



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:26 AM UTC

PDB ID : 6WIH / pdb_00006wih
Title : N-terminal mutation of ISCU2 (L35H36) traps Nfs1 Cys loop in the active site of ISCU2 without metal present. Structure of human mitochondrial complex Nfs1-ISCU2(L35H36)-ISD11 with E.coli ACP1 at 1.9 Å resolution (NIAU)2.
Authors : Boniecki, M.T.; Cygler, M.
Deposited on : 2020-04-09
Resolution : 1.90 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

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

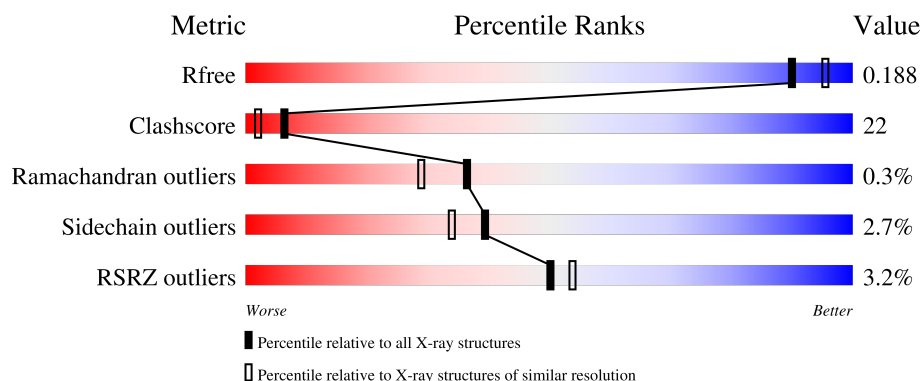
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	2% 75% 20% ...
2	B	91	5% 75% 14% 8%
3	C	77	6% 82% 12% ...
4	D	143	3% 65% 22% 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	P15	A	523	-	-	X	-
11	PGE	A	528	-	-	X	-
11	PGE	A	536	-	-	X	-
12	EDT	B	411	-	-	X	-
14	1PE	D	202	-	-	X	-
6	PEG	B	405	-	-	X	-
6	PEG	B	407	-	-	X	-
7	GOL	B	402	-	-	X	-
8	EDO	A	507	-	-	X	-
8	EDO	A	509	-	-	X	-
8	EDO	A	513	-	-	X	-
8	EDO	A	520	-	-	X	-
8	EDO	A	521	-	-	X	-
8	EDO	A	526	-	-	X	-

2 Entry composition

There are 15 unique types of molecules in this entry. The entry contains 6194 atoms, of which 6 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	401	Total	C	N	O	S	0	9	0
			3134	1977	542	594	21			

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	84	Total	C	N	O	S	0	10	0
			741	469	148	123	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
			550	343	85	121	1			

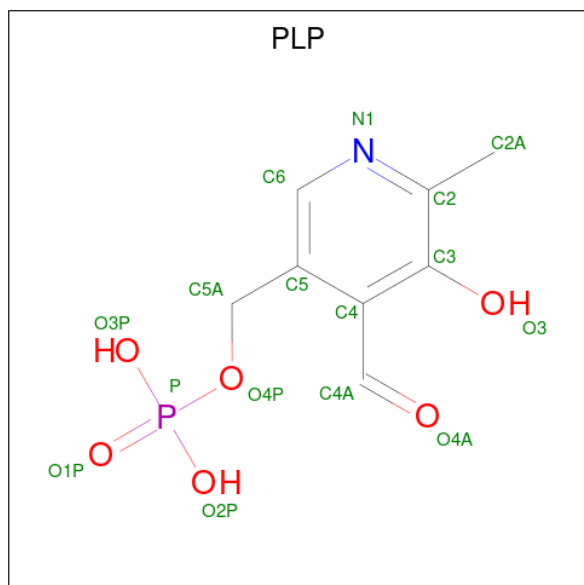
- Molecule 4 is a protein called Iron-sulfur cluster assembly enzyme ISCU, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	125	Total	C	N	O	S	0	1	0
			926	583	156	180	7			

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	36	HIS	SER	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: $C_8H_{10}NO_6P$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	C	H	N	O	P	0	0
			21	8	6	1	5	1		

- Molecule 6 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



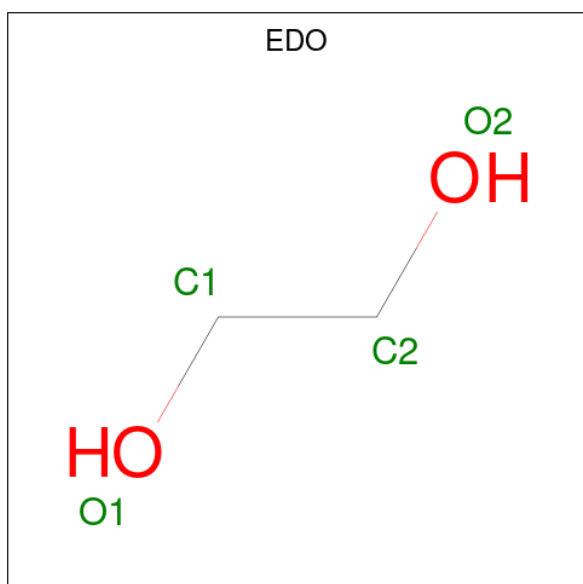
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 8 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



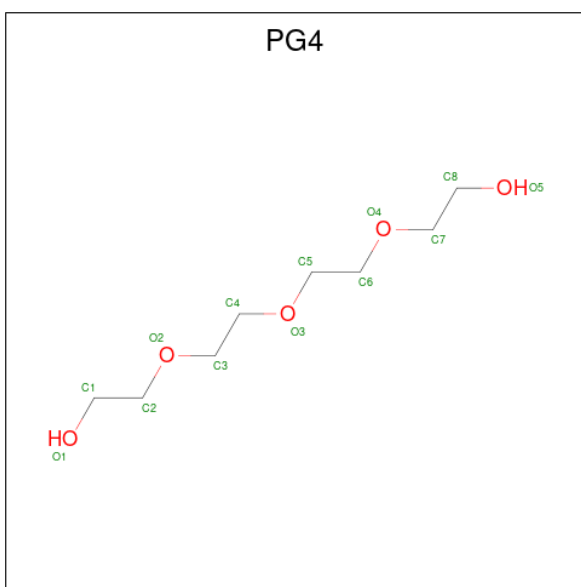
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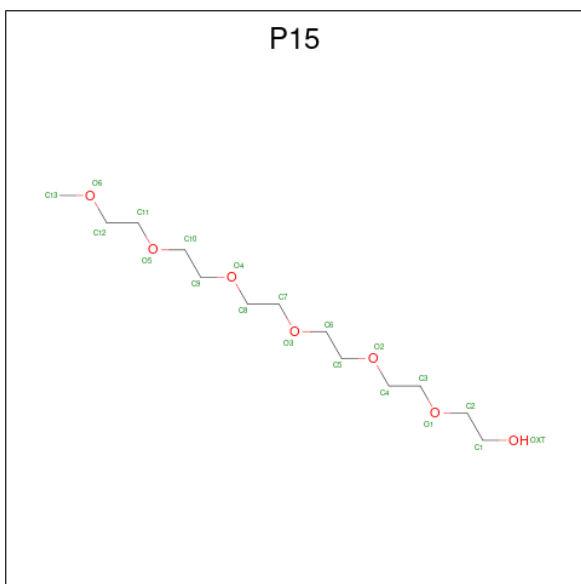
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total 4	C 2	O 2	0	0
8	A	1	Total 4	C 2	O 2	0	0
8	A	1	Total 4	C 2	O 2	0	0
8	B	1	Total 4	C 2	O 2	0	0
8	B	1	Total 4	C 2	O 2	0	0
8	B	1	Total 4	C 2	O 2	0	0
8	B	1	Total 4	C 2	O 2	0	0
8	B	1	Total 4	C 2	O 2	0	0
8	B	1	Total 4	C 2	O 2	0	0
8	B	1	Total 4	C 2	O 2	0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

- Molecule 9 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



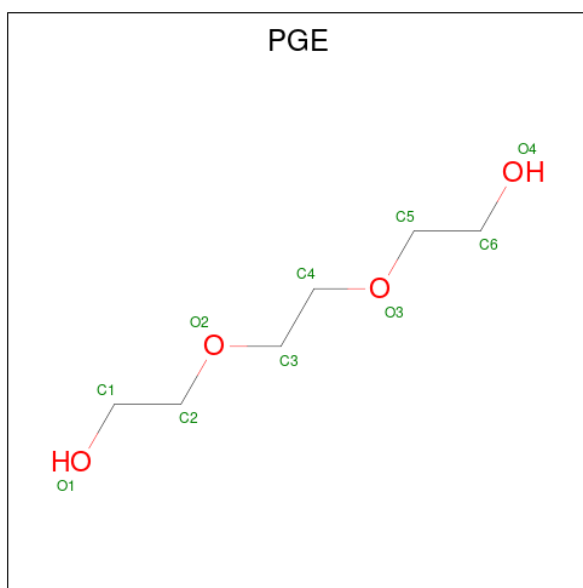
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			13	8	5		
9	A	1	Total	C	O	0	0
			13	8	5		

- Molecule 10 is 2,5,8,11,14,17-HEXAOXANONADECAN-19-OL (CCD ID: P15) (formula: $C_{13}H_{28}O_7$).



