



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 10, 2026 – 10:08 pm BST

PDB ID : 9TPB / pdb_00009tpb
Title : Canilysin from *Capnocytophaga canimorsus* (Strain 5)
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Deposited on : 2025-12-17
Resolution : 2.50 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	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

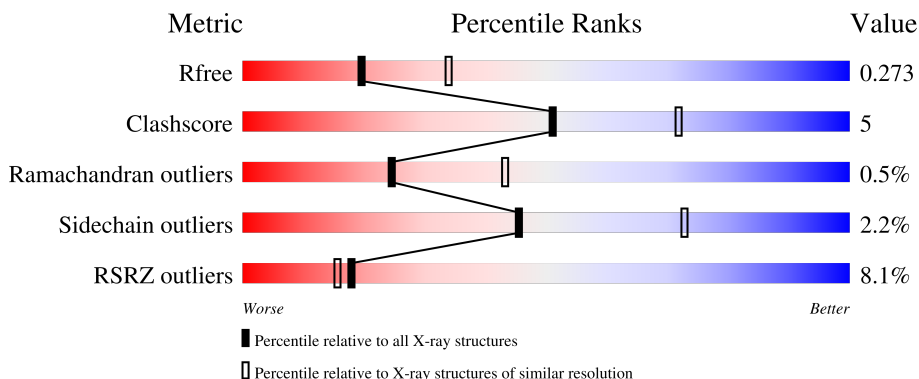
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	356	
1	B	356	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 5234 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 Metallopeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	304	Total	C	N	O	S	0	0	0
			2467	1602	399	459	7			
1	B	304	Total	C	N	O	S	0	0	0
			2467	1602	399	459	7			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	10	MET	-	initiating methionine	UNP F9YTK6
A	11	GLY	-	expression tag	UNP F9YTK6
A	12	SER	-	expression tag	UNP F9YTK6
A	13	SER	-	expression tag	UNP F9YTK6
A	14	HIS	-	expression tag	UNP F9YTK6
A	15	HIS	-	expression tag	UNP F9YTK6
A	16	HIS	-	expression tag	UNP F9YTK6
A	17	HIS	-	expression tag	UNP F9YTK6
A	18	HIS	-	expression tag	UNP F9YTK6
A	19	HIS	-	expression tag	UNP F9YTK6
A	20	SER	-	expression tag	UNP F9YTK6
A	21	SER	-	expression tag	UNP F9YTK6
A	22	GLY	-	expression tag	UNP F9YTK6
A	23	GLU	-	expression tag	UNP F9YTK6
A	24	ASN	-	expression tag	UNP F9YTK6
A	25	LEU	-	expression tag	UNP F9YTK6
A	26	TYR	-	expression tag	UNP F9YTK6
A	27	PHE	-	expression tag	UNP F9YTK6
A	28	GLN	-	expression tag	UNP F9YTK6
A	29	GLY	-	expression tag	UNP F9YTK6
A	30	HIS	-	expression tag	UNP F9YTK6
A	31	MET	-	expression tag	UNP F9YTK6
B	10	MET	-	initiating methionine	UNP F9YTK6
B	11	GLY	-	expression tag	UNP F9YTK6
B	12	SER	-	expression tag	UNP F9YTK6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	13	SER	-	expression tag	UNP F9YTK6
B	14	HIS	-	expression tag	UNP F9YTK6
B	15	HIS	-	expression tag	UNP F9YTK6
B	16	HIS	-	expression tag	UNP F9YTK6
B	17	HIS	-	expression tag	UNP F9YTK6
B	18	HIS	-	expression tag	UNP F9YTK6
B	19	HIS	-	expression tag	UNP F9YTK6
B	20	SER	-	expression tag	UNP F9YTK6
B	21	SER	-	expression tag	UNP F9YTK6
B	22	GLY	-	expression tag	UNP F9YTK6
B	23	GLU	-	expression tag	UNP F9YTK6
B	24	ASN	-	expression tag	UNP F9YTK6
B	25	LEU	-	expression tag	UNP F9YTK6
B	26	TYR	-	expression tag	UNP F9YTK6
B	27	PHE	-	expression tag	UNP F9YTK6
B	28	GLN	-	expression tag	UNP F9YTK6
B	29	GLY	-	expression tag	UNP F9YTK6
B	30	HIS	-	expression tag	UNP F9YTK6
B	31	MET	-	expression tag	UNP F9YTK6

