



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 15, 2024 – 03:13 pm GMT

PDB ID : 6YX8  
Title : The structure of allophycocyanin from cyanobacterium Nostoc sp. WR13, the C2221 crystal form.  
Authors : Patel, H.M.; Roszak, A.W.; Madamwar, D.; Cogdell, R.J.  
Deposited on : 2020-04-30  
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 : 4.02b-467  
Mogul : 1.8.4, CSD as541be (2020)  
Xtrriage (Phenix) : 1.13  
EDS : 2.36  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

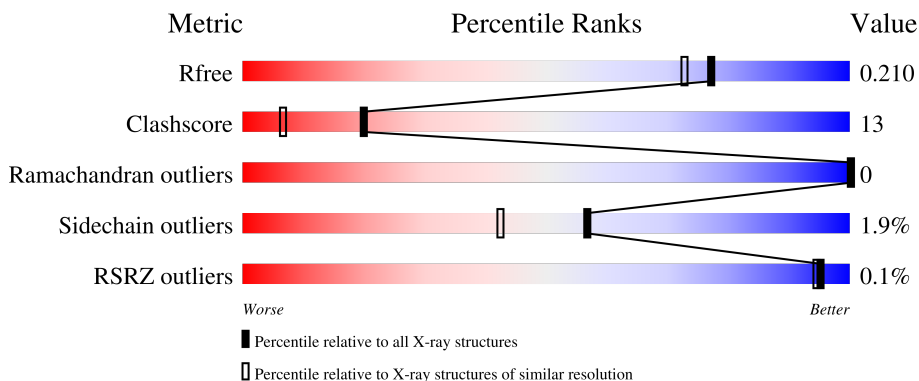
# 1 Overall quality at a glance

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))
$R_{free}$	130704	4003 (1.86-1.82)
Clashscore	141614	4233 (1.86-1.82)
Ramachandran outliers	138981	4185 (1.86-1.82)
Sidechain outliers	138945	4186 (1.86-1.82)
RSRZ outliers	127900	3957 (1.86-1.82)

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	AAA	160	89% 10% .
1	CCC	160	88% 11% .
1	EEE	160	90% 9% .
2	BBB	161	92% 8% .
2	DDD	161	90% 8% .

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Mol	Chain	Length	Quality of chain
2	FFF	161	 86% 13%

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
10	PG4	DDD	207	-	-	-	X
10	PG4	EEE	201	-	-	-	X
4	MPD	BBB	204	-	-	-	X
4	MPD	DDD	204	-	-	-	X
4	MPD	FFF	203	-	-	X	-
4	MPD	FFF	204	-	-	-	X
6	1PE	BBB	217	-	-	-	X
7	PGE	EEE	219	-	-	X	-
7	PGE	FFF	210	-	-	-	X
8	PEG	AAA	206[B]	-	-	X	-
8	PEG	AAA	207	-	-	X	X
8	PEG	AAA	208	-	-	X	-
8	PEG	BBB	210	-	-	-	X
8	PEG	CCC	206	-	-	X	-
8	PEG	CCC	211	-	-	-	X
8	PEG	DDD	210	-	-	X	X
8	PEG	EEE	207	-	-	X	X
8	PEG	FFF	211	-	-	-	X
8	PEG	FFF	212	-	-	-	X
8	PEG	FFF	214	-	-	-	X
9	EDO	BBB	212	-	-	-	X
9	EDO	DDD	213	-	-	X	-
9	EDO	EEE	210	-	-	-	X
9	EDO	EEE	211	-	-	-	X
9	EDO	EEE	212	-	-	-	X
9	EDO	EEE	214	-	-	X	-
9	EDO	EEE	216	-	-	-	X
9	EDO	FFF	215	-	-	X	-

## 2 Entry composition i

There are 12 unique types of molecules in this entry. The entry contains 9412 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.

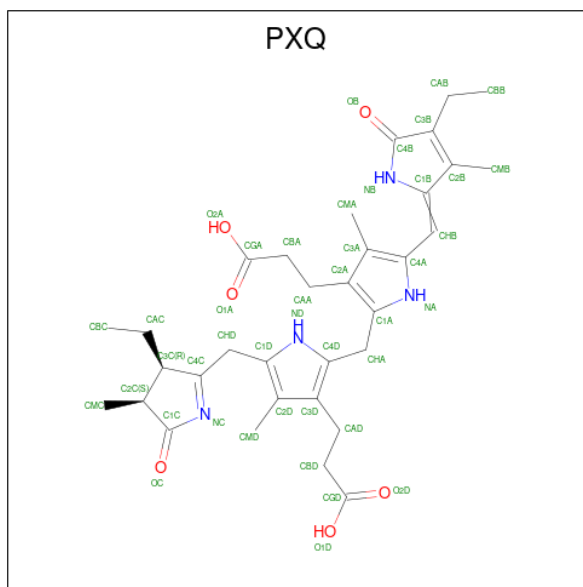
- Molecule 1 is a protein called ALPHA SUBUNIT OF CYANOBACTERIAL ALLOPHYCO-CYANIN PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	AAA	160	Total 1248	C 784	N 217	O 241	S 6	0	9	0
1	CCC	160	Total 1243	C 781	N 215	O 241	S 6	0	8	0
1	EEE	160	Total 1254	C 791	N 220	O 237	S 6	0	10	0

- Molecule 2 is a protein called BETA SUBUNIT OF CYANOBACTERIAL ALLOPHYCO-CYANIN PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	BBB	161	Total 1239	C 784	N 208	O 241	S 6	0	3	0
2	DDD	161	Total 1254	C 793	N 212	O 243	S 6	0	6	0
2	FFF	161	Total 1280	C 812	N 208	O 254	S 6	0	13	0

- Molecule 3 is 3-[5-[[[3 {R},4 {R}]-3-ethyl-4-methyl-5-oxidanylidene-3,4-dihydropyrrol-2-yl]methyl]-2-[[5-[( {Z})-(4-ethyl-3-methyl-5-oxidanylidene-pyrrol-2-ylidene)methyl]-3-(3-hydroxy-3-oxopropyl)-4-methyl-1 {H}-pyrrol-2-yl]methyl]-4-methyl-1 {H}-pyrrol-3-yl]propanoic acid (three-letter code: PXQ) (formula: C<sub>33</sub>H<sub>42</sub>N<sub>4</sub>O<sub>6</sub>) (labeled as "Ligand of Interest" by depositor).



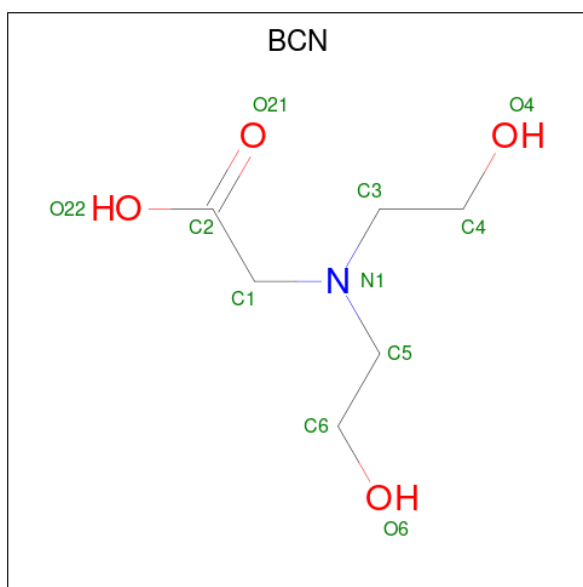
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	AAA	1	43	33	4	6	0	0
3	BBB	1	43	33	4	6	0	0
3	CCC	1	43	33	4	6	0	0
3	DDD	1	43	33	4	6	0	0
3	EEE	1	43	33	4	6	0	0
3	FFF	1	43	33	4	6	0	0

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>2</sub>).



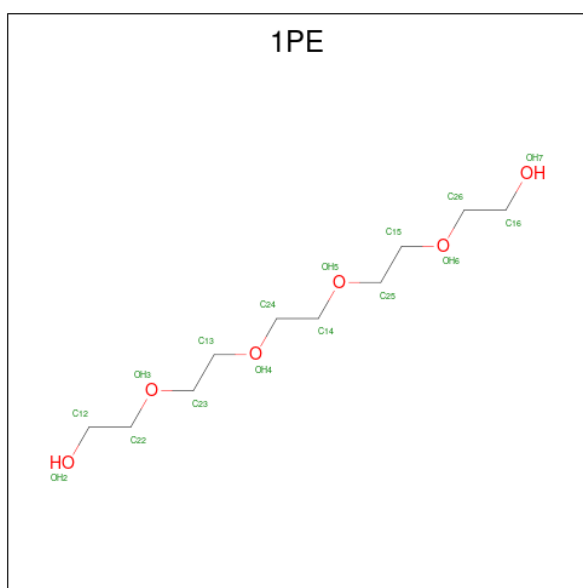
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	AAA	1	Total C O 8 6 2	0	0
4	BBB	1	Total C O 8 6 2	0	0
4	BBB	1	Total C O 8 6 2	0	0
4	BBB	1	Total C O 8 6 2	0	0
4	CCC	1	Total C O 8 6 2	0	0
4	DDD	1	Total C O 8 6 2	0	0
4	DDD	1	Total C O 8 6 2	0	0
4	DDD	1	Total C O 8 6 2	0	0
4	FFF	1	Total C O 8 6 2	0	0
4	FFF	1	Total C O 8 6 2	0	0
4	FFF	1	Total C O 8 6 2	0	0
4	FFF	1	Total C O 8 6 2	0	0

- Molecule 5 is BICINE (three-letter code: BCN) (formula: C<sub>6</sub>H<sub>13</sub>NO<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
5	AAA	1	Total	C	N	O	0	0
			11	6	1	4		
5	DDD	1	Total	C	N	O	0	0
			11	6	1	4		

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



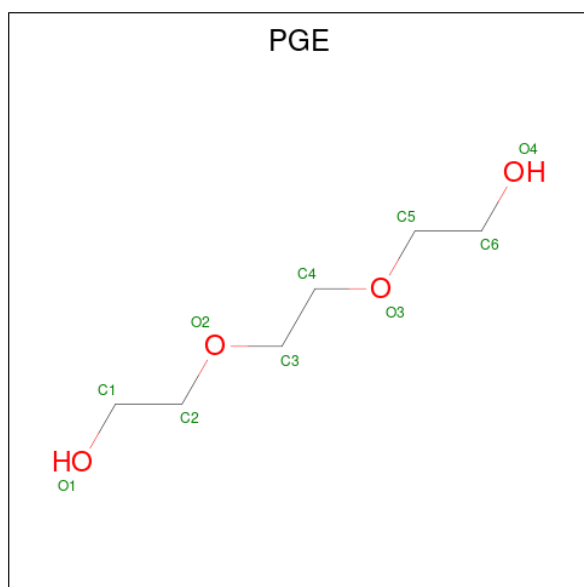
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	AAA	1	Total	C	O	0	0
			16	10	6		
6	BBB	1	Total	C	O	0	0
			16	10	6		

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

- Molecule 7 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	AAA	1	Total	C	O	0	0
			10	6	4		
7	AAA	1	Total	C	O	0	0
			10	6	4		
7	BBB	1	Total	C	O	0	0
			10	6	4		
7	BBB	1	Total	C	O	0	0
			10	6	4		
7	CCC	1	Total	C	O	0	0
			10	6	4		
7	EEE	1	Total	C	O	0	0
			10	6	4		
7	EEE	1	Total	C	O	0	0
			10	6	4		
7	FFF	1	Total	C	O	0	0
			10	6	4		

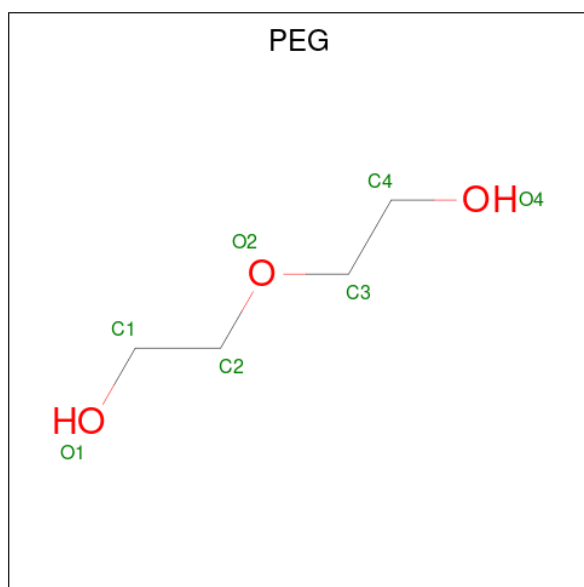
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	FFF	1	Total	C	O	0	0
			10	6	4		
7	FFF	1	Total	C	O	0	0
			10	6	4		
7	FFF	1	Total	C	O	0	0
			10	6	4		

- Molecule 8 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).



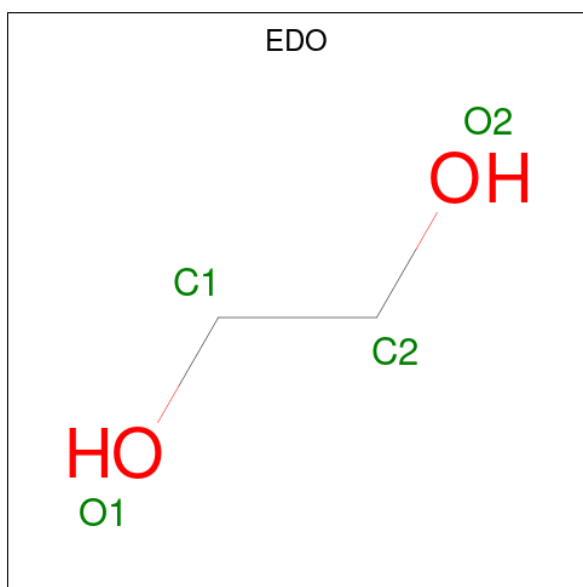
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	AAA	1	Total	C	O	0	1
			14	8	6		
8	AAA	1	Total	C	O	0	0
			7	4	3		
8	AAA	1	Total	C	O	0	0
			7	4	3		
8	AAA	1	Total	C	O	0	0
			7	4	3		
8	BBB	1	Total	C	O	0	0
			7	4	3		
8	BBB	1	Total	C	O	0	0
			7	4	3		
8	BBB	1	Total	C	O	0	0
			7	4	3		
8	CCC	1	Total	C	O	0	1
			14	8	6		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	CCC	1	Total	C	O	0	0
			7	4	3		
8	CCC	1	Total	C	O	0	0
			7	4	3		
8	CCC	1	Total	C	O	0	0
			7	4	3		
8	CCC	1	Total	C	O	0	0
			7	4	3		
8	DDD	1	Total	C	O	0	0
			7	4	3		
8	DDD	1	Total	C	O	0	0
			7	4	3		
8	DDD	1	Total	C	O	0	0
			7	4	3		
8	EEE	1	Total	C	O	0	0
			7	4	3		
8	EEE	1	Total	C	O	0	0
			7	4	3		
8	EEE	1	Total	C	O	0	0
			7	4	3		
8	EEE	1	Total	C	O	0	0
			7	4	3		
8	FFF	1	Total	C	O	0	0
			7	4	3		
8	FFF	1	Total	C	O	0	0
			7	4	3		
8	FFF	1	Total	C	O	0	0
			7	4	3		
8	FFF	1	Total	C	O	0	0
			7	4	3		

- Molecule 9 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	AAA	1	Total C O 8 4 4	0	1
9	AAA	1	Total C O 4 2 2	0	0
9	AAA	1	Total C O 4 2 2	0	0
9	AAA	1	Total C O 4 2 2	0	0
9	AAA	1	Total C O 4 2 2	0	0
9	AAA	1	Total C O 4 2 2	0	0
9	BBB	1	Total C O 4 2 2	0	0
9	BBB	1	Total C O 4 2 2	0	0
9	BBB	1	Total C O 4 2 2	0	0
9	BBB	1	Total C O 4 2 2	0	0
9	BBB	1	Total C O 4 2 2	0	0
9	CCC	1	Total C O 4 2 2	0	0
9	CCC	1	Total C O 4 2 2	0	0
9	CCC	1	Total C O 4 2 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	DDD	1	Total	C	O	0	0
			4	2	2		
9	DDD	1	Total	C	O	0	0
			4	2	2		
9	DDD	1	Total	C	O	0	0
			4	2	2		
9	DDD	1	Total	C	O	0	1
			8	4	4		
9	DDD	1	Total	C	O	0	0
			4	2	2		
9	DDD	1	Total	C	O	0	1
			8	4	4		
9	DDD	1	Total	C	O	0	0
			4	2	2		
9	EEE	1	Total	C	O	0	0
			4	2	2		
9	EEE	1	Total	C	O	0	0
			4	2	2		
9	EEE	1	Total	C	O	0	0
			4	2	2		
9	EEE	1	Total	C	O	0	0
			4	2	2		
9	EEE	1	Total	C	O	0	0
			4	2	2		
9	EEE	1	Total	C	O	0	1
			8	4	4		
9	EEE	1	Total	C	O	0	0
			4	2	2		
9	EEE	1	Total	C	O	0	0
			4	2	2		
9	FFF	1	Total	C	O	0	0
			4	2	2		
9	FFF	1	Total	C	O	0	0
			4	2	2		
9	FFF	1	Total	C	O	0	0
			4	2	2		
9	FFF	1	Total	C	O	0	0
			4	2	2		