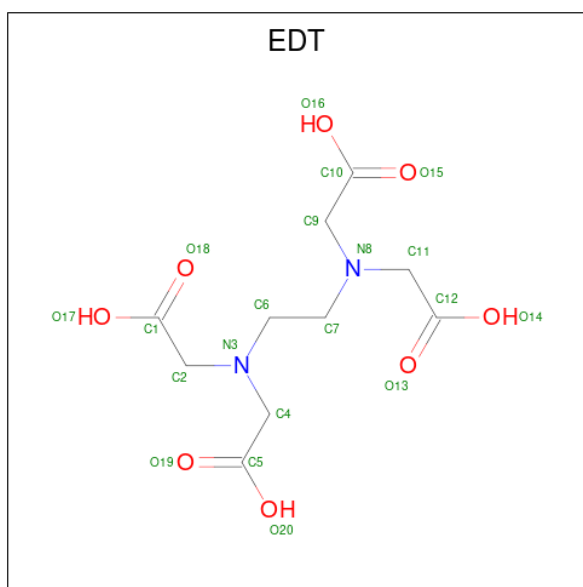
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			20	13	7		

- Molecule 11 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



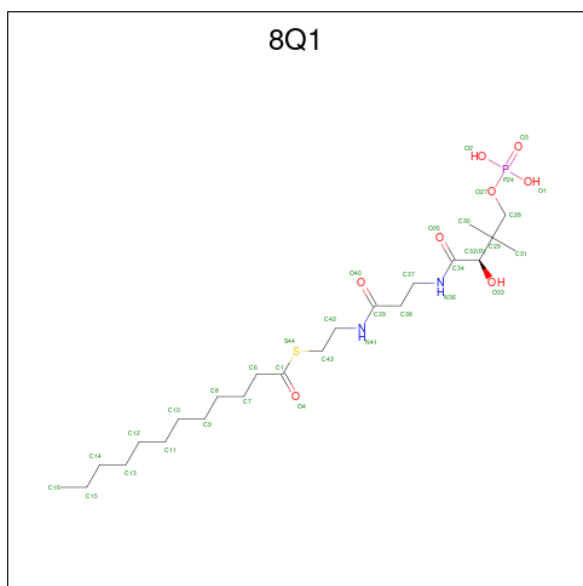
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	A	1	Total	C	O	0	0
			10	6	4		
11	A	1	Total	C	O	0	0
			10	6	4		
11	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 12 is {[-(BIS-CARBOXYMETHYL-AMINO)-ETHYL]-CARBOXYMETHYL-AMINO}-ACETIC ACID (CCD ID: EDT) (formula: $C_{10}H_{16}N_2O_8$).



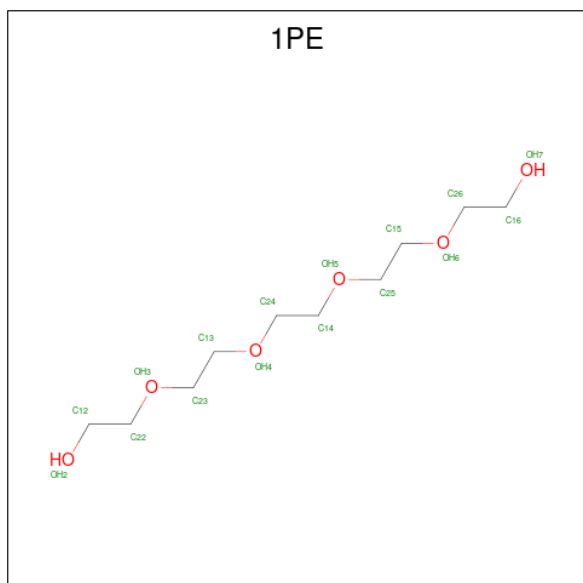
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
12	B	1	Total	C	N	O	0	0
			20	10	2	8		

- Molecule 13 is S-[2-({N-[(2R)-2-hydroxy-3,3-dimethyl-4-(phosphonooxy)butanoyl]-beta-alanyl}amino)ethyl] dodecanethioate (CCD ID: 8Q1) (formula: $C_{23}H_{45}N_2O_8PS$).



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

- Molecule 14 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).



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

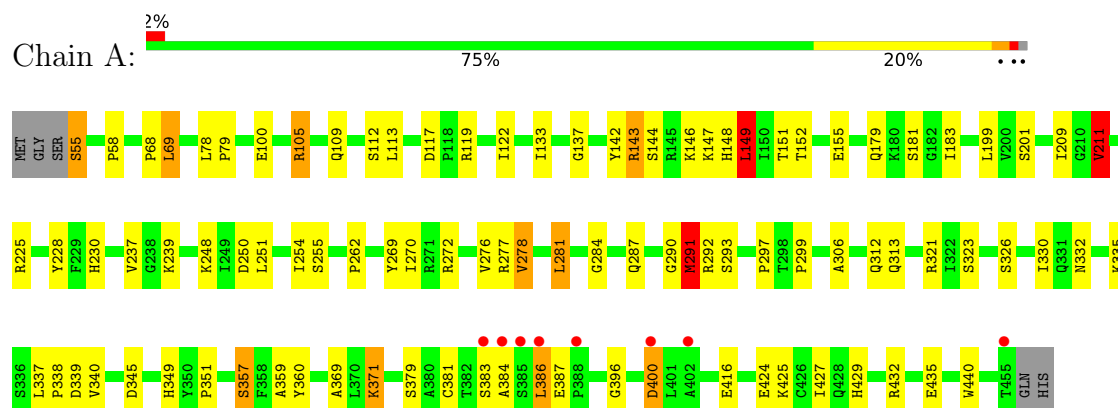
- Molecule 15 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	A	303	Total	O	0	0
			303	303		
15	B	60	Total	O	0	0
			60	60		
15	C	19	Total	O	0	0
			19	19		
15	D	49	Total	O	0	0
			49	49		

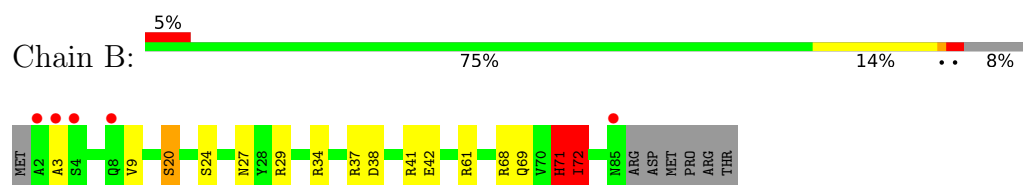
3 Residue-property plots

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

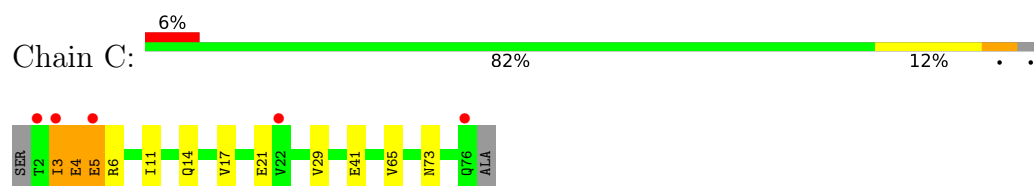
- Molecule 1: Cysteine desulfurase, mitochondrial



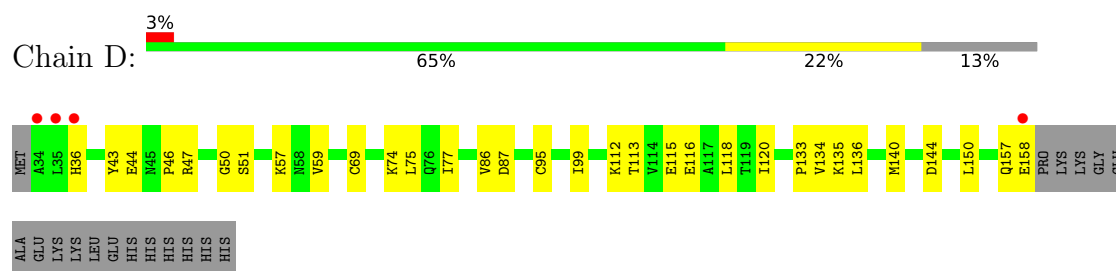
- Molecule 2: LYR motif-containing protein 4



- 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	86.35Å 86.35Å 245.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.97 – 1.90 48.97 – 1.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.97-1.90) 100.0 (48.97-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.33 (at 1.90Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.151 , 0.186 0.163 , 0.188	Depositor DCC
R_{free} test set	3717 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	28.1	Xtriage
Anisotropy	0.189	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 54.4	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	6194	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.29% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, 1PE, EDO, PG4, GOL, 8Q1, PLP, P15, PGE, EDT

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	1.38	25/3218 (0.8%)	1.13	14/4356 (0.3%)
2	B	1.49	8/780 (1.0%)	1.33	12/1045 (1.1%)
3	C	0.58	0/554	0.82	2/754 (0.3%)
4	D	0.82	2/941 (0.2%)	0.82	0/1273
All	All	1.26	35/5493 (0.6%)	1.09	28/7428 (0.4%)

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	8
2	B	0	6
All	All	0	14

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	357[A]	SER	CA-C	12.16	1.67	1.52
1	A	357[B]	SER	CA-C	12.16	1.67	1.52
1	A	211[A]	VAL	CA-C	11.64	1.66	1.52
1	A	211[B]	VAL	CA-C	11.64	1.66	1.52
2	B	24[A]	SER	CA-C	11.53	1.68	1.52
2	B	24[B]	SER	CA-C	11.53	1.68	1.52
1	A	149[A]	LEU	N-CA	8.51	1.56	1.45
1	A	149[B]	LEU	N-CA	8.51	1.56	1.45
2	B	71[A]	HIS	CA-C	8.34	1.64	1.52
2	B	71[B]	HIS	CA-C	8.34	1.64	1.52
2	B	72[A]	ILE	CA-C	8.04	1.64	1.52
2	B	72[B]	ILE	CA-C	8.04	1.64	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	69[A]	LEU	CA-C	7.81	1.63	1.52
1	A	69[B]	LEU	CA-C	7.81	1.63	1.52
1	A	69[A]	LEU	N-CA	7.40	1.55	1.46
1	A	69[B]	LEU	N-CA	7.40	1.55	1.46
2	B	20[A]	SER	N-CA	6.54	1.54	1.46
2	B	20[B]	SER	N-CA	6.54	1.54	1.46
1	A	262	PRO	C-O	-6.36	1.16	1.23
1	A	79	PRO	C-O	-6.05	1.16	1.24
1	A	58	PRO	C-O	-6.00	1.16	1.23
1	A	299	PRO	C-O	-5.88	1.16	1.24
4	D	46	PRO	C-O	-5.86	1.17	1.23
1	A	113	LEU	C-O	-5.41	1.17	1.24
4	D	99	ILE	C-O	-5.33	1.18	1.24
1	A	105[A]	ARG	CA-C	5.33	1.60	1.52
1	A	105[B]	ARG	CA-C	5.33	1.60	1.52
1	A	68	PRO	C-O	-5.32	1.17	1.23
1	A	78	LEU	C-O	-5.29	1.19	1.24
1	A	105[A]	ARG	N-CA	5.19	1.52	1.46
1	A	105[B]	ARG	N-CA	5.19	1.52	1.46
1	A	313	GLN	C-O	-5.10	1.17	1.24
1	A	297	PRO	C-O	-5.09	1.17	1.24
1	A	255	SER	C-O	-5.08	1.18	1.24
1	A	416	GLU	C-O	-5.03	1.18	1.24

