



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 01:03 PM UTC

PDB ID : 3HHL / pdb_00003hhl
Title : Crystal structure of methylated RPA0582 protein
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Deposited on : 2009-05-15
Resolution : 2.65 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

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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

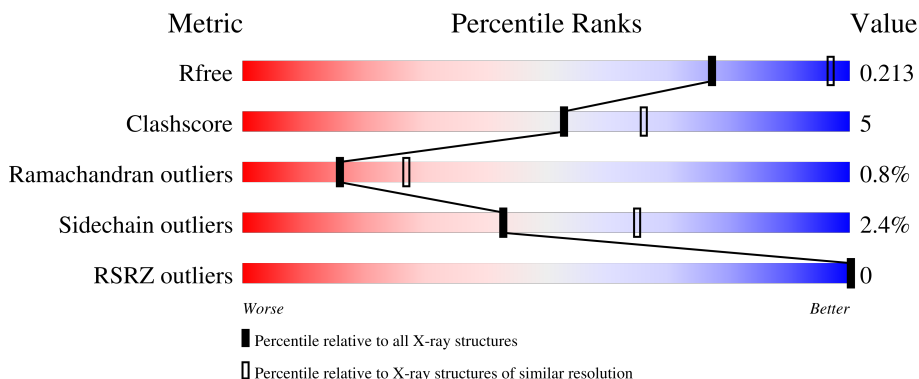
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

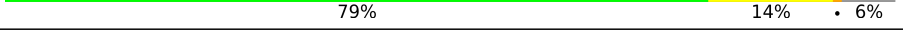
The reported resolution of this entry is 2.65 Å.

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}	180053	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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 ≥ 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	143	 85% 8% • 6%
2	B	143	 79% 14% • 6%
3	C	143	 82% 10% • 6%
4	D	143	 79% 14% • 6%

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 4944 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 RPA0582.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	135	Total	C	N	O	Se	0	2	0
			1098	696	203	195	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q6NC90
A	61	SER	PHE	SEE REMARK 999	UNP Q6NC90

- Molecule 2 is a protein called RPA0582.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	135	Total	C	N	O	Se	0	3	0
			1103	698	208	194	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	GLY	-	expression tag	UNP Q6NC90
B	61	SER	PHE	SEE REMARK 999	UNP Q6NC90

- Molecule 3 is a protein called RPA0582.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	135	Total	C	N	O	Se	0	5	0
			1119	710	210	195	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	GLY	-	expression tag	UNP Q6NC90
C	61	SER	PHE	SEE REMARK 999	UNP Q6NC90

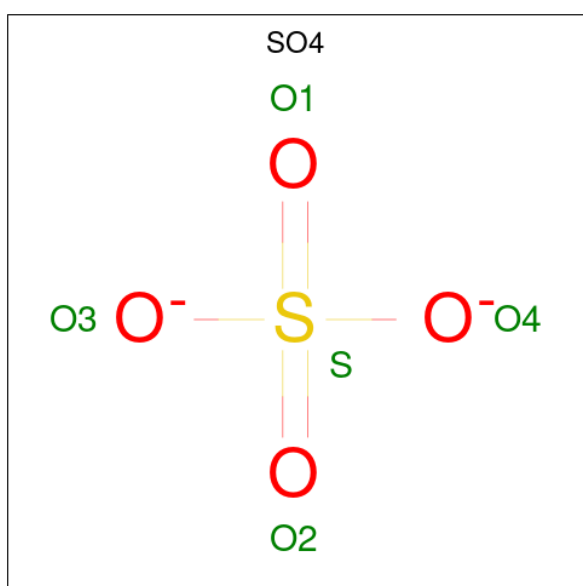
- Molecule 4 is a protein called RPA0582.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	135	Total	C	N	O	Se	0	3	0
			1103	697	205	198	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	0	GLY	-	expression tag	UNP Q6NC90
D	61	SER	PHE	SEE REMARK 999	UNP Q6NC90

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

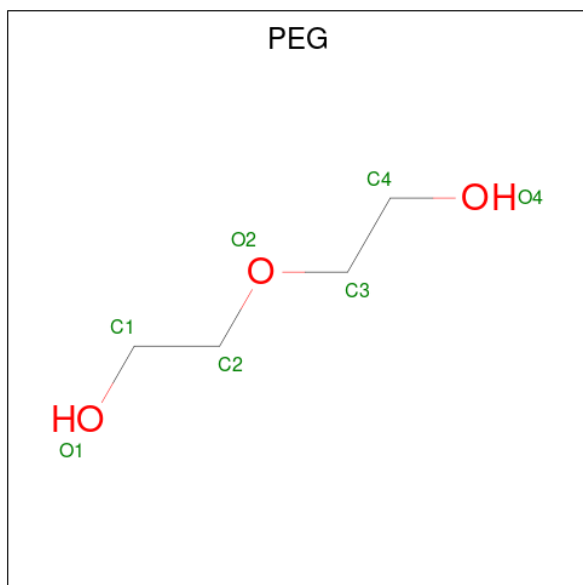
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Cl	0	0
			1	1		

