



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 10:33 AM UTC

PDB ID : 6UXE / pdb_00006uxe
Title : Structure of the human mitochondrial desulfurase complex Nfs1-ISCU2(M140I)-ISD11 with E.coli ACP1 at 1.57 Å resolution showing flexibility of N terminal end of ISCU2
Authors : Boniecki, M.T.; Cygler, M.
Deposited on : 2019-11-07
Resolution : 1.57 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

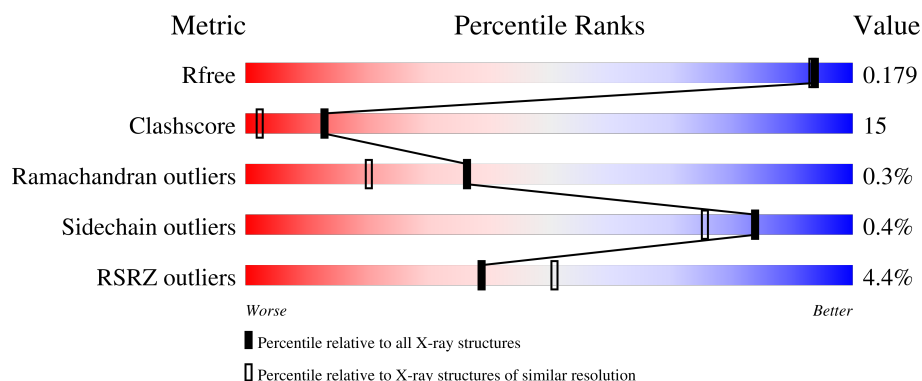
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.57 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	1094 (1.58-1.58)
Clashscore	190562	1105 (1.58-1.58)
Ramachandran outliers	187476	1082 (1.58-1.58)
Sidechain outliers	187428	1081 (1.58-1.58)
RSRZ outliers	180081	1094 (1.58-1.58)

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	406	3% 72% 23% • •
2	B	91	5% 79% 12% • 7%
3	C	77	10% 90% 5% • •
4	D	143	4% 73% 15% 12%

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	P15	A	526	-	-	X	-
11	DTT	A	533	-	-	X	-
12	PGE	A	534	-	-	X	-
12	PGE	A	543	-	-	X	-
7	EDO	A	507	-	-	X	-
7	EDO	A	511	-	-	X	-
7	EDO	A	514	-	-	X	-
7	EDO	A	554	-	-	X	-
7	EDO	B	107	-	-	X	-
8	PEG	A	531	-	-	X	-
8	PEG	B	104	-	-	X	-
8	PEG	B	106	-	-	X	-

2 Entry composition [i](#)

There are 17 unique types of molecules in this entry. The entry contains 6265 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 Cysteine desulfurase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	395	Total	C	N	O	S	0	22	0
			3141	1993	539	587	22			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	52	MET	-	initiating methionine	UNP Q9Y697
A	53	GLY	-	expression tag	UNP Q9Y697
A	54	SER	-	expression tag	UNP Q9Y697
A	55	SER	-	expression tag	UNP Q9Y697

- Molecule 2 is a protein called LYR motif-containing protein 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	85	Total	C	N	O	S	0	8	0
			728	461	142	124	1			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	11	ALA	SER	variant	UNP Q9HD34

- Molecule 3 is a protein called Acyl carrier protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	75	Total	C	N	O	S	0	0	0
			535	335	84	115	1			

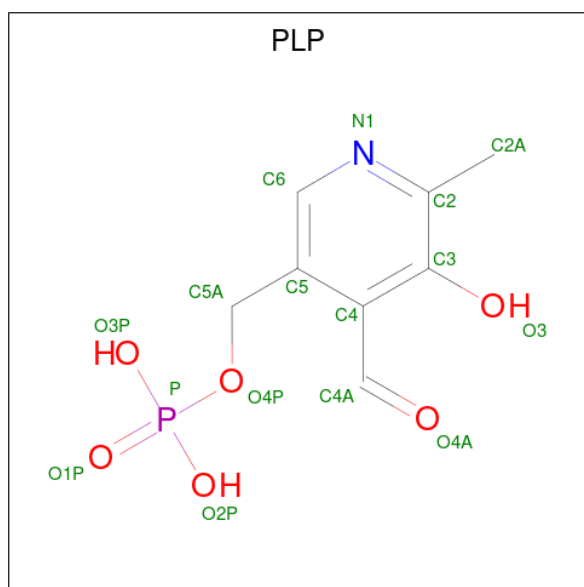
- Molecule 4 is a protein called Iron-sulfur cluster assembly enzyme ISCU, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	126	Total	C	N	O	S	0	7	0
			977	621	164	186	6			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	33	MET	-	initiating methionine	UNP Q9H1K1
D	34	ALA	-	expression tag	UNP Q9H1K1
D	140	ILE	MET	engineered mutation	UNP Q9H1K1
D	168	LEU	-	expression tag	UNP Q9H1K1
D	169	GLU	-	expression tag	UNP Q9H1K1
D	170	HIS	-	expression tag	UNP Q9H1K1
D	171	HIS	-	expression tag	UNP Q9H1K1
D	172	HIS	-	expression tag	UNP Q9H1K1
D	173	HIS	-	expression tag	UNP Q9H1K1
D	174	HIS	-	expression tag	UNP Q9H1K1
D	175	HIS	-	expression tag	UNP Q9H1K1

- Molecule 5 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 7 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

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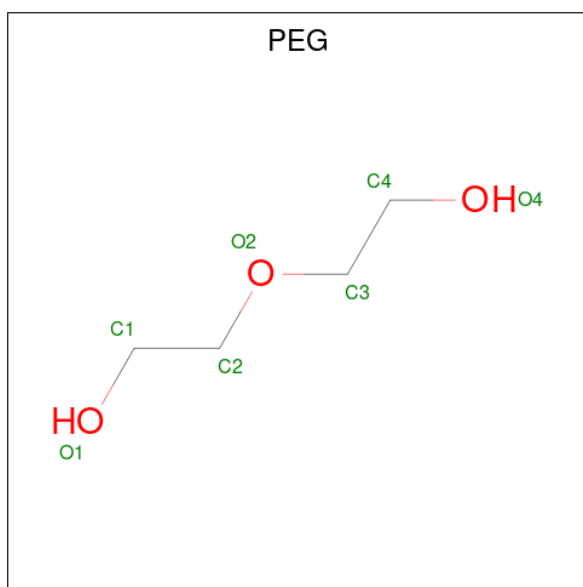
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total 4	C 2	O 2	0	0
7	A	1	Total 4	C 2	O 2	0	0
7	A	1	Total 4	C 2	O 2	0	0
7	A	1	Total 4	C 2	O 2	0	0
7	A	1	Total 4	C 2	O 2	0	0
7	A	1	Total 4	C 2	O 2	0	0
7	A	1	Total 4	C 2	O 2	0	0
7	A	1	Total 4	C 2	O 2	0	0
7	A	1	Total 4	C 2	O 2	0	0
7	A	1	Total 4	C 2	O 2	0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

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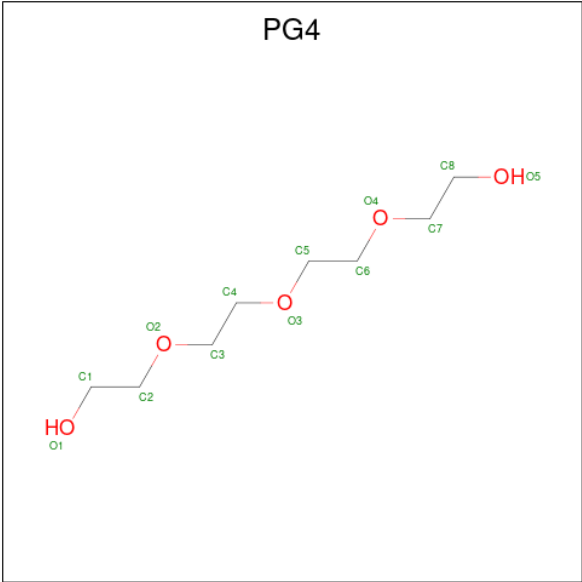
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			4	2	2		
7	B	1	Total	C	O	0	0
			4	2	2		
7	B	1	Total	C	O	0	0
			4	2	2		
7	B	1	Total	C	O	0	0
			4	2	2		
7	B	1	Total	C	O	0	0
			4	2	2		
7	B	1	Total	C	O	0	0
			4	2	2		
7	C	1	Total	C	O	0	0
			4	2	2		
7	C	1	Total	C	O	0	0
			4	2	2		
7	C	1	Total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