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	6	ARG	N-CA-C	-11.29	98.73	111.71
1	A	291[A]	MET	N-CA-C	10.84	125.76	112.54
1	A	291[B]	MET	N-CA-C	10.84	125.76	112.54
2	B	71[A]	HIS	N-CA-C	10.21	123.68	111.33
2	B	71[B]	HIS	N-CA-C	10.21	123.68	111.33
2	B	24[A]	SER	N-CA-C	9.48	122.61	111.71
2	B	24[B]	SER	N-CA-C	9.48	122.61	111.71
1	A	211[A]	VAL	CA-C-O	-9.23	110.32	120.85
1	A	211[B]	VAL	CA-C-O	-9.23	110.32	120.85
2	B	20[A]	SER	N-CA-C	7.31	121.30	112.38
2	B	20[B]	SER	N-CA-C	7.31	121.30	112.38
2	B	72[A]	ILE	N-CA-C	6.74	122.24	113.07
2	B	72[B]	ILE	N-CA-C	6.74	122.24	113.07
1	A	211[A]	VAL	N-CA-C	6.45	117.40	108.89
1	A	211[B]	VAL	N-CA-C	6.45	117.40	108.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	5	GLU	CB-CA-C	6.40	122.05	110.72
2	B	71[A]	HIS	CB-CA-C	-5.88	99.76	110.63
2	B	71[B]	HIS	CB-CA-C	-5.88	99.76	110.63
1	A	149[A]	LEU	CA-C-O	-5.80	114.55	121.11
1	A	149[B]	LEU	CA-C-O	-5.80	114.55	121.11
1	A	149[A]	LEU	N-CA-C	5.65	118.76	109.72
1	A	149[B]	LEU	N-CA-C	5.65	118.76	109.72
2	B	72[A]	ILE	CA-C-O	-5.52	114.51	120.47
2	B	72[B]	ILE	CA-C-O	-5.52	114.51	120.47
1	A	69[A]	LEU	CB-CA-C	-5.47	101.13	109.89
1	A	69[B]	LEU	CB-CA-C	-5.47	101.13	109.89
1	A	105[A]	ARG	N-CA-C	5.21	117.36	111.11
1	A	105[B]	ARG	N-CA-C	5.21	117.36	111.11

There are no chirality outliers.

All (14) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	149[A]	LEU	Mainchain
1	A	149[B]	LEU	Mainchain
1	A	211[A]	VAL	Mainchain
1	A	211[B]	VAL	Mainchain
1	A	357[A]	SER	Mainchain
1	A	357[B]	SER	Mainchain
1	A	69[A]	LEU	Mainchain
1	A	69[B]	LEU	Mainchain
2	B	20[A]	SER	Mainchain
2	B	20[B]	SER	Mainchain
2	B	71[A]	HIS	Mainchain
2	B	71[B]	HIS	Mainchain
2	B	72[A]	ILE	Mainchain
2	B	72[B]	ILE	Mainchain

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	3134	0	3147	138	0
2	B	741	0	790	31	0
3	C	550	0	496	10	0
4	D	926	0	934	38	0
5	A	15	6	7	0	0
6	A	49	0	70	7	0
6	B	14	0	20	20	0
7	A	18	0	24	6	0
7	B	6	0	8	5	0
7	D	6	0	8	0	0
8	A	100	0	146	65	0
8	B	28	0	42	5	0
8	C	12	0	18	3	0
8	D	12	0	18	4	0
9	A	26	0	36	8	0
10	A	20	0	28	15	0
11	A	30	0	42	31	0
12	B	20	0	12	9	0
13	C	34	0	0	0	0
14	D	16	0	22	9	0
15	A	303	0	0	32	0
15	B	60	0	0	6	0
15	C	19	0	0	0	0
15	D	49	0	0	7	0
All	All	6188	6	5868	256	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 22.