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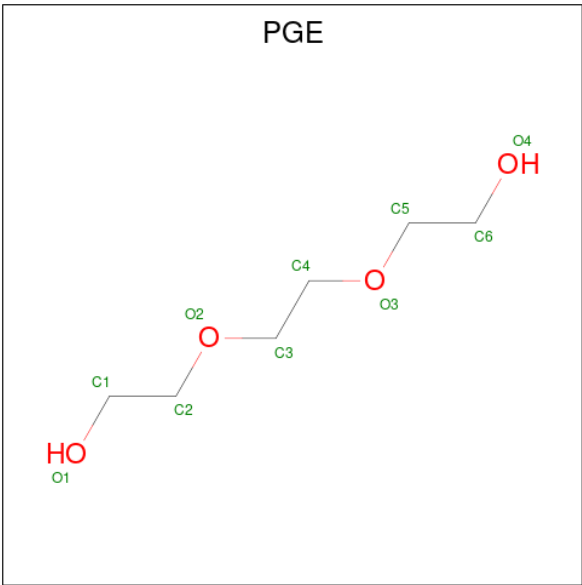
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Cl	0	0
			1	1		
6	C	1	Total	Cl	0	0
			1	1		
6	D	1	Total	Cl	0	0
			1	1		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



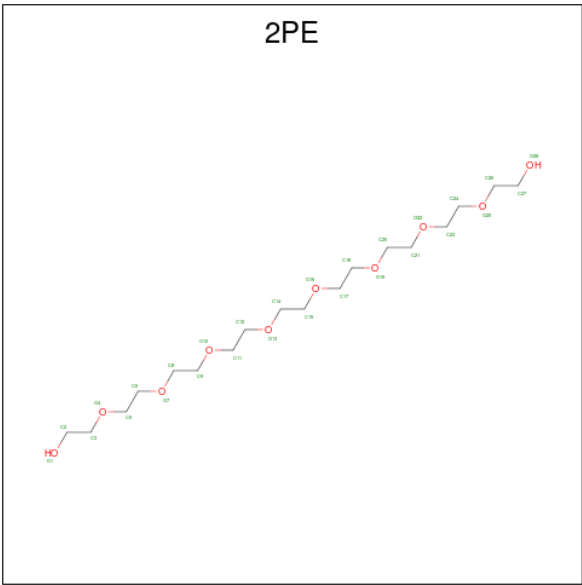
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			7	4	3		
7	A	1	Total	C	O	0	0
			7	4	3		
7	B	1	Total	C	O	0	0
			7	4	3		
7	B	1	Total	C	O	0	0
			7	4	3		
7	B	1	Total	C	O	0	0
			7	4	3		
7	C	1	Total	C	O	0	0
			7	4	3		
7	D	1	Total	C	O	0	0
			7	4	3		

- Molecule 8 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			10	6	4		
8	B	1	Total	C	O	0	0
			10	6	4		
8	C	1	Total	C	O	0	0
			10	6	4		

- Molecule 9 is NONAETHYLENE GLYCOL (CCD ID: 2PE) (formula: C₁₈H₃₈O₁₀).



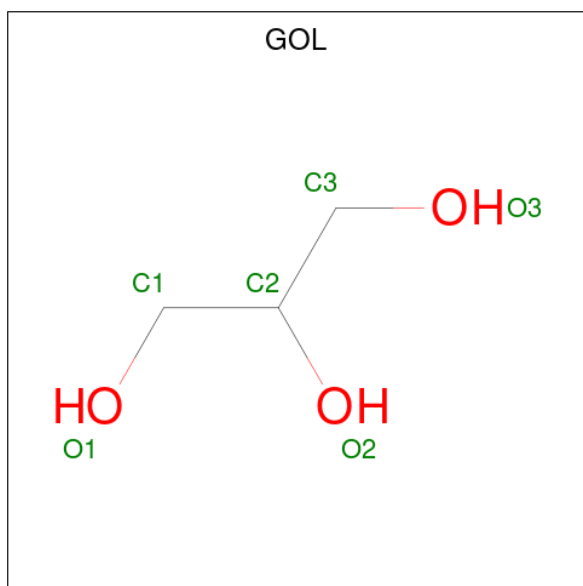
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			28	18	10		