- Molecule 8 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



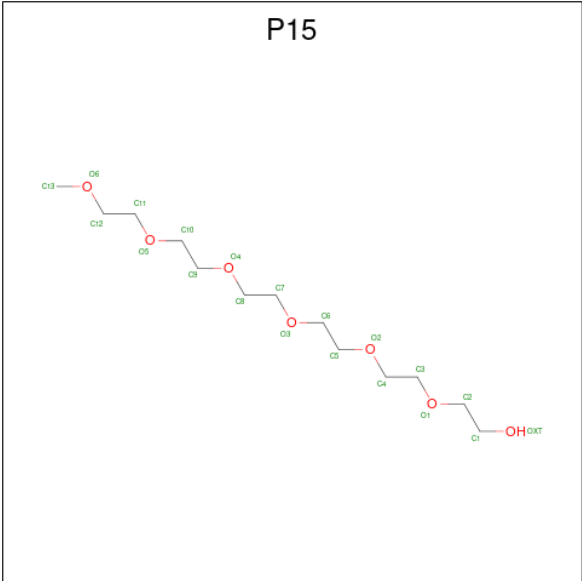
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			7	4	3		
8	A	1	Total	C	O	0	0
			7	4	3		
8	A	1	Total	C	O	0	0
			7	4	3		
8	A	1	Total	C	O	0	0
			7	4	3		
8	A	1	Total	C	O	0	0
			7	4	3		
8	A	1	Total	C	O	0	0
			7	4	3		
8	B	1	Total	C	O	0	0
			7	4	3		
8	B	1	Total	C	O	0	0
			7	4	3		
8	C	1	Total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

- Molecule 9 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



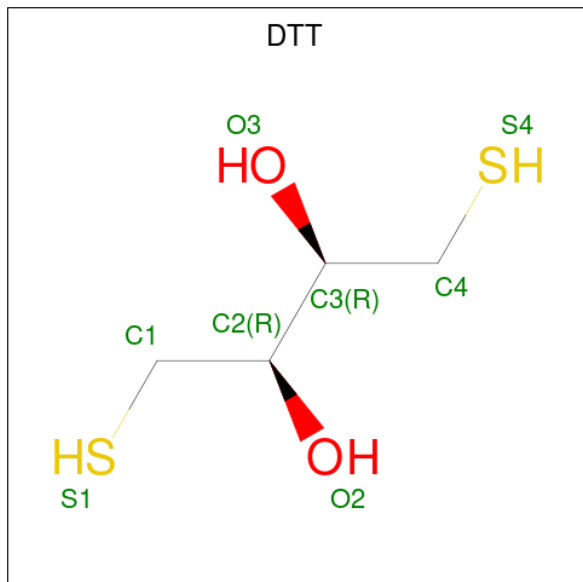
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			13	8	5		
9	A	1	Total	C	O	0	0
			13	8	5		

- Molecule 10 is 2,5,8,11,14,17-HEXAOXANONADECAN-19-OL (CCD ID: P15) (formula: C₁₃H₂₈O₇).



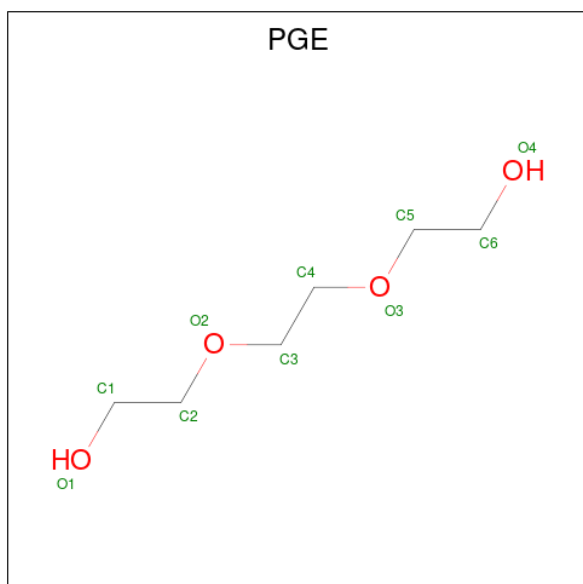
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			20	13	7		

- Molecule 11 is 2,3-DIHYDROXY-1,4-DITHIOBUTANE (CCD ID: DTT) (formula: $C_4H_{10}O_2S_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	A	1	Total	C	O	S	0	0
			8	4	2	2		

- Molecule 12 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



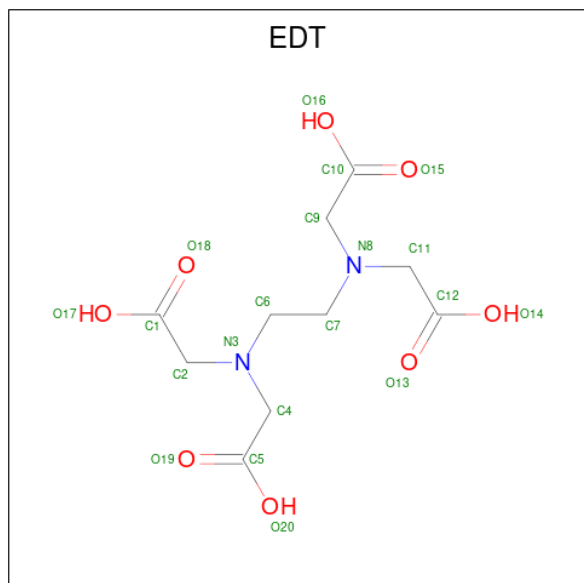
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	A	1	Total	C	O	0	0
			10	6	4		

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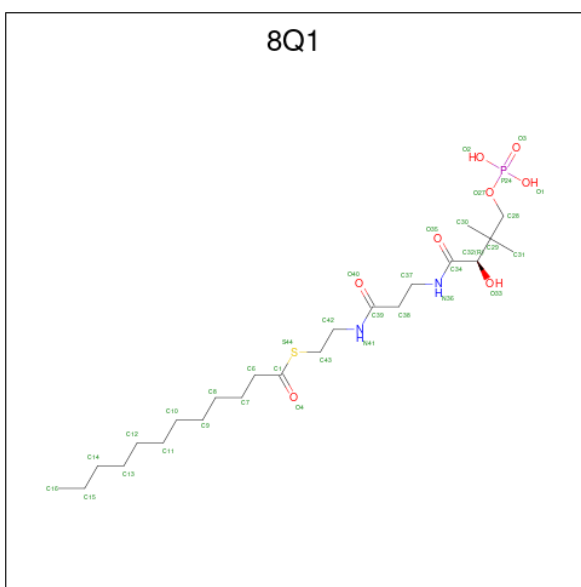
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	A	1	Total	C	O	0	0
			10	6	4		
12	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 13 is {[-(BIS-CARBOXYMETHYL-AMINO)-ETHYL]-CARBOXYMETHYL-AMINO}-ACETIC ACID (CCD ID: EDT) (formula: C₁₀H₁₆N₂O₈).



