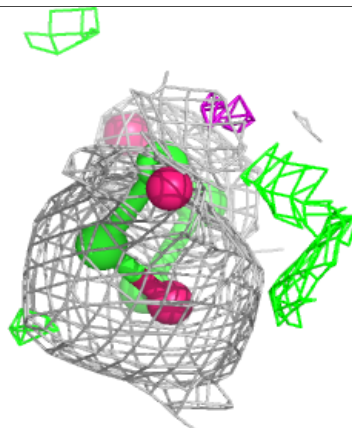
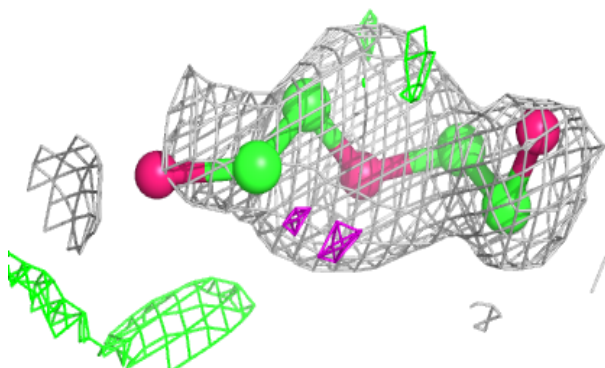
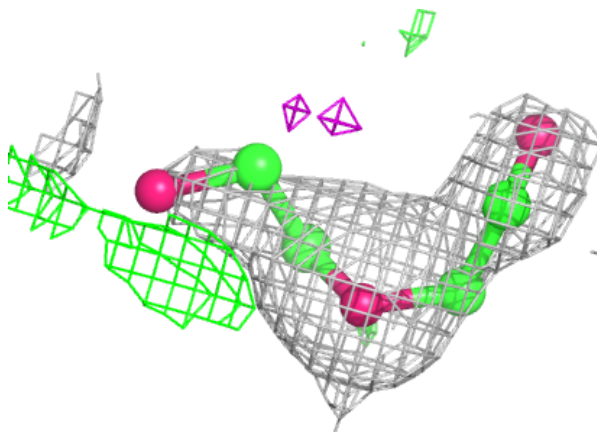


**Electron density around PEG A 201:**

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

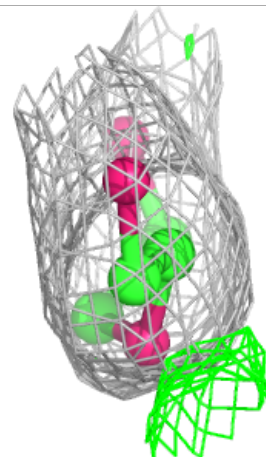
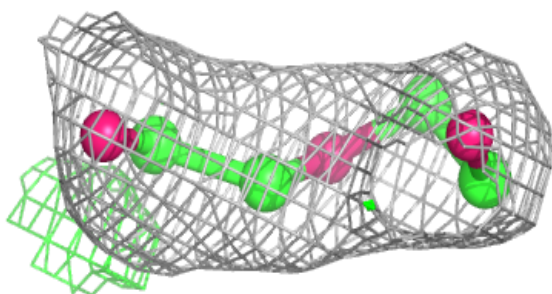
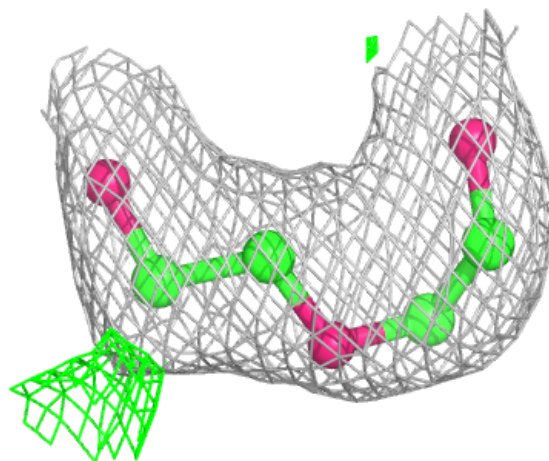
**Electron density around PEG A 205:**