- Molecule 10 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
10	DDD	1	Total	C	O	0	0
			13	8	5		
10	DDD	1	Total	C	O	0	0
			13	8	5		
10	DDD	1	Total	C	O	0	0
			13	8	5		
10	EEE	1	Total	C	O	0	0
			13	8	5		
10	EEE	1	Total	C	O	0	0
			13	8	5		

- Molecule 11 is (4R)-2-METHYLPENTANE-2,4-DIOL (three-letter code: MRD) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	EEE	1	Total	C O	0	0
			8	6 2		

- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	AAA	161	Total	O	0	0
			161	161		
12	BBB	220	Total	O	0	0
			220	220		
12	CCC	121	Total	O	0	0
			121	121		
12	DDD	109	Total	O	0	0
			109	109		
12	EEE	155	Total	O	0	0
			155	155		
12	FFF	162	Total	O	0	0
			162	162		

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


These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: ALPHA SUBUNIT OF CYANOBACTERIAL ALLOPHYCOCYANIN PROTEIN

Chain AAA:  89% 10%



- Molecule 1: ALPHA SUBUNIT OF CYANOBACTERIAL ALLOPHYCOCYANIN PROTEIN

Chain CCC:  88% 11%



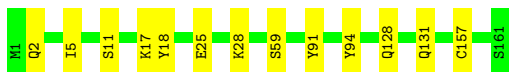
- Molecule 1: ALPHA SUBUNIT OF CYANOBACTERIAL ALLOPHYCOCYANIN PROTEIN

Chain EEE:  90% 9%



- Molecule 2: BETA SUBUNIT OF CYANOBACTERIAL ALLOPHYCOCYANIN PROTEIN

Chain BBB:  92% 8%



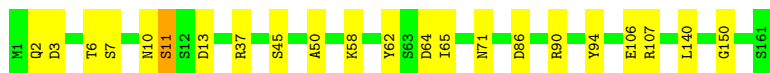
- Molecule 2: BETA SUBUNIT OF CYANOBACTERIAL ALLOPHYCOCYANIN PROTEIN

Chain DDD:  90% 8%



- Molecule 2: BETA SUBUNIT OF CYANOBACTERIAL ALLOPHYCOCYANIN PROTEIN

Chain FFF:  86% 13%





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	104.68Å 177.34Å 122.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	90.14 – 1.83 90.14 – 1.83	Depositor EDS
% Data completeness (in resolution range)	87.3 (90.14-1.83) 87.3 (90.14-1.83)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.62 (at 1.83Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.160 , 0.198 0.174 , 0.210	Depositor DCC
$R_{free}$ test set	4339 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	29.1	Xtrriage
Anisotropy	0.012	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 57.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.006 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.015 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	9412	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% 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: BCN, MRD, MPD, MEN, PEG, EDO, PG4, 1PE, PXQ, PGE

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	AAA	0.79	0/1290	0.80	0/1741
1	CCC	0.74	0/1282	0.82	1/1731 (0.1%)
1	EEE	0.75	0/1299	0.84	1/1751 (0.1%)
2	BBB	0.77	0/1258	0.82	1/1699 (0.1%)
2	DDD	0.74	0/1279	0.94	4/1728 (0.2%)
2	FFF	0.74	0/1329	0.80	3/1799 (0.2%)
All	All	0.75	0/7737	0.84	10/10449 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	DDD	77[A]	ARG	NE-CZ-NH2	-10.23	115.19	120.30
2	DDD	77[B]	ARG	NE-CZ-NH2	-10.23	115.19	120.30
2	DDD	77[A]	ARG	NE-CZ-NH1	7.92	124.26	120.30
2	DDD	77[B]	ARG	NE-CZ-NH1	7.92	124.26	120.30
2	FFF	94	TYR	CB-CG-CD1	5.38	124.23	121.00
1	EEE	86	TYR	CB-CG-CD1	5.33	124.20	121.00
2	FFF	94	TYR	CB-CG-CD2	-5.30	117.82	121.00
2	BBB	94	TYR	CB-CG-CD2	-5.13	117.92	121.00
1	CCC	82	ARG	CG-CD-NE	5.03	122.37	111.80
2	FFF	37	ARG	NE-CZ-NH2	-5.00	117.80	120.30

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	AAA	1248	0	1280	22	0
1	CCC	1243	0	1270	22	1
1	EEE	1254	0	1306	31	1
2	BBB	1239	0	1268	18	0
2	DDD	1254	0	1285	29	0
2	FFF	1280	0	1319	34	0
3	AAA	43	0	0	1	0
3	BBB	43	0	0	0	0
3	CCC	43	0	0	0	0
3	DDD	43	0	0	1	0
3	EEE	43	0	0	0	0
3	FFF	43	0	0	0	0
4	AAA	8	0	14	0	0
4	BBB	24	0	42	3	0
4	CCC	8	0	14	0	0
4	DDD	24	0	42	5	0
4	FFF	32	0	56	8	0
5	AAA	11	0	12	2	0
5	DDD	11	0	11	5	0
6	AAA	16	0	22	0	0
6	BBB	48	0	66	4	0
6	FFF	16	0	22	0	0
7	AAA	20	0	28	3	0
7	BBB	20	0	28	4	0
7	CCC	10	0	14	0	0
7	EEE	20	0	28	7	0
7	FFF	40	0	56	11	0
8	AAA	35	0	50	15	0
8	BBB	21	0	30	1	0
8	CCC	42	0	60	7	0
8	DDD	21	0	30	11	0
8	EEE	28	0	40	4	0
8	FFF	28	0	40	3	0
9	AAA	28	0	42	0	0
9	BBB	20	0	30	2	0
9	CCC	12	0	18	2	0
9	DDD	36	0	54	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	EEE	40	0	59	15	0
9	FFF	16	0	24	15	0
10	DDD	39	0	54	10	0
10	EEE	26	0	36	6	0
11	EEE	8	0	14	0	0
12	AAA	161	0	0	11	0
12	BBB	220	0	0	13	0
12	CCC	121	0	0	2	0
12	DDD	109	0	0	8	0
12	EEE	155	0	0	8	0
12	FFF	162	0	0	8	0
All	All	9412	0	8764	218	1

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 13.

