



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2026 – 02:32 AM UTC

PDB ID : 9B2L / pdb_00009b2l
Title : Structure of the quorum quenching lactonase GcL D122N mutant - bimetallic center
Authors : Corbella, M.; Bravo, J.A.; Demkiv, A.O.; Calixto, A.R.; Sompiyachoke, K.; Bergonzi, C.; Kamerlin, S.C.L.; Elias, M.
Deposited on : 2024-03-15
Resolution : 2.20 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
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

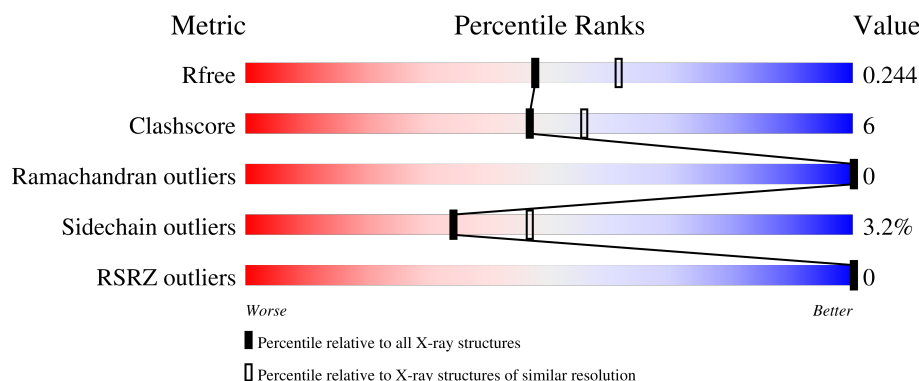
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	275	 86% 13% .
1	B	275	 87% 11% .

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 4669 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 quorum-quenching N-acyl-homoserine lactonase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	2	0
			2251	1431	395	411	14			
1	B	275	Total	C	N	O	S	0	2	0
			2252	1432	396	410	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	122	ASN	ASP	engineered mutation	UNP A0A023DFE8
B	122	ASN	ASP	engineered mutation	UNP A0A023DFE8

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



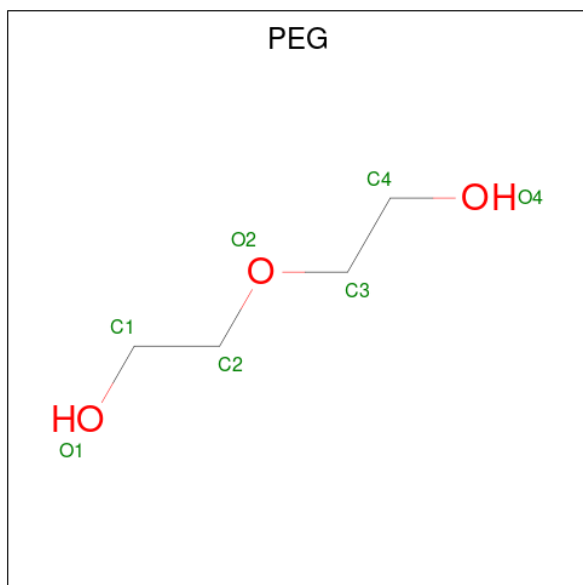
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	O S	0	0
			5	4 1		

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

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

- Molecule 4 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



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

- 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	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is FE (III) ION (CCD ID: FE) (formula: Fe) (labeled as "Ligand of Interest" by

depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total 1	Fe 1	0	0
6	B	1	Total 1	Fe 1	0	0

- Molecule 7 is COBALT (II) ION (CCD ID: CO) (formula: Co) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total 1	Co 1	0	0
7	B	1	Total 1	Co 1	0	0

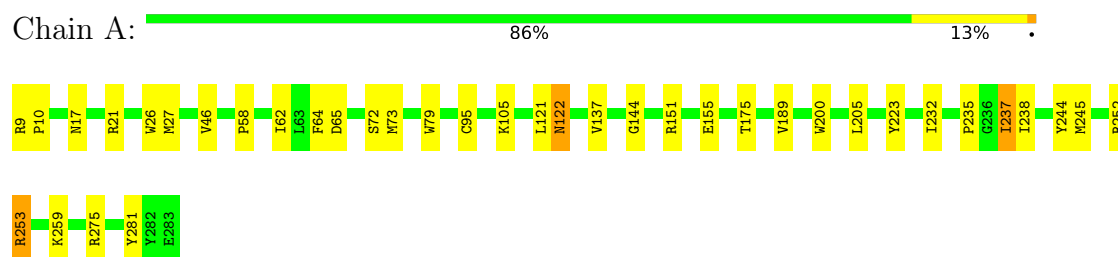
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	53	Total 53	O 53	0	0
8	B	57	Total 57	O 57	0	0