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

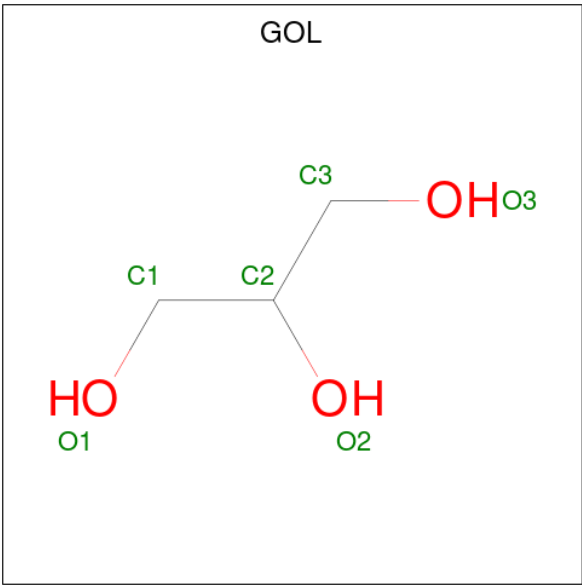
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



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

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

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

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).

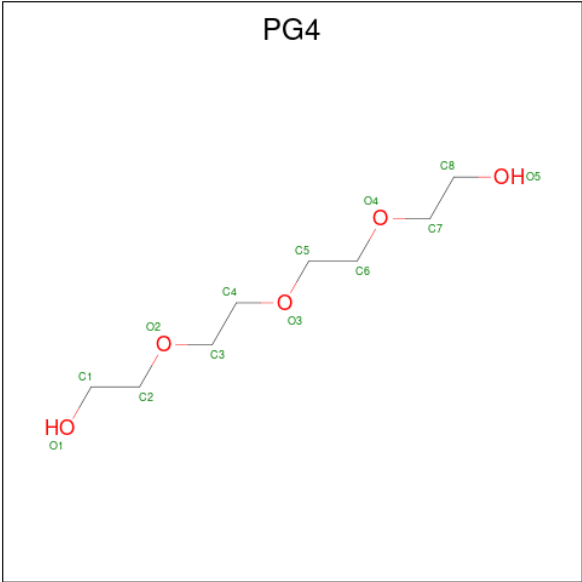


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Ca	0	0
			1	1		

- Molecule 7 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	120	Total	O	0	0
			120	120		
8	B	104	Total	O	0	0
			104	104		

- Molecule 1: Metallopeptidase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	123.25Å 94.56Å 83.96Å 90.00° 128.43° 90.00°	Depositor
Resolution (Å)	62.68 – 2.50 62.68 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (62.68-2.50) 99.8 (62.68-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.20_4459	Depositor
R, R_{free}	0.212 , 0.273 0.212 , 0.273	Depositor DCC
R_{free} test set	595 reflections (2.27%)	wwPDB-VP
Wilson B-factor (Å ²)	47.9	Xtriage
Anisotropy	0.557	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5234	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.53% 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: ZN, PGE, EDO, CA, PG4, GOL

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.36	0/2529	0.58	0/3421
1	B	0.38	0/2529	0.59	0/3421
All	All	0.37	0/5058	0.58	0/6842

There are no bond length outliers.

There are no bond angle outliers.