All (218) 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)
2:FFF:6[A]:THR:HG22	12:FFF:322:HOH:O	1.13	1.31
2:DDD:39[A]:ARG:NH2	8:DDD:210:PEG:H41	1.48	1.28
8:AAA:206[B]:PEG:H32	12:AAA:302:HOH:O	1.28	1.27
4:DDD:205:MPD:H52	4:DDD:205:MPD:CM	1.51	1.24
4:DDD:205:MPD:HM3	4:DDD:205:MPD:C5	1.56	1.24
8:CCC:206:PEG:H12	8:CCC:206:PEG:H42	1.29	1.14
8:EEE:207:PEG:O1	8:EEE:207:PEG:C3	1.95	1.12
6:BBB:206:1PE:H162	12:BBB:339:HOH:O	1.47	1.11
8:EEE:207:PEG:O1	8:EEE:207:PEG:H31	1.40	1.11
2:DDD:39[A]:ARG:HH22	8:DDD:210:PEG:H41	0.96	1.08
2:FFF:86[B]:ASP:OD1	12:FFF:301:HOH:O	1.73	1.06
8:CCC:206:PEG:H12	8:CCC:206:PEG:C4	1.79	1.05
8:AAA:207:PEG:C1	8:AAA:207:PEG:H42	1.85	1.04
2:BBB:128:GLN:HG2	7:BBB:208:PGE:H6	1.41	1.02
8:AAA:207:PEG:H12	8:AAA:207:PEG:C4	1.89	0.98
2:DDD:39[A]:ARG:HH22	8:DDD:210:PEG:C4	1.79	0.95
2:DDD:39[B]:ARG:HD3	8:DDD:210:PEG:H12	1.48	0.94
8:AAA:206[B]:PEG:H22	8:AAA:206[B]:PEG:O4	1.65	0.94
2:BBB:17[A]:LYS:NZ	12:BBB:302:HOH:O	2.00	0.94
2:DDD:39[A]:ARG:NH2	8:DDD:210:PEG:C4	2.33	0.92
7:FFF:210:PGE:H32	7:FFF:210:PGE:H62	1.53	0.91
2:BBB:17[A]:LYS:CE	12:BBB:302:HOH:O	2.22	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:EEE:122:ILE:H	9:EEE:214:EDO:H11	1.40	0.87
2:DDD:36:LEU:O	2:DDD:39[A]:ARG:HG3	1.75	0.86
1:EEE:16:TYR:OH	12:FFF:301:HOH:O	1.93	0.86
12:EEE:321:HOH:O	2:FFF:6[B]:THR:HG21	1.78	0.83
1:EEE:76[B]:MET:SD	12:EEE:377:HOH:O	2.37	0.83
4:FFF:203:MPD:H52	4:FFF:203:MPD:H11	1.62	0.81
1:EEE:73:GLY:H	10:EEE:204:PG4:C1	1.94	0.80
1:CCC:34[A]:GLN:NE2	1:CCC:147[A]:GLU:OE1	2.16	0.78
2:BBB:17[A]:LYS:HE2	12:BBB:302:HOH:O	1.83	0.78
7:FFF:210:PGE:H32	7:FFF:210:PGE:C6	2.14	0.78
1:CCC:5:LYS:NZ	1:CCC:99:ASP:OD1	2.17	0.77
8:CCC:206:PEG:C4	8:CCC:206:PEG:C1	2.54	0.77
10:DDD:208:PG4:H42	10:DDD:208:PG4:H11	1.67	0.76
7:EEE:219:PGE:H32	12:EEE:301:HOH:O	1.84	0.76
2:FFF:45[A]:SER:OG	12:FFF:302:HOH:O	2.03	0.76
8:AAA:208:PEG:H41	12:AAA:301:HOH:O	1.86	0.75
8:EEE:207:PEG:H31	8:EEE:207:PEG:HO1	1.49	0.75
9:FFF:215:EDO:H11	12:FFF:327:HOH:O	1.85	0.75
2:DDD:71:MEN:HB3	12:DDD:377:HOH:O	1.87	0.74
5:DDD:206:BCN:H42	12:DDD:308:HOH:O	1.87	0.74
1:EEE:122:ILE:N	9:EEE:214:EDO:H11	2.02	0.74
8:AAA:207:PEG:H42	8:AAA:207:PEG:H12	1.50	0.73
1:EEE:122:ILE:H	9:EEE:214:EDO:C1	2.03	0.72
2:FFF:86[B]:ASP:OD2	9:FFF:215:EDO:H22	1.90	0.71
2:FFF:86[B]:ASP:OD2	9:FFF:215:EDO:O1	2.07	0.71
8:AAA:208:PEG:O2	12:AAA:301:HOH:O	2.09	0.70
4:FFF:203:MPD:H11	4:FFF:203:MPD:C5	2.23	0.69
2:BBB:157:CYS:SG	12:BBB:444:HOH:O	2.50	0.69
1:CCC:28:PHE:CE1	1:CCC:35[B]:ARG:NH2	2.60	0.68
10:DDD:208:PG4:C6	10:DDD:208:PG4:H31	2.24	0.68
1:EEE:73:GLY:H	10:EEE:204:PG4:H12	1.60	0.67
1:EEE:40:GLN:HG2	7:EEE:219:PGE:H5	1.77	0.66
8:AAA:206[B]:PEG:C3	12:AAA:302:HOH:O	2.05	0.65
2:DDD:127:VAL:HG11	9:DDD:217[A]:EDO:H22	1.79	0.64
2:BBB:131:GLN:HG2	7:BBB:208:PGE:H52	1.80	0.64
1:EEE:73:GLY:H	10:EEE:204:PG4:H11	1.62	0.64
1:EEE:76[A]:MET:HG2	12:EEE:377:HOH:O	1.95	0.64
2:BBB:91:TYR:OH	12:BBB:303:HOH:O	2.15	0.64
1:EEE:116:LYS:O	9:EEE:213:EDO:H12	1.98	0.64
8:AAA:206[A]:PEG:O4	12:AAA:302:HOH:O	2.14	0.63
1:AAA:76[B]:MET:SD	12:AAA:361:HOH:O	2.55	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BBB:203:MPD:H11	12:BBB:399:HOH:O	1.98	0.63
8:FFF:213:PEG:O4	8:FFF:214:PEG:H21	1.98	0.63
2:FFF:58:LYS:HD2	7:FFF:209:PGE:O4	1.99	0.63
9:FFF:215:EDO:C1	12:FFF:327:HOH:O	2.46	0.62
1:EEE:120:THR:O	9:EEE:214:EDO:C1	2.47	0.62
7:EEE:219:PGE:C3	12:EEE:301:HOH:O	2.45	0.62
1:EEE:28:PHE:O	1:EEE:35[B]:ARG:NH2	2.33	0.61
2:FFF:106[B]:GLU:HG2	2:FFF:107:ARG:HG3	1.81	0.61
2:FFF:3:ASP:H	2:FFF:6[B]:THR:HG22	1.66	0.61
2:FFF:106[B]:GLU:CG	2:FFF:107:ARG:HG3	2.30	0.61
10:DDD:208:PG4:C6	10:DDD:208:PG4:C3	2.78	0.60
2:DDD:39[B]:ARG:HH21	8:DDD:210:PEG:H21	1.65	0.60
1:EEE:120:THR:O	9:EEE:214:EDO:H12	2.01	0.60
8:AAA:207:PEG:H42	8:AAA:207:PEG:H11	1.80	0.60
1:CCC:40[A]:GLN:NE2	10:DDD:207:PG4:H31	2.17	0.59
4:BBB:204:MPD:HM3	12:BBB:315:HOH:O	2.03	0.59
1:CCC:40[A]:GLN:HE22	10:DDD:207:PG4:H31	1.67	0.59
1:CCC:160:GLN:NE2	1:CCC:160:GLN:HA	2.18	0.59
2:FFF:90:ARG:HE	9:FFF:215:EDO:H21	1.68	0.59
1:CCC:11:ASP:OD2	2:DDD:107[B]:ARG:HD3	2.02	0.59
1:EEE:122:ILE:H	9:EEE:214:EDO:C2	2.16	0.59
2:FFF:140:LEU:HD21	4:FFF:203:MPD:H32	1.85	0.58
1:AAA:101:THR:HG21	8:AAA:208:PEG:H31	1.85	0.58
12:DDD:302:HOH:O	9:FFF:218:EDO:C2	2.51	0.58
4:FFF:203:MPD:HO2	7:FFF:210:PGE:HO4	1.52	0.58
2:FFF:7[B]:SER:OG	7:FFF:207:PGE:H12	2.04	0.57
8:AAA:206[B]:PEG:H22	8:AAA:206[B]:PEG:HO4	1.69	0.57
1:CCC:28:PHE:CE1	1:CCC:35[B]:ARG:CZ	2.88	0.57
1:CCC:160:GLN:HA	1:CCC:160:GLN:HE21	1.69	0.57
1:EEE:34:GLN:OE1	1:EEE:37[A]:ARG:NH2	2.36	0.57
1:AAA:13:GLU:OE1	7:AAA:205:PGE:H12	2.04	0.57
2:FFF:3:ASP:H	2:FFF:6[B]:THR:CG2	2.17	0.57
1:AAA:105:GLU:OE2	8:AAA:208:PEG:H42	2.05	0.56
2:DDD:36:LEU:HA	2:DDD:39[A]:ARG:HG2	1.86	0.56
2:FFF:86[B]:ASP:OD2	9:FFF:215:EDO:C1	2.54	0.56
1:EEE:24[A]:ARG:NE	1:EEE:24[A]:ARG:HA	2.20	0.56
1:EEE:122:ILE:HG22	9:EEE:214:EDO:H21	1.88	0.55
8:EEE:207:PEG:O1	8:EEE:207:PEG:H32	2.00	0.55
2:DDD:39[A]:ARG:NH2	8:DDD:210:PEG:C3	2.69	0.55
2:FFF:90:ARG:HE	9:FFF:215:EDO:C1	2.18	0.55
1:AAA:76[B]:MET:HG3	12:AAA:393:HOH:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:35[B]:ARG:NH1	1:AAA:147[B]:GLU:OE1	2.40	0.55
1:CCC:28:PHE:CZ	1:CCC:35[B]:ARG:NH2	2.75	0.54
1:EEE:24[B]:ARG:NH1	12:EEE:307:HOH:O	2.39	0.54
1:EEE:37[A]:ARG:HG3	7:EEE:219:PGE:H62	1.88	0.54
1:EEE:122:ILE:H	9:EEE:214:EDO:H22	1.72	0.54
2:FFF:86[B]:ASP:OD2	9:FFF:215:EDO:C2	2.56	0.54
10:DDD:208:PG4:H31	10:DDD:208:PG4:H62	1.89	0.54
1:CCC:5:LYS:NZ	1:CCC:99:ASP:CG	2.62	0.53
10:DDD:208:PG4:C3	10:DDD:208:PG4:H62	2.39	0.53
2:DDD:39[A]:ARG:HD2	12:DDD:359:HOH:O	2.10	0.52
2:DDD:39[B]:ARG:NH2	8:DDD:210:PEG:H21	2.25	0.52
2:FFF:62:TYR:O	9:FFF:216:EDO:H21	2.09	0.52
2:BBB:59:SER:HA	7:BBB:208:PGE:H1	1.91	0.52
2:FFF:6[A]:THR:CG2	12:FFF:322:HOH:O	1.97	0.52
2:BBB:25:GLU:OE1	6:BBB:206:1PE:OH2	2.16	0.52
1:AAA:5:LYS:NZ	12:AAA:308:HOH:O	2.42	0.52
9:EEE:210:EDO:O1	12:EEE:301:HOH:O	2.19	0.51
1:AAA:27:SER:OG	7:AAA:215:PGE:H22	2.10	0.51
1:AAA:35[B]:ARG:NH2	1:AAA:147[B]:GLU:OE1	2.44	0.51
4:BBB:203:MPD:O4	4:BBB:203:MPD:O2	2.14	0.51
1:AAA:62:PRO:HB3	5:AAA:203:BCN:H51	1.91	0.51
1:AAA:112:ARG:NH1	1:AAA:160:GLN:O	2.44	0.51
2:BBB:17[A]:LYS:HE3	12:BBB:301:HOH:O	2.11	0.51
2:BBB:17[B]:LYS:NZ	12:BBB:302:HOH:O	2.44	0.51
1:CCC:5:LYS:HZ1	1:CCC:99:ASP:CG	2.13	0.50
2:DDD:77[B]:ARG:NH1	9:DDD:216:EDO:H11	2.27	0.50
1:CCC:40[B]:GLN:NE2	12:CCC:303:HOH:O	2.33	0.50
1:AAA:35[B]:ARG:CZ	1:AAA:147[B]:GLU:OE1	2.59	0.50
2:DDD:76:ARG:HG3	8:DDD:211:PEG:H31	1.94	0.49
2:FFF:2:GLN:HB2	2:FFF:6[B]:THR:HG23	1.92	0.49
2:FFF:90:ARG:HE	9:FFF:215:EDO:C2	2.24	0.49
8:BBB:210:PEG:H41	9:BBB:212:EDO:O2	2.13	0.49
1:AAA:76[A]:MET:HG2	12:AAA:393:HOH:O	2.12	0.49
1:EEE:40:GLN:CB	7:EEE:219:PGE:H5	2.43	0.49
1:AAA:13:GLU:CD	7:AAA:205:PGE:H12	2.34	0.48
4:FFF:203:MPD:O2	7:FFF:210:PGE:O4	2.29	0.48
7:FFF:210:PGE:H62	7:FFF:210:PGE:C3	2.35	0.48
1:EEE:73:GLY:N	10:EEE:204:PG4:H11	2.27	0.48
1:CCC:104:GLU:HA	1:CCC:108:ILE:HB	1.95	0.48
2:FFF:50:ALA:HB1	4:FFF:203:MPD:H52	1.96	0.48
1:EEE:40:GLN:CG	7:EEE:219:PGE:H5	2.41	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CCC:40[A]:GLN:NE2	10:DDD:207:PG4:C3	2.77	0.48
2:BBB:17[C]:LYS:HE2	12:BBB:302:HOH:O	2.14	0.47
2:DDD:39[B]:ARG:HD3	8:DDD:210:PEG:C1	2.32	0.47
4:DDD:205:MPD:H52	4:DDD:205:MPD:HM3	0.63	0.47
2:FFF:13:ASP:OD2	9:FFF:218:EDO:C2	2.62	0.47
4:DDD:205:MPD:CM	4:DDD:205:MPD:C5	2.29	0.47
2:BBB:2:GLN:HE22	9:BBB:215:EDO:H21	1.79	0.47
2:BBB:59:SER:CB	7:BBB:208:PGE:H1	2.44	0.47
12:BBB:437:HOH:O	5:DDD:206:BCN:H41	2.14	0.47
2:FFF:90:ARG:HE	9:FFF:215:EDO:H12	1.80	0.47
1:AAA:36:LEU:CD1	2:BBB:28:LYS:HG3	2.45	0.46
8:CCC:206:PEG:H42	8:CCC:206:PEG:C1	2.18	0.46
2:DDD:127:VAL:CG1	9:DDD:217[A]:EDO:H22	2.45	0.46
2:FFF:64:ASP:HB3	7:FFF:208:PGE:H4	1.96	0.46
1:CCC:35[B]:ARG:NH2	2:DDD:31:PHE:CZ	2.84	0.46
2:DDD:25:GLU:OE2	10:DDD:207:PG4:O1	2.34	0.46
1:CCC:28:PHE:HA	9:CCC:210:EDO:C2	2.46	0.46
1:EEE:73:GLY:N	10:EEE:204:PG4:C1	2.71	0.46
8:CCC:211:PEG:H22	2:DDD:18:TYR:HB2	1.97	0.46
5:AAA:203:BCN:H11	5:AAA:203:BCN:H61	1.40	0.46
10:DDD:208:PG4:H11	10:DDD:208:PG4:C4	2.43	0.45
1:EEE:120:THR:O	9:EEE:214:EDO:H11	2.16	0.45
8:CCC:211:PEG:O1	8:CCC:211:PEG:H32	2.16	0.45
4:FFF:204:MPD:H53	4:FFF:204:MPD:O2	2.17	0.45
8:AAA:209:PEG:H41	8:AAA:209:PEG:H21	1.76	0.45
6:BBB:217:1PE:H222	6:BBB:217:1PE:H131	1.25	0.45
2:DDD:47:ASN:ND2	9:DDD:213:EDO:C2	2.80	0.45
1:AAA:46[A]:ARG:HD2	2:BBB:18:TYR:CE1	2.52	0.44
1:AAA:76[B]:MET:CE	3:AAA:201:PXQ:CGD	2.95	0.44
9:DDD:213:EDO:C1	12:DDD:360:HOH:O	2.64	0.44
4:DDD:203:MPD:HM3	4:DDD:203:MPD:O4	2.17	0.44
9:DDD:213:EDO:H12	12:DDD:360:HOH:O	2.16	0.44
2:FFF:10[A]:ASN:ND2	12:FFF:305:HOH:O	2.31	0.44
2:DDD:36:LEU:HA	2:DDD:39[A]:ARG:CG	2.47	0.44
2:FFF:13:ASP:OD2	9:FFF:218:EDO:C1	2.66	0.44
1:AAA:97:ALA:HA	2:BBB:5:ILE:HG21	2.00	0.43
2:FFF:11[B]:SER:OG	7:FFF:207:PGE:H3	2.18	0.43
6:BBB:217:1PE:H131	6:BBB:217:1PE:H151	1.99	0.43
2:FFF:90:ARG:NE	9:FFF:215:EDO:H12	2.32	0.43
8:CCC:205[A]:PEG:H11	8:CCC:205[A]:PEG:H31	1.61	0.43
8:DDD:212:PEG:C4	8:DDD:212:PEG:C1	2.93	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:FFF:65[A]:ILE:HG12	7:FFF:208:PGE:H2	1.99	0.43
2:FFF:106[A]:GLU:OE2	8:FFF:211:PEG:H21	2.18	0.43
2:FFF:150:GLY:HA3	8:FFF:212:PEG:H12	2.00	0.42
2:DDD:13:ASP:OD2	5:DDD:206:BCN:H61	2.19	0.42
2:DDD:47:ASN:HD22	9:DDD:213:EDO:C2	2.32	0.42
2:DDD:77[B]:ARG:NE	3:DDD:202:PXQ:O1D	2.52	0.42
5:DDD:206:BCN:H62	5:DDD:206:BCN:H31	1.84	0.42
4:FFF:203:MPD:C5	4:FFF:203:MPD:C1	2.91	0.42
1:AAA:28:PHE:CE1	1:AAA:35[B]:ARG:NE	2.88	0.42
8:AAA:206[A]:PEG:C3	12:AAA:302:HOH:O	2.67	0.42
12:BBB:437:HOH:O	5:DDD:206:BCN:C4	2.68	0.42
2:DDD:77[A]:ARG:NH2	12:DDD:311:HOH:O	2.52	0.41
1:EEE:104:GLU:HA	1:EEE:108:ILE:HB	2.02	0.41
2:DDD:47:ASN:ND2	9:DDD:213:EDO:O2	2.51	0.41
1:CCC:9:ASN:ND2	12:CCC:313:HOH:O	2.52	0.41
1:CCC:113:GLU:HG2	9:CCC:201:EDO:H12	2.02	0.41
1:EEE:116:LYS:O	9:EEE:213:EDO:C1	2.67	0.41
9:EEE:213:EDO:H11	9:EEE:217:EDO:H12	2.02	0.41
1:EEE:73:GLY:N	10:EEE:204:PG4:H12	2.30	0.41
7:EEE:205:PGE:H12	12:EEE:385:HOH:O	2.20	0.41
1:AAA:35[B]:ARG:HD3	1:AAA:95:ILE:O	2.20	0.41
1:CCC:24[A]:ARG:HD2	1:CCC:24[A]:ARG:HA	1.83	0.41
2:DDD:39[A]:ARG:CD	12:DDD:359:HOH:O	2.69	0.41
1:AAA:76[B]:MET:CG	12:AAA:393:HOH:O	2.66	0.41
1:AAA:101:THR:N	1:AAA:102:PRO:CD	2.83	0.41
1:EEE:35[B]:ARG:NH1	9:EEE:212:EDO:C1	2.84	0.41
2:FFF:106[B]:GLU:HG3	2:FFF:107:ARG:HG3	2.03	0.41
2:BBB:17[B]:LYS:HB2	2:BBB:17[B]:LYS:HE2	1.70	0.40
1:CCC:160:GLN:NE2	1:CCC:160:GLN:CA	2.84	0.40
1:EEE:122:ILE:HG22	9:EEE:214:EDO:C2	2.51	0.40
2:FFF:7[B]:SER:OG	7:FFF:207:PGE:C1	2.69	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CCC:23:ASP:OD2	1:EEE:24[B]:ARG:NH2[3_555]	2.02	0.18

## 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	AAA	167/160 (104%)	167 (100%)	0	0	100	100
1	CCC	166/160 (104%)	166 (100%)	0	0	100	100
1	EEE	168/160 (105%)	167 (99%)	1 (1%)	0	100	100
2	BBB	162/161 (101%)	159 (98%)	3 (2%)	0	100	100
2	DDD	164/161 (102%)	161 (98%)	3 (2%)	0	100	100
2	FFF	172/161 (107%)	171 (99%)	1 (1%)	0	100	100
All	All	999/963 (104%)	991 (99%)	8 (1%)	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	AAA	131/122 (107%)	130 (99%)	1 (1%)	81	75
1	CCC	130/122 (107%)	124 (95%)	6 (5%)	27	10
1	EEE	132/122 (108%)	129 (98%)	3 (2%)	50	34
2	BBB	129/125 (103%)	128 (99%)	1 (1%)	81	75
2	DDD	131/125 (105%)	125 (95%)	6 (5%)	27	10
2	FFF	139/125 (111%)	137 (99%)	2 (1%)	67	55
All	All	792/741 (107%)	773 (98%)	19 (2%)	57	32

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	160	GLN
2	BBB	11	SER
1	CCC	34[A]	GLN
1	CCC	34[B]	GLN
1	CCC	36	LEU
1	CCC	60	LYS
1	CCC	76[A]	MET
1	CCC	76[B]	MET
2	DDD	2	GLN
2	DDD	11	SER
2	DDD	24	LEU
2	DDD	39[A]	ARG
2	DDD	39[B]	ARG
2	DDD	58	LYS
1	EEE	24[A]	ARG
1	EEE	24[B]	ARG
1	EEE	46	ARG
2	FFF	11[A]	SER
2	FFF	11[B]	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

3 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	$\# Z  > 2$
2	MEN	BBB	71	2	7,8,9	0.58	0	6,9,11	0.47	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	MEN	FFF	71	2	7,8,9	0.80	0	6,9,11	1.11	1 (16%)
2	MEN	DDD	71	2	7,8,9	1.73	1 (14%)	6,9,11	2.03	1 (16%)

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	MEN	BBB	71	2	-	2/7/8/10	-
2	MEN	FFF	71	2	-	2/7/8/10	-
2	MEN	DDD	71	2	-	2/7/8/10	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	DDD	71	MEN	CB-CG	3.91	1.60	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	DDD	71	MEN	OD1-CG-CB	4.52	128.11	121.50
2	FFF	71	MEN	CB-CG-ND2	2.52	118.88	115.48

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	BBB	71	MEN	CA-CB-CG-OD1
2	FFF	71	MEN	CA-CB-CG-OD1
2	BBB	71	MEN	CA-CB-CG-ND2
2	DDD	71	MEN	CA-CB-CG-ND2
2	FFF	71	MEN	CA-CB-CG-ND2
2	DDD	71	MEN	CA-CB-CG-OD1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	DDD	71	MEN	1	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