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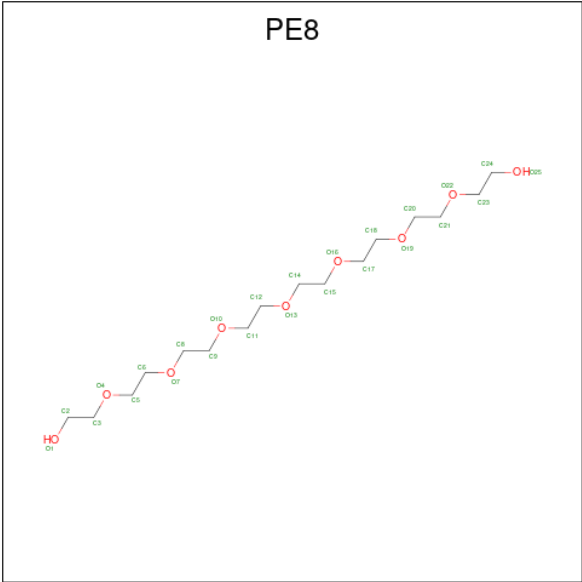
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	D	1	Total	C	O	0	0
			28	18	10		

- Molecule 10 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



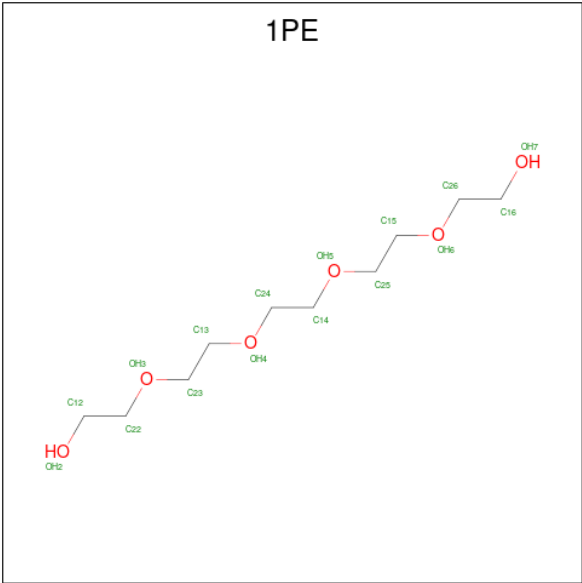
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			6	3	3		
10	B	1	Total	C	O	0	0
			6	3	3		
10	C	1	Total	C	O	0	0
			6	3	3		
10	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 11 is 3,6,9,12,15,18,21-HEPTAOXATRICOSANE-1,23-DIOL (CCD ID: PE8) (formula: $C_{16}H_{34}O_9$).



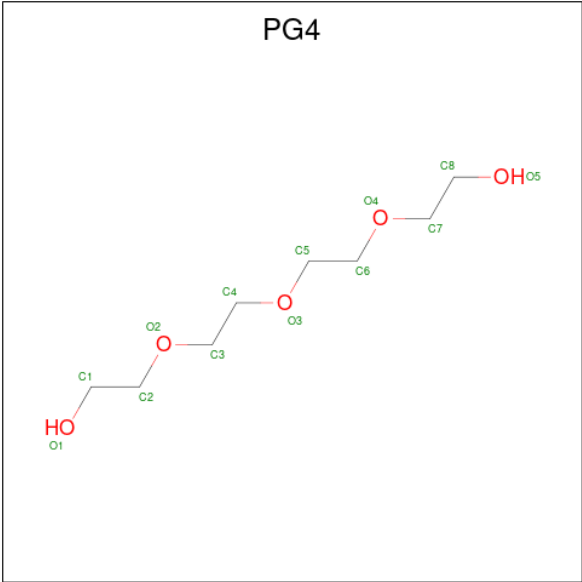
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	B	1	Total	C	O	0	0
			25	16	9		
11	C	1	Total	C	O	0	0
			25	16	9		

- Molecule 12 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: C₁₀H₂₂O₆).



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

- Molecule 13 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



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

- Molecule 14 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	A	57	Total	O	0	0
			57	57		
14	B	66	Total	O	0	0
			66	66		
14	C	70	Total	O	0	0
			70	70		
14	D	58	Total	O	0	0
			58	58		

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: RPA0582

Chain A: 



- Molecule 2: RPA0582

Chain B: 



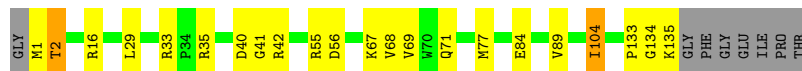
- Molecule 3: RPA0582

Chain C: 