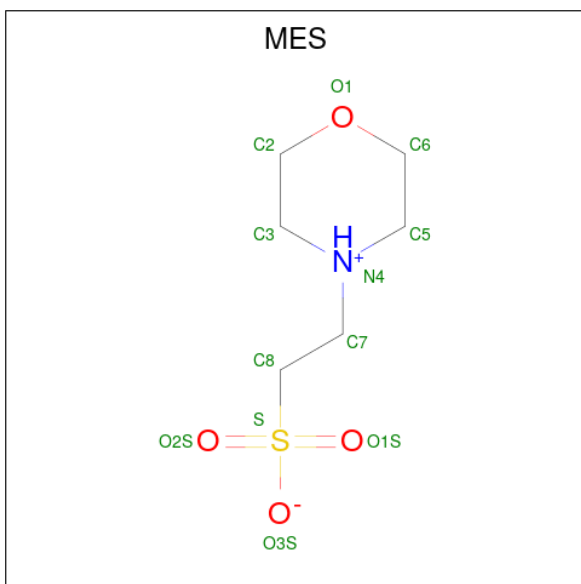
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
13	B	1	Total	C	N	O	0	0
			20	10	2	8		

- Molecule 14 is S-[2-({N-[(2R)-2-hydroxy-3,3-dimethyl-4-(phosphonooxy)butanoyl]-beta-alanyl}amino)ethyl] dodecanethioate (CCD ID: 8Q1) (formula: C₂₃H₄₅N₂O₈PS) (labeled as "Ligand of Interest" by depositor).



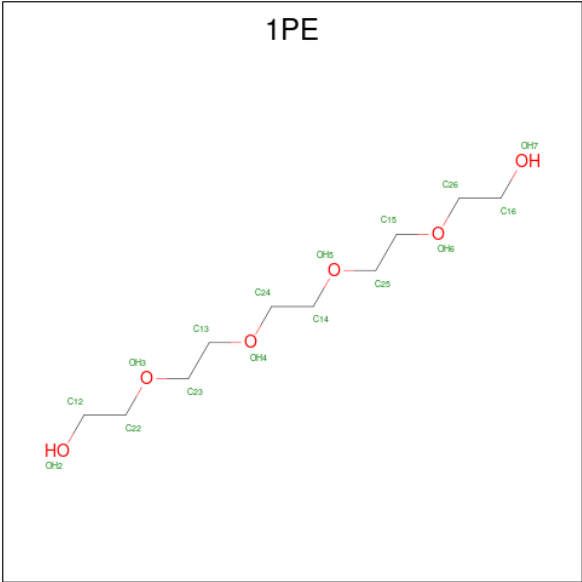
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
14	C	1	Total	C	N	O	P	S	0	0
			34	23	2	7	1	1		

- Molecule 15 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
15	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 16 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $\text{C}_{10}\text{H}_{22}\text{O}_6$).



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

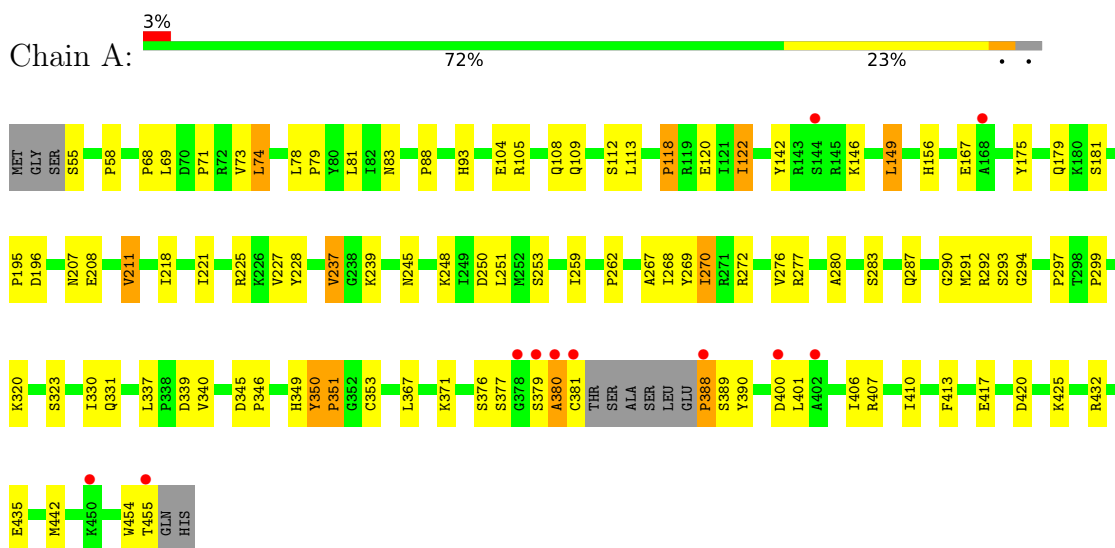
- Molecule 17 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
17	A	260	Total	O	0	0
			260	260		
17	B	71	Total	O	0	0
			71	71		
17	C	18	Total	O	0	0
			18	18		
17	D	41	Total	O	0	0
			41	41		

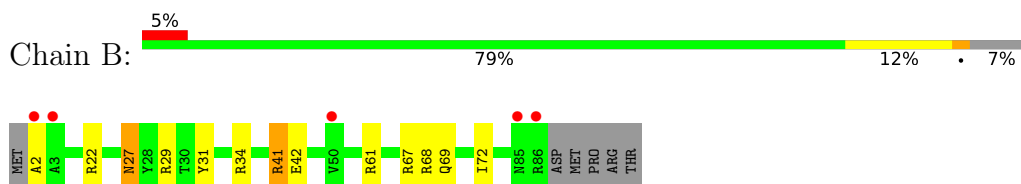
3 Residue-property plots [i](#)

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.

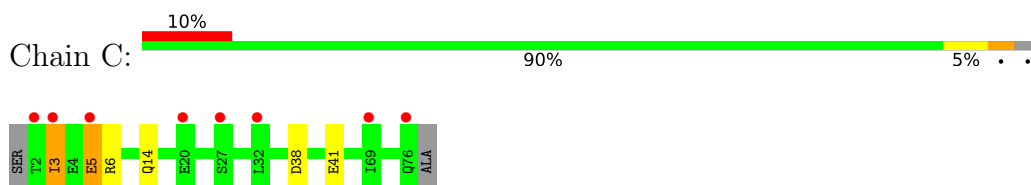
- Molecule 1: Cysteine desulfurase, mitochondrial



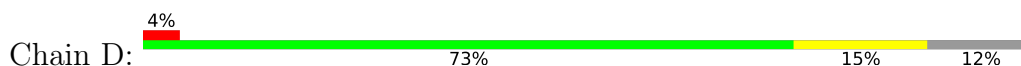
- Molecule 2: LYR motif-containing protein 4

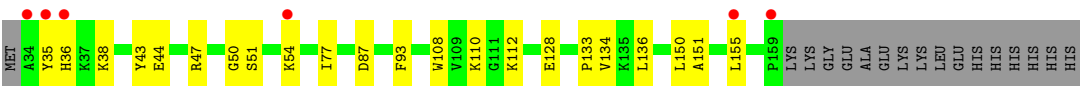


- Molecule 3: Acyl carrier protein



- Molecule 4: Iron-sulfur cluster assembly enzyme ISCU, mitochondrial





4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	86.41Å 86.41Å 245.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.98 – 1.57 48.98 – 1.57	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.98-1.57) 100.0 (48.98-1.57)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 1.57Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.156 , 0.180 (Not available) , 0.179	Depositor DCC
R_{free} test set	6521 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	21.6	Xtriage
Anisotropy	0.157	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6265	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.13% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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, PEG, GOL, 8Q1, P15, 1PE, PG4, MES, EDT, EDO, PLP, DTT

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	1.42	40/3260 (1.2%)	1.12	14/4410 (0.3%)
2	B	1.08	5/761 (0.7%)	0.92	0/1021
3	C	0.48	0/539	0.64	0/735
4	D	0.68	1/1002 (0.1%)	0.74	0/1357
All	All	1.20	46/5562 (0.8%)	1.00	14/7523 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
3	C	0	1
All	All	0	5

The worst 5 of 46 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	237[A]	VAL	CA-C	12.60	1.68	1.52
1	A	237[B]	VAL	CA-C	12.60	1.68	1.52
1	A	211[A]	VAL	CA-C	10.78	1.65	1.52
1	A	211[B]	VAL	CA-C	10.78	1.65	1.52
1	A	353	CYS	C-O	-9.36	1.13	1.24

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	237[A]	VAL	CA-C-O	-7.39	112.01	120.95
1	A	237[B]	VAL	CA-C-O	-7.39	112.01	120.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	388	PRO	CA-N-CD	-6.43	103.00	112.00
1	A	149[A]	LEU	CA-C-O	-6.39	113.89	121.11
1	A	149[B]	LEU	CA-C-O	-6.39	113.89	121.11