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	2467	0	2437	23	0
1	B	2467	0	2437	30	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	10	0	14	0	0
3	B	20	0	28	0	0
4	A	6	0	8	0	0
4	B	12	0	15	1	0
5	A	12	0	18	1	0
6	B	1	0	0	0	0
7	B	13	0	18	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	120	0	0	2	0
8	B	104	0	0	2	0
All	All	5234	0	4975	52	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 52 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:ILE:HD11	1:B:311:VAL:HA	1.62	0.81
1:B:296:GLU:HB2	1:B:354:LEU:HD21	1.65	0.78
1:B:112:ASP:HB2	1:B:165:ASP:HA	1.73	0.70
1:B:51:ASN:OD1	1:B:57:LYS:NZ	2.29	0.66
1:A:213:PRO:HG3	1:A:364:PRO:HG3	1.79	0.64

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	300/356 (84%)	287 (96%)	11 (4%)	2 (1%)	18	34
1	B	300/356 (84%)	288 (96%)	11 (4%)	1 (0%)	36	55
All	All	600/712 (84%)	575 (96%)	22 (4%)	3 (0%)	24	43

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	256	THR
1	B	256	THR

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Mol	Chain	Res	Type
1	A	258	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	268/316 (85%)	264 (98%)	4 (2%)	57	80
1	B	268/316 (85%)	260 (97%)	8 (3%)	36	64
All	All	536/632 (85%)	524 (98%)	12 (2%)	45	73

5 of 12 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	62	ILE
1	B	84	ILE
1	B	160	ASN
1	B	114	PHE
1	A	295	LYS

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	167	ASN
1	B	100	HIS
1	B	167	ASN
1	B	260	GLN
1	B	300	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 3 are monoatomic - leaving 10 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	$\# Z > 2$
5	EDO	A	405	-	3,3,3	0.62	0	2,2,2	0.10	0
4	GOL	A	403	-	5,5,5	1.03	0	5,5,5	0.87	0
7	PG4	B	403	-	12,12,12	0.14	0	11,11,11	0.55	0
3	PGE	A	402	-	9,9,9	0.28	0	8,8,8	0.54	0
4	GOL	B	407	-	5,5,5	0.97	0	5,5,5	1.14	0
5	EDO	A	404	-	3,3,3	0.50	0	2,2,2	0.30	0
3	PGE	B	404	-	9,9,9	0.35	0	8,8,8	0.36	0
3	PGE	B	405	-	9,9,9	0.35	0	8,8,8	0.36	0
4	GOL	B	406	-	5,5,5	1.22	1 (20%)	5,5,5	0.87	0
5	EDO	A	406	-	3,3,3	0.74	0	2,2,2	0.29	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
5	EDO	A	405	-	-	0/1/1/1	-
4	GOL	A	403	-	-	2/4/4/4	-
7	PG4	B	403	-	-	5/10/10/10	-
3	PGE	A	402	-	-	1/7/7/7	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	407	-	-	2/4/4/4	-
5	EDO	A	404	-	-	0/1/1/1	-
3	PGE	B	404	-	-	2/7/7/7	-
3	PGE	B	405	-	-	3/7/7/7	-
4	GOL	B	406	-	-	0/4/4/4	-
5	EDO	A	406	-	-	0/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	406	GOL	O2-C2	-2.42	1.36	1.43

There are no bond angle outliers.

There are no chirality outliers.

5 of 15 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	403	GOL	C1-C2-C3-O3
3	B	405	PGE	C6-C5-O3-C4
7	B	403	PG4	O2-C3-C4-O3
3	A	402	PGE	O3-C5-C6-O4
3	B	404	PGE	O3-C5-C6-O4

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	405	EDO	1	0
7	B	403	PG4	2	0
4	B	406	GOL	1	0

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	304/356 (85%)	0.67	22 (7%) 21 19	39, 51, 95, 179	0
1	B	304/356 (85%)	0.64	27 (8%) 15 13	38, 50, 100, 172	0
All	All	608/712 (85%)	0.66	49 (8%) 18 15	38, 50, 96, 179	0

The worst 5 of 49 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	138	PHE	5.6
1	B	161	ILE	5.2
1	B	138	PHE	4.9
1	A	161	ILE	4.8
1	A	114	PHE	4.8

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](#)