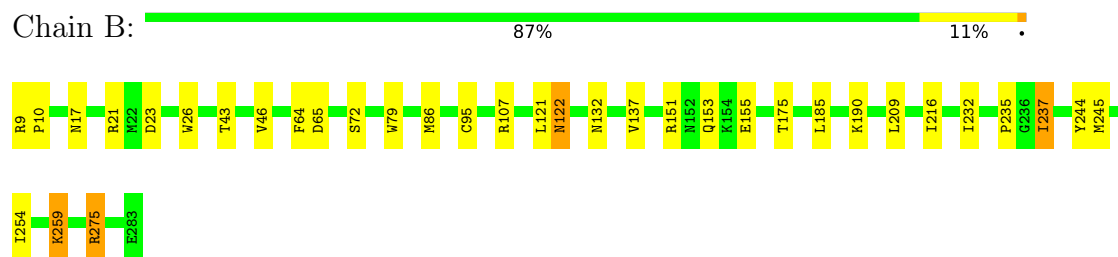
3 Residue-property plots

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: quorum-quenching N-acyl-homoserine lactonase



- Molecule 1: quorum-quenching N-acyl-homoserine lactonase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	109.22Å 109.22Å 221.66Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	73.89 – 2.20 73.89 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (73.89-2.20) 99.8 (73.89-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0419	Depositor
R, R_{free}	0.214 , 0.241 0.218 , 0.244	Depositor DCC
R_{free} test set	2498 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	39.4	Xtriage
Anisotropy	0.651	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 33.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.459 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4669	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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: SO4, PEG, EDO, FE, CO, PGE

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.41	0/2315	0.79	0/3142
1	B	0.41	0/2317	0.78	0/3145
All	All	0.41	0/4632	0.79	0/6287

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	3
1	B	0	1
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	151	ARG	Sidechain
1	A	252	ARG	Sidechain
1	A	253	ARG	Sidechain
1	B	275	ARG	Sidechain

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	2251	0	2161	25	0
1	B	2252	0	2160	26	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
3	A	7	0	10	0	0
3	B	7	0	10	1	0
4	A	10	0	14	1	0
5	A	4	0	6	0	0
5	B	4	0	6	1	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
8	A	53	0	0	0	0
8	B	57	0	0	0	0
All	All	4669	0	4367	51	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 6.