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	211[A]	VAL	Mainchain
1	A	211[B]	VAL	Mainchain
1	A	237[A]	VAL	Mainchain
1	A	237[B]	VAL	Mainchain
3	C	5	GLU	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3141	0	3188	112	0
2	B	728	0	766	19	0
3	C	535	0	472	7	0
4	D	977	0	984	19	0
5	A	15	0	7	0	0
6	A	30	0	40	8	0
6	D	6	0	8	0	0
7	A	132	0	193	46	0
7	B	36	0	53	8	0
7	C	12	0	18	3	0
7	D	20	0	30	4	0
8	A	56	0	80	10	0
8	B	14	0	20	12	0
8	C	7	0	10	1	0
9	A	26	0	36	4	0
10	A	20	0	28	9	0
11	A	8	0	10	5	0
12	A	30	0	42	16	0
13	B	20	0	12	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
14	C	34	0	0	0	0
15	C	12	0	13	1	0
16	D	16	0	22	3	0
17	A	260	0	0	14	0
17	B	71	0	0	6	0
17	C	18	0	0	1	0
17	D	41	0	0	1	0
All	All	6265	0	6032	179	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 15.

The worst 5 of 179 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:68:ARG:HH11	8:B:106:PEG:H22	1.14	1.04
6:A:503:GOL:H12	2:B:34[A]:ARG:HH21	1.25	1.02
1:A:245:ASN:HD21	11:A:533:DTT:H2	1.28	0.97
1:A:208[A]:GLU:HG2	1:A:379:SER:O	1.67	0.94
1:A:112:SER:HA	9:A:519:PG4:H11	1.53	0.89

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	413/406 (102%)	403 (98%)	10 (2%)	0	100	100
2	B	91/91 (100%)	89 (98%)	2 (2%)	0	100	100
3	C	73/77 (95%)	70 (96%)	2 (3%)	1 (1%)	9	1
4	D	130/143 (91%)	128 (98%)	0	2 (2%)	8	1

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	707/717 (99%)	690 (98%)	14 (2%)	3 (0%)	36 12

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	35[A]	TYR
4	D	35[B]	TYR
3	C	3	ILE

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	345/346 (100%)	345 (100%)	0	100 100
2	B	76/80 (95%)	75 (99%)	1 (1%)	61 37
3	C	49/66 (74%)	49 (100%)	0	100 100
4	D	103/118 (87%)	102 (99%)	1 (1%)	68 49
All	All	573/610 (94%)	571 (100%)	2 (0%)	84 79

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	41	ARG
4	D	47	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	GLN
1	A	172	GLN
1	A	179	GLN
1	A	207	ASN
1	A	444	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