All (256) 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:34[A]:ARG:HH21	7:B:402:GOL:H12	1.06	1.16
4:D:112:LYS:HG2	14:D:202:1PE:H222	1.37	1.06
1:A:381:CYS:SG	4:D:135:LYS:NZ	2.32	1.01
2:B:68:ARG:HH11	6:B:407:PEG:H22	1.28	0.97
9:A:517:PG4:H12	15:A:691:HOH:O	1.63	0.97
2:B:34[A]:ARG:NH2	7:B:402:GOL:H12	1.80	0.94
2:B:29:ARG:HH21	12:B:411:EDT:H061	1.35	0.91
1:A:337:LEU:O	10:A:523:P15:H131	1.73	0.88
2:B:41[B]:ARG:HG3	15:B:522:HOH:O	1.75	0.87
1:A:278:VAL:HG22	1:A:291[B]:MET:HE1	1.56	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:386:LEU:HB2	1:A:387:GLU:HG3	1.59	0.85
2:B:68:ARG:HD3	6:B:407:PEG:H22	1.57	0.84
8:A:520:EDO:H22	10:A:523:P15:H42	1.60	0.83
1:A:112:SER:HA	9:A:517:PG4:H11	1.61	0.83
4:D:112:LYS:CG	14:D:202:1PE:H222	2.10	0.82
1:A:425:LYS:HZ1	8:A:526:EDO:H22	1.41	0.82
7:A:503:GOL:O2	8:A:509:EDO:H21	1.81	0.81
1:A:351:PRO:HA	11:A:528:PGE:H2	1.62	0.80
6:A:540:PEG:H42	15:A:737:HOH:O	1.83	0.79
1:A:270:ILE:HD12	1:A:291[A]:MET:HE3	1.64	0.79
1:A:332:ASN:HD21	9:A:518:PG4:H32	1.48	0.79
3:C:29:VAL:HG21	8:C:302:EDO:H11	1.65	0.78
8:A:514:EDO:H21	6:B:405:PEG:H12	1.64	0.78
2:B:42:GLU:HB2	6:B:405:PEG:H11	1.66	0.78
15:A:664:HOH:O	4:D:74:LYS:HE3	1.83	0.77
1:A:425:LYS:NZ	8:A:526:EDO:H22	1.99	0.77
4:D:51:SER:H	8:D:201:EDO:H11	1.50	0.77
1:A:149[A]:LEU:HD12	1:A:199:LEU:HB3	1.65	0.77
1:A:272:ARG:HH12	8:A:521:EDO:H21	1.50	0.76
4:D:95:CYS:SG	4:D:135:LYS:NZ	2.56	0.76
1:A:142:TYR:OH	1:A:228:TYR:OH	1.58	0.76
1:A:248:LYS:HB3	8:A:512:EDO:H21	1.67	0.76
1:A:292:ARG:HH22	8:A:513:EDO:H11	1.48	0.76
4:D:112:LYS:NZ	14:D:202:1PE:H142	2.02	0.75
1:A:371:LYS:HD3	8:A:507:EDO:H22	1.69	0.75
1:A:109:GLN:HE22	11:A:536:PGE:C6	2.01	0.74
4:D:112:LYS:CE	14:D:202:1PE:H162	2.17	0.74
11:A:528:PGE:H6	15:A:606:HOH:O	1.87	0.74
4:D:115:GLU:HG2	15:D:337:HOH:O	1.88	0.74
8:D:205:EDO:H12	15:D:327:HOH:O	1.88	0.74
1:A:321[A]:ARG:NH1	7:B:402:GOL:O2	2.21	0.73
2:B:41[B]:ARG:NH1	3:C:41:GLU:OE2	2.20	0.73
1:A:143:ARG:HD2	8:A:535:EDO:H22	1.71	0.72
4:D:69:CYS:HB2	15:D:338:HOH:O	1.90	0.72
1:A:340:VAL:H	10:A:523:P15:C13	2.03	0.71
2:B:34[B]:ARG:NE	15:B:501:HOH:O	2.23	0.71
2:B:29:ARG:NH2	12:B:411:EDT:H041	2.05	0.71
8:A:507:EDO:H11	4:D:44:GLU:HG2	1.71	0.71
1:A:371:LYS:CD	8:A:507:EDO:H22	2.21	0.70
3:C:4:GLU:HG2	3:C:73:ASN:OD1	1.92	0.70
2:B:68:ARG:HH11	6:B:407:PEG:C2	2.02	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:34[A]:ARG:HH21	7:B:402:GOL:C1	1.96	0.69
1:A:225:ARG:NH1	7:A:505:GOL:O1	2.20	0.69
1:A:292:ARG:HH22	8:A:513:EDO:C1	2.05	0.69
3:C:4:GLU:HA	3:C:4:GLU:OE1	1.92	0.69
1:A:287:GLN:HG3	8:A:539:EDO:H11	1.73	0.68
8:A:520:EDO:H22	10:A:523:P15:C4	2.24	0.68
1:A:432[A]:ARG:NH2	15:A:604:HOH:O	2.25	0.68
9:A:518:PG4:H71	15:A:603:HOH:O	1.92	0.68
2:B:29:ARG:HH21	12:B:411:EDT:H041	1.57	0.68
1:A:332:ASN:ND2	9:A:518:PG4:H32	2.08	0.68
2:B:42:GLU:CB	6:B:405:PEG:H11	2.24	0.68
4:D:77:ILE:CD1	4:D:150:LEU:HD23	2.24	0.67
4:D:50:GLY:HA3	8:D:201:EDO:H12	1.77	0.67
1:A:400:ASP:OD2	10:A:523:P15:H41	1.95	0.67
1:A:351:PRO:HB3	11:A:528:PGE:H3	1.77	0.67
15:A:791:HOH:O	7:B:402:GOL:H2	1.95	0.66
1:A:425:LYS:NZ	8:A:526:EDO:C2	2.59	0.66
8:A:520:EDO:C2	10:A:523:P15:H42	2.25	0.66
1:A:340:VAL:HB	10:A:523:P15:H133	1.76	0.65
1:A:432[A]:ARG:HH12	8:A:506:EDO:H11	1.62	0.65
4:D:51:SER:H	8:D:201:EDO:C1	2.09	0.65
1:A:349:HIS:CE1	11:A:528:PGE:H22	2.30	0.65
4:D:118:LEU:HD13	4:D:150:LEU:HD13	1.78	0.65
4:D:77:ILE:HD13	4:D:150:LEU:HD23	1.80	0.64
4:D:140[B]:MET:HE3	4:D:144:ASP:OD1	1.97	0.64
4:D:112:LYS:NZ	14:D:202:1PE:H162	2.12	0.64
1:A:278:VAL:HG22	1:A:291[B]:MET:CE	2.25	0.63
1:A:400:ASP:OD2	10:A:523:P15:H22	1.99	0.63
1:A:383:SER:O	1:A:386:LEU:HD22	1.99	0.63
1:A:312:GLN:HE21	8:A:522:EDO:H11	1.63	0.63
6:B:407:PEG:O2	8:B:408:EDO:H22	1.98	0.63
4:D:112:LYS:HE3	14:D:202:1PE:H252	1.81	0.62
1:A:119:ARG:HH12	8:A:521:EDO:H12	1.64	0.62
8:A:514:EDO:C2	6:B:405:PEG:H12	2.28	0.62
3:C:17:VAL:HG13	3:C:21:GLU:HB2	1.82	0.61
4:D:112:LYS:CE	14:D:202:1PE:H142	2.30	0.61
1:A:133:ILE:HA	1:A:281:LEU:HD22	1.82	0.61
1:A:109:GLN:HE22	11:A:536:PGE:H62	1.64	0.61
1:A:105[A]:ARG:NH2	11:A:536:PGE:H6	2.17	0.60
1:A:340:VAL:H	10:A:523:P15:H133	1.65	0.60
1:A:119:ARG:HH22	8:A:521:EDO:H12	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:GLN:HE22	11:A:536:PGE:H5	1.66	0.59
1:A:384:ALA:HB3	1:A:386:LEU:HD21	1.84	0.59
1:A:122:ILE:HD11	1:A:291[A]:MET:CE	2.32	0.59
1:A:225:ARG:HH12	7:A:505:GOL:HO1	1.49	0.59
1:A:292:ARG:NH1	8:A:513:EDO:H12	2.17	0.59
1:A:109:GLN:HE22	11:A:536:PGE:C5	2.16	0.59
2:B:29:ARG:NH2	12:B:411:EDT:H061	2.14	0.58
8:A:531:EDO:H11	15:A:773:HOH:O	2.03	0.58
8:A:542:EDO:H22	15:A:606:HOH:O	2.03	0.58
1:A:349:HIS:ND1	11:A:528:PGE:C2	2.66	0.58
1:A:396:GLY:CA	8:A:532:EDO:H22	2.34	0.58
1:A:142:TYR:CE1	1:A:228:TYR:HE2	2.21	0.58
3:C:41:GLU:HB2	8:C:303:EDO:H11	1.86	0.58
1:A:272:ARG:NH1	8:A:521:EDO:H21	2.18	0.58
1:A:371:LYS:HD3	8:A:507:EDO:C2	2.33	0.57
9:A:517:PG4:H22	15:A:630:HOH:O	2.03	0.57
2:B:68:ARG:NH1	6:B:407:PEG:H22	2.09	0.57
1:A:239:LYS:HE2	15:A:726:HOH:O	2.03	0.57
12:B:411:EDT:O18	12:B:411:EDT:H042	2.01	0.57
3:C:17:VAL:CG1	3:C:21:GLU:HB2	2.33	0.57
11:A:537:PGE:H5	15:A:850:HOH:O	2.03	0.57
1:A:292:ARG:NH2	8:A:513:EDO:H11	2.20	0.57
1:A:349:HIS:ND1	11:A:528:PGE:H22	2.18	0.57
1:A:351:PRO:HG3	11:A:528:PGE:H62	1.87	0.56
4:D:118:LEU:HD13	4:D:150:LEU:CD1	2.35	0.56
1:A:105[A]:ARG:NE	15:A:607:HOH:O	2.29	0.56
1:A:183:ILE:HD12	8:A:515:EDO:C2	2.35	0.56
1:A:384:ALA:HB3	1:A:386:LEU:CD2	2.36	0.56
6:A:516:PEG:H22	15:A:879:HOH:O	2.04	0.56
4:D:134:VAL:HG23	15:D:340:HOH:O	2.05	0.55
1:A:323:SER:OG	11:A:528:PGE:H62	2.05	0.55
4:D:112:LYS:HZ1	14:D:202:1PE:H142	1.71	0.55
1:A:292:ARG:NH2	8:A:513:EDO:C1	2.69	0.55
4:D:57:LYS:O	4:D:157:GLN:NE2	2.37	0.55
1:A:326:SER:OG	11:A:528:PGE:H1	2.08	0.54
10:A:523:P15:H61	15:A:720:HOH:O	2.06	0.54
1:A:183:ILE:HD12	8:A:515:EDO:H21	1.88	0.54
1:A:340:VAL:H	10:A:523:P15:H132	1.71	0.54
1:A:424:GLU:OE2	8:A:530:EDO:H12	2.08	0.54
1:A:142:TYR:CE1	1:A:228:TYR:CE2	2.95	0.54
6:A:524:PEG:H32	15:A:714:HOH:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:B:411:EDT:O15	12:B:411:EDT:H111	2.07	0.53
2:B:27[B]:ASN:ND2	8:B:403:EDO:O2	2.30	0.52
2:B:69:GLN:OE1	2:B:72[B]:ILE:HD11	2.10	0.52
7:A:503:GOL:HO2	8:A:509:EDO:H21	1.72	0.51
1:A:340:VAL:O	10:A:523:P15:H122	2.10	0.51
1:A:147:LYS:HD2	15:A:858:HOH:O	2.10	0.51
1:A:429:HIS:NE2	8:A:526:EDO:H12	2.25	0.51
1:A:146:LYS:HE3	15:A:749:HOH:O	2.10	0.50
1:A:396:GLY:HA2	8:A:532:EDO:H22	1.91	0.50
8:A:509:EDO:H12	15:A:752:HOH:O	2.11	0.50
1:A:292:ARG:HH12	8:A:513:EDO:H12	1.76	0.50
1:A:432[A]:ARG:NH1	8:A:506:EDO:H11	2.26	0.50
1:A:425:LYS:HZ2	8:A:526:EDO:C2	2.23	0.50
1:A:55:SER:HB3	15:A:776:HOH:O	2.11	0.49
1:A:384:ALA:HB1	1:A:386:LEU:HD11	1.93	0.49
2:B:38:ASP:HB3	6:B:405:PEG:C2	2.43	0.49
8:A:507:EDO:H12	4:D:43:TYR:CE2	2.47	0.49
1:A:105[B]:ARG:CD	11:A:537:PGE:H1	2.43	0.49
9:A:517:PG4:O1	6:A:541:PEG:H12	2.12	0.49
6:B:407:PEG:C2	8:B:408:EDO:H22	2.43	0.49
3:C:14:GLN:O	8:C:303:EDO:H21	2.13	0.49
1:A:181:SER:CB	8:A:515:EDO:H21	2.42	0.49
1:A:122:ILE:CD1	1:A:291[A]:MET:HE3	2.43	0.48
1:A:312:GLN:NE2	8:A:522:EDO:H11	2.28	0.48
1:A:386:LEU:HD13	1:A:386:LEU:N	2.28	0.48
10:A:523:P15:H111	15:A:741:HOH:O	2.11	0.48
2:B:68:ARG:HG3	6:B:407:PEG:H31	1.95	0.48
1:A:384:ALA:CB	1:A:386:LEU:HD11	2.43	0.48
1:A:144:SER:HB3	8:A:535:EDO:H21	1.95	0.48
1:A:105[B]:ARG:HD2	11:A:537:PGE:H1	1.95	0.48
1:A:137:GLY:HA3	1:A:278:VAL:HG23	1.94	0.48
1:A:312:GLN:HE21	8:A:522:EDO:C1	2.26	0.47
11:A:528:PGE:H5	15:A:601:HOH:O	2.14	0.47
1:A:359:ALA:HA	8:A:520:EDO:O1	2.14	0.47
8:A:507:EDO:H22	4:D:44:GLU:OE2	2.14	0.47
2:B:42:GLU:CG	6:B:405:PEG:H11	2.44	0.47
1:A:432[A]:ARG:HH12	8:A:506:EDO:C1	2.26	0.47
9:A:517:PG4:H31	9:A:517:PG4:H52	1.39	0.47
12:B:411:EDT:O15	12:B:411:EDT:C11	2.63	0.47
1:A:181:SER:HB2	1:A:345:ASP:HB2	1.97	0.47
1:A:292:ARG:CZ	8:A:513:EDO:H12	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:38:ASP:HB3	6:B:405:PEG:H22	1.96	0.47
1:A:250:ASP:HB3	1:A:276:VAL:HG21	1.96	0.47
8:B:406:EDO:H11	6:B:407:PEG:C4	2.44	0.47
1:A:360:TYR:CD2	1:A:440:TRP:HE3	2.33	0.46
11:A:536:PGE:H12	2:B:61[B]:ARG:NH1	2.30	0.46
1:A:435:GLU:HG3	7:A:504:GOL:H32	1.98	0.45
1:A:371:LYS:HD2	8:A:507:EDO:H22	1.95	0.45
1:A:432[B]:ARG:HE	8:A:534:EDO:H21	1.81	0.45
11:A:536:PGE:H2	15:B:549:HOH:O	2.16	0.45
4:D:113:THR:OG1	4:D:116:GLU:HG3	2.16	0.45
4:D:134:VAL:HG12	4:D:134:VAL:O	2.17	0.45
1:A:209:ILE:HG13	1:A:211[B]:VAL:HG13	1.99	0.45
1:A:306:ALA:CB	11:A:536:PGE:H52	2.46	0.45
2:B:29:ARG:HG2	12:B:411:EDT:H091	1.99	0.45
4:D:75:LEU:HD23	4:D:75:LEU:C	2.42	0.45
4:D:133:PRO:HA	4:D:136:LEU:HG	1.98	0.45
1:A:335:LYS:NZ	15:A:603:HOH:O	2.49	0.45
11:A:536:PGE:H62	15:A:687:HOH:O	2.16	0.45
1:A:321[B]:ARG:CZ	2:B:34[B]:ARG:NH2	2.80	0.44
1:A:338:PRO:HD3	8:A:508:EDO:H11	1.99	0.44
1:A:424:GLU:CD	8:A:530:EDO:H12	2.41	0.44
2:B:34[B]:ARG:HG2	15:B:501:HOH:O	2.17	0.44
1:A:117:ASP:HA	6:A:541:PEG:H31	2.00	0.44
1:A:349:HIS:CE1	11:A:528:PGE:C2	2.99	0.44
1:A:151:THR:OG1	1:A:155:GLU:HG3	2.17	0.44
1:A:105[A]:ARG:CD	11:A:537:PGE:H1	2.48	0.44
1:A:148:HIS:C	1:A:149[A]:LEU:HD22	2.41	0.43
1:A:381:CYS:SG	4:D:135:LYS:CE	3.06	0.43
12:B:411:EDT:H062	12:B:411:EDT:H092	1.85	0.43
11:A:528:PGE:O1	15:A:601:HOH:O	2.21	0.43
2:B:37[B]:ARG:NH2	15:B:509:HOH:O	2.51	0.43
1:A:277:ARG:NH2	15:A:612:HOH:O	2.48	0.43
1:A:292:ARG:NH2	8:A:513:EDO:H12	2.33	0.43
1:A:330:ILE:CD1	11:A:528:PGE:H1	2.49	0.43
1:A:119:ARG:NH2	8:A:521:EDO:H12	2.33	0.43
1:A:371:LYS:HB3	8:A:507:EDO:H21	1.99	0.43
1:A:369:ALA:HA	8:A:507:EDO:H12	2.01	0.43
4:D:120:ILE:HG12	14:D:202:1PE:C25	2.49	0.43
1:A:142:TYR:HE1	1:A:228:TYR:HE2	1.65	0.43
2:B:68:ARG:CG	6:B:407:PEG:H31	2.49	0.43
4:D:133:PRO:HD2	15:D:340:HOH:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:118:LEU:CD1	4:D:150:LEU:HD12	2.49	0.43
1:A:330:ILE:HD12	11:A:528:PGE:H1	2.00	0.42
1:A:332:ASN:OD1	1:A:335:LYS:HE3	2.19	0.42
6:B:407:PEG:O2	8:B:408:EDO:C2	2.65	0.42
1:A:290:GLY:O	6:A:540:PEG:H21	2.18	0.42
6:B:407:PEG:H12	6:B:407:PEG:H32	1.34	0.42
1:A:119:ARG:NH1	8:A:521:EDO:H12	2.31	0.42
1:A:122:ILE:CD1	1:A:291[A]:MET:CE	2.97	0.42
11:A:536:PGE:H5	11:A:536:PGE:H32	1.84	0.42
1:A:105[A]:ARG:HH22	11:A:536:PGE:H6	1.84	0.42
1:A:201:SER:HA	1:A:230:HIS:O	2.20	0.42
1:A:237:VAL:HG21	1:A:254:ILE:HG12	2.01	0.42
1:A:277:ARG:NE	15:A:612:HOH:O	2.33	0.42
3:C:3:ILE:HG23	3:C:5:GLU:H	1.85	0.42
1:A:100:GLU:HG2	8:A:511:EDO:O2	2.20	0.42
1:A:122:ILE:N	1:A:122:ILE:HD12	2.35	0.41
4:D:36:HIS:CD2	15:D:334:HOH:O	2.72	0.41
4:D:87:ASP:OD2	15:D:301:HOH:O	2.22	0.41
1:A:179:GLN:NE2	15:A:610:HOH:O	2.33	0.41
1:A:337:LEU:HD11	1:A:427:ILE:HA	2.03	0.41
2:B:68:ARG:HD3	6:B:407:PEG:C2	2.40	0.41
1:A:384:ALA:C	1:A:386:LEU:HD13	2.45	0.41
1:A:284:GLY:HA3	1:A:293:SER:HB3	2.02	0.41
2:B:71[B]:HIS:CE1	15:B:510:HOH:O	2.74	0.41
1:A:251:LEU:HA	1:A:269:TYR:O	2.20	0.41
8:A:538:EDO:H22	15:A:710:HOH:O	2.20	0.41
1:A:109:GLN:NE2	11:A:536:PGE:H5	2.34	0.41
1:A:330:ILE:HD12	11:A:528:PGE:C1	2.50	0.41
1:A:340:VAL:N	10:A:523:P15:H133	2.34	0.41
8:A:531:EDO:C2	15:A:773:HOH:O	2.69	0.41
6:A:540:PEG:H22	15:A:822:HOH:O	2.20	0.41
15:A:664:HOH:O	4:D:74:LYS:CE	2.55	0.41
1:A:119:ARG:HH22	8:A:521:EDO:C1	2.32	0.40
8:A:514:EDO:O1	6:B:405:PEG:H31	2.22	0.40
3:C:11:ILE:HD13	3:C:65:VAL:HG22	2.02	0.40
1:A:339:ASP:N	10:A:523:P15:H132	2.36	0.40
1:A:360:TYR:N	8:A:520:EDO:O1	2.39	0.40
7:A:503:GOL:O2	8:A:509:EDO:C2	2.61	0.40
4:D:59:VAL:HA	4:D:77:ILE:O	2.22	0.40