105 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	$\# Z  > 2$
8	PEG	CCC	208	-	6,6,6	0.11	0	5,5,5	0.15	0
8	PEG	BBB	211	-	6,6,6	0.13	0	5,5,5	0.15	0
9	EDO	EEE	214	-	3,3,3	0.86	0	2,2,2	0.22	0
8	PEG	AAA	206[B]	-	6,6,6	0.24	0	5,5,5	0.14	0
7	PGE	FFF	208	-	9,9,9	0.12	0	8,8,8	0.12	0
8	PEG	CCC	206	-	6,6,6	0.14	0	5,5,5	0.07	0
7	PGE	AAA	205	-	9,9,9	0.25	0	8,8,8	0.23	0
11	MRD	EEE	203	-	7,7,7	0.30	0	9,10,10	0.59	0
9	EDO	DDD	214	-	3,3,3	0.19	0	2,2,2	0.24	0
8	PEG	DDD	210	-	6,6,6	0.14	0	5,5,5	0.11	0
4	MPD	DDD	204	-	7,7,7	0.15	0	9,10,10	0.44	0
9	EDO	DDD	216	-	3,3,3	0.23	0	2,2,2	0.43	0
9	EDO	DDD	201	-	3,3,3	0.21	0	2,2,2	0.04	0
8	PEG	CCC	211	-	6,6,6	0.14	0	5,5,5	0.14	0
4	MPD	DDD	205	-	7,7,7	0.36	0	9,10,10	0.28	0
9	EDO	EEE	216	-	3,3,3	0.09	0	2,2,2	0.21	0
3	PXQ	FFF	201	2	38,46,46	3.95	6 (15%)	40,67,67	1.29	4 (10%)
8	PEG	FFF	212	-	6,6,6	0.32	0	5,5,5	0.35	0
9	EDO	EEE	215[B]	-	3,3,3	0.25	0	2,2,2	0.29	0
7	PGE	BBB	207	-	9,9,9	0.19	0	8,8,8	0.14	0
9	EDO	CCC	210	-	3,3,3	0.10	0	2,2,2	0.41	0
9	EDO	DDD	218	-	3,3,3	0.30	0	2,2,2	0.42	0
9	EDO	BBB	214	-	3,3,3	0.13	0	2,2,2	0.26	0
9	EDO	EEE	218	-	3,3,3	0.06	0	2,2,2	0.23	0
9	EDO	FFF	217	-	3,3,3	0.22	0	2,2,2	0.38	0
9	EDO	AAA	211	-	3,3,3	0.40	0	2,2,2	0.47	0
8	PEG	AAA	208	-	6,6,6	0.23	0	5,5,5	0.23	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
9	EDO	FFF	215	-	3,3,3	0.51	0	2,2,2	0.36	0
9	EDO	EEE	210	-	3,3,3	0.06	0	2,2,2	0.24	0
9	EDO	CCC	201	-	3,3,3	0.12	0	2,2,2	0.27	0
9	EDO	CCC	209	-	3,3,3	0.09	0	2,2,2	0.22	0
9	EDO	FFF	216	-	3,3,3	0.15	0	2,2,2	0.33	0
8	PEG	BBB	209	-	6,6,6	0.22	0	5,5,5	0.22	0
6	1PE	AAA	204	-	15,15,15	0.51	0	14,14,14	0.60	0
8	PEG	CCC	205[B]	-	6,6,6	0.22	0	5,5,5	0.11	0
7	PGE	AAA	215	-	9,9,9	0.14	0	8,8,8	0.10	0
9	EDO	FFF	218	-	3,3,3	0.39	0	2,2,2	0.69	0
7	PGE	BBB	208	-	9,9,9	0.37	0	8,8,8	0.30	0
9	EDO	DDD	213	-	3,3,3	0.24	0	2,2,2	0.48	0
10	PG4	EEE	201	-	12,12,12	0.16	0	11,11,11	0.16	0
9	EDO	AAA	210[A]	-	3,3,3	0.12	0	2,2,2	0.27	0
9	EDO	DDD	217[B]	-	3,3,3	0.13	0	2,2,2	0.34	0
8	PEG	AAA	207	-	6,6,6	0.12	0	5,5,5	0.33	0
6	1PE	FFF	206	-	15,15,15	0.40	0	14,14,14	0.45	0
7	PGE	EEE	205	-	9,9,9	0.25	0	8,8,8	0.13	0
6	1PE	BBB	217	-	15,15,15	0.54	0	14,14,14	0.16	0
9	EDO	DDD	215[B]	-	3,3,3	0.15	0	2,2,2	0.28	0
7	PGE	EEE	219	-	9,9,9	0.22	0	8,8,8	0.16	0
6	1PE	BBB	205	-	15,15,15	0.42	0	14,14,14	0.60	0
8	PEG	AAA	206[A]	-	6,6,6	0.19	0	5,5,5	0.11	0
9	EDO	BBB	216	-	3,3,3	0.12	0	2,2,2	0.26	0
9	EDO	BBB	212	-	3,3,3	0.07	0	2,2,2	0.24	0
4	MPD	FFF	203	-	7,7,7	0.17	0	9,10,10	0.58	0
3	PXQ	BBB	201	2	38,46,46	4.09	5 (13%)	40,67,67	1.39	3 (7%)
3	PXQ	AAA	201	1	38,46,46	3.71	5 (13%)	40,67,67	1.46	5 (12%)
9	EDO	AAA	214	-	3,3,3	0.14	0	2,2,2	0.33	0
7	PGE	CCC	204	-	9,9,9	0.14	0	8,8,8	0.09	0
9	EDO	EEE	215[A]	-	3,3,3	0.35	0	2,2,2	0.61	0
10	PG4	DDD	209	-	12,12,12	0.24	0	11,11,11	0.22	0
4	MPD	BBB	204	-	7,7,7	0.37	0	9,10,10	0.59	0
3	PXQ	CCC	202	1	38,46,46	4.27	7 (18%)	40,67,67	1.48	4 (10%)
8	PEG	AAA	209	-	6,6,6	0.15	0	5,5,5	0.06	0
7	PGE	FFF	209	-	9,9,9	0.21	0	8,8,8	0.22	0
4	MPD	DDD	203	-	7,7,7	0.31	0	9,10,10	0.53	0
4	MPD	BBB	202	-	7,7,7	0.28	0	9,10,10	0.53	0
9	EDO	AAA	213	-	3,3,3	0.08	0	2,2,2	0.17	0
10	PG4	DDD	207	-	12,12,12	0.16	0	11,11,11	0.21	0
4	MPD	AAA	202	-	7,7,7	0.33	0	9,10,10	0.49	0
8	PEG	BBB	210	-	6,6,6	0.11	0	5,5,5	0.07	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	MPD	CCC	203	-	7,7,7	0.27	0	9,10,10	0.45	0
8	PEG	EEE	208	-	6,6,6	0.17	0	5,5,5	0.17	0
3	PXQ	DDD	202	2	38,46,46	4.09	5 (13%)	40,67,67	1.21	6 (15%)
8	PEG	CCC	207	-	6,6,6	0.29	0	5,5,5	0.14	0
8	PEG	CCC	205[A]	-	6,6,6	0.21	0	5,5,5	0.17	0
8	PEG	EEE	209	-	6,6,6	0.14	0	5,5,5	0.08	0
3	PXQ	EEE	202	1	38,46,46	3.88	7 (18%)	40,67,67	1.23	4 (10%)
4	MPD	FFF	202	-	7,7,7	0.45	0	9,10,10	0.46	0
8	PEG	EEE	206	-	6,6,6	0.12	0	5,5,5	0.11	0
8	PEG	FFF	213	-	6,6,6	0.16	0	5,5,5	0.08	0
8	PEG	EEE	207	-	6,6,6	0.23	0	5,5,5	0.17	0
9	EDO	DDD	217[A]	-	3,3,3	0.37	0	2,2,2	0.60	0
8	PEG	DDD	211	-	6,6,6	0.23	0	5,5,5	0.20	0
9	EDO	EEE	212	-	3,3,3	0.07	0	2,2,2	0.11	0
10	PG4	EEE	204	-	12,12,12	0.27	0	11,11,11	0.20	0
9	EDO	EEE	217	-	3,3,3	0.11	0	2,2,2	0.23	0
8	PEG	DDD	212	-	6,6,6	0.09	0	5,5,5	0.30	0
9	EDO	BBB	213	-	3,3,3	0.17	0	2,2,2	0.13	0
9	EDO	DDD	215[A]	-	3,3,3	0.11	0	2,2,2	0.24	0
9	EDO	AAA	210[B]	-	3,3,3	0.10	0	2,2,2	0.26	0
6	1PE	BBB	206	-	15,15,15	0.50	0	14,14,14	0.68	0
10	PG4	DDD	208	-	12,12,12	0.16	0	11,11,11	0.15	0
8	PEG	FFF	214	-	6,6,6	0.13	0	5,5,5	0.19	0
4	MPD	BBB	203	-	7,7,7	0.23	0	9,10,10	0.50	0
5	BCN	DDD	206	-	10,10,10	0.90	1 (10%)	11,11,11	0.98	0
9	EDO	AAA	212	-	3,3,3	0.09	0	2,2,2	0.24	0
9	EDO	AAA	216	-	3,3,3	0.06	0	2,2,2	0.04	0
5	BCN	AAA	203	-	10,10,10	0.62	0	11,11,11	0.84	0
9	EDO	EEE	211	-	3,3,3	0.07	0	2,2,2	0.23	0
9	EDO	EEE	213	-	3,3,3	0.26	0	2,2,2	0.47	0
4	MPD	FFF	204	-	7,7,7	0.29	0	9,10,10	0.62	0
8	PEG	FFF	211	-	6,6,6	0.16	0	5,5,5	0.10	0
9	EDO	BBB	215	-	3,3,3	0.11	0	2,2,2	0.22	0
7	PGE	FFF	207	-	9,9,9	0.16	0	8,8,8	0.22	0
7	PGE	FFF	210	-	9,9,9	0.23	0	8,8,8	0.13	0
4	MPD	FFF	205	-	7,7,7	0.28	0	9,10,10	0.57	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	PEG	CCC	208	-	-	2/4/4/4	-
8	PEG	BBB	211	-	-	2/4/4/4	-
9	EDO	EEE	214	-	-	0/1/1/1	-
8	PEG	AAA	206[B]	-	-	3/4/4/4	-
7	PGE	FFF	208	-	-	4/7/7/7	-
8	PEG	CCC	206	-	-	4/4/4/4	-
7	PGE	AAA	205	-	-	2/7/7/7	-
11	MRD	EEE	203	-	-	0/5/5/5	-
9	EDO	DDD	214	-	-	1/1/1/1	-
8	PEG	DDD	210	-	-	0/4/4/4	-
4	MPD	DDD	204	-	-	1/5/5/5	-
9	EDO	DDD	216	-	-	1/1/1/1	-
9	EDO	DDD	201	-	-	1/1/1/1	-
8	PEG	CCC	211	-	-	1/4/4/4	-
4	MPD	DDD	205	-	-	5/5/5/5	-
9	EDO	EEE	216	-	-	1/1/1/1	-
3	PXQ	FFF	201	2	-	7/23/58/58	0/4/4/4
8	PEG	FFF	212	-	-	0/4/4/4	-
9	EDO	EEE	215[B]	-	-	0/1/1/1	-
7	PGE	BBB	207	-	-	3/7/7/7	-
9	EDO	CCC	210	-	-	1/1/1/1	-
9	EDO	DDD	218	-	-	0/1/1/1	-
9	EDO	BBB	214	-	-	0/1/1/1	-
9	EDO	EEE	218	-	-	0/1/1/1	-
9	EDO	FFF	217	-	-	1/1/1/1	-
9	EDO	AAA	211	-	-	1/1/1/1	-
8	PEG	AAA	208	-	-	3/4/4/4	-
9	EDO	FFF	215	-	-	0/1/1/1	-
9	EDO	EEE	210	-	-	1/1/1/1	-
9	EDO	CCC	201	-	-	1/1/1/1	-
9	EDO	CCC	209	-	-	1/1/1/1	-
9	EDO	FFF	216	-	-	0/1/1/1	-
8	PEG	BBB	209	-	-	3/4/4/4	-
6	1PE	AAA	204	-	-	2/13/13/13	-
8	PEG	CCC	205[B]	-	-	3/4/4/4	-
7	PGE	AAA	215	-	-	1/7/7/7	-
9	EDO	FFF	218	-	-	1/1/1/1	-
7	PGE	BBB	208	-	-	5/7/7/7	-
9	EDO	DDD	213	-	-	1/1/1/1	-
10	PG4	EEE	201	-	-	5/10/10/10	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	EDO	AAA	210[A]	-	-	1/1/1/1	-
9	EDO	DDD	217[B]	-	-	1/1/1/1	-
8	PEG	AAA	207	-	-	3/4/4/4	-
6	1PE	FFF	206	-	-	6/13/13/13	-
7	PGE	EEE	205	-	-	4/7/7/7	-
6	1PE	BBB	217	-	-	10/13/13/13	-
9	EDO	DDD	215[B]	-	-	1/1/1/1	-
7	PGE	EEE	219	-	-	4/7/7/7	-
6	1PE	BBB	205	-	-	9/13/13/13	-
8	PEG	AAA	206[A]	-	-	4/4/4/4	-
9	EDO	BBB	216	-	-	1/1/1/1	-
9	EDO	BBB	212	-	-	1/1/1/1	-
4	MPD	FFF	203	-	-	2/5/5/5	-
3	PXQ	BBB	201	2	-	8/23/58/58	0/4/4/4
3	PXQ	AAA	201	1	-	7/23/58/58	0/4/4/4
9	EDO	AAA	214	-	-	0/1/1/1	-
7	PGE	CCC	204	-	-	4/7/7/7	-
9	EDO	EEE	215[A]	-	-	1/1/1/1	-
10	PG4	DDD	209	-	-	7/10/10/10	-
4	MPD	BBB	204	-	-	4/5/5/5	-
3	PXQ	CCC	202	1	-	7/23/58/58	0/4/4/4
8	PEG	AAA	209	-	-	2/4/4/4	-
7	PGE	FFF	209	-	-	5/7/7/7	-
4	MPD	DDD	203	-	-	0/5/5/5	-
4	MPD	BBB	202	-	-	0/5/5/5	-
9	EDO	AAA	213	-	-	0/1/1/1	-
10	PG4	DDD	207	-	-	6/10/10/10	-
4	MPD	AAA	202	-	-	2/5/5/5	-
8	PEG	BBB	210	-	-	2/4/4/4	-
4	MPD	CCC	203	-	-	0/5/5/5	-
8	PEG	EEE	208	-	-	3/4/4/4	-
3	PXQ	DDD	202	2	-	9/23/58/58	0/4/4/4
8	PEG	CCC	207	-	-	0/4/4/4	-
8	PEG	CCC	205[A]	-	-	3/4/4/4	-
8	PEG	EEE	209	-	-	2/4/4/4	-
3	PXQ	EEE	202	1	-	9/23/58/58	0/4/4/4
4	MPD	FFF	202	-	-	1/5/5/5	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	PEG	EEE	206	-	-	1/4/4/4	-
8	PEG	FFF	213	-	-	2/4/4/4	-
8	PEG	EEE	207	-	-	2/4/4/4	-
9	EDO	DDD	217[A]	-	-	1/1/1/1	-
8	PEG	DDD	211	-	-	3/4/4/4	-
9	EDO	EEE	212	-	-	0/1/1/1	-
10	PG4	EEE	204	-	-	5/10/10/10	-
9	EDO	EEE	217	-	-	0/1/1/1	-
8	PEG	DDD	212	-	-	2/4/4/4	-
9	EDO	BBB	213	-	-	1/1/1/1	-
9	EDO	DDD	215[A]	-	-	1/1/1/1	-
9	EDO	AAA	210[B]	-	-	1/1/1/1	-
6	1PE	BBB	206	-	-	6/13/13/13	-
10	PG4	DDD	208	-	-	7/10/10/10	-
8	PEG	FFF	214	-	-	4/4/4/4	-
4	MPD	BBB	203	-	-	2/5/5/5	-
5	BCN	DDD	206	-	-	4/10/10/10	-
9	EDO	AAA	212	-	-	1/1/1/1	-
9	EDO	AAA	216	-	-	0/1/1/1	-
5	BCN	AAA	203	-	-	5/10/10/10	-
9	EDO	EEE	211	-	-	1/1/1/1	-
9	EDO	EEE	213	-	-	0/1/1/1	-
4	MPD	FFF	204	-	-	3/5/5/5	-
8	PEG	FFF	211	-	-	0/4/4/4	-
9	EDO	BBB	215	-	-	1/1/1/1	-
7	PGE	FFF	207	-	-	5/7/7/7	-
7	PGE	FFF	210	-	-	3/7/7/7	-
4	MPD	FFF	205	-	-	2/5/5/5	-

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	DDD	202	PXQ	CHA-C1A	-15.86	1.39	1.51
3	CCC	202	PXQ	CHA-C4D	-15.47	1.39	1.51
3	CCC	202	PXQ	C4C-NC	15.21	1.40	1.29
3	BBB	201	PXQ	CHA-C4D	-15.19	1.40	1.51
3	DDD	202	PXQ	CHA-C4D	-15.18	1.40	1.51
3	BBB	201	PXQ	CHA-C1A	-14.20	1.40	1.51
3	EEE	202	PXQ	CHA-C1A	-14.07	1.40	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	FFF	201	PXQ	C4C-NC	13.65	1.39	1.29
3	FFF	201	PXQ	CHA-C1A	-13.61	1.41	1.51
3	EEE	202	PXQ	C4C-NC	13.32	1.39	1.29
3	AAA	201	PXQ	CHA-C1A	-13.06	1.41	1.51
3	FFF	201	PXQ	CHA-C4D	-12.85	1.41	1.51
3	CCC	202	PXQ	CHA-C1A	-12.65	1.42	1.51
3	AAA	201	PXQ	CHA-C4D	-12.31	1.42	1.51
3	BBB	201	PXQ	C4C-NC	12.13	1.38	1.29
3	AAA	201	PXQ	C4C-NC	12.01	1.38	1.29
3	EEE	202	PXQ	CHA-C4D	-11.99	1.42	1.51
3	DDD	202	PXQ	C4C-NC	10.45	1.37	1.29
3	FFF	201	PXQ	C2A-C1A	4.89	1.46	1.39
3	BBB	201	PXQ	C2A-C1A	4.72	1.46	1.39
3	AAA	201	PXQ	C2A-C1A	4.52	1.46	1.39
3	DDD	202	PXQ	C2A-C1A	4.49	1.45	1.39
3	EEE	202	PXQ	C2A-C1A	4.38	1.45	1.39
3	CCC	202	PXQ	C2A-C1A	4.26	1.45	1.39
3	CCC	202	PXQ	C3D-C4D	3.45	1.44	1.39
3	AAA	201	PXQ	C3D-C4D	2.87	1.43	1.39
3	EEE	202	PXQ	C3D-C4D	2.81	1.43	1.39
3	CCC	202	PXQ	CHB-C1B	2.79	1.40	1.34
3	FFF	201	PXQ	C3D-C4D	2.78	1.43	1.39
3	BBB	201	PXQ	O2A-CGA	-2.76	1.21	1.30
3	FFF	201	PXQ	O1D-CGD	-2.59	1.22	1.30
3	EEE	202	PXQ	CHD-C4C	-2.52	1.36	1.48
3	DDD	202	PXQ	O1D-CGD	-2.50	1.22	1.30
5	DDD	206	BCN	O22-C2	-2.50	1.22	1.30
3	CCC	202	PXQ	CHD-C4C	-2.31	1.37	1.48
3	EEE	202	PXQ	O1D-CGD	-2.01	1.24	1.30