- Molecule 4: RPA0582

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	193.01Å 193.01Å 117.97Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	32.17 – 2.65 32.17 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.6 (32.17-2.65) 99.2 (32.17-2.65)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 2.64Å)	Xtriage
Refinement program	REFMAC 5.5.0089	Depositor
R, R_{free}	0.178 , 0.206 0.186 , 0.213	Depositor DCC
R_{free} test set	2400 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	53.8	Xtriage
Anisotropy	0.011	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 31.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.015 for -2/3*h-1/3*k-4/3*l,-1/3*h-2/3*k+4/3*l,-1/3*h+1/3*k+1/3*l 0.012 for -h,1/3*h-1/3*k-4/3*l,-1/3*h-2/3*k+1/3*l 0.014 for -1/3*h+1/3*k+4/3*l,-k,2/3*h+1/3*k+1/3*l 0.470 for -h,2/3*h+1/3*k+4/3*l,1/3*h+2/3*k-1/3*l 0.477 for -1/3*h-2/3*k+4/3*l,-2/3*h-1/3*k-4/3*l,1/3*h-1/3*k-1/3*l 0.470 for 1/3*h+2/3*k-4/3*l,-k,-2/3*h-1/3*k-1/3*l 0.016 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4944	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² 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: PGE, 2PE, SO4, PEG, PG4, PE8, 1PE, GOL, MLZ, MLY, CL

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	1.02	0/1090	0.92	2/1471 (0.1%)
2	B	1.02	0/1112	0.90	1/1500 (0.1%)
3	C	0.98	0/1109	0.90	0/1494
4	D	1.01	2/1105 (0.2%)	0.91	1/1494 (0.1%)
All	All	1.01	2/4416 (0.0%)	0.91	4/5959 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	69	VAL	C-O	-5.59	1.18	1.24
4	D	2	THR	CA-CB	5.21	1.60	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	40	ASP	N-CA-C	-5.25	103.22	110.35
1	A	38	TYR	CA-C-N	5.24	124.68	119.24
1	A	38	TYR	C-N-CA	5.24	124.68	119.24
2	B	116	ARG	N-CA-C	-5.08	105.74	111.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1098	0	1077	9	0
2	B	1103	0	1089	20	0
3	C	1119	0	1113	9	0
4	D	1103	0	1072	13	0
5	A	10	0	0	0	0
5	C	5	0	0	0	0
6	A	1	0	0	0	0
6	B	1	0	0	1	0
6	C	1	0	0	1	0
6	D	1	0	0	0	0
7	A	14	0	20	0	0
7	B	21	0	30	0	0
7	C	7	0	10	0	0
7	D	7	0	10	0	0
8	A	10	0	14	0	0
8	B	10	0	14	0	0
8	C	10	0	14	0	0
9	A	28	0	38	1	0
9	D	28	0	38	1	0
10	A	6	0	8	0	0
10	B	6	0	8	0	0
10	C	6	0	8	0	0
10	D	6	0	8	0	0
11	B	25	0	34	2	0
11	C	25	0	34	1	0
12	B	16	0	22	0	0
13	C	13	0	18	0	0
13	D	13	0	18	0	0
14	A	57	0	0	1	0
14	B	66	0	0	0	0
14	C	70	0	0	1	0
14	D	58	0	0	1	0
All	All	4944	0	4697	47	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 5.

The worst 5 of 47 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:B:16[B]:ARG:HG2	2:B:16[B]:ARG:HH11	1.09	1.15
4:D:16[B]:ARG:HH11	4:D:16[B]:ARG:HG2	1.20	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:16[B]:ARG:HG2	2:B:16[B]:ARG:NH1	1.90	0.85
4:D:16[B]:ARG:HG2	4:D:16[B]:ARG:NH1	1.95	0.76
2:B:16[B]:ARG:HH11	2:B:16[B]:ARG:CG	1.95	0.75

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	133/143 (93%)	129 (97%)	3 (2%)	1 (1%)	16	27
2	B	135/143 (94%)	131 (97%)	3 (2%)	1 (1%)	18	30
3	C	135/143 (94%)	130 (96%)	4 (3%)	1 (1%)	18	30
4	D	135/143 (94%)	130 (96%)	4 (3%)	1 (1%)	18	30
All	All	538/572 (94%)	520 (97%)	14 (3%)	4 (1%)	16	30

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	133	PRO
3	C	133	PRO
4	D	133	PRO
1	A	133	PRO

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	107/108 (99%)	105 (98%)	2 (2%)	50	71
2	B	109/109 (100%)	107 (98%)	2 (2%)	51	72
3	C	109/107 (102%)	105 (96%)	4 (4%)	30	50
4	D	108/109 (99%)	106 (98%)	2 (2%)	50	71
All	All	433/433 (100%)	423 (98%)	10 (2%)	43	66

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	104	ILE
4	D	2	THR
4	D	104	ILE
2	B	104	ILE
3	C	1	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	HIS
4	D	4	HIS
4	D	71	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