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	408/406 (100%)	399 (98%)	9 (2%)	0	100	100
2	B	92/91 (101%)	89 (97%)	2 (2%)	1 (1%)	11	4
3	C	73/77 (95%)	69 (94%)	3 (4%)	1 (1%)	9	3
4	D	124/143 (87%)	123 (99%)	1 (1%)	0	100	100
All	All	697/717 (97%)	680 (98%)	15 (2%)	2 (0%)	36	29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	3	ALA
3	C	3	ILE

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	342/346 (99%)	329 (96%)	13 (4%)	29	21
2	B	78/80 (98%)	78 (100%)	0	100	100
3	C	54/66 (82%)	53 (98%)	1 (2%)	50	47
4	D	98/118 (83%)	95 (97%)	3 (3%)	35	29
All	All	572/610 (94%)	555 (97%)	17 (3%)	39	30

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	SER
1	A	143	ARG
1	A	149[A]	LEU
1	A	149[B]	LEU
1	A	152	THR
1	A	278	VAL
1	A	281	LEU
1	A	291[A]	MET
1	A	291[B]	MET
1	A	371	LYS
1	A	379	SER
1	A	386	LEU
1	A	400	ASP
3	C	4	GLU
4	D	47	ARG
4	D	86	VAL
4	D	158	GLU

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	194	GLN
1	A	312	GLN
1	A	343	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