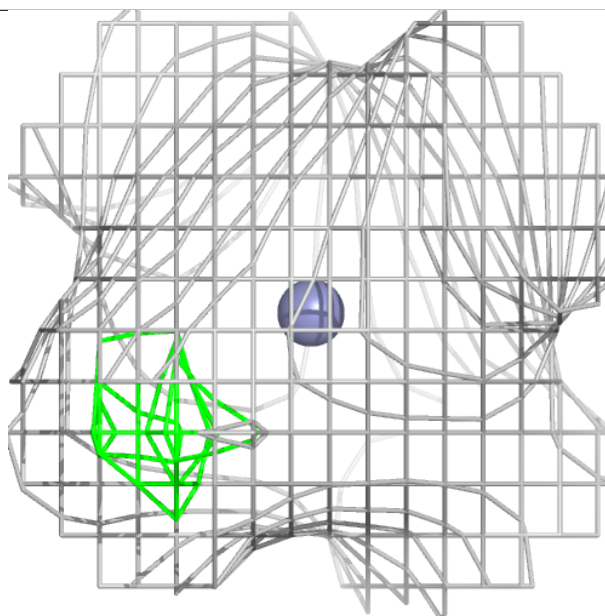
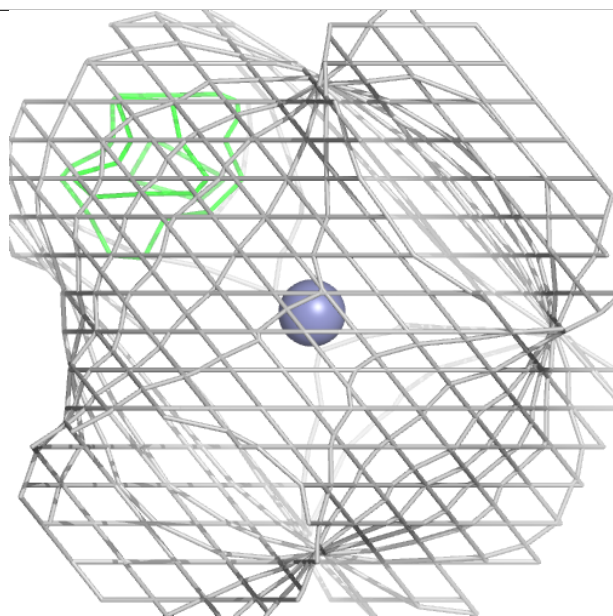
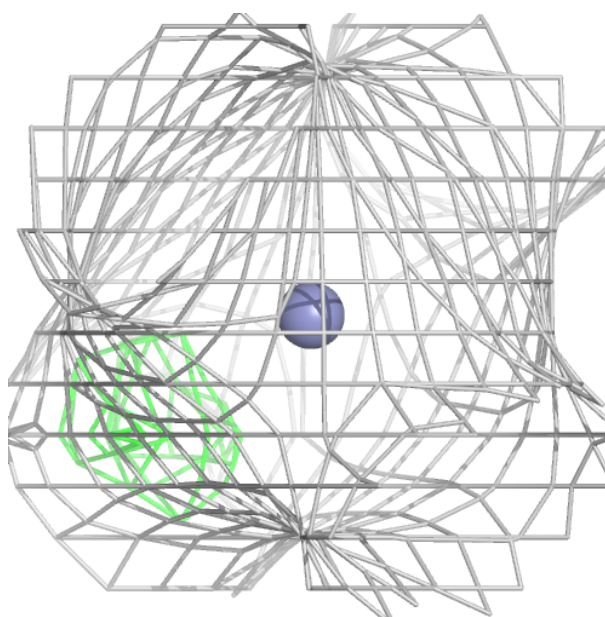
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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
5	EDO	A	406	4/4	0.65	0.20	50,52,61,62	0
5	EDO	A	405	4/4	0.69	0.20	52,53,58,59	0
4	GOL	B	407	6/6	0.77	0.20	61,65,69,69	0
4	GOL	B	406	6/6	0.79	0.14	58,61,64,65	0
3	PGE	B	405	10/10	0.79	0.17	59,70,73,77	0
5	EDO	A	404	4/4	0.80	0.17	65,68,73,74	0
3	PGE	A	402	10/10	0.81	0.19	56,63,67,70	0
4	GOL	A	403	6/6	0.82	0.14	64,71,75,80	0
3	PGE	B	404	10/10	0.88	0.17	62,71,75,79	0
7	PG4	B	403	13/13	0.88	0.14	59,65,70,72	0
6	CA	B	402	1/1	0.94	0.14	87,87,87,87	0
2	ZN	B	401	1/1	0.99	0.03	44,44,44,44	0
2	ZN	A	401	1/1	1.00	0.02	44,44,44,44	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.