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



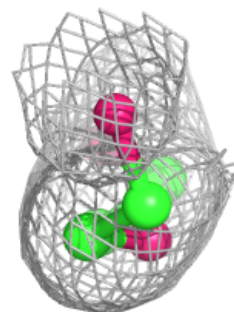
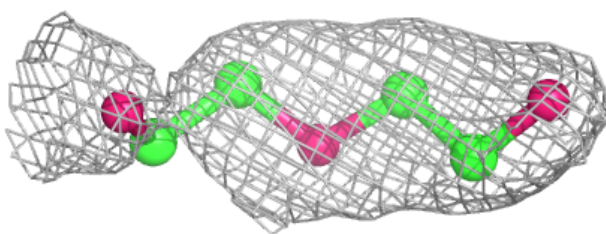
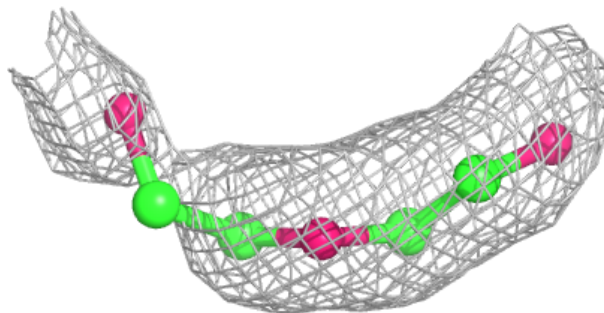
**Electron density around PEG C 201:**

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



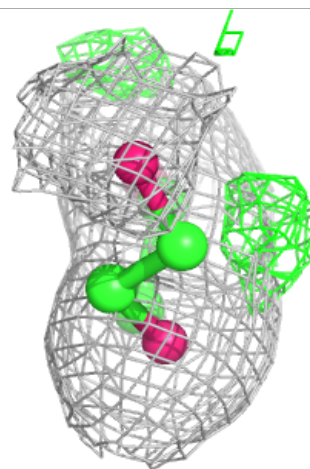
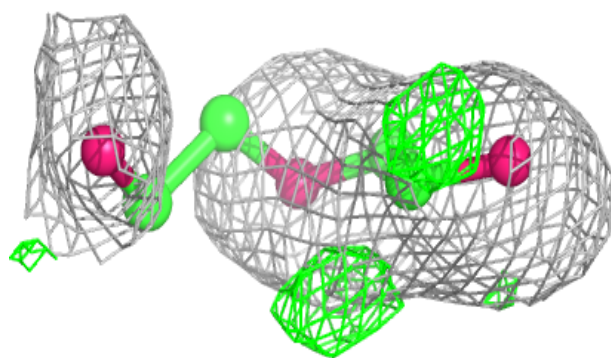
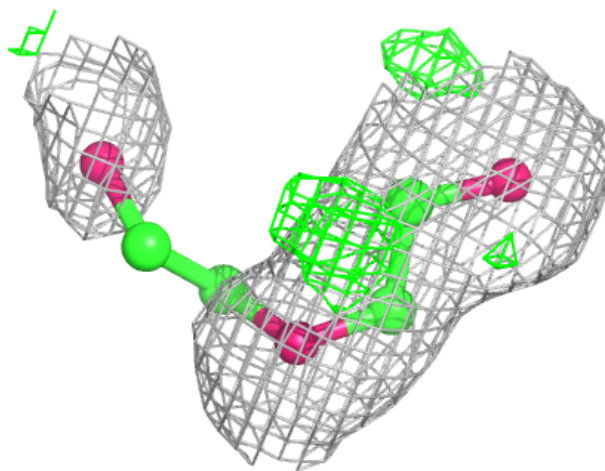
**Electron density around PEG A 202:**

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



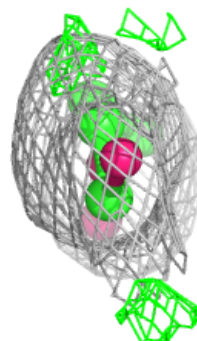
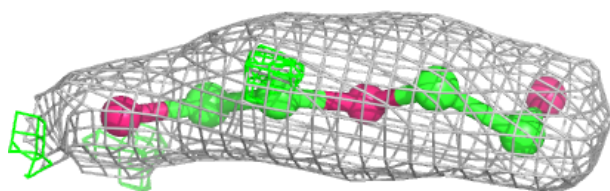
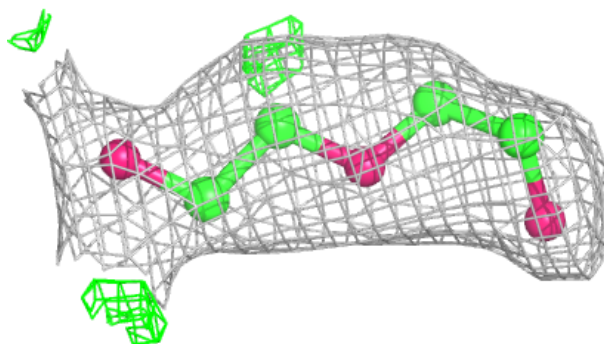
**Electron density around PEG B 202:**