11 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
3	MLZ	C	114	3	8,9,10	0.96	0	4,9,11	0.60	0
4	MLY	D	67	4	9,10,11	0.67	0	6,11,13	1.09	1 (16%)
1	MLZ	A	129	1	8,9,10	1.68	3 (37%)	4,9,11	1.14	0
1	MLY	A	67	1	9,10,11	0.61	0	6,11,13	0.53	0
3	MLY	C	67	3	9,10,11	0.65	0	6,11,13	0.39	0
1	MLY	A	135	1	9,10,11	1.13	1 (11%)	6,11,13	0.36	0
3	MLZ	C	129	3	8,9,10	1.07	1 (12%)	4,9,11	0.58	0
3	MLY	C	135	3	9,10,11	1.05	1 (11%)	6,11,13	0.74	0
4	MLY	D	135	4	9,10,11	0.88	0	6,11,13	0.42	0
2	MLY	B	135	2	9,10,11	0.53	0	6,11,13	0.43	0
2	MLZ	B	67	2	8,9,10	1.34	1 (12%)	4,9,11	0.68	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MLZ	C	114	3	-	0/7/8/10	-
4	MLY	D	67	4	-	2/8/9/11	-
1	MLZ	A	129	1	-	3/7/8/10	-
1	MLY	A	67	1	-	3/8/9/11	-
3	MLY	C	67	3	-	2/8/9/11	-
1	MLY	A	135	1	-	1/8/9/11	-
3	MLZ	C	129	3	-	3/7/8/10	-
3	MLY	C	135	3	-	4/8/9/11	-
4	MLY	D	135	4	-	2/8/9/11	-
2	MLY	B	135	2	-	4/8/9/11	-
2	MLZ	B	67	2	-	1/7/8/10	-

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	129	MLZ	CD-CE	-2.83	1.40	1.51
2	B	67	MLZ	CB-CA	-2.47	1.49	1.53
1	A	129	MLZ	CB-CA	-2.37	1.50	1.53
1	A	135	MLY	CB-CA	2.30	1.57	1.53
3	C	135	MLY	CB-CA	2.19	1.56	1.53

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	67	MLY	CD-CE-NZ	2.48	120.12	113.71

There are no chirality outliers.

5 of 25 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	129	MLZ	CD-CE-NZ-CM
2	B	67	MLZ	CD-CE-NZ-CM
2	B	135	MLY	O-C-CA-CB
3	C	135	MLY	C-CA-CB-CG
4	D	67	MLY	O-C-CA-CB

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	135	MLY	1	0
3	C	135	MLY	1	0
4	D	135	MLY	1	0
2	B	135	MLY	3	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 28 ligands modelled in this entry, 4 are monoatomic - leaving 24 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
13	PG4	D	144	-	12,12,12	0.71	0	11,11,11	0.55	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	PGE	A	147	-	9,9,9	0.67	0	8,8,8	0.41	0
11	PE8	B	148	-	24,24,24	0.70	0	23,23,23	0.54	0
5	SO4	C	143	-	4,4,4	0.25	0	6,6,6	0.21	0
13	PG4	C	145	-	12,12,12	0.60	0	11,11,11	0.34	0
9	2PE	D	146	-	27,27,27	0.61	0	26,26,26	0.43	0
7	PEG	B	146	-	6,6,6	0.41	0	5,5,5	0.44	0
7	PEG	B	144	-	6,6,6	0.54	0	5,5,5	0.23	0
10	GOL	D	147	-	5,5,5	0.27	0	5,5,5	0.35	0
8	PGE	B	147	-	9,9,9	0.59	0	8,8,8	0.67	0
11	PE8	C	148	-	24,24,24	0.75	0	23,23,23	0.72	1 (4%)
10	GOL	C	149	-	5,5,5	0.31	0	5,5,5	0.40	0
12	1PE	B	150	-	15,15,15	0.60	0	14,14,14	0.39	0
10	GOL	B	149	-	5,5,5	0.52	0	5,5,5	0.75	0
8	PGE	C	147	-	9,9,9	0.61	0	8,8,8	0.52	0
10	GOL	A	150	-	5,5,5	0.34	0	5,5,5	0.16	0
7	PEG	B	145	-	6,6,6	0.47	0	5,5,5	0.41	0
7	PEG	C	146	-	6,6,6	0.49	0	5,5,5	0.19	0
5	SO4	A	144	-	4,4,4	0.25	0	6,6,6	0.13	0
7	PEG	A	149	-	6,6,6	0.41	0	5,5,5	0.41	0
7	PEG	D	145	-	6,6,6	0.51	0	5,5,5	0.23	0
7	PEG	A	146	-	6,6,6	0.47	0	5,5,5	0.23	0
5	SO4	A	143	-	4,4,4	0.30	0	6,6,6	0.28	0
9	2PE	A	148	-	27,27,27	0.66	0	26,26,26	0.48	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
13	PG4	D	144	-	-	6/10/10/10	-
8	PGE	A	147	-	-	3/7/7/7	-
11	PE8	B	148	-	-	12/22/22/22	-
13	PG4	C	145	-	-	2/10/10/10	-
9	2PE	D	146	-	-	13/25/25/25	-
7	PEG	B	146	-	-	3/4/4/4	-
7	PEG	B	144	-	-	2/4/4/4	-
10	GOL	D	147	-	-	0/4/4/4	-
8	PGE	B	147	-	-	4/7/7/7	-
11	PE8	C	148	-	-	13/22/22/22	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	GOL	C	149	-	-	0/4/4/4	-
12	1PE	B	150	-	-	8/13/13/13	-
10	GOL	B	149	-	-	3/4/4/4	-
8	PGE	C	147	-	-	2/7/7/7	-
10	GOL	A	150	-	-	4/4/4/4	-
7	PEG	B	145	-	-	3/4/4/4	-
7	PEG	C	146	-	-	2/4/4/4	-
7	PEG	A	149	-	-	3/4/4/4	-
7	PEG	D	145	-	-	2/4/4/4	-
7	PEG	A	146	-	-	3/4/4/4	-
9	2PE	A	148	-	-	13/25/25/25	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	C	148	PE8	O19-C18-C17	2.16	120.21	110.35