79 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	$\# Z > 2$
6	GOL	A	502	-	5,5,5	0.41	0	5,5,5	1.35	1 (20%)
7	EDO	A	522	-	3,3,3	0.74	0	2,2,2	0.78	0
7	EDO	A	553	-	3,3,3	0.41	0	2,2,2	0.35	0
8	PEG	A	531	-	6,6,6	0.55	0	5,5,5	0.85	0
7	EDO	A	538	-	3,3,3	0.47	0	2,2,2	0.44	0
7	EDO	A	514	-	3,3,3	1.30	0	2,2,2	0.81	0
7	EDO	A	523	-	3,3,3	0.46	0	2,2,2	0.53	0
7	EDO	A	529	-	3,3,3	0.45	0	2,2,2	0.33	0
7	EDO	A	508	-	3,3,3	0.43	0	2,2,2	0.31	0
7	EDO	A	512	-	3,3,3	0.40	0	2,2,2	0.77	0
7	EDO	A	517	-	3,3,3	0.07	0	2,2,2	0.42	0
7	EDO	A	528	-	3,3,3	0.49	0	2,2,2	0.20	0
12	PGE	A	534	-	9,9,9	0.47	0	8,8,8	0.99	1 (12%)
7	EDO	A	536	-	3,3,3	0.39	0	2,2,2	0.57	0
7	EDO	A	552	-	3,3,3	0.42	0	2,2,2	0.72	0
7	EDO	A	547	-	3,3,3	0.41	0	2,2,2	0.50	0
12	PGE	A	544	-	9,9,9	0.38	0	8,8,8	0.23	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	EDO	A	541	-	3,3,3	0.39	0	2,2,2	0.23	0
7	EDO	B	102	-	3,3,3	0.47	0	2,2,2	0.75	0
7	EDO	B	111	-	3,3,3	0.37	0	2,2,2	0.33	0
8	PEG	A	518	-	6,6,6	0.59	0	5,5,5	0.62	0
13	EDT	B	110	-	19,19,19	1.35	2 (10%)	24,24,24	1.39	3 (12%)
7	EDO	A	510	-	3,3,3	0.41	0	2,2,2	0.68	0
7	EDO	A	554	-	3,3,3	0.47	0	2,2,2	0.44	0
16	1PE	D	203	-	15,15,15	0.54	0	14,14,14	0.25	0
7	EDO	B	108	-	3,3,3	0.42	0	2,2,2	0.43	0
7	EDO	D	202	-	3,3,3	0.45	0	2,2,2	0.43	0
7	EDO	D	204	-	3,3,3	0.44	0	2,2,2	0.33	0
7	EDO	A	524	-	3,3,3	0.42	0	2,2,2	0.52	0
7	EDO	A	525	-	3,3,3	0.43	0	2,2,2	0.45	0
9	PG4	A	520	-	12,12,12	0.56	0	11,11,11	0.40	0
8	PEG	A	527	-	6,6,6	0.50	0	5,5,5	0.35	0
7	EDO	A	542	-	3,3,3	0.48	0	2,2,2	0.35	0
7	EDO	A	506	-	3,3,3	0.36	0	2,2,2	0.25	0
7	EDO	D	206	-	3,3,3	0.46	0	2,2,2	0.52	0
7	EDO	A	516	-	3,3,3	0.52	0	2,2,2	0.24	0
12	PGE	A	543	-	9,9,9	0.27	0	8,8,8	0.58	0
7	EDO	A	530	-	3,3,3	0.49	0	2,2,2	0.32	0
14	8Q1	C	301	3	28,33,34	2.43	9 (32%)	32,40,43	1.21	2 (6%)
7	EDO	B	107	-	3,3,3	0.46	0	2,2,2	0.17	0
7	EDO	A	546	-	3,3,3	0.73	0	2,2,2	1.35	0
7	EDO	A	515	-	3,3,3	0.47	0	2,2,2	0.14	0
6	GOL	A	551	-	5,5,5	0.49	0	5,5,5	0.53	0
6	GOL	A	504	-	5,5,5	0.57	0	5,5,5	0.29	0
10	P15	A	526	-	19,19,19	0.65	0	18,18,18	0.75	0
8	PEG	C	305	-	6,6,6	0.50	0	5,5,5	0.24	0
8	PEG	A	549	-	6,6,6	0.49	0	5,5,5	0.33	0
8	PEG	B	106	-	6,6,6	0.64	0	5,5,5	0.82	0
5	PLP	A	501	1	15,15,16	1.11	1 (6%)	21,22,23	1.06	0
7	EDO	C	304	-	3,3,3	0.51	0	2,2,2	0.58	0
7	EDO	D	201	-	3,3,3	0.44	0	2,2,2	0.27	0
6	GOL	A	505	-	5,5,5	0.87	0	5,5,5	0.65	0
8	PEG	B	104	-	6,6,6	0.55	0	5,5,5	0.26	0
7	EDO	D	207	-	3,3,3	0.42	0	2,2,2	0.51	0
8	PEG	A	532	-	6,6,6	0.51	0	5,5,5	0.60	0
7	EDO	C	303	-	3,3,3	0.45	0	2,2,2	0.48	0
6	GOL	D	205	-	5,5,5	0.56	0	5,5,5	0.32	0
7	EDO	B	112	-	3,3,3	0.44	0	2,2,2	0.21	0
7	EDO	B	109	-	3,3,3	0.46	0	2,2,2	0.49	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	PEG	A	521	-	6,6,6	0.49	0	5,5,5	0.94	0
7	EDO	A	540	-	3,3,3	0.43	0	2,2,2	0.47	0
7	EDO	A	513	-	3,3,3	0.40	0	2,2,2	0.51	0
7	EDO	A	511	-	3,3,3	0.53	0	2,2,2	0.64	0
7	EDO	A	539	-	3,3,3	0.42	0	2,2,2	0.34	0
7	EDO	A	535	-	3,3,3	0.40	0	2,2,2	0.47	0
8	PEG	A	548	-	6,6,6	0.50	0	5,5,5	0.35	0
8	PEG	A	550	-	6,6,6	0.49	0	5,5,5	0.29	0
11	DTT	A	533	-	7,7,7	0.64	0	4,8,8	0.96	0
7	EDO	B	103	-	3,3,3	0.41	0	2,2,2	0.37	0
7	EDO	A	507	-	3,3,3	1.99	1 (33%)	2,2,2	1.59	1 (50%)
7	EDO	B	101	-	3,3,3	0.86	0	2,2,2	1.15	0
7	EDO	B	105	-	3,3,3	0.41	0	2,2,2	0.39	0
7	EDO	A	537	-	3,3,3	0.41	0	2,2,2	0.32	0
7	EDO	A	545	-	3,3,3	0.43	0	2,2,2	0.29	0
15	MES	C	302	-	12,12,12	1.96	5 (41%)	15,16,16	1.94	3 (20%)
7	EDO	C	306	-	3,3,3	0.43	0	2,2,2	0.32	0
6	GOL	A	503	-	5,5,5	0.46	0	5,5,5	0.53	0
9	PG4	A	519	-	12,12,12	0.54	0	11,11,11	0.48	0
7	EDO	A	509	-	3,3,3	0.65	0	2,2,2	0.99	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
6	GOL	A	502	-	-	0/4/4/4	-
7	EDO	A	522	-	-	0/1/1/1	-
7	EDO	A	553	-	-	1/1/1/1	-
8	PEG	A	531	-	-	2/4/4/4	-
7	EDO	A	538	-	-	0/1/1/1	-
7	EDO	A	514	-	-	1/1/1/1	-
7	EDO	A	523	-	-	1/1/1/1	-
7	EDO	A	529	-	-	1/1/1/1	-
7	EDO	A	508	-	-	0/1/1/1	-
7	EDO	A	512	-	-	1/1/1/1	-
7	EDO	A	517	-	-	1/1/1/1	-
7	EDO	A	528	-	-	1/1/1/1	-
12	PGE	A	534	-	-	5/7/7/7	-
7	EDO	A	536	-	-	1/1/1/1	-
7	EDO	A	552	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	EDO	A	547	-	-	1/1/1/1	-
12	PGE	A	544	-	-	3/7/7/7	-
7	EDO	A	541	-	-	1/1/1/1	-
7	EDO	B	102	-	-	1/1/1/1	-
7	EDO	B	111	-	-	1/1/1/1	-
8	PEG	A	518	-	-	2/4/4/4	-
13	EDT	B	110	-	-	12/21/21/21	-
7	EDO	A	510	-	-	1/1/1/1	-
7	EDO	A	554	-	-	1/1/1/1	-
16	1PE	D	203	-	-	7/13/13/13	-
7	EDO	B	108	-	-	1/1/1/1	-
7	EDO	D	202	-	-	1/1/1/1	-
7	EDO	D	204	-	-	0/1/1/1	-
7	EDO	A	524	-	-	1/1/1/1	-
7	EDO	A	525	-	-	1/1/1/1	-
9	PG4	A	520	-	-	5/10/10/10	-
8	PEG	A	527	-	-	1/4/4/4	-
7	EDO	A	542	-	-	0/1/1/1	-
7	EDO	A	506	-	-	1/1/1/1	-
7	EDO	D	206	-	-	1/1/1/1	-
7	EDO	A	516	-	-	1/1/1/1	-
12	PGE	A	543	-	-	3/7/7/7	-
7	EDO	A	530	-	-	1/1/1/1	-
14	8Q1	C	301	3	-	2/38/40/41	-
7	EDO	B	107	-	-	0/1/1/1	-
7	EDO	A	546	-	-	1/1/1/1	-
7	EDO	A	515	-	-	1/1/1/1	-
6	GOL	A	551	-	-	4/4/4/4	-
6	GOL	A	504	-	-	4/4/4/4	-
10	P15	A	526	-	-	10/17/17/17	-
8	PEG	C	305	-	-	4/4/4/4	-
8	PEG	A	549	-	-	2/4/4/4	-
8	PEG	B	106	-	-	3/4/4/4	-
5	PLP	A	501	1	-	0/6/6/8	0/1/1/1
7	EDO	C	304	-	-	1/1/1/1	-
7	EDO	D	201	-	-	0/1/1/1	-
6	GOL	A	505	-	-	2/4/4/4	-
8	PEG	B	104	-	-	3/4/4/4	-
7	EDO	D	207	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	PEG	A	532	-	-	0/4/4/4	-
7	EDO	C	303	-	-	1/1/1/1	-
6	GOL	D	205	-	-	1/4/4/4	-
7	EDO	B	112	-	-	1/1/1/1	-
7	EDO	B	109	-	-	0/1/1/1	-
8	PEG	A	521	-	-	2/4/4/4	-
7	EDO	A	540	-	-	1/1/1/1	-
7	EDO	A	513	-	-	1/1/1/1	-
7	EDO	A	511	-	-	0/1/1/1	-
7	EDO	A	539	-	-	0/1/1/1	-
7	EDO	A	535	-	-	1/1/1/1	-
8	PEG	A	548	-	-	4/4/4/4	-
8	PEG	A	550	-	-	2/4/4/4	-
11	DTT	A	533	-	-	2/8/8/8	-
7	EDO	B	103	-	-	0/1/1/1	-
7	EDO	A	507	-	-	0/1/1/1	-
7	EDO	B	101	-	-	0/1/1/1	-
7	EDO	B	105	-	-	1/1/1/1	-
7	EDO	A	537	-	-	0/1/1/1	-
7	EDO	A	545	-	-	1/1/1/1	-
15	MES	C	302	-	-	3/6/14/14	0/1/1/1
7	EDO	C	306	-	-	1/1/1/1	-
6	GOL	A	503	-	-	4/4/4/4	-
9	PG4	A	519	-	-	5/10/10/10	-
7	EDO	A	509	-	-	0/1/1/1	-

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	C	301	8Q1	O4-C1	-6.28	1.11	1.21
14	C	301	8Q1	O40-C39	-4.78	1.13	1.23
14	C	301	8Q1	C39-N41	4.48	1.44	1.33
14	C	301	8Q1	O35-C34	-4.23	1.15	1.23
15	C	302	MES	C8-S	4.08	1.83	1.77

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	C	302	MES	O1S-S-C8	4.52	113.56	106.73
15	C	302	MES	O3S-S-O2S	-3.86	101.73	111.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	B	110	EDT	C11-N8-C9	3.75	120.45	111.66
13	B	110	EDT	C11-N8-C7	2.80	118.66	111.91
14	C	301	8Q1	C42-N41-C39	-2.58	118.03	122.82

There are no chirality outliers.

5 of 130 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	503	GOL	O1-C1-C2-C3
6	A	503	GOL	C1-C2-C3-O3
6	A	504	GOL	C1-C2-C3-O3
6	A	551	GOL	C1-C2-C3-O3
6	A	551	GOL	O2-C2-C3-O3

There are no ring outliers.