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

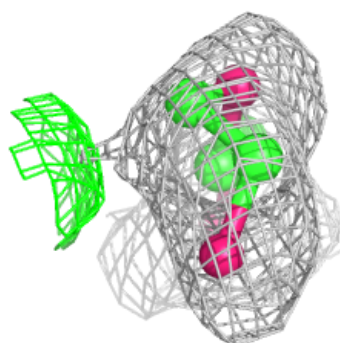
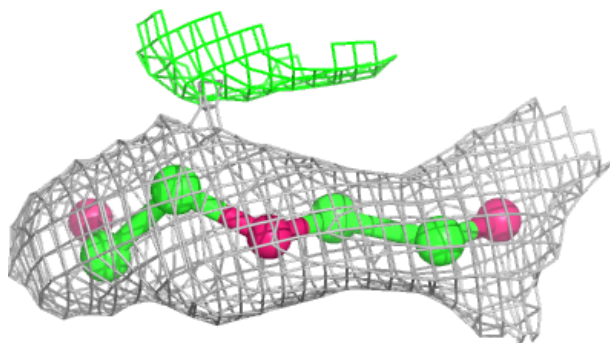
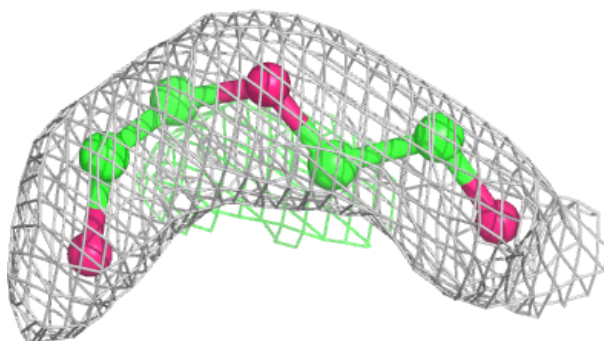


**Electron density around PEG C 203:**

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

**Electron density around PEG A 203:**

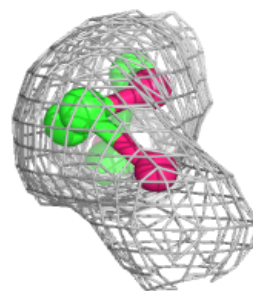
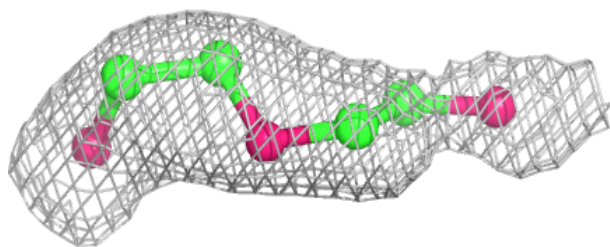
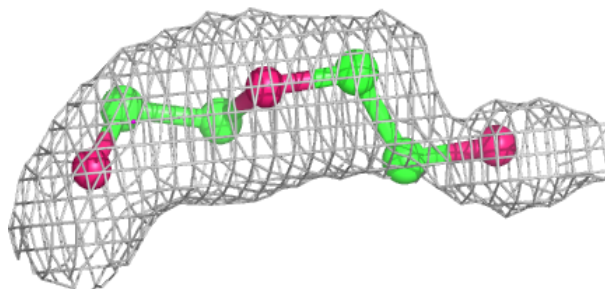
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



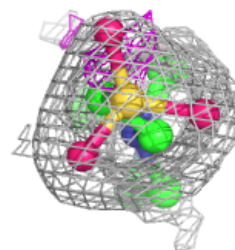
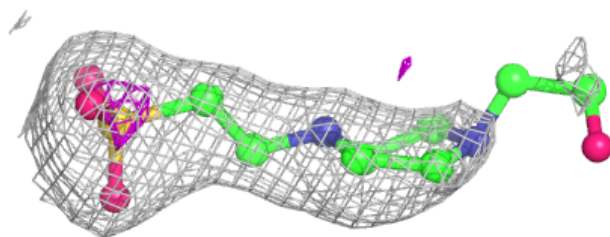
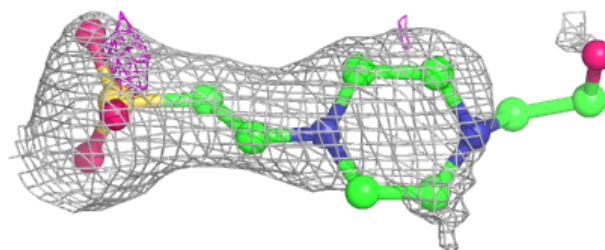