There are no chirality outliers.

5 of 101 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	A	150	GOL	C1-C2-C3-O3
10	B	149	GOL	C1-C2-C3-O3
11	B	148	PE8	O4-C5-C6-O7
11	C	148	PE8	O7-C8-C9-O10
9	A	148	2PE	O13-C14-C15-O16

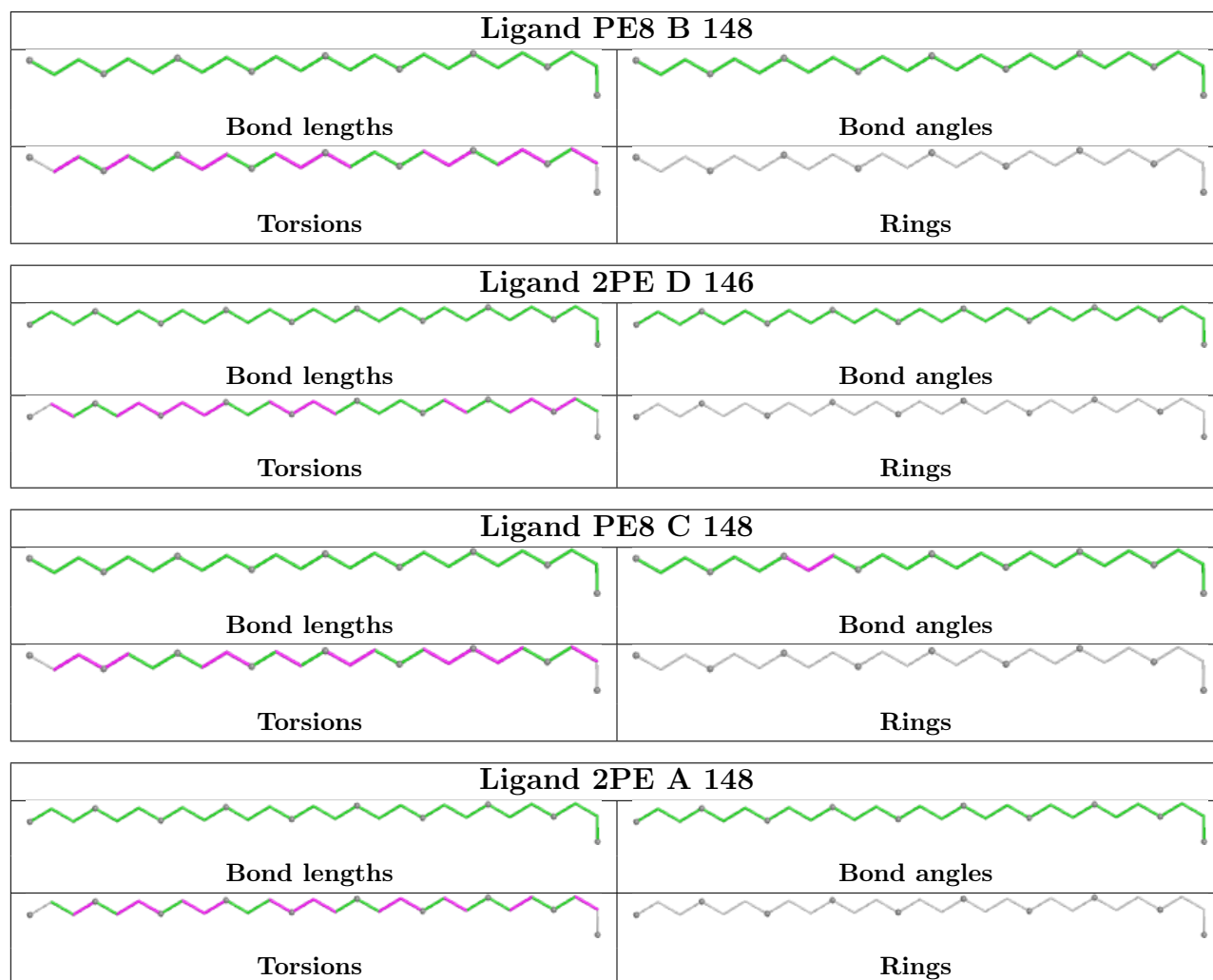
There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	B	148	PE8	2	0
9	D	146	2PE	1	0
11	C	148	PE8	1	0
9	A	148	2PE	1	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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, 95th 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 ²)	Q<0.9
1	A	128/143 (89%)	-1.49	0 100 100	16, 37, 55, 79	2 (1%)
2	B	129/143 (90%)	-1.51	0 100 100	17, 36, 55, 79	3 (2%)
3	C	127/143 (88%)	-1.49	0 100 100	18, 36, 53, 77	5 (3%)
4	D	129/143 (90%)	-1.49	0 100 100	15, 37, 55, 79	3 (2%)
All	All	513/572 (89%)	-1.49	0 100 100	15, 37, 55, 79	13 (2%)