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	AAA	201	PXQ	CAD-C3D-C4D	-5.26	123.60	127.30
3	CCC	202	PXQ	CAA-C2A-C1A	-5.22	123.63	127.30
3	BBB	201	PXQ	CAD-C3D-C4D	-4.96	123.81	127.30
3	FFF	201	PXQ	CAD-C3D-C4D	-4.40	124.21	127.30
3	CCC	202	PXQ	CAD-C3D-C4D	-3.98	124.50	127.30
3	BBB	201	PXQ	CMB-C2B-C1B	3.90	129.03	124.17
3	EEE	202	PXQ	CMB-C2B-C1B	3.33	128.32	124.17
3	EEE	202	PXQ	CHB-C1B-C2B	-3.24	120.57	126.97
3	AAA	201	PXQ	CHB-C1B-C2B	-3.13	120.79	126.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	AAA	201	PXQ	CAA-C2A-C1A	-3.12	125.11	127.30
3	DDD	202	PXQ	CAD-C3D-C4D	-3.01	125.18	127.30
3	AAA	201	PXQ	CMB-C2B-C1B	2.93	127.83	124.17
3	CCC	202	PXQ	CHB-C1B-C2B	-2.85	121.33	126.97
3	CCC	202	PXQ	CAB-C3B-C2B	2.83	132.36	127.53
3	EEE	202	PXQ	CAD-C3D-C4D	-2.79	125.34	127.30
3	BBB	201	PXQ	CAA-C2A-C1A	-2.73	125.38	127.30
3	FFF	201	PXQ	CHD-C4C-NC	2.64	127.85	124.81
3	DDD	202	PXQ	CAA-C2A-C1A	-2.50	125.54	127.30
3	EEE	202	PXQ	CAA-C2A-C1A	-2.50	125.54	127.30
3	FFF	201	PXQ	CAA-C2A-C1A	-2.49	125.55	127.30
3	DDD	202	PXQ	CAB-C3B-C2B	2.31	131.48	127.53
3	FFF	201	PXQ	CMB-C2B-C1B	2.24	126.96	124.17
3	DDD	202	PXQ	CBD-CAD-C3D	-2.14	108.96	112.62
3	DDD	202	PXQ	O1A-CGA-CBA	-2.13	116.25	123.08
3	AAA	201	PXQ	C3B-C4B-NB	-2.12	105.06	106.78
3	DDD	202	PXQ	CMB-C2B-C1B	2.09	126.78	124.17

There are no chirality outliers.

All (260) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	AAA	201	PXQ	C2A-C1A-CHA-C4D
3	AAA	201	PXQ	C3D-C4D-CHA-C1A
3	AAA	201	PXQ	NA-C4A-CHB-C1B
3	AAA	201	PXQ	C3A-C4A-CHB-C1B
3	AAA	201	PXQ	C2D-C1D-CHD-C4C
3	BBB	201	PXQ	C2A-C1A-CHA-C4D
3	BBB	201	PXQ	C3D-C4D-CHA-C1A
3	BBB	201	PXQ	NA-C4A-CHB-C1B
3	BBB	201	PXQ	C2D-C1D-CHD-C4C
3	CCC	202	PXQ	C2A-C1A-CHA-C4D
3	CCC	202	PXQ	NA-C4A-CHB-C1B
3	CCC	202	PXQ	C3A-C4A-CHB-C1B
3	CCC	202	PXQ	C2D-C1D-CHD-C4C
3	DDD	202	PXQ	C2A-C1A-CHA-C4D
3	DDD	202	PXQ	NA-C4A-CHB-C1B
3	DDD	202	PXQ	C3A-C4A-CHB-C1B
3	DDD	202	PXQ	C2D-C1D-CHD-C4C
3	EEE	202	PXQ	C2A-C1A-CHA-C4D
3	EEE	202	PXQ	NA-C4A-CHB-C1B
3	EEE	202	PXQ	C3A-C4A-CHB-C1B

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Mol	Chain	Res	Type	Atoms
3	EEE	202	PXQ	C2D-C1D-CHD-C4C
3	FFF	201	PXQ	C3D-C4D-CHA-C1A
3	FFF	201	PXQ	C2D-C1D-CHD-C4C
4	DDD	205	MPD	O2-C2-C3-C4
4	DDD	205	MPD	CM-C2-C3-C4
4	DDD	205	MPD	C2-C3-C4-O4
4	DDD	205	MPD	C2-C3-C4-C5
5	AAA	203	BCN	N1-C1-C2-O21
5	AAA	203	BCN	N1-C1-C2-O22
5	DDD	206	BCN	C4-C3-N1-C1
6	BBB	217	1PE	C13-C23-OH3-C22
8	CCC	205[B]	PEG	C1-C2-O2-C3
8	CCC	206	PEG	C1-C2-O2-C3
8	DDD	212	PEG	C1-C2-O2-C3
7	EEE	205	PGE	C1-C2-O2-C3
8	AAA	207	PEG	C1-C2-O2-C3
8	CCC	205[A]	PEG	C1-C2-O2-C3
8	EEE	207	PEG	C1-C2-O2-C3
7	FFF	209	PGE	O2-C3-C4-O3
7	FFF	209	PGE	C3-C4-O3-C5
7	FFF	208	PGE	O2-C3-C4-O3
7	AAA	205	PGE	O2-C3-C4-O3
8	AAA	206[B]	PEG	C4-C3-O2-C2
10	DDD	209	PG4	O2-C3-C4-O3
6	BBB	217	1PE	OH5-C14-C24-OH4
7	EEE	219	PGE	O2-C3-C4-O3
10	EEE	204	PG4	O2-C3-C4-O3
6	BBB	217	1PE	OH2-C12-C22-OH3
8	EEE	206	PEG	O2-C3-C4-O4
8	FFF	214	PEG	O1-C1-C2-O2
10	DDD	208	PG4	O3-C5-C6-O4
6	BBB	205	1PE	OH5-C14-C24-OH4
7	BBB	208	PGE	O2-C3-C4-O3
7	BBB	208	PGE	C4-C3-O2-C2
6	FFF	206	1PE	OH7-C16-C26-OH6
7	BBB	207	PGE	O1-C1-C2-O2
7	BBB	208	PGE	O1-C1-C2-O2
7	CCC	204	PGE	O1-C1-C2-O2
7	EEE	205	PGE	O3-C5-C6-O4
7	FFF	210	PGE	O3-C5-C6-O4
8	AAA	208	PEG	O2-C3-C4-O4
8	BBB	209	PEG	O2-C3-C4-O4

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Mol	Chain	Res	Type	Atoms
8	CCC	205[B]	PEG	O2-C3-C4-O4
8	CCC	206	PEG	O2-C3-C4-O4
8	CCC	208	PEG	O1-C1-C2-O2
8	DDD	212	PEG	O1-C1-C2-O2
8	EEE	207	PEG	O1-C1-C2-O2
8	EEE	208	PEG	O1-C1-C2-O2
8	AAA	207	PEG	C4-C3-O2-C2
5	AAA	203	BCN	C6-C5-N1-C1
10	EEE	201	PG4	O2-C3-C4-O3
8	CCC	206	PEG	C4-C3-O2-C2
7	BBB	207	PGE	O3-C5-C6-O4
7	FFF	207	PGE	O3-C5-C6-O4
8	CCC	211	PEG	O2-C3-C4-O4
8	FFF	214	PEG	O2-C3-C4-O4
10	EEE	204	PG4	O4-C7-C8-O5
10	DDD	208	PG4	C4-C3-O2-C2
10	DDD	208	PG4	O2-C3-C4-O3
10	DDD	208	PG4	C3-C4-O3-C5
8	FFF	214	PEG	C1-C2-O2-C3
7	EEE	219	PGE	O3-C5-C6-O4
7	FFF	209	PGE	O3-C5-C6-O4
8	AAA	207	PEG	O1-C1-C2-O2
8	BBB	210	PEG	O1-C1-C2-O2
8	BBB	211	PEG	O2-C3-C4-O4
10	EEE	201	PG4	O4-C7-C8-O5
6	BBB	217	1PE	OH4-C13-C23-OH3
8	CCC	205[A]	PEG	O2-C3-C4-O4
9	AAA	212	EDO	O1-C1-C2-O2
9	BBB	213	EDO	O1-C1-C2-O2
9	BBB	215	EDO	O1-C1-C2-O2
9	CCC	210	EDO	O1-C1-C2-O2
9	DDD	213	EDO	O1-C1-C2-O2
9	DDD	214	EDO	O1-C1-C2-O2
9	DDD	215[B]	EDO	O1-C1-C2-O2
9	DDD	217[B]	EDO	O1-C1-C2-O2
9	EEE	216	EDO	O1-C1-C2-O2
9	FFF	217	EDO	O1-C1-C2-O2
5	DDD	206	BCN	N1-C1-C2-O22
6	BBB	205	1PE	OH6-C15-C25-OH5
6	BBB	206	1PE	OH6-C15-C25-OH5
7	FFF	208	PGE	O1-C1-C2-O2
6	AAA	204	1PE	OH5-C14-C24-OH4

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Mol	Chain	Res	Type	Atoms
6	BBB	206	1PE	OH7-C16-C26-OH6
7	AAA	215	PGE	O1-C1-C2-O2
7	BBB	208	PGE	O3-C5-C6-O4
7	FFF	210	PGE	O1-C1-C2-O2
8	AAA	206[A]	PEG	O1-C1-C2-O2
8	AAA	206[B]	PEG	O1-C1-C2-O2
8	DDD	211	PEG	O1-C1-C2-O2
10	DDD	207	PG4	O3-C5-C6-O4
6	FFF	206	1PE	OH5-C14-C24-OH4
7	FFF	208	PGE	O3-C5-C6-O4
10	DDD	208	PG4	O4-C7-C8-O5
10	DDD	209	PG4	O4-C7-C8-O5
9	EEE	215[A]	EDO	O1-C1-C2-O2
6	BBB	205	1PE	OH4-C13-C23-OH3
6	BBB	206	1PE	OH4-C13-C23-OH3
10	DDD	207	PG4	O4-C7-C8-O5
7	FFF	210	PGE	O2-C3-C4-O3
8	AAA	206[A]	PEG	O2-C3-C4-O4
7	FFF	207	PGE	O2-C3-C4-O3
8	BBB	210	PEG	O2-C3-C4-O4
8	CCC	208	PEG	O2-C3-C4-O4
9	AAA	210[B]	EDO	O1-C1-C2-O2
9	BBB	212	EDO	O1-C1-C2-O2
9	BBB	216	EDO	O1-C1-C2-O2
9	CCC	201	EDO	O1-C1-C2-O2
9	DDD	215[A]	EDO	O1-C1-C2-O2
9	DDD	217[A]	EDO	O1-C1-C2-O2
9	EEE	210	EDO	O1-C1-C2-O2
9	EEE	211	EDO	O1-C1-C2-O2
9	FFF	218	EDO	O1-C1-C2-O2
10	DDD	207	PG4	O2-C3-C4-O3
6	BBB	205	1PE	C12-C22-OH3-C23
6	BBB	205	1PE	OH7-C16-C26-OH6
7	FFF	207	PGE	O1-C1-C2-O2
6	BBB	205	1PE	C16-C26-OH6-C15
8	AAA	206[A]	PEG	C4-C3-O2-C2
10	DDD	209	PG4	C3-C4-O3-C5
5	DDD	206	BCN	N1-C1-C2-O21
6	BBB	217	1PE	C12-C22-OH3-C23
6	BBB	217	1PE	C24-C14-OH5-C25
7	CCC	204	PGE	C3-C4-O3-C5
8	BBB	209	PEG	C4-C3-O2-C2

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Mol	Chain	Res	Type	Atoms
10	DDD	208	PG4	C5-C6-O4-C7
6	BBB	206	1PE	C12-C22-OH3-C23
7	FFF	209	PGE	C6-C5-O3-C4
6	BBB	217	1PE	C23-C13-OH4-C24
10	EEE	204	PG4	C3-C4-O3-C5
8	DDD	211	PEG	O2-C3-C4-O4
8	FFF	213	PEG	O2-C3-C4-O4
7	CCC	204	PGE	C6-C5-O3-C4
8	DDD	211	PEG	C4-C3-O2-C2
10	DDD	209	PG4	O3-C5-C6-O4
10	EEE	201	PG4	O3-C5-C6-O4
6	BBB	205	1PE	C14-C24-OH4-C13
7	EEE	205	PGE	C6-C5-O3-C4
8	EEE	208	PEG	C4-C3-O2-C2
10	DDD	209	PG4	C1-C2-O2-C3
4	BBB	203	MPD	C2-C3-C4-C5
4	FFF	202	MPD	C2-C3-C4-C5
4	FFF	203	MPD	C2-C3-C4-C5
10	DDD	207	PG4	C1-C2-O2-C3
6	FFF	206	1PE	C12-C22-OH3-C23
6	FFF	206	1PE	C23-C13-OH4-C24
8	AAA	208	PEG	O1-C1-C2-O2
7	EEE	205	PGE	C4-C3-O2-C2
5	AAA	203	BCN	C4-C3-N1-C5
9	DDD	216	EDO	O1-C1-C2-O2
8	AAA	206[B]	PEG	C1-C2-O2-C3
4	AAA	202	MPD	C2-C3-C4-O4
4	BBB	203	MPD	C2-C3-C4-O4
4	FFF	203	MPD	C2-C3-C4-O4
10	EEE	201	PG4	C6-C5-O3-C4
6	BBB	205	1PE	C23-C13-OH4-C24
4	BBB	204	MPD	C1-C2-C3-C4
4	BBB	204	MPD	CM-C2-C3-C4
4	DDD	205	MPD	C1-C2-C3-C4
4	FFF	204	MPD	C1-C2-C3-C4
10	DDD	207	PG4	C3-C4-O3-C5
8	CCC	205[A]	PEG	C4-C3-O2-C2
8	BBB	211	PEG	O1-C1-C2-O2
7	FFF	207	PGE	C1-C2-O2-C3
10	EEE	204	PG4	C1-C2-O2-C3
8	EEE	209	PEG	C1-C2-O2-C3
8	FFF	213	PEG	C4-C3-O2-C2