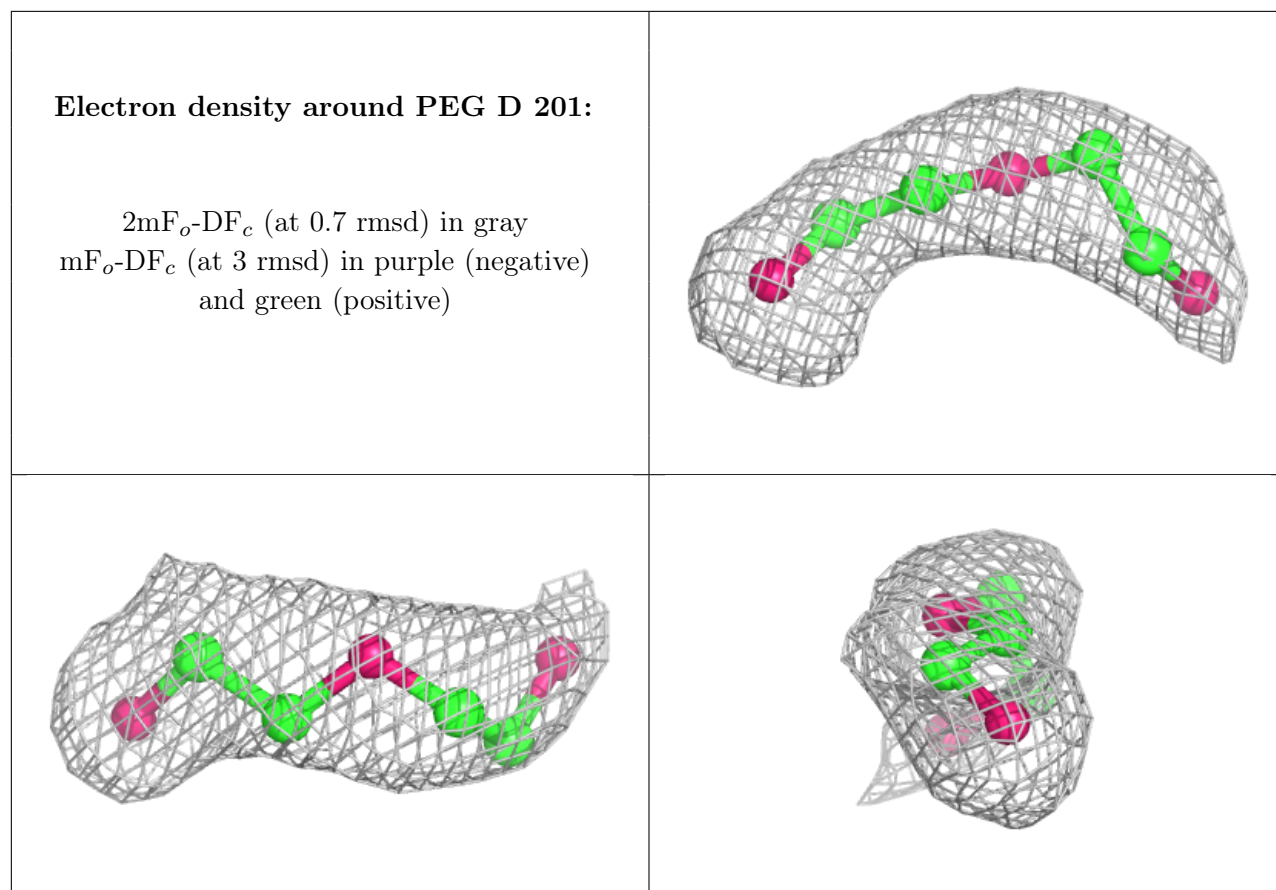
**Electron density around PEG C 202:**

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

**Electron density around EPE D 203:**

$2mF_o-DF_c$  (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.