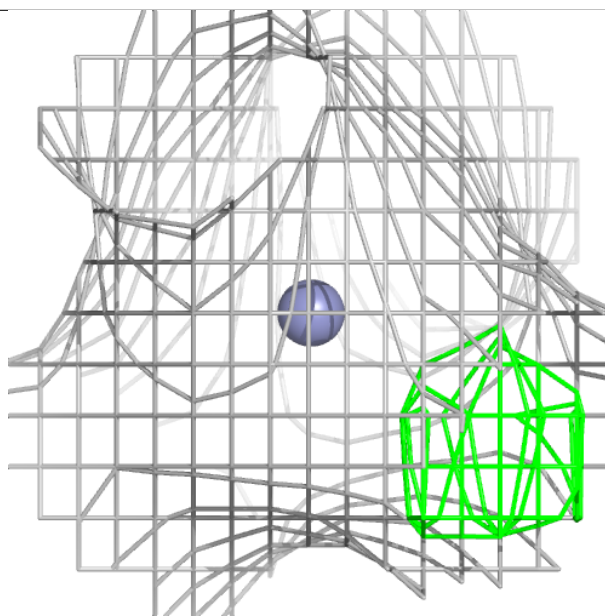
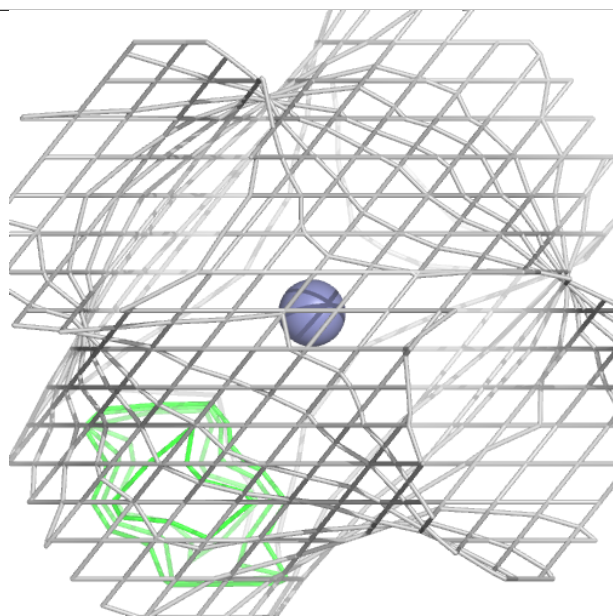
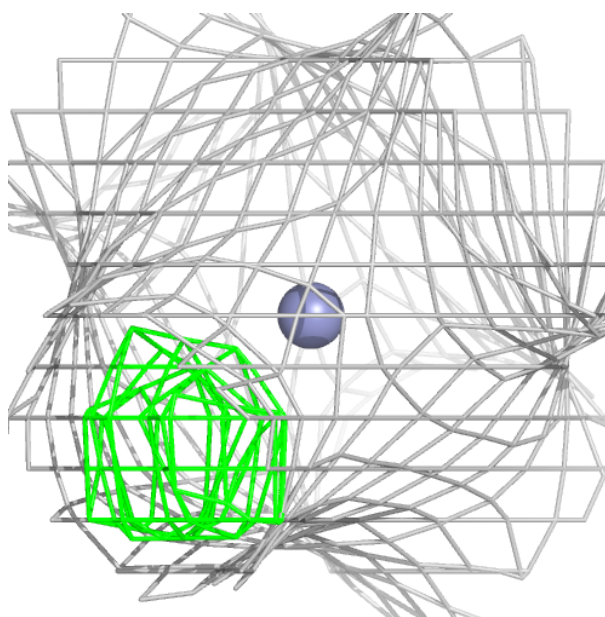
Electron density around ZN B 401:

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



Electron density around ZN A 401:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.