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Mol	Chain	Res	Type	Atoms
5	AAA	203	BCN	N1-C3-C4-O4
8	AAA	209	PEG	O1-C1-C2-O2
8	EEE	208	PEG	O2-C3-C4-O4
6	BBB	217	1PE	C16-C26-OH6-C15
7	EEE	219	PGE	C3-C4-O3-C5
6	BBB	206	1PE	C16-C26-OH6-C15
10	DDD	209	PG4	O1-C1-C2-O2
8	AAA	209	PEG	C4-C3-O2-C2
8	AAA	206[A]	PEG	C1-C2-O2-C3
8	CCC	205[B]	PEG	C4-C3-O2-C2
8	FFF	214	PEG	C4-C3-O2-C2
8	AAA	208	PEG	C4-C3-O2-C2
6	BBB	217	1PE	OH7-C16-C26-OH6
8	BBB	209	PEG	O1-C1-C2-O2
8	CCC	206	PEG	O1-C1-C2-O2
3	BBB	201	PXQ	CAD-CBD-CGD-O1D
9	AAA	210[A]	EDO	O1-C1-C2-O2
7	BBB	208	PGE	C6-C5-O3-C4
10	EEE	201	PG4	C1-C2-O2-C3
3	BBB	201	PXQ	CAA-CBA-CGA-O2A
3	BBB	201	PXQ	CAD-CBD-CGD-O2D
3	BBB	201	PXQ	CAA-CBA-CGA-O1A
3	FFF	201	PXQ	CAA-CBA-CGA-O1A
5	DDD	206	BCN	C4-C3-N1-C5
3	DDD	202	PXQ	CAA-CBA-CGA-O1A
6	FFF	206	1PE	C15-C25-OH5-C14
6	BBB	217	1PE	C15-C25-OH5-C14
10	EEE	204	PG4	C5-C6-O4-C7
3	EEE	202	PXQ	CAA-CBA-CGA-O1A
6	AAA	204	1PE	OH6-C15-C25-OH5
3	FFF	201	PXQ	CAA-CBA-CGA-O2A
10	DDD	207	PG4	C5-C6-O4-C7
3	AAA	201	PXQ	CAA-CBA-CGA-O2A
3	AAA	201	PXQ	CAA-CBA-CGA-O1A
3	DDD	202	PXQ	CAA-CBA-CGA-O2A
6	BBB	206	1PE	C24-C14-OH5-C25
4	BBB	204	MPD	O2-C2-C3-C4
4	FFF	204	MPD	O2-C2-C3-C4
4	FFF	205	MPD	O2-C2-C3-C4
7	EEE	219	PGE	C4-C3-O2-C2
6	FFF	206	1PE	OH2-C12-C22-OH3
3	EEE	202	PXQ	CAA-CBA-CGA-O2A

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Mol	Chain	Res	Type	Atoms
3	CCC	202	PXQ	CAA-CBA-CGA-O1A
3	CCC	202	PXQ	CAA-CBA-CGA-O2A
3	FFF	201	PXQ	CAD-CBD-CGD-O1D
7	FFF	208	PGE	C1-C2-O2-C3
3	DDD	202	PXQ	CAD-CBD-CGD-O1D
7	FFF	209	PGE	O1-C1-C2-O2
8	EEE	209	PEG	O1-C1-C2-O2
9	AAA	211	EDO	O1-C1-C2-O2
9	CCC	209	EDO	O1-C1-C2-O2
9	DDD	201	EDO	O1-C1-C2-O2
3	DDD	202	PXQ	CAD-CBD-CGD-O2D
3	FFF	201	PXQ	CAD-CBD-CGD-O2D
10	DDD	209	PG4	C6-C5-O3-C4
7	CCC	204	PGE	O2-C3-C4-O3
6	BBB	205	1PE	OH2-C12-C22-OH3
7	AAA	205	PGE	O3-C5-C6-O4
3	CCC	202	PXQ	C3D-C4D-CHA-C1A
3	DDD	202	PXQ	C3D-C4D-CHA-C1A
3	EEE	202	PXQ	C3D-C4D-CHA-C1A
3	FFF	201	PXQ	C2A-C1A-CHA-C4D
4	AAA	202	MPD	C2-C3-C4-C5
4	BBB	204	MPD	C2-C3-C4-C5
7	FFF	207	PGE	C4-C3-O2-C2
10	DDD	208	PG4	C1-C2-O2-C3
7	BBB	207	PGE	O2-C3-C4-O3
3	EEE	202	PXQ	CAD-CBD-CGD-O1D
4	DDD	204	MPD	C2-C3-C4-O4
4	FFF	204	MPD	C2-C3-C4-O4
4	FFF	205	MPD	C2-C3-C4-O4
3	EEE	202	PXQ	CAD-CBD-CGD-O2D

There are no ring outliers.

56 monomers are involved in 150 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	EEE	214	EDO	10	0
8	AAA	206[B]	PEG	4	0
7	FFF	208	PGE	2	0
8	CCC	206	PEG	4	0
7	AAA	205	PGE	2	0
8	DDD	210	PEG	9	0
9	DDD	216	EDO	1	0

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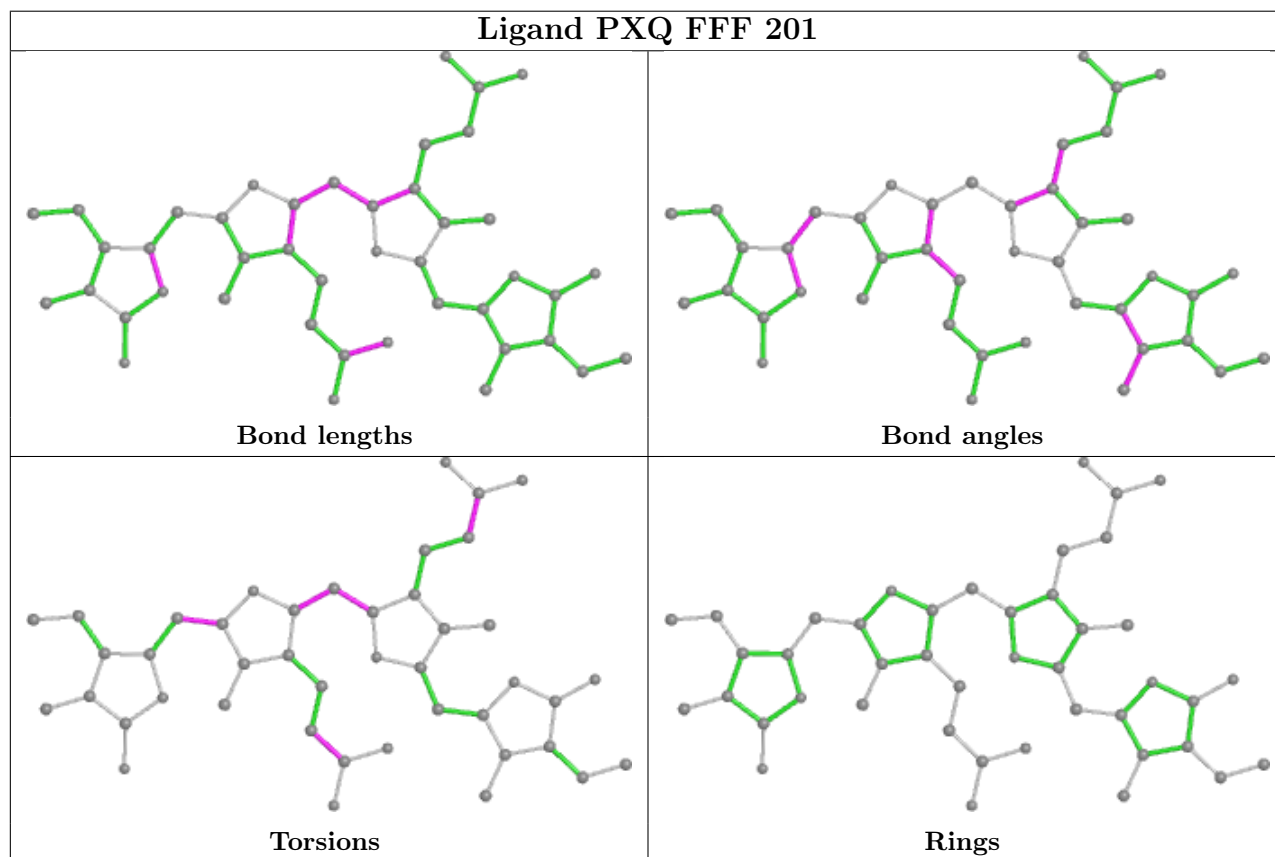
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	CCC	211	PEG	2	0
4	DDD	205	MPD	4	0
8	FFF	212	PEG	1	0
9	CCC	210	EDO	1	0
8	AAA	208	PEG	4	0
9	FFF	215	EDO	11	0
9	EEE	210	EDO	1	0
9	CCC	201	EDO	1	0
9	FFF	216	EDO	1	0
7	AAA	215	PGE	1	0
9	FFF	218	EDO	3	0
7	BBB	208	PGE	4	0
9	DDD	213	EDO	5	0
8	AAA	207	PEG	4	0
7	EEE	205	PGE	1	0
6	BBB	217	1PE	2	0
7	EEE	219	PGE	6	0
8	AAA	206[A]	PEG	2	0
9	BBB	212	EDO	1	0
4	FFF	203	MPD	7	0
3	AAA	201	PXQ	1	0
4	BBB	204	MPD	1	0
8	AAA	209	PEG	1	0
7	FFF	209	PGE	1	0
4	DDD	203	MPD	1	0
10	DDD	207	PG4	4	0
8	BBB	210	PEG	1	0
3	DDD	202	PXQ	1	0
8	CCC	205[A]	PEG	1	0
8	FFF	213	PEG	1	0
8	EEE	207	PEG	4	0
9	DDD	217[A]	EDO	2	0
8	DDD	211	PEG	1	0
9	EEE	212	EDO	1	0
10	EEE	204	PG4	6	0
9	EEE	217	EDO	1	0
8	DDD	212	PEG	1	0
6	BBB	206	1PE	2	0
10	DDD	208	PG4	6	0
8	FFF	214	PEG	1	0
4	BBB	203	MPD	2	0
5	DDD	206	BCN	5	0