62 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	PEG	A	541	-	6,6,6	0.49	0	5,5,5	0.47	0
8	EDO	A	520	-	3,3,3	0.63	0	2,2,2	1.09	0
13	8Q1	C	301	3	28,33,34	2.18	8 (28%)	32,40,43	1.71	7 (21%)
8	EDO	A	529	-	3,3,3	0.49	0	2,2,2	1.00	0
11	PGE	A	528	-	9,9,9	0.50	0	8,8,8	0.71	0
8	EDO	D	205	-	3,3,3	0.44	0	2,2,2	0.09	0
8	EDO	A	522	-	3,3,3	0.43	0	2,2,2	0.24	0
8	EDO	A	525	-	3,3,3	0.44	0	2,2,2	0.31	0
8	EDO	A	506	-	3,3,3	0.47	0	2,2,2	0.14	0
8	EDO	A	507	-	3,3,3	1.53	0	2,2,2	1.60	1 (50%)
7	GOL	B	402	-	5,5,5	0.53	0	5,5,5	0.87	0
8	EDO	A	510	-	3,3,3	0.43	0	2,2,2	0.16	0
10	P15	A	523	-	19,19,19	0.63	0	18,18,18	0.47	0
8	EDO	A	542	-	3,3,3	0.50	0	2,2,2	0.23	0
11	PGE	A	537	-	9,9,9	0.38	0	8,8,8	0.26	0
6	PEG	B	407	-	6,6,6	0.73	0	5,5,5	0.77	0
6	PEG	A	540	-	6,6,6	0.53	0	5,5,5	0.36	0
8	EDO	B	410	-	3,3,3	0.47	0	2,2,2	0.60	0
7	GOL	A	505	-	5,5,5	0.14	0	5,5,5	0.44	0
8	EDO	B	409	-	3,3,3	0.42	0	2,2,2	0.26	0
8	EDO	A	513	-	3,3,3	1.06	0	2,2,2	0.86	0
6	PEG	A	524	-	6,6,6	0.53	0	5,5,5	0.52	0
7	GOL	D	204	-	5,5,5	0.63	0	5,5,5	0.32	0
8	EDO	B	408	-	3,3,3	0.42	0	2,2,2	0.75	0
8	EDO	A	534	-	3,3,3	0.46	0	2,2,2	0.10	0
6	PEG	B	405	-	6,6,6	0.57	0	5,5,5	0.55	0
6	PEG	A	519	-	6,6,6	0.50	0	5,5,5	0.63	0
8	EDO	D	203	-	3,3,3	0.46	0	2,2,2	0.34	0
8	EDO	A	539	-	3,3,3	0.46	0	2,2,2	0.46	0
11	PGE	A	536	-	9,9,9	0.33	0	8,8,8	0.48	0
6	PEG	A	516	-	6,6,6	0.51	0	5,5,5	0.70	0
6	PEG	A	502	-	6,6,6	0.61	0	5,5,5	0.78	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	EDO	A	521	-	3,3,3	0.42	0	2,2,2	0.62	0
8	EDO	A	509	-	3,3,3	0.51	0	2,2,2	0.41	0
7	GOL	A	503	-	5,5,5	0.63	0	5,5,5	1.15	0
5	PLP	A	501	-	15,15,16	1.46	1 (6%)	21,22,23	0.91	1 (4%)
8	EDO	C	303	-	3,3,3	0.41	0	2,2,2	0.19	0
8	EDO	C	304	-	3,3,3	0.43	0	2,2,2	0.33	0
8	EDO	A	515	-	3,3,3	0.72	0	2,2,2	0.94	0
8	EDO	D	201	-	3,3,3	0.46	0	2,2,2	0.49	0
9	PG4	A	517	-	12,12,12	0.56	0	11,11,11	0.27	0
14	1PE	D	202	-	15,15,15	0.53	0	14,14,14	0.30	0
8	EDO	A	526	-	3,3,3	0.45	0	2,2,2	0.36	0
8	EDO	A	531	-	3,3,3	0.48	0	2,2,2	0.29	0
8	EDO	B	406	-	3,3,3	0.47	0	2,2,2	0.24	0
8	EDO	A	530	-	3,3,3	0.42	0	2,2,2	0.19	0
8	EDO	B	401	-	3,3,3	0.42	0	2,2,2	0.66	0
6	PEG	A	527	-	6,6,6	0.51	0	5,5,5	0.22	0
8	EDO	A	533	-	3,3,3	0.44	0	2,2,2	0.36	0
9	PG4	A	518	-	12,12,12	0.57	0	11,11,11	0.37	0
7	GOL	A	504	-	5,5,5	0.56	0	5,5,5	0.25	0
8	EDO	B	403	-	3,3,3	0.73	0	2,2,2	1.07	0
8	EDO	A	508	-	3,3,3	0.47	0	2,2,2	0.32	0
8	EDO	A	538	-	3,3,3	0.46	0	2,2,2	0.41	0
8	EDO	A	511	-	3,3,3	0.29	0	2,2,2	0.40	0
8	EDO	A	535	-	3,3,3	0.43	0	2,2,2	0.36	0
8	EDO	B	404	-	3,3,3	0.51	0	2,2,2	0.62	0
8	EDO	C	302	-	3,3,3	0.49	0	2,2,2	0.33	0
8	EDO	A	512	-	3,3,3	0.44	0	2,2,2	0.32	0
8	EDO	A	532	-	3,3,3	0.42	0	2,2,2	0.14	0
12	EDT	B	411	-	19,19,19	1.17	0	24,24,24	1.26	3 (12%)
8	EDO	A	514	-	3,3,3	0.51	0	2,2,2	0.17	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	PEG	A	541	-	-	2/4/4/4	-
8	EDO	A	520	-	-	0/1/1/1	-
13	8Q1	C	301	3	-	1/38/40/41	-
8	EDO	A	529	-	-	1/1/1/1	-
11	PGE	A	528	-	-	5/7/7/7	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	EDO	D	205	-	-	1/1/1/1	-
8	EDO	A	522	-	-	1/1/1/1	-
8	EDO	A	525	-	-	1/1/1/1	-
8	EDO	A	506	-	-	0/1/1/1	-
8	EDO	A	507	-	-	0/1/1/1	-
7	GOL	B	402	-	-	4/4/4/4	-
8	EDO	A	510	-	-	0/1/1/1	-
10	P15	A	523	-	-	9/17/17/17	-
8	EDO	A	542	-	-	1/1/1/1	-
11	PGE	A	537	-	-	4/7/7/7	-
6	PEG	B	407	-	-	3/4/4/4	-
6	PEG	A	540	-	-	4/4/4/4	-
8	EDO	B	410	-	-	0/1/1/1	-
7	GOL	A	505	-	-	2/4/4/4	-
8	EDO	B	409	-	-	0/1/1/1	-
8	EDO	A	513	-	-	1/1/1/1	-
6	PEG	A	524	-	-	1/4/4/4	-
7	GOL	D	204	-	-	0/4/4/4	-
8	EDO	B	408	-	-	1/1/1/1	-
8	EDO	A	534	-	-	1/1/1/1	-
6	PEG	B	405	-	-	1/4/4/4	-
6	PEG	A	519	-	-	2/4/4/4	-
8	EDO	D	203	-	-	0/1/1/1	-
8	EDO	A	539	-	-	0/1/1/1	-
11	PGE	A	536	-	-	3/7/7/7	-
6	PEG	A	516	-	-	1/4/4/4	-
6	PEG	A	502	-	-	2/4/4/4	-
8	EDO	A	521	-	-	1/1/1/1	-
8	EDO	A	509	-	-	1/1/1/1	-
7	GOL	A	503	-	-	1/4/4/4	-
5	PLP	A	501	-	-	0/6/6/8	0/1/1/1
8	EDO	C	303	-	-	1/1/1/1	-
8	EDO	C	304	-	-	0/1/1/1	-
8	EDO	A	515	-	-	1/1/1/1	-
8	EDO	D	201	-	-	0/1/1/1	-
9	PG4	A	517	-	-	6/10/10/10	-
14	1PE	D	202	-	-	9/13/13/13	-
8	EDO	A	526	-	-	0/1/1/1	-
8	EDO	A	531	-	-	1/1/1/1	-
8	EDO	B	406	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	EDO	A	530	-	-	1/1/1/1	-
8	EDO	B	401	-	-	1/1/1/1	-
6	PEG	A	527	-	-	3/4/4/4	-
8	EDO	A	533	-	-	0/1/1/1	-
9	PG4	A	518	-	-	5/10/10/10	-
7	GOL	A	504	-	-	4/4/4/4	-
8	EDO	B	403	-	-	1/1/1/1	-
8	EDO	A	508	-	-	0/1/1/1	-
8	EDO	A	538	-	-	1/1/1/1	-
8	EDO	A	511	-	-	1/1/1/1	-
8	EDO	A	535	-	-	0/1/1/1	-
8	EDO	B	404	-	-	0/1/1/1	-
8	EDO	C	302	-	-	1/1/1/1	-
8	EDO	A	512	-	-	1/1/1/1	-
8	EDO	A	532	-	-	0/1/1/1	-
12	EDT	B	411	-	-	11/21/21/21	-
8	EDO	A	514	-	-	0/1/1/1	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
13	C	301	8Q1	C39-N41	5.41	1.46	1.33
13	C	301	8Q1	C34-N36	4.99	1.45	1.33
13	C	301	8Q1	C6-C1	3.61	1.54	1.50
5	A	501	PLP	C3-C2	-3.40	1.37	1.41
13	C	301	8Q1	C12-C13	-2.87	1.37	1.51
13	C	301	8Q1	C15-C14	-2.61	1.35	1.51
13	C	301	8Q1	C28-C29	2.57	1.56	1.52
13	C	301	8Q1	C38-C39	2.39	1.56	1.51
13	C	301	8Q1	O40-C39	-2.23	1.18	1.23

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	C	301	8Q1	C37-N36-C34	-4.48	114.49	122.55
13	C	301	8Q1	C42-N41-C39	-3.95	115.47	122.82
12	B	411	EDT	C11-N8-C9	2.96	118.59	111.66
13	C	301	8Q1	C38-C37-N36	-2.91	105.81	112.00
13	C	301	8Q1	O40-C39-N41	-2.62	117.89	123.03
13	C	301	8Q1	C38-C39-N41	2.17	120.30	116.34
13	C	301	8Q1	O4-C1-S44	2.12	125.38	122.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	C	301	8Q1	C8-C7-C6	-2.10	105.40	113.13
12	B	411	EDT	C7-C6-N3	-2.07	108.07	112.98
12	B	411	EDT	C11-N8-C7	2.05	116.86	111.91
8	A	507	EDO	O2-C2-C1	-2.04	96.82	112.39
5	A	501	PLP	C2A-C2-C3	-2.04	118.41	120.80

There are no chirality outliers.

All (103) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	504	GOL	O1-C1-C2-O2
7	A	504	GOL	O1-C1-C2-C3
7	A	504	GOL	C1-C2-C3-O3
7	A	504	GOL	O2-C2-C3-O3
7	A	505	GOL	C1-C2-C3-O3
7	B	402	GOL	C1-C2-C3-O3
12	B	411	EDT	C10-C9-N8-C11
12	B	411	EDT	N8-C11-C12-O13
12	B	411	EDT	N8-C11-C12-O14
9	A	517	PG4	C3-C4-O3-C5
6	A	502	PEG	C4-C3-O2-C2
6	B	407	PEG	C1-C2-O2-C3
12	B	411	EDT	C1-C2-N3-C4
9	A	517	PG4	O2-C3-C4-O3
10	A	523	P15	O5-C10-C9-O4
10	A	523	P15	O5-C11-C12-O6
9	A	517	PG4	O3-C5-C6-O4
14	D	202	1PE	OH4-C13-C23-OH3
10	A	523	P15	O2-C5-C6-O3
10	A	523	P15	C3-C4-O2-C5
9	A	518	PG4	O2-C3-C4-O3
9	A	518	PG4	O3-C5-C6-O4
10	A	523	P15	O3-C7-C8-O4
11	A	528	PGE	O2-C3-C4-O3
7	B	402	GOL	O1-C1-C2-O2
6	A	502	PEG	O1-C1-C2-O2
6	A	540	PEG	O1-C1-C2-O2
11	A	536	PGE	C3-C4-O3-C5
6	A	516	PEG	O1-C1-C2-O2
6	A	519	PEG	O2-C3-C4-O4
6	A	527	PEG	O2-C3-C4-O4
6	A	540	PEG	O2-C3-C4-O4

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Mol	Chain	Res	Type	Atoms
6	A	541	PEG	O1-C1-C2-O2
6	B	407	PEG	O1-C1-C2-O2
7	B	402	GOL	O1-C1-C2-C3
11	A	528	PGE	O3-C5-C6-O4
7	A	505	GOL	O2-C2-C3-O3
14	D	202	1PE	OH6-C15-C25-OH5
9	A	517	PG4	O4-C7-C8-O5
6	A	527	PEG	O1-C1-C2-O2
9	A	518	PG4	O1-C1-C2-O2
7	B	402	GOL	O2-C2-C3-O3
11	A	528	PGE	C3-C4-O3-C5
10	A	523	P15	OXT-C1-C2-O1
8	A	509	EDO	O1-C1-C2-O2
8	A	512	EDO	O1-C1-C2-O2
8	A	522	EDO	O1-C1-C2-O2
8	A	529	EDO	O1-C1-C2-O2
8	A	530	EDO	O1-C1-C2-O2
8	A	538	EDO	O1-C1-C2-O2
8	C	303	EDO	O1-C1-C2-O2
11	A	536	PGE	O2-C3-C4-O3
7	A	503	GOL	O2-C2-C3-O3
12	B	411	EDT	C7-C6-N3-C2
8	A	511	EDO	O1-C1-C2-O2
6	A	524	PEG	C1-C2-O2-C3
6	A	540	PEG	C1-C2-O2-C3
6	A	527	PEG	C1-C2-O2-C3
10	A	523	P15	C10-C9-O4-C8
9	A	518	PG4	C1-C2-O2-C3
6	A	540	PEG	C4-C3-O2-C2
9	A	518	PG4	C5-C6-O4-C7
11	A	537	PGE	O3-C5-C6-O4
12	B	411	EDT	C7-C6-N3-C4
11	A	528	PGE	C1-C2-O2-C3
14	D	202	1PE	C25-C15-OH6-C26
10	A	523	P15	C8-C7-O3-C6
14	D	202	1PE	C15-C25-OH5-C14
14	D	202	1PE	C16-C26-OH6-C15
14	D	202	1PE	C14-C24-OH4-C13
14	D	202	1PE	C13-C23-OH3-C22
8	A	515	EDO	O1-C1-C2-O2
8	A	531	EDO	O1-C1-C2-O2
8	B	403	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
9	A	517	PG4	C8-C7-O4-C6
12	B	411	EDT	C10-C9-N8-C7
11	A	536	PGE	O1-C1-C2-O2
6	B	405	PEG	C1-C2-O2-C3
9	A	517	PG4	C4-C3-O2-C2
8	B	408	EDO	O1-C1-C2-O2
10	A	523	P15	O1-C3-C4-O2
6	B	407	PEG	C4-C3-O2-C2
12	B	411	EDT	O15-C10-C9-N8
11	A	528	PGE	C6-C5-O3-C4
8	A	513	EDO	O1-C1-C2-O2
8	A	525	EDO	O1-C1-C2-O2
8	A	534	EDO	O1-C1-C2-O2
8	A	542	EDO	O1-C1-C2-O2
8	B	401	EDO	O1-C1-C2-O2
8	C	302	EDO	O1-C1-C2-O2
8	D	205	EDO	O1-C1-C2-O2
12	B	411	EDT	C6-C7-N8-C9
14	D	202	1PE	C24-C14-OH5-C25
6	A	541	PEG	C1-C2-O2-C3
12	B	411	EDT	O16-C10-C9-N8
11	A	537	PGE	C1-C2-O2-C3
11	A	537	PGE	C3-C4-O3-C5
14	D	202	1PE	OH5-C14-C24-OH4
8	A	521	EDO	O1-C1-C2-O2
6	A	519	PEG	O1-C1-C2-O2
12	B	411	EDT	O18-C1-C2-N3
13	C	301	8Q1	C6-C7-C8-C9
11	A	537	PGE	O2-C3-C4-O3