47 monomers are involved in 127 short contacts:

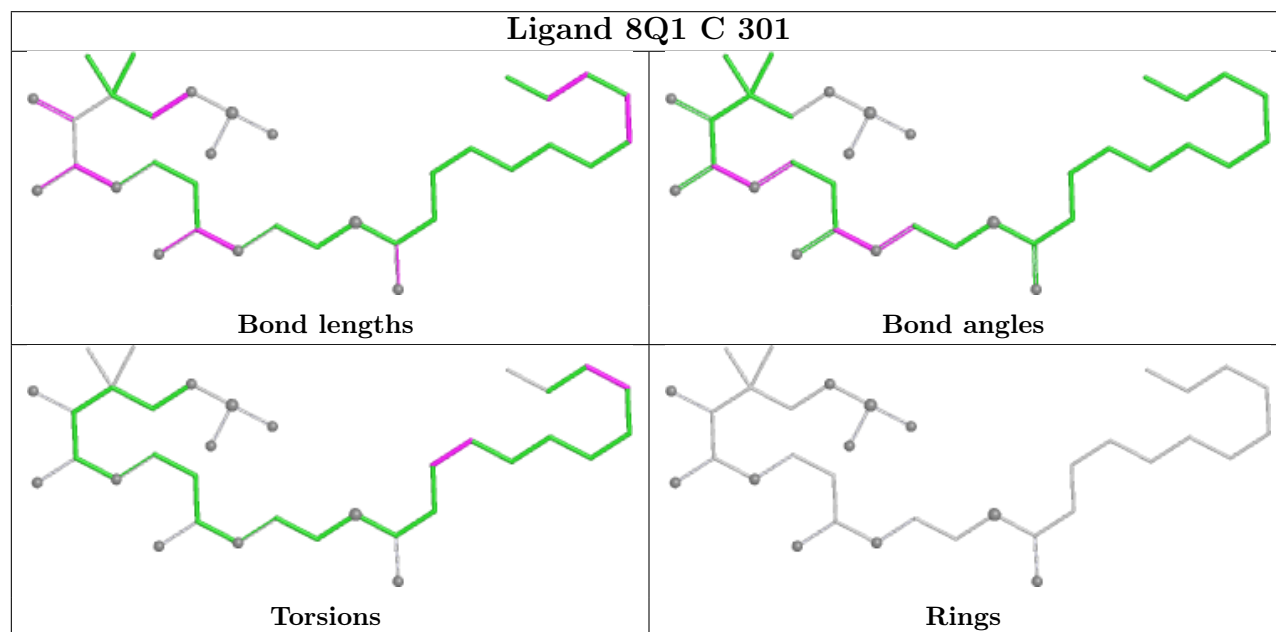
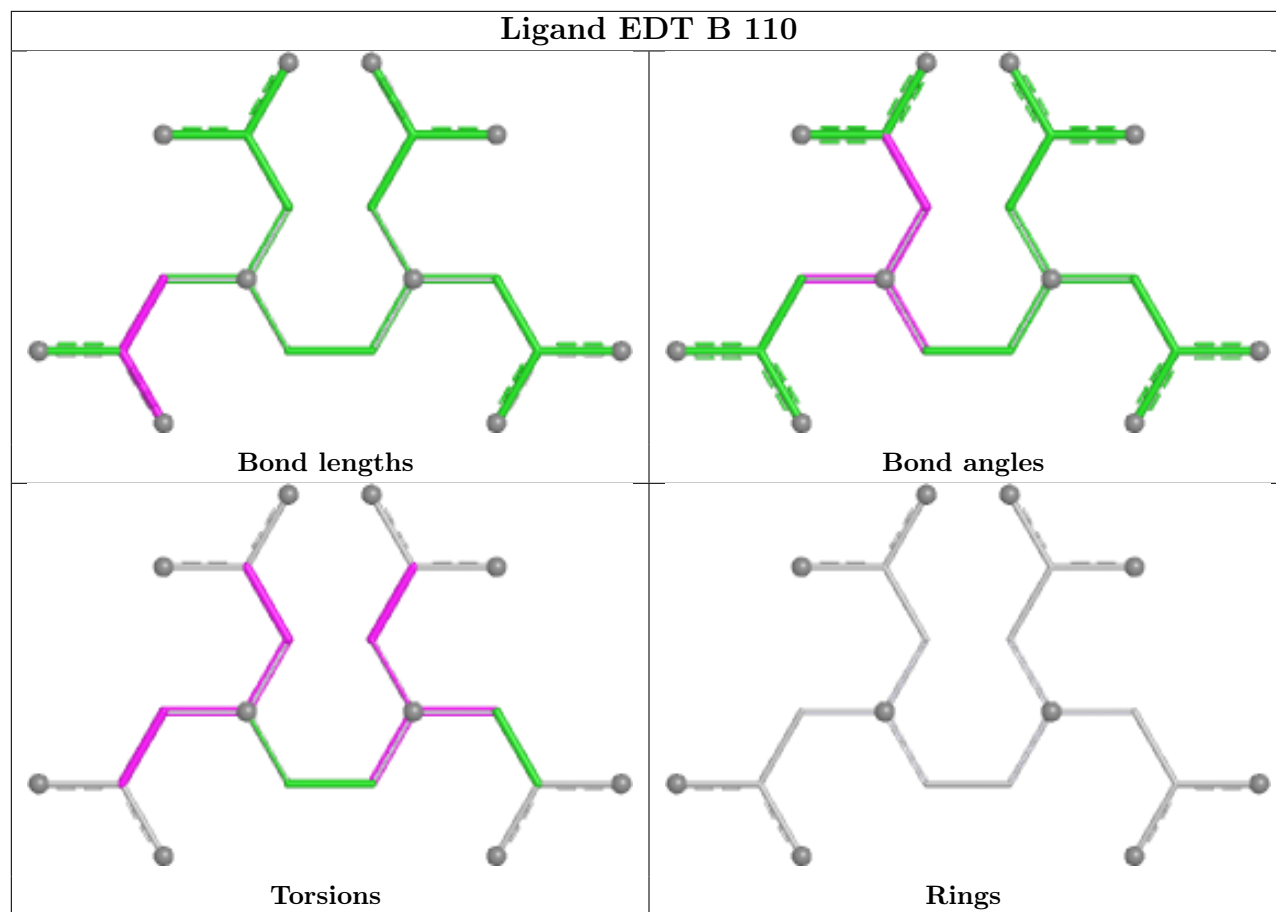
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	502	GOL	1	0
7	A	522	EDO	1	0
8	A	531	PEG	5	0
7	A	514	EDO	7	0
7	A	523	EDO	1	0
7	A	517	EDO	3	0
12	A	534	PGE	9	0
7	A	552	EDO	3	0
7	A	547	EDO	1	0
7	A	541	EDO	1	0
7	B	111	EDO	1	0
8	A	518	PEG	1	0
13	B	110	EDT	4	0
7	A	510	EDO	1	0
7	A	554	EDO	8	0
16	D	203	1PE	3	0
7	D	202	EDO	2	0
7	A	524	EDO	1	0
7	A	506	EDO	1	0
12	A	543	PGE	7	0
7	A	530	EDO	1	0
7	B	107	EDO	4	0
7	A	546	EDO	2	0