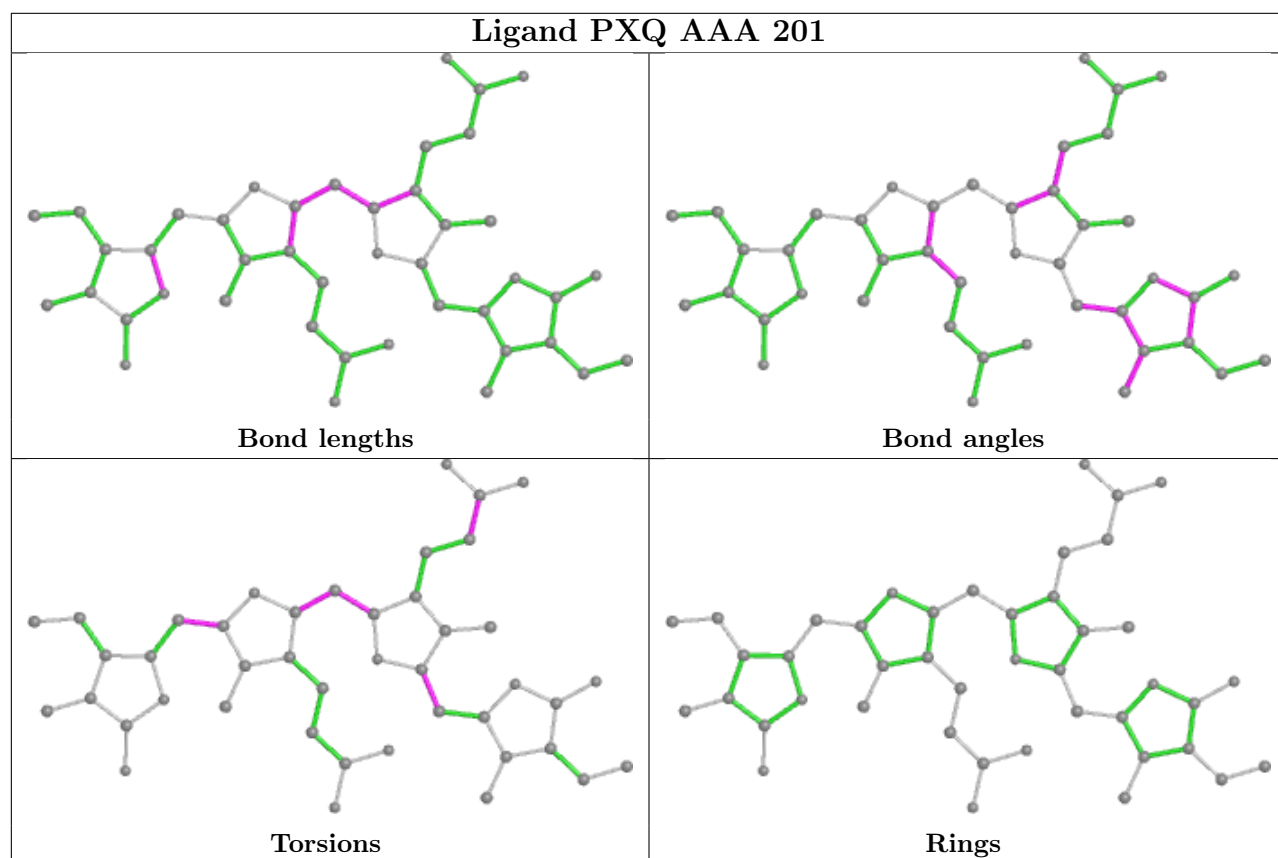
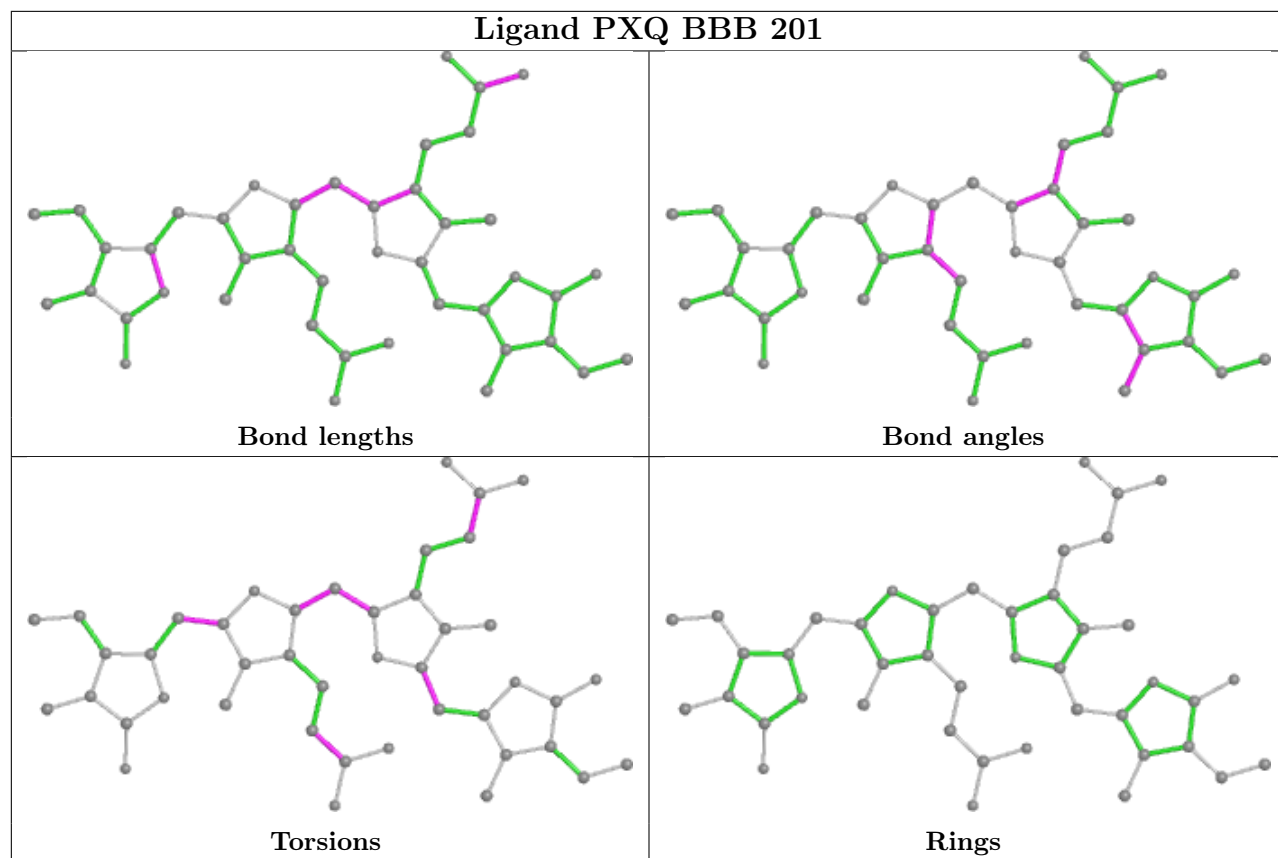
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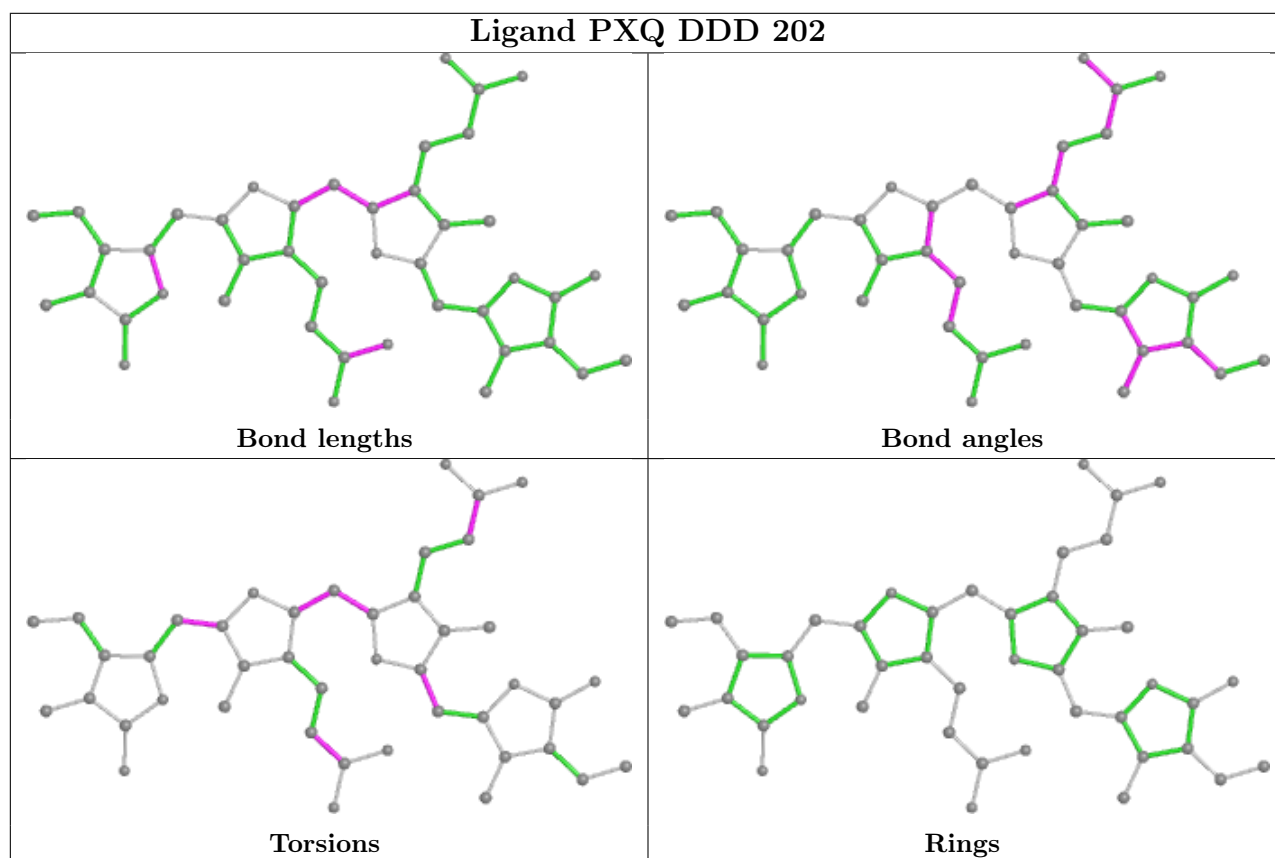
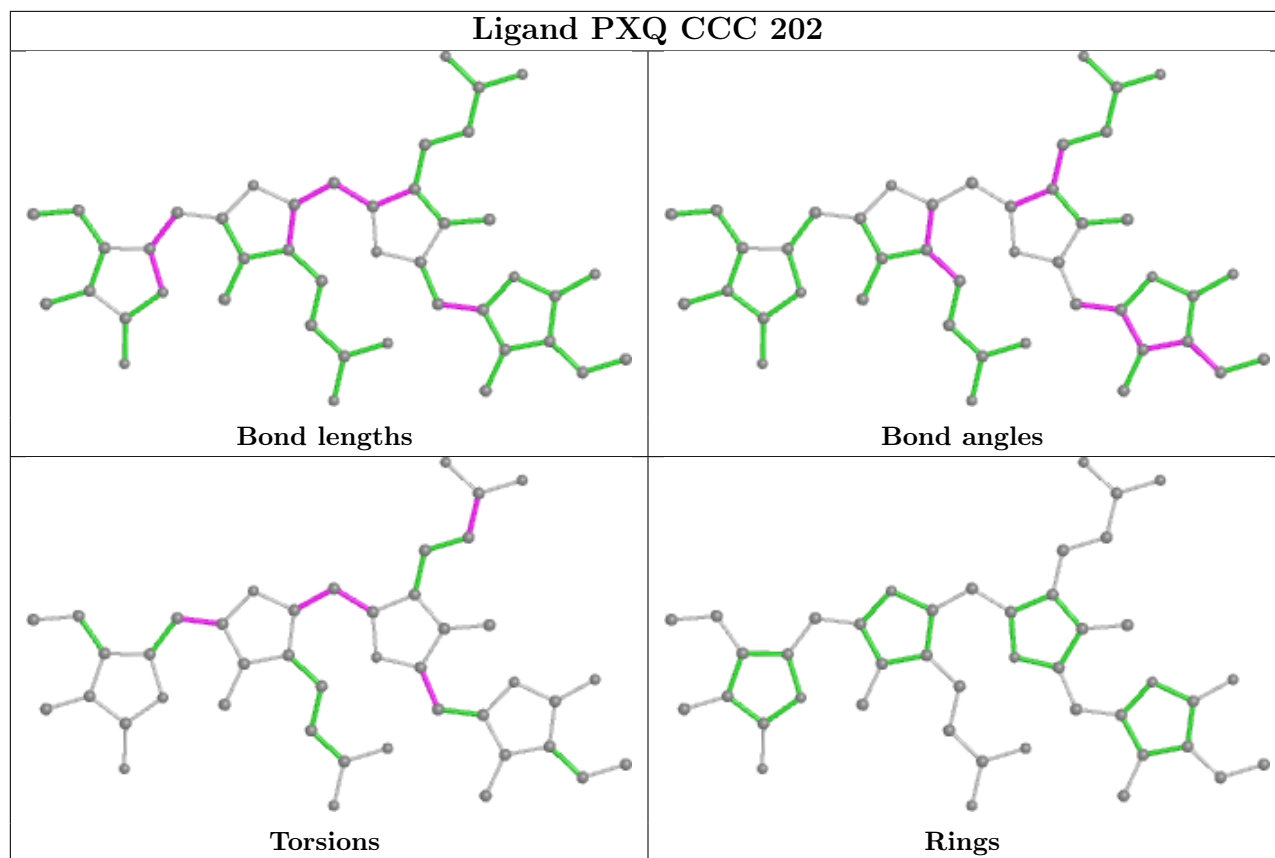
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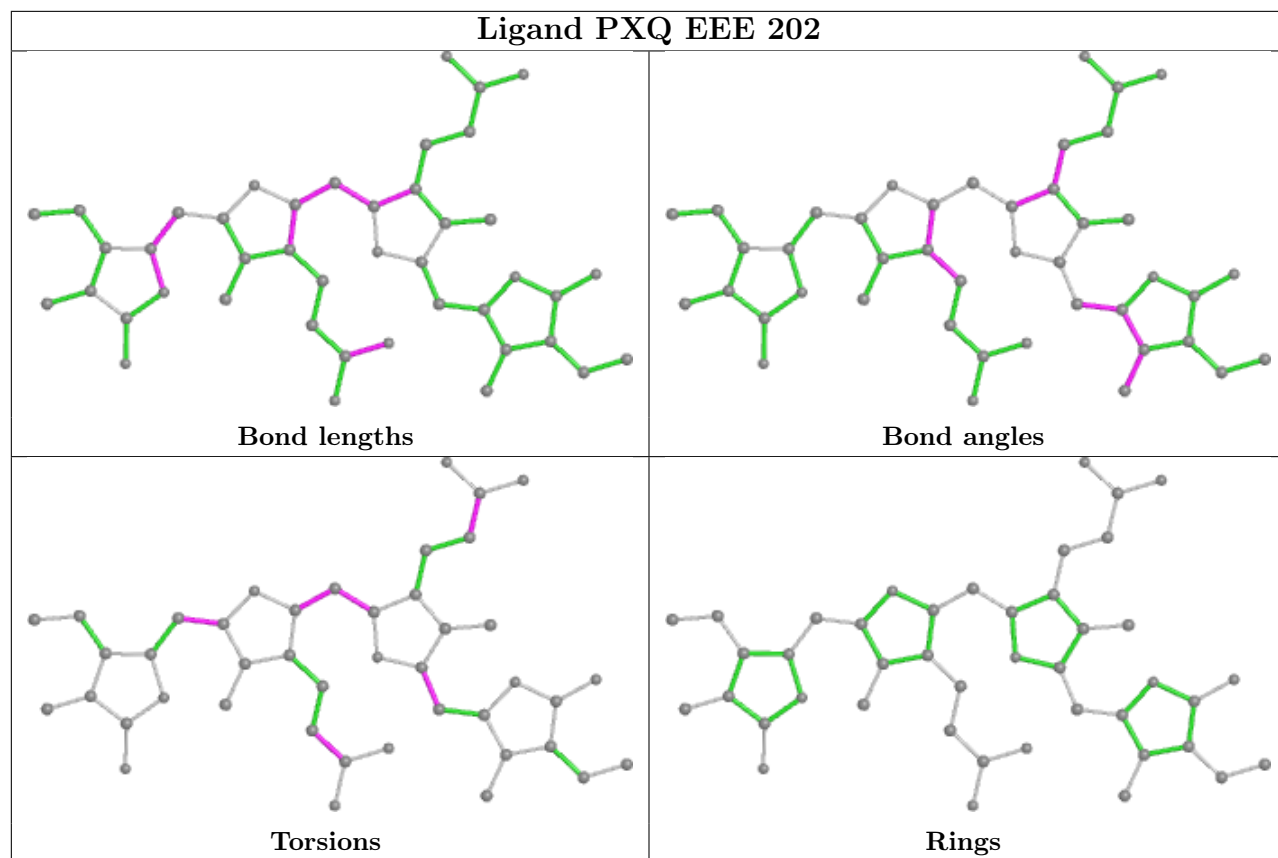
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	AAA	203	BCN	2	0
9	EEE	213	EDO	3	0
4	FFF	204	MPD	1	0
8	FFF	211	PEG	1	0
9	BBB	215	EDO	1	0
7	FFF	207	PGE	3	0
7	FFF	210	PGE	5	0

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	AAA	160/160 (100%)	-0.58	0 100   100	21, 29, 42, 73	0
1	CCC	160/160 (100%)	-0.19	1 (0%) 89   89	23, 38, 60, 70	0
1	EEE	160/160 (100%)	-0.45	0 100   100	21, 31, 47, 58	0
2	BBB	160/161 (99%)	-0.60	0 100   100	22, 28, 38, 58	0
2	DDD	160/161 (99%)	-0.38	0 100   100	21, 33, 50, 70	0
2	FFF	160/161 (99%)	-0.52	0 100   100	25, 30, 43, 64	0
All	All	960/963 (99%)	-0.45	1 (0%) 95   94	21, 31, 49, 73	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	CCC	12	ALA	2.1

### 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
2	MEN	DDD	71	9/10	0.96	0.08	23,25,29,31	0
2	MEN	BBB	71	9/10	0.97	0.06	22,23,24,25	0
2	MEN	FFF	71	9/10	0.97	0.07	25,28,32,34	0

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.



## 6.4 Ligands

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
9	EDO	FFF	217	4/4	0.44	0.37	44,46,47,48	4
7	PGE	FFF	210	10/10	0.47	0.48	42,47,48,48	10
9	EDO	BBB	215	4/4	0.48	0.32	43,46,47,50	4
9	EDO	CCC	201	4/4	0.54	0.32	43,44,46,46	4
9	EDO	DDD	214	4/4	0.56	0.40	46,47,48,48	4
8	PEG	CCC	207	7/7	0.56	0.33	46,47,48,49	7
9	EDO	EEE	211	4/4	0.60	0.41	43,43,43,44	4
8	PEG	BBB	210	7/7	0.60	0.76	47,48,50,50	7
8	PEG	AAA	208	7/7	0.61	0.29	43,44,45,46	7
8	PEG	CCC	206	7/7	0.62	0.34	47,48,50,50	7
9	EDO	DDD	218	4/4	0.62	0.31	40,42,42,42	4
8	PEG	AAA	209	7/7	0.65	0.29	42,43,46,47	7
6	1PE	BBB	217	16/16	0.65	0.51	42,46,46,47	16
9	EDO	DDD	215[B]	4/4	0.66	0.27	45,46,46,47	4
9	EDO	DDD	215[A]	4/4	0.66	0.27	45,45,46,47	4
8	PEG	BBB	209	7/7	0.68	0.25	36,42,47,48	7
8	PEG	DDD	211	7/7	0.68	0.35	46,47,48,49	7
8	PEG	AAA	207	7/7	0.69	0.46	38,39,43,43	7
4	MPD	DDD	204	8/8	0.69	0.49	44,46,47,48	8
8	PEG	EEE	207	7/7	0.69	0.50	41,44,47,47	7
8	PEG	FFF	211	7/7	0.69	0.43	43,43,46,48	7
8	PEG	FFF	214	7/7	0.70	0.41	45,46,47,48	7
9	EDO	CCC	209	4/4	0.70	0.35	43,43,43,45	4
10	PG4	DDD	207	13/13	0.70	0.44	41,42,45,45	13
5	BCN	AAA	203	11/11	0.71	0.28	45,46,51,53	11
10	PG4	DDD	209	13/13	0.71	0.31	43,47,49,49	13
7	PGE	EEE	205	10/10	0.72	0.22	36,43,46,46	10
9	EDO	AAA	216	4/4	0.73	0.34	43,46,46,50	4
7	PGE	CCC	204	10/10	0.73	0.40	45,46,48,49	10
7	PGE	AAA	215	10/10	0.73	0.26	41,54,56,57	10
6	1PE	BBB	205	16/16	0.74	0.40	32,38,50,52	16
8	PEG	CCC	211	7/7	0.74	0.46	43,44,45,46	7
8	PEG	DDD	210	7/7	0.74	0.50	51,51,51,52	7
7	PGE	AAA	205	10/10	0.74	0.30	35,41,46,46	10
8	PEG	DDD	212	7/7	0.74	0.30	39,44,48,48	7
7	PGE	FFF	209	10/10	0.75	0.34	44,46,47,49	10

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
9	EDO	EEE	212	4/4	0.75	0.72	45,45,46,46	4
10	PG4	DDD	208	13/13	0.75	0.35	40,46,49,49	13
9	EDO	EEE	218	4/4	0.75	0.36	43,44,44,45	4
9	EDO	EEE	216	4/4	0.76	0.50	44,44,44,46	4
9	EDO	AAA	210[A]	4/4	0.76	0.38	42,43,43,43	4
9	EDO	AAA	210[B]	4/4	0.76	0.38	57,57,58,58	4
10	PG4	EEE	201	13/13	0.76	0.46	41,44,50,50	13
10	PG4	EEE	204	13/13	0.76	0.21	31,51,72,72	13
11	MRD	EEE	203	8/8	0.76	0.18	42,43,46,48	8
4	MPD	BBB	204	8/8	0.77	0.57	40,41,42,44	8
9	EDO	EEE	217	4/4	0.77	0.26	46,47,48,48	4
8	PEG	BBB	211	7/7	0.77	0.24	42,45,47,48	7
7	PGE	EEE	219	10/10	0.77	0.31	39,47,48,49	10
7	PGE	BBB	207	10/10	0.77	0.32	41,43,44,46	10
9	EDO	DDD	217[A]	4/4	0.77	0.19	39,44,44,44	4
9	EDO	DDD	217[B]	4/4	0.77	0.19	43,44,44,45	4
9	EDO	BBB	212	4/4	0.77	0.46	43,44,45,47	4
4	MPD	AAA	202	8/8	0.77	0.36	37,39,40,41	8
8	PEG	FFF	212	7/7	0.77	0.40	28,36,45,47	7
4	MPD	BBB	203	8/8	0.78	0.35	42,46,48,52	8
4	MPD	FFF	204	8/8	0.78	0.61	45,46,51,52	8
9	EDO	CCC	210	4/4	0.78	0.24	43,44,44,44	4
5	BCN	DDD	206	11/11	0.79	0.21	35,40,53,55	11
9	EDO	EEE	210	4/4	0.80	0.56	46,46,46,46	4
4	MPD	FFF	205	8/8	0.80	0.20	37,44,47,49	8
9	EDO	FFF	215	4/4	0.80	0.25	36,40,40,43	4
8	PEG	CCC	208	7/7	0.80	0.49	44,45,47,48	7
6	1PE	BBB	206	16/16	0.80	0.21	38,44,48,48	16
8	PEG	EEE	208	7/7	0.82	0.48	45,46,47,47	7
8	PEG	EEE	209	7/7	0.82	0.30	41,44,45,48	7
4	MPD	FFF	203	8/8	0.82	0.49	38,40,41,41	8
9	EDO	BBB	214	4/4	0.82	0.72	44,47,47,48	4
9	EDO	EEE	215[B]	4/4	0.83	0.23	33,35,36,37	4
9	EDO	DDD	213	4/4	0.83	0.20	35,41,43,45	4
9	EDO	DDD	216	4/4	0.83	0.21	43,44,44,44	4
9	EDO	BBB	216	4/4	0.83	0.55	42,42,42,43	4
8	PEG	FFF	213	7/7	0.83	0.18	41,43,45,45	7
9	EDO	EEE	215[A]	4/4	0.83	0.23	37,37,38,39	4
7	PGE	BBB	208	10/10	0.84	0.17	38,41,44,45	10
4	MPD	DDD	205	8/8	0.84	0.38	41,43,44,45	8
9	EDO	EEE	213	4/4	0.84	0.21	47,47,49,50	4
9	EDO	DDD	201	4/4	0.84	0.35	40,42,42,44	4