There are no ring outliers.

46 monomers are involved in 173 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	541	PEG	2	0
8	A	520	EDO	5	0
11	A	528	PGE	15	0
8	D	205	EDO	1	0
8	A	522	EDO	3	0
8	A	506	EDO	3	0
8	A	507	EDO	9	0
7	B	402	GOL	5	0

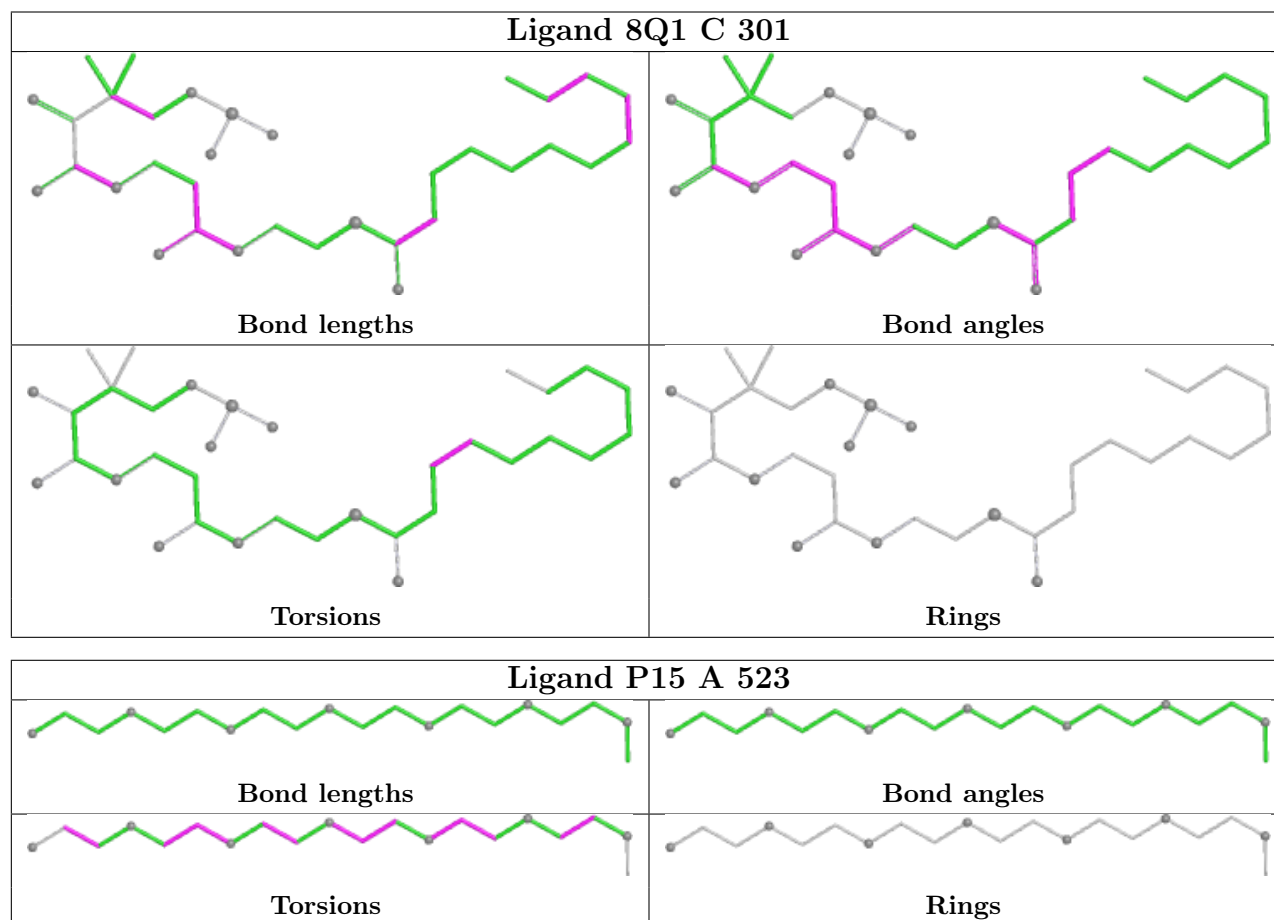
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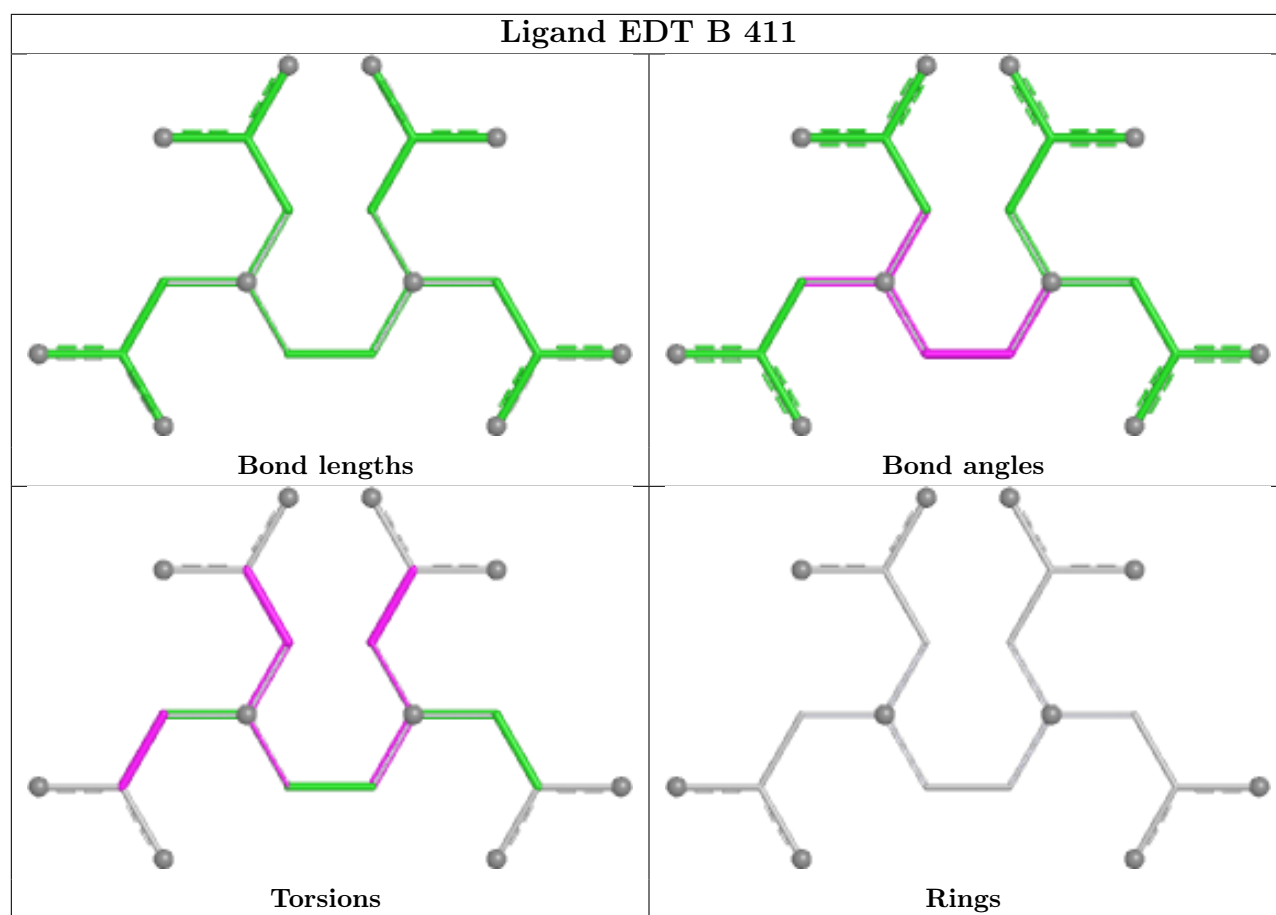
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	A	523	P15	15	0
8	A	542	EDO	1	0
11	A	537	PGE	4	0
6	B	407	PEG	12	0
6	A	540	PEG	3	0
7	A	505	GOL	2	0
8	A	513	EDO	8	0
6	A	524	PEG	1	0
8	B	408	EDO	3	0
8	A	534	EDO	1	0
6	B	405	PEG	8	0
8	A	539	EDO	1	0
11	A	536	PGE	12	0
6	A	516	PEG	1	0
8	A	521	EDO	7	0
8	A	509	EDO	4	0
7	A	503	GOL	3	0
8	C	303	EDO	2	0
8	A	515	EDO	3	0
8	D	201	EDO	3	0
9	A	517	PG4	5	0
14	D	202	1PE	9	0
8	A	526	EDO	5	0
8	A	531	EDO	2	0
8	B	406	EDO	1	0
8	A	530	EDO	2	0
9	A	518	PG4	3	0
7	A	504	GOL	1	0
8	B	403	EDO	1	0
8	A	508	EDO	1	0
8	A	538	EDO	1	0
8	A	511	EDO	1	0
8	A	535	EDO	2	0
8	C	302	EDO	1	0
8	A	512	EDO	1	0
8	A	532	EDO	2	0
12	B	411	EDT	9	0
8	A	514	EDO	3	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