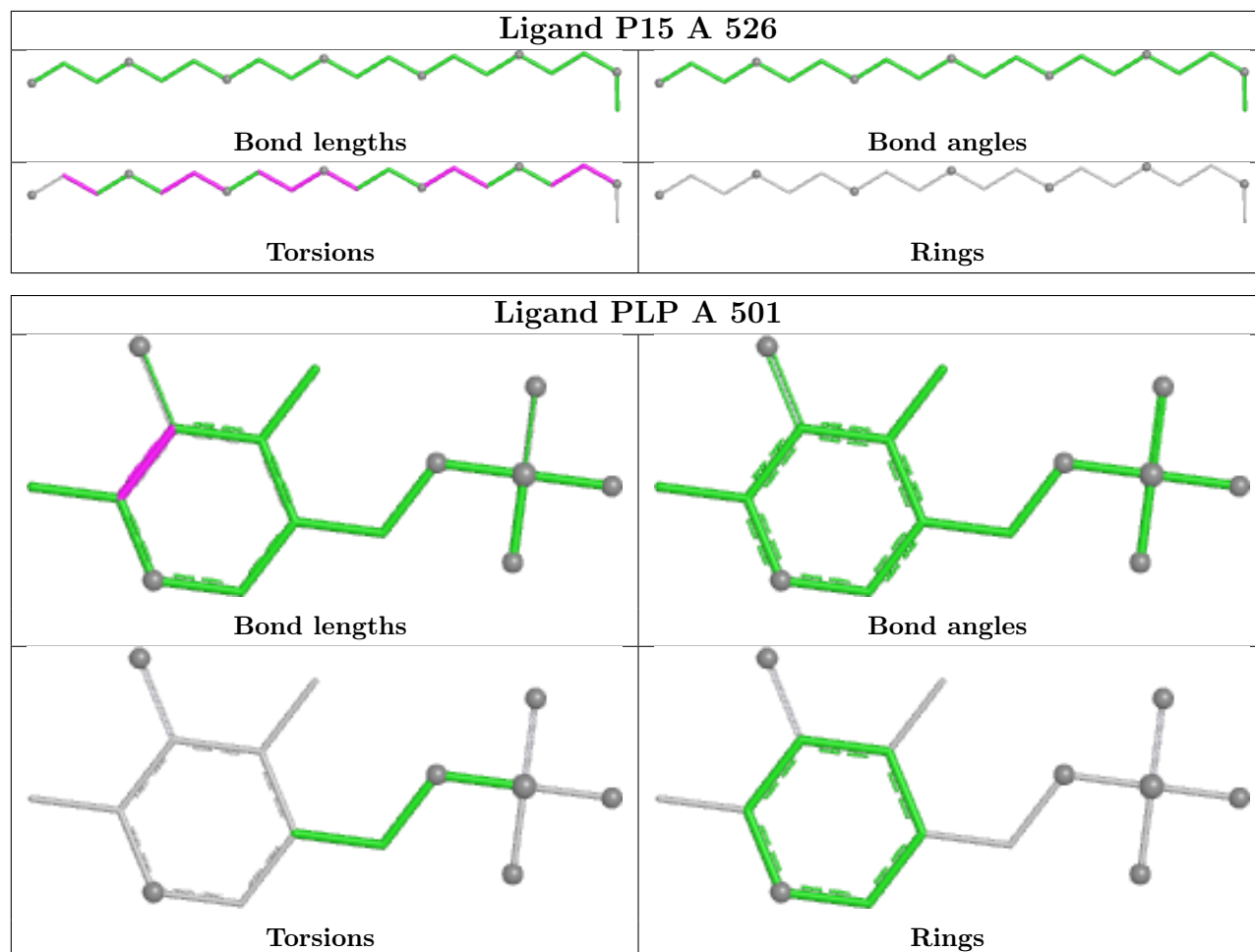
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	515	EDO	2	0
6	A	504	GOL	1	0
10	A	526	P15	9	0
8	C	305	PEG	1	0
8	A	549	PEG	1	0
8	B	106	PEG	8	0
7	C	304	EDO	2	0
7	D	201	EDO	2	0
6	A	505	GOL	3	0
8	B	104	PEG	4	0
8	A	532	PEG	1	0
7	C	303	EDO	1	0
7	A	511	EDO	4	0
7	A	535	EDO	3	0
8	A	548	PEG	1	0
8	A	550	PEG	1	0
11	A	533	DTT	5	0
7	B	103	EDO	1	0
7	A	507	EDO	6	0
7	B	101	EDO	2	0
15	C	302	MES	1	0
6	A	503	GOL	3	0
9	A	519	PG4	4	0
7	A	509	EDO	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 [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, 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	395/406 (97%)	-0.08	11 (2%) 55 67	10, 20, 44, 89	26 (6%)
2	B	85/91 (93%)	0.12	5 (5%) 28 38	10, 23, 46, 88	8 (9%)
3	C	75/77 (97%)	0.98	8 (10%) 11 17	28, 50, 84, 93	0
4	D	126/143 (88%)	0.29	6 (4%) 35 47	11, 32, 56, 97	7 (5%)
All	All	681/717 (94%)	0.13	30 (4%) 39 51	10, 24, 61, 97	41 (6%)

The worst 5 of 30 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	380	ALA	11.2
4	D	35[A]	TYR	7.9
1	A	378	GLY	7.7
4	D	34[A]	ALA	6.7
2	B	2	ALA	5.4

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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($\text{\AA}^2$)	Q<0.9
7	EDO	A	524	4/4	0.59	0.20	57,59,63,64	4
7	EDO	A	553	4/4	0.59	0.32	87,88,90,91	0
8	PEG	C	305	7/7	0.59	0.25	57,71,76,76	0
7	EDO	A	540	4/4	0.66	0.25	76,77,78,80	0
7	EDO	A	529	4/4	0.68	0.31	84,85,87,87	0
7	EDO	A	530	4/4	0.69	0.34	46,51,56,59	4
8	PEG	A	532	7/7	0.70	0.27	69,72,77,83	0
7	EDO	A	542	4/4	0.70	0.21	42,53,63,69	4
7	EDO	B	111	4/4	0.71	0.27	73,74,78,84	0
8	PEG	A	550	7/7	0.72	0.22	74,75,77,79	0
7	EDO	A	525	4/4	0.72	0.23	73,74,75,77	0
7	EDO	D	201	4/4	0.73	0.31	83,87,88,88	0
6	GOL	A	551	6/6	0.73	0.27	36,71,92,98	0
9	PG4	A	520	13/13	0.73	0.22	33,65,81,82	0
7	EDO	D	204	4/4	0.75	0.20	63,66,72,75	0
8	PEG	B	104	7/7	0.75	0.25	36,58,79,80	0
7	EDO	D	207	4/4	0.75	0.22	70,71,76,82	0
7	EDO	C	306	4/4	0.75	0.18	53,67,67,68	0
13	EDT	B	110	20/20	0.75	0.24	35,65,91,96	20
7	EDO	B	109	4/4	0.76	0.25	64,67,73,79	0
7	EDO	A	515	4/4	0.77	0.23	58,59,60,63	0
12	PGE	A	544	10/10	0.77	0.20	84,91,95,95	0
9	PG4	A	519	13/13	0.77	0.23	36,61,71,75	13
7	EDO	A	539	4/4	0.78	0.21	65,66,67,69	0
7	EDO	B	103	4/4	0.78	0.17	65,67,74,81	0
8	PEG	A	549	7/7	0.79	0.19	43,73,80,80	0
7	EDO	A	547	4/4	0.79	0.19	43,56,68,76	0
7	EDO	A	552	4/4	0.79	0.16	38,43,46,54	0
7	EDO	A	512	4/4	0.79	0.23	37,37,39,43	4
7	EDO	A	510	4/4	0.79	0.18	47,56,57,59	0
7	EDO	D	206	4/4	0.79	0.17	55,61,61,70	4
7	EDO	B	108	4/4	0.79	0.17	88,90,91,92	0
7	EDO	A	535	4/4	0.79	0.21	43,50,53,56	4
8	PEG	A	548	7/7	0.80	0.26	79,82,87,87	0
10	P15	A	526	20/20	0.80	0.24	20,74,88,90	0
8	PEG	A	518	7/7	0.80	0.18	33,41,54,61	0
7	EDO	B	107	4/4	0.80	0.21	18,24,29,41	4
11	DTT	A	533	8/8	0.81	0.22	47,75,81,85	8
7	EDO	A	508	4/4	0.81	0.15	47,51,55,59	0
7	EDO	A	537	4/4	0.81	0.19	55,59,71,79	0
7	EDO	A	509	4/4	0.82	0.17	31,36,38,47	4
7	EDO	A	528	4/4	0.82	0.20	32,44,45,46	4
8	PEG	A	527	7/7	0.82	0.21	61,71,77,80	0