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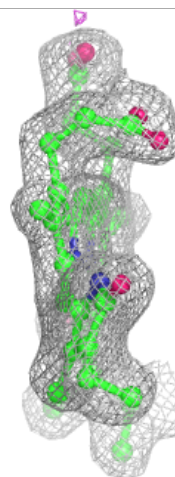
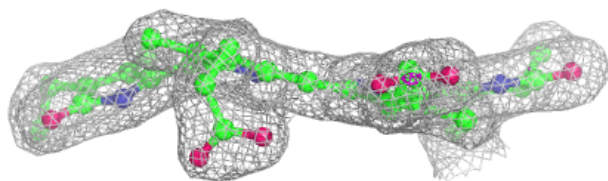
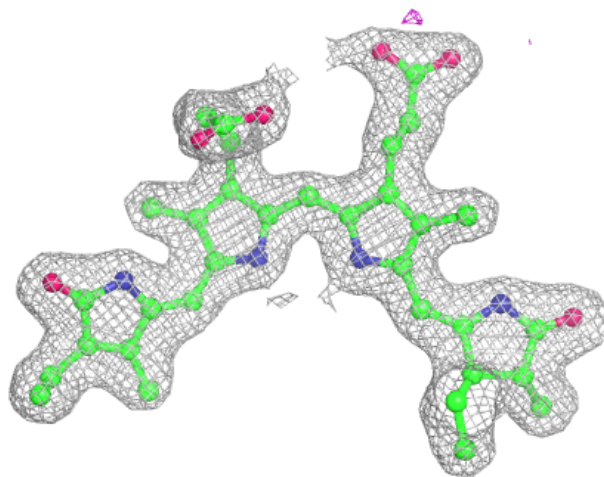
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
9	EDO	FFF	216	4/4	0.84	0.22	39,41,42,43	4
9	EDO	AAA	213	4/4	0.84	0.36	44,45,45,45	4
6	1PE	FFF	206	16/16	0.85	0.21	39,44,49,52	16
7	PGE	FFF	207	10/10	0.85	0.22	41,45,52,53	10
7	PGE	FFF	208	10/10	0.85	0.38	43,44,48,48	10
8	PEG	AAA	206[B]	7/7	0.86	0.16	27,35,42,42	7
8	PEG	AAA	206[A]	7/7	0.86	0.16	45,46,47,48	7
9	EDO	BBB	213	4/4	0.86	0.30	43,43,45,46	4
9	EDO	AAA	214	4/4	0.86	0.25	39,39,40,40	4
8	PEG	EEE	206	7/7	0.87	0.82	47,48,49,50	7
8	PEG	CCC	205[B]	7/7	0.87	0.21	42,43,44,44	7
8	PEG	CCC	205[A]	7/7	0.87	0.21	38,43,46,47	7
9	EDO	AAA	212	4/4	0.88	0.38	41,42,42,43	4
9	EDO	EEE	214	4/4	0.89	0.35	37,38,42,45	4
9	EDO	AAA	211	4/4	0.89	0.16	31,36,40,47	4
4	MPD	DDD	203	8/8	0.89	0.33	39,40,42,43	8
6	1PE	AAA	204	16/16	0.93	0.11	42,49,73,76	0
4	MPD	FFF	202	8/8	0.93	0.16	38,43,47,48	8
9	EDO	FFF	218	4/4	0.94	0.29	38,40,41,41	4
3	PXQ	FFF	201	43/43	0.95	0.08	28,35,41,49	0
3	PXQ	AAA	201	43/43	0.95	0.08	22,25,34,44	0
4	MPD	BBB	202	8/8	0.96	0.10	33,33,39,39	0
3	PXQ	EEE	202	43/43	0.97	0.07	19,22,27,33	0
3	PXQ	BBB	201	43/43	0.97	0.08	23,27,37,43	0
3	PXQ	CCC	202	43/43	0.97	0.07	22,26,30,35	0
3	PXQ	DDD	202	43/43	0.97	0.07	23,28,38,43	0
4	MPD	CCC	203	8/8	0.98	0.07	26,30,30,31	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.

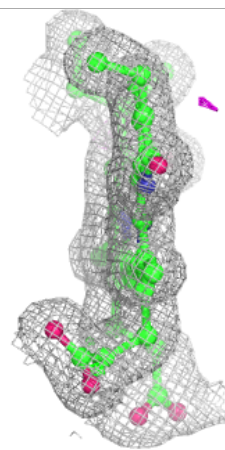
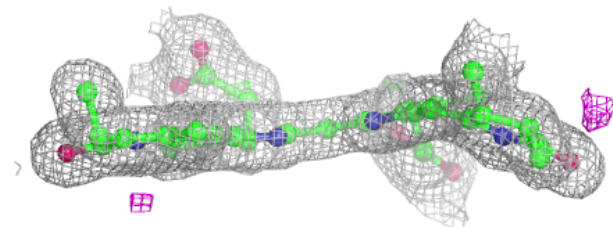
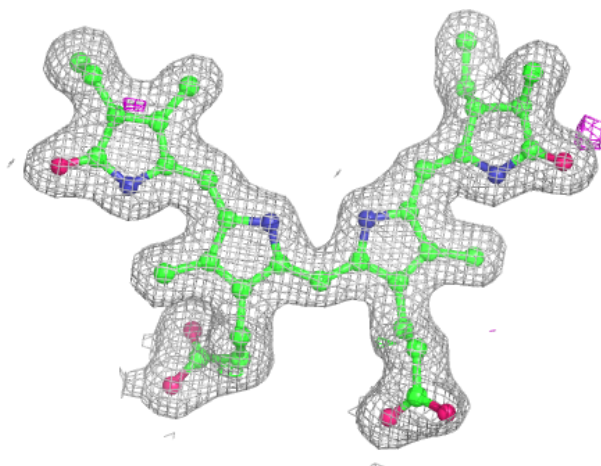
**Electron density around PXQ FFF 201:**

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



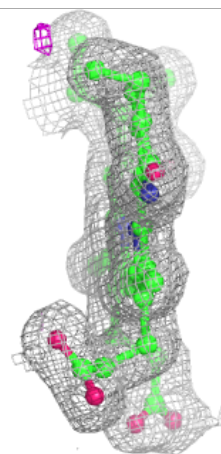
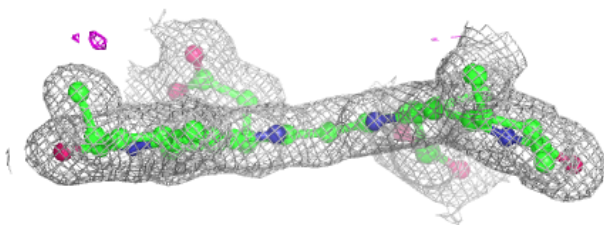
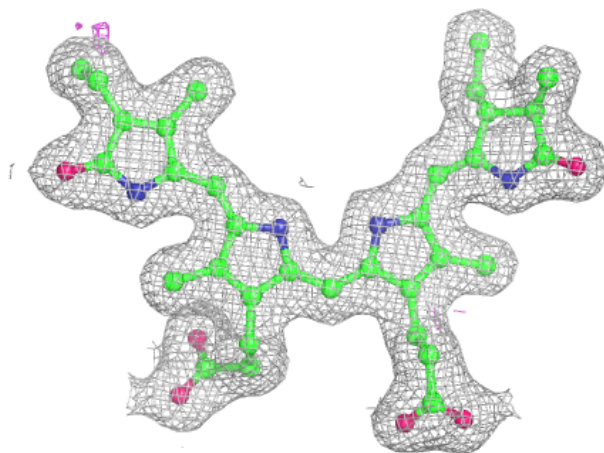
**Electron density around PXQ AAA 201:**

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



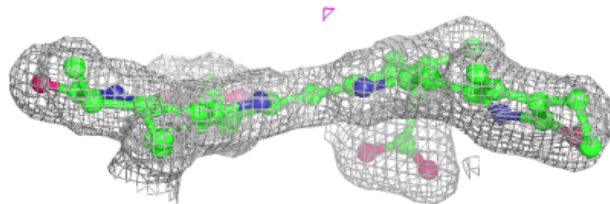
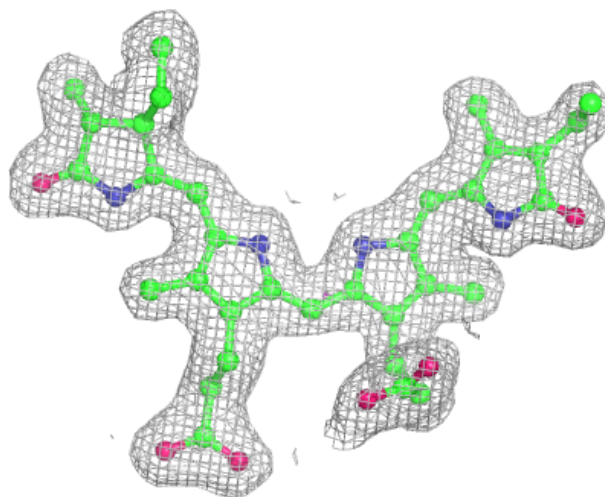
**Electron density around PXQ EEE 202:**

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



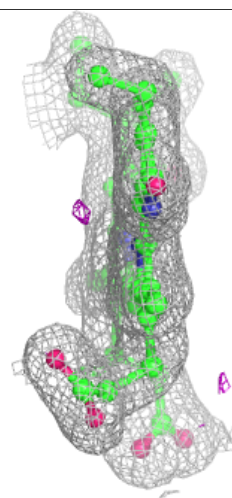
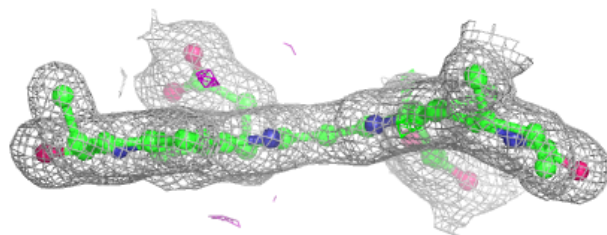
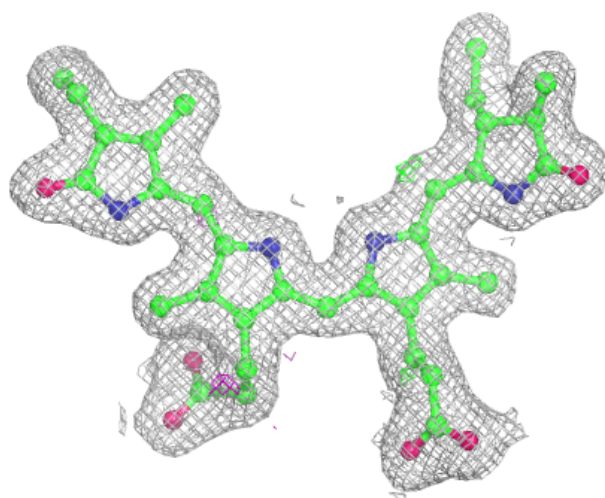
**Electron density around PXQ BBB 201:**

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

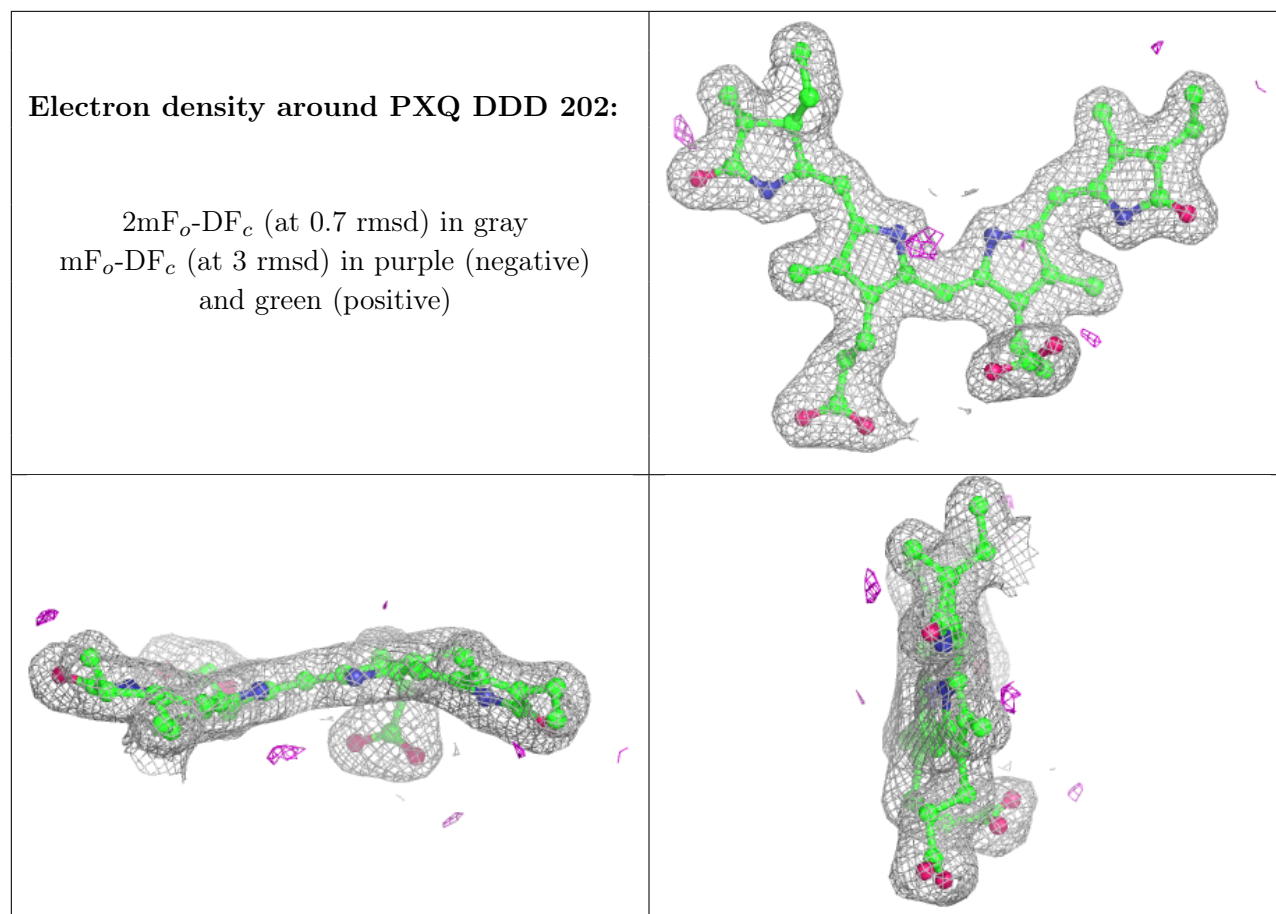


**Electron density around PXQ CCC 202:**

$2mF_o-DF_c$  (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.