There are no RSRZ outliers to report.

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th 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 ²)	Q<0.9
1	MLY	A	135	11/12	0.95	0.20	74,80,82,82	0
2	MLY	B	135	11/12	0.95	0.16	74,78,81,81	0
3	MLY	C	135	11/12	0.95	0.22	73,77,79,79	0
4	MLY	D	135	11/12	0.97	0.18	72,79,82,82	0
1	MLY	A	67	11/12	0.98	0.07	37,38,69,69	0
3	MLY	C	67	11/12	0.98	0.06	36,38,62,64	0
3	MLZ	C	114	10/11	0.99	0.03	35,37,47,50	0
3	MLZ	C	129	10/11	0.99	0.05	33,34,59,61	0
1	MLZ	A	129	10/11	0.99	0.05	33,34,62,62	0
4	MLY	D	67	11/12	0.99	0.07	38,39,67,68	0
2	MLZ	B	67	10/11	0.99	0.05	38,38,60,60	0

6.3 Carbohydrates ⓘ

There are no oligosaccharides 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, 95th 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.

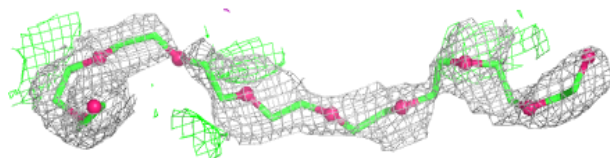
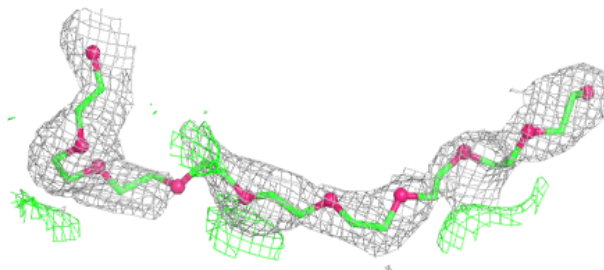
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	PEG	B	144	7/7	0.97	0.11	127,128,129,130	0
7	PEG	D	145	7/7	0.97	0.09	110,111,111,112	0
8	PGE	B	147	10/10	0.97	0.10	89,96,100,100	0
8	PGE	C	147	10/10	0.97	0.09	88,89,92,92	0
13	PG4	C	145	13/13	0.97	0.12	110,114,118,118	0
7	PEG	A	146	7/7	0.98	0.07	77,78,81,82	0
8	PGE	A	147	10/10	0.98	0.08	84,85,87,88	0
7	PEG	A	149	7/7	0.98	0.07	105,107,108,109	0
5	SO4	A	144	5/5	0.98	0.07	125,125,125,126	0
9	2PE	D	146	28/28	0.98	0.09	99,105,116,117	0
10	GOL	B	149	6/6	0.98	0.07	73,76,80,82	0
12	1PE	B	150	16/16	0.98	0.10	108,111,116,116	0
7	PEG	C	146	7/7	0.98	0.07	76,77,79,80	0
13	PG4	D	144	13/13	0.98	0.11	110,113,116,116	0
6	CL	A	145	1/1	0.99	0.06	54,54,54,54	0
10	GOL	A	150	6/6	0.99	0.08	77,79,80,80	0
7	PEG	B	145	7/7	0.99	0.07	76,76,78,78	0
10	GOL	C	149	6/6	0.99	0.07	77,78,79,79	0
10	GOL	D	147	6/6	0.99	0.06	82,86,87,88	0
11	PE8	B	148	25/25	0.99	0.09	93,100,107,108	0
11	PE8	C	148	25/25	0.99	0.09	97,104,111,112	0
7	PEG	B	146	7/7	0.99	0.06	80,81,84,85	0
6	CL	B	143	1/1	0.99	0.05	54,54,54,54	0
9	2PE	A	148	28/28	0.99	0.09	102,107,116,116	0
5	SO4	C	143	5/5	1.00	0.03	51,52,53,53	5
6	CL	C	144	1/1	1.00	0.03	55,55,55,55	0
6	CL	D	143	1/1	1.00	0.06	54,54,54,54	0
5	SO4	A	143	5/5	1.00	0.03	55,57,57,58	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.