6.1 Protein, DNA and RNA chains

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	401/406 (98%)	-0.32	8 (1%) 65 69	12, 24, 51, 95	19 (4%)
2	B	84/91 (92%)	0.02	5 (5%) 27 29	12, 26, 53, 108	10 (11%)
3	C	75/77 (97%)	0.65	5 (6%) 24 25	34, 62, 98, 117	0
4	D	125/143 (87%)	0.03	4 (3%) 50 54	23, 41, 68, 107	2 (1%)
All	All	685/717 (95%)	-0.11	22 (3%) 50 54	12, 29, 72, 117	31 (4%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	386	LEU	8.0
1	A	384	ALA	7.0
2	B	2	ALA	6.5
2	B	3	ALA	6.3
1	A	455	THR	5.3
4	D	35	LEU	5.2
1	A	383	SER	4.6
3	C	3	ILE	4.2
1	A	385	SER	4.1
3	C	2	THR	3.8
3	C	76	GLN	3.7
4	D	34	ALA	3.6
1	A	402	ALA	2.6
2	B	85	ASN	2.6
3	C	5	GLU	2.5
4	D	158	GLU	2.5
4	D	36	HIS	2.3
3	C	22	VAL	2.3
1	A	388	PRO	2.2
2	B	4	SER	2.2
1	A	400	ASP	2.2
2	B	8	GLN	2.1

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
8	EDO	A	521	4/4	0.50	0.23	58,59,60,63	4
8	EDO	B	401	4/4	0.66	0.20	83,84,84,85	0
8	EDO	A	522	4/4	0.69	0.25	85,86,87,87	0
8	EDO	A	535	4/4	0.71	0.20	59,62,67,72	4
8	EDO	A	533	4/4	0.71	0.19	79,79,79,79	0
8	EDO	B	404	4/4	0.73	0.20	58,63,63,66	0
9	PG4	A	518	13/13	0.73	0.20	81,87,91,91	0
12	EDT	B	411	20/20	0.74	0.23	53,63,78,79	20
6	PEG	B	405	7/7	0.75	0.22	48,60,74,76	0
8	EDO	A	534	4/4	0.76	0.21	70,73,75,75	0
8	EDO	A	525	4/4	0.77	0.24	91,92,92,94	0
8	EDO	D	205	4/4	0.77	0.19	75,77,77,77	0
8	EDO	A	539	4/4	0.77	0.18	54,60,63,63	0
8	EDO	A	526	4/4	0.77	0.21	67,68,70,71	0
6	PEG	A	541	7/7	0.78	0.22	73,82,88,90	0
9	PG4	A	517	13/13	0.78	0.18	51,59,63,63	13
7	GOL	D	204	6/6	0.79	0.17	41,49,54,55	6
6	PEG	A	524	7/7	0.79	0.21	62,66,67,70	0
7	GOL	B	402	6/6	0.79	0.16	37,61,62,65	0
8	EDO	B	408	4/4	0.80	0.18	29,32,33,45	4
8	EDO	D	203	4/4	0.80	0.19	74,77,78,80	0
6	PEG	A	502	7/7	0.80	0.19	41,53,66,67	0
8	EDO	A	538	4/4	0.81	0.21	67,74,77,79	0
8	EDO	A	514	4/4	0.81	0.20	61,63,66,66	0
8	EDO	B	410	4/4	0.81	0.18	66,67,68,70	0
11	PGE	A	537	10/10	0.81	0.17	81,86,88,88	0
8	EDO	A	511	4/4	0.81	0.18	33,33,34,37	4

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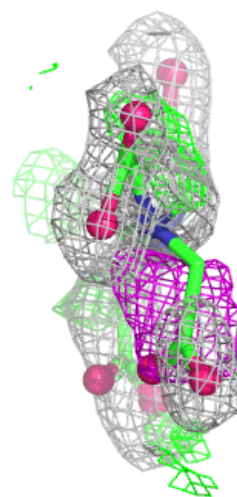
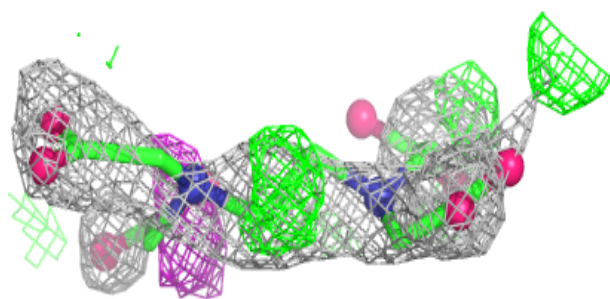
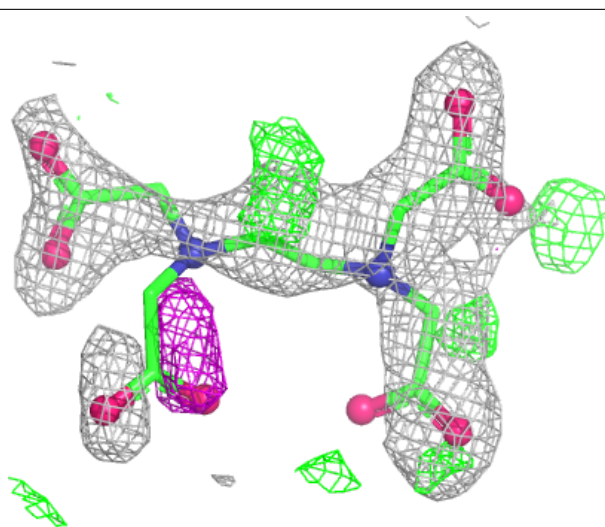
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	PEG	B	407	7/7	0.82	0.17	32,36,48,51	7
8	EDO	A	542	4/4	0.82	0.15	52,55,59,60	0
7	GOL	A	504	6/6	0.82	0.16	59,74,76,79	0
8	EDO	A	530	4/4	0.83	0.16	75,76,76,78	0
8	EDO	B	406	4/4	0.83	0.21	65,69,69,70	0
8	EDO	A	532	4/4	0.84	0.15	73,73,73,73	0
8	EDO	B	409	4/4	0.84	0.12	74,74,75,77	0
8	EDO	A	506	4/4	0.84	0.16	51,55,59,61	0
8	EDO	A	529	4/4	0.85	0.18	34,38,38,40	4
10	P15	A	523	20/20	0.85	0.19	28,77,81,82	20
8	EDO	C	303	4/4	0.85	0.21	38,45,45,48	4
8	EDO	C	304	4/4	0.85	0.16	66,71,73,74	0
6	PEG	A	540	7/7	0.86	0.14	45,62,66,67	0
6	PEG	A	527	7/7	0.86	0.15	64,70,72,72	0
8	EDO	D	201	4/4	0.86	0.13	48,50,52,53	0
11	PGE	A	536	10/10	0.86	0.17	44,57,64,64	10
8	EDO	A	531	4/4	0.86	0.21	72,75,77,77	0
8	EDO	A	508	4/4	0.86	0.13	51,55,57,58	0
6	PEG	A	516	7/7	0.87	0.14	49,53,55,59	0
7	GOL	A	505	6/6	0.87	0.15	32,36,40,40	6
6	PEG	A	519	7/7	0.87	0.14	54,55,58,61	0
8	EDO	A	510	4/4	0.88	0.15	62,64,64,66	0
8	EDO	A	520	4/4	0.88	0.20	26,31,33,39	4
8	EDO	A	509	4/4	0.88	0.18	36,49,52,62	0
14	1PE	D	202	16/16	0.89	0.13	59,71,75,77	0
11	PGE	A	528	10/10	0.90	0.15	25,40,47,48	10
8	EDO	A	512	4/4	0.91	0.13	50,56,61,66	0
7	GOL	A	503	6/6	0.92	0.11	16,23,30,32	6
8	EDO	A	515	4/4	0.92	0.12	23,25,29,30	4
8	EDO	B	403	4/4	0.92	0.12	29,34,38,46	4
8	EDO	C	302	4/4	0.93	0.13	41,45,50,50	4
13	8Q1	C	301	34/35	0.96	0.07	27,34,43,47	0
8	EDO	A	513	4/4	0.97	0.10	17,20,24,29	4
8	EDO	A	507	4/4	0.97	0.10	10,11,16,20	4
5	PLP	A	501	15/16	0.99	0.04	17,20,25,27	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.

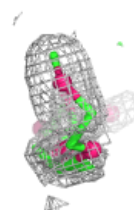
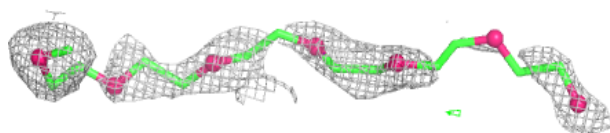
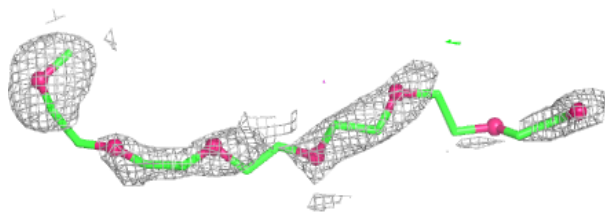
Electron density around EDT B 411:

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

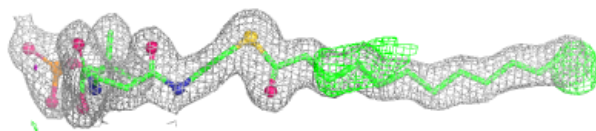
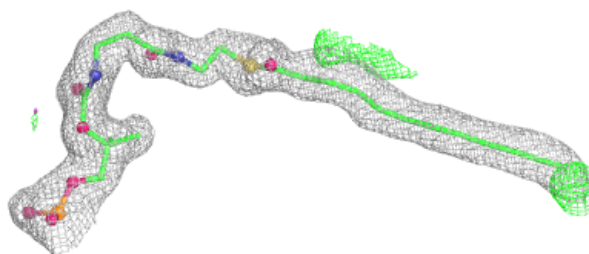


Electron density around P15 A 523:

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

**Electron density around 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.