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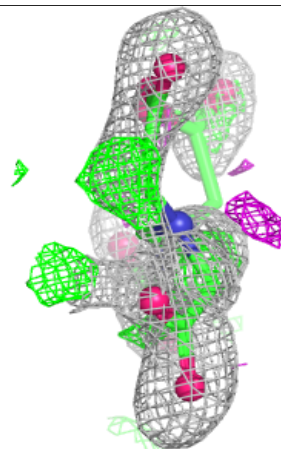
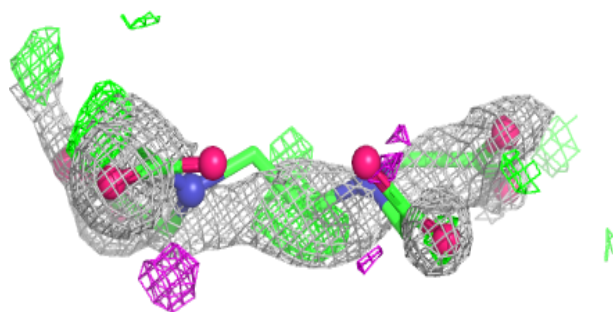
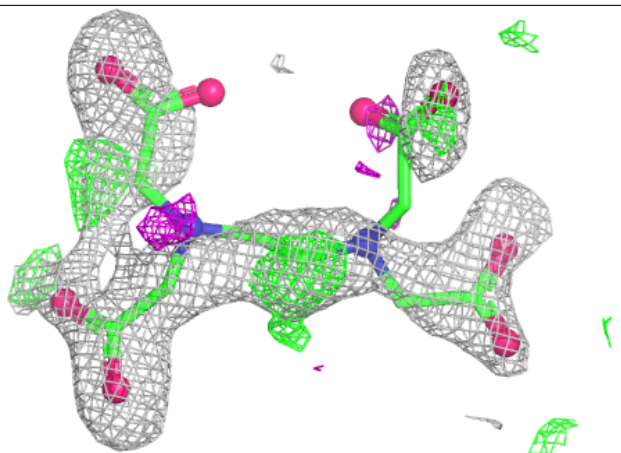
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	EDO	A	541	4/4	0.82	0.20	42,60,66,70	0
7	EDO	B	112	4/4	0.82	0.16	57,61,69,72	0
7	EDO	B	102	4/4	0.83	0.18	48,49,58,59	0
7	EDO	A	536	4/4	0.83	0.20	41,51,64,65	0
7	EDO	B	105	4/4	0.83	0.20	63,66,67,69	0
6	GOL	A	503	6/6	0.83	0.18	28,65,66,67	0
8	PEG	A	521	7/7	0.83	0.17	44,45,54,69	0
7	EDO	A	523	4/4	0.83	0.18	40,45,47,49	0
12	PGE	A	543	10/10	0.84	0.20	25,45,58,61	10
7	EDO	A	554	4/4	0.84	0.19	58,64,74,81	0
12	PGE	A	534	10/10	0.84	0.17	27,42,48,49	10
7	EDO	A	545	4/4	0.85	0.17	54,63,69,77	0
7	EDO	A	538	4/4	0.85	0.21	61,63,74,75	0
8	PEG	B	106	7/7	0.85	0.18	18,40,51,56	7
7	EDO	A	506	4/4	0.86	0.17	45,47,52,54	0
6	GOL	A	504	6/6	0.86	0.15	35,49,56,66	6
15	MES	C	302	12/12	0.86	0.15	52,60,66,69	0
6	GOL	A	505	6/6	0.87	0.16	34,47,48,55	6
7	EDO	A	546	4/4	0.87	0.15	26,28,32,39	4
16	1PE	D	203	16/16	0.87	0.15	54,63,71,76	0
6	GOL	D	205	6/6	0.88	0.14	35,39,43,47	6
6	GOL	A	502	6/6	0.88	0.15	13,25,25,30	6
7	EDO	A	513	4/4	0.88	0.16	40,56,57,60	0
8	PEG	A	531	7/7	0.89	0.17	23,38,66,70	7
7	EDO	A	511	4/4	0.89	0.15	36,36,37,37	4
7	EDO	A	517	4/4	0.89	0.16	13,20,34,35	4
7	EDO	B	101	4/4	0.90	0.13	23,29,31,42	4
7	EDO	C	303	4/4	0.90	0.15	39,39,42,43	4
7	EDO	D	202	4/4	0.90	0.15	29,35,50,59	4
7	EDO	C	304	4/4	0.91	0.13	33,45,48,49	4
7	EDO	A	516	4/4	0.91	0.16	23,24,33,40	4
7	EDO	A	514	4/4	0.94	0.14	22,25,26,34	4
7	EDO	A	522	4/4	0.94	0.13	22,31,37,38	4
7	EDO	A	507	4/4	0.95	0.15	18,19,20,33	4
14	8Q1	C	301	34/35	0.96	0.08	23,29,41,42	0
5	PLP	A	501	15/16	0.98	0.05	14,16,25,26	0

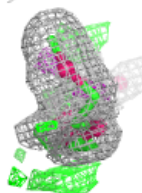
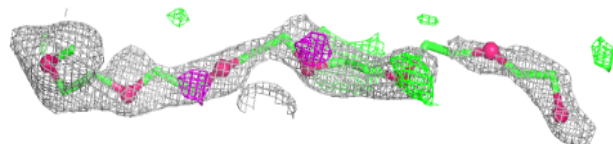
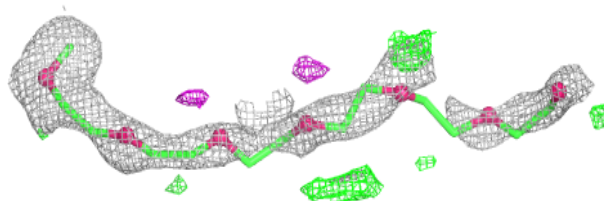
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around EDT B 110:

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

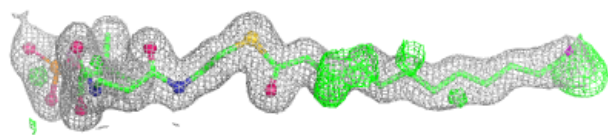
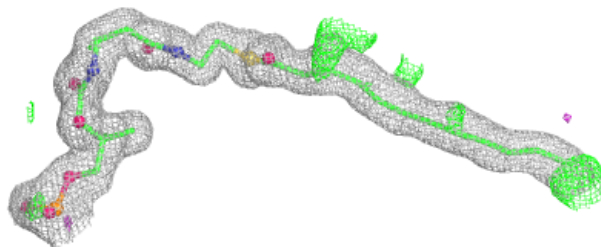
**Electron density around P15 A 526:**

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and green (positive)

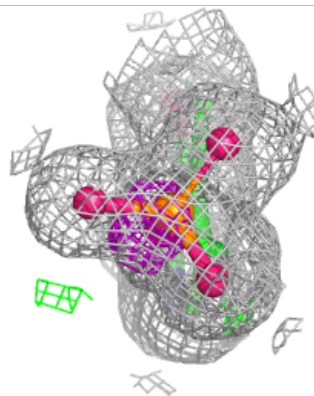
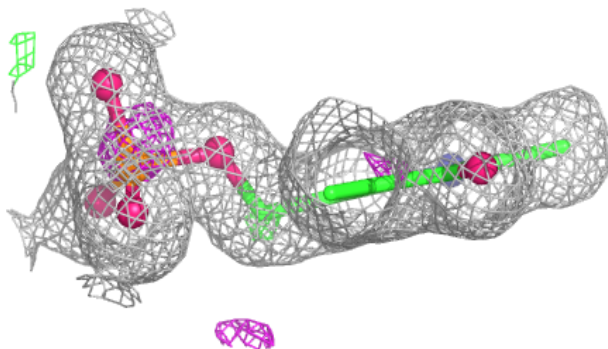
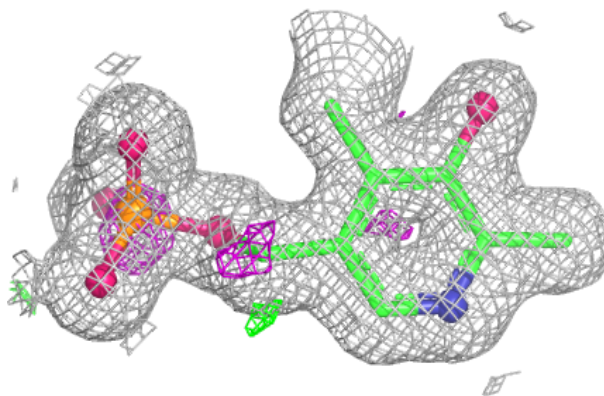


Electron density around 8Q1 C 301:

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

**Electron density around PLP A 501:**

$2mF_o-DF_c$ (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.