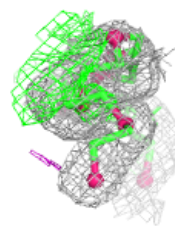
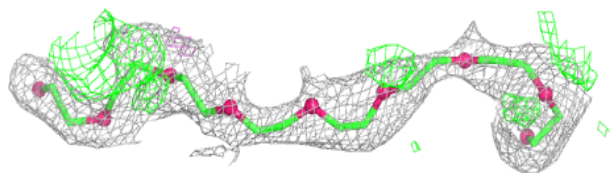
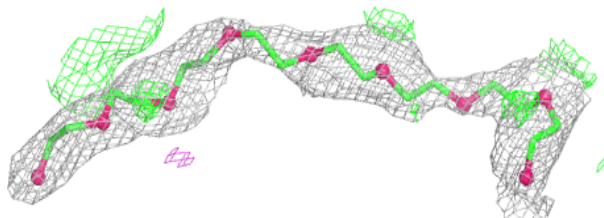
Electron density around 2PE D 146:

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



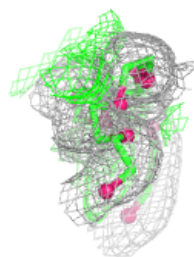
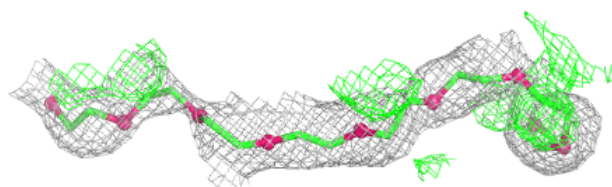
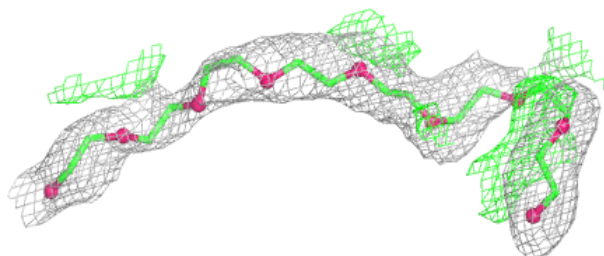
Electron density around PE8 B 148:

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

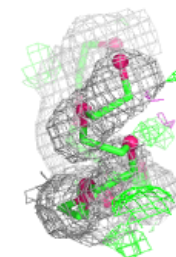
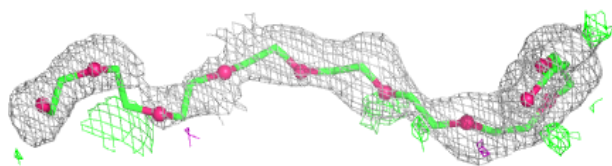
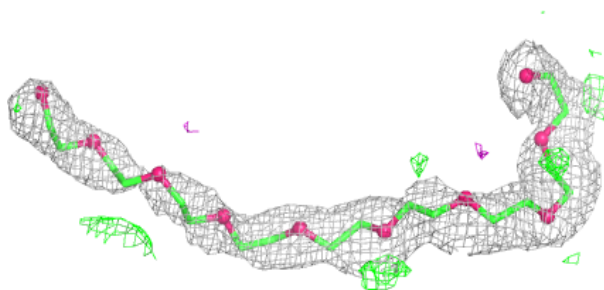


Electron density around PE8 C 148:

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

**Electron density around 2PE A 148:**

$2mF_o-DF_c$ (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.