All (51) 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:B:9:ARG:HB3	1:B:10:PRO:HD3	1.35	1.08
1:A:9:ARG:HB3	1:A:10:PRO:HD3	1.49	0.95
1:A:122:ASN:HD22	1:A:122:ASN:H	1.31	0.78
1:B:122:ASN:HD22	1:B:122:ASN:H	1.35	0.71
1:B:9:ARG:HB3	1:B:10:PRO:CD	2.19	0.70
1:A:9:ARG:HB3	1:A:10:PRO:CD	2.25	0.65
1:A:9:ARG:CB	1:A:10:PRO:HD3	2.25	0.63
1:B:275:ARG:HE	5:B:403:EDO:H22	1.65	0.61
1:B:26:TRP:CE3	1:B:237:ILE:HG12	2.41	0.56
1:A:17:ASN:HA	1:A:95:CYS:O	2.07	0.55
1:A:26:TRP:CE3	1:A:237:ILE:HG12	2.42	0.55
1:B:122:ASN:H	1:B:122:ASN:ND2	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:ASN:HD22	1:A:122:ASN:N	2.03	0.54
1:A:21:ARG:HA	1:A:46:VAL:O	2.08	0.53
1:B:121:LEU:H	1:B:121:LEU:HD23	1.73	0.53
1:B:122:ASN:HD22	1:B:122:ASN:N	2.06	0.53
1:B:72:SER:HB3	1:B:79:TRP:CE2	2.43	0.53
1:A:235:PRO:HD3	1:A:244:TYR:CZ	2.44	0.53
1:B:235:PRO:HD3	1:B:244:TYR:CZ	2.44	0.52
1:A:121:LEU:H	1:A:121:LEU:HD23	1.74	0.52
1:A:121:LEU:HD23	1:A:121:LEU:N	2.26	0.49
1:A:64:PHE:O	1:A:65:ASP:HB2	2.13	0.49
1:A:121:LEU:H	1:A:121:LEU:CD2	2.26	0.49
1:B:121:LEU:HD23	1:B:121:LEU:N	2.28	0.49
1:B:21:ARG:HA	1:B:46:VAL:O	2.13	0.49
1:B:121:LEU:H	1:B:121:LEU:CD2	2.27	0.48
1:B:9:ARG:CB	1:B:10:PRO:HD3	2.22	0.48
1:B:64:PHE:O	1:B:65:ASP:HB2	2.13	0.48
1:A:72:SER:HB3	1:A:79:TRP:CE2	2.49	0.47
1:B:17:ASN:HA	1:B:95:CYS:O	2.14	0.47
1:A:144:GLY:HA3	1:A:200:TRP:CE3	2.50	0.47
1:A:9:ARG:HG2	1:A:58:PRO:HG3	1.98	0.45
1:B:26:TRP:HE3	1:B:237:ILE:HG12	1.79	0.45
1:A:73:MET:HE3	1:A:73:MET:HB3	1.86	0.44
1:B:216:ILE:HD11	1:B:254:ILE:HG22	2.00	0.44
1:B:151:ARG:HD2	1:B:153:GLN:HE21	1.82	0.43
1:B:137:VAL:O	1:B:175:THR:HA	2.19	0.43
1:B:232:ILE:HB	1:B:245:MET:HE1	2.01	0.43
1:B:185:LEU:HD12	1:B:185:LEU:HA	1.89	0.43
1:B:10:PRO:HG2	1:B:209:LEU:HD23	2.01	0.43
1:B:132:ASN:HA	3:B:404:PEG:H21	2.01	0.43
1:A:122:ASN:H	1:A:122:ASN:ND2	2.07	0.42
1:A:275:ARG:NH2	1:A:281:TYR:O	2.49	0.42
1:A:26:TRP:HE3	1:A:237:ILE:HG12	1.82	0.41
1:A:137:VAL:O	1:A:175:THR:HA	2.21	0.41
1:B:259:LYS:HB3	1:B:259:LYS:HE3	1.86	0.41
1:A:232:ILE:HB	1:A:245:MET:HE1	2.01	0.41
1:A:235:PRO:HD2	1:A:238:ILE:HD12	2.03	0.40
1:A:105:LYS:HG3	4:A:303:PGE:H3	2.04	0.40
1:A:62:ILE:HD13	1:A:189:VAL:HG11	2.02	0.40
1:B:23:ASP:OD1	1:B:43:THR:OG1	2.35	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	275/275 (100%)	267 (97%)	8 (3%)	0	100	100
1	B	275/275 (100%)	267 (97%)	8 (3%)	0	100	100
All	All	550/550 (100%)	534 (97%)	16 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	238/236 (101%)	230 (97%)	8 (3%)	32	44
1	B	238/236 (101%)	231 (97%)	7 (3%)	37	51
All	All	476/472 (101%)	461 (97%)	15 (3%)	34	47

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	MET
1	A	122	ASN
1	A	155	GLU
1	A	205	LEU
1	A	223	TYR
1	A	237	ILE
1	A	253	ARG
1	A	259	LYS

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Mol	Chain	Res	Type
1	B	86	MET
1	B	107	ARG
1	B	122	ASN
1	B	155	GLU
1	B	190	LYS
1	B	237	ILE
1	B	259	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	31	HIS
1	A	38	ASN
1	A	69	ASN
1	A	71	ASN
1	A	122	ASN
1	A	132	ASN
1	A	179	HIS
1	A	182	ASN
1	B	17	ASN
1	B	25	ASN
1	B	31	HIS
1	B	38	ASN
1	B	69	ASN
1	B	85	GLN
1	B	122	ASN
1	B	132	ASN
1	B	153	GLN
1	B	182	ASN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 4 are monoatomic - leaving 9 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	$\# Z > 2$
2	SO4	A	301	-	4,4,4	0.34	0	6,6,6	0.08	0
2	SO4	A	304	-	4,4,4	0.33	0	6,6,6	0.07	0
5	EDO	B	403	-	3,3,3	0.06	0	2,2,2	0.04	0
2	SO4	B	401	-	4,4,4	0.33	0	6,6,6	0.08	0
5	EDO	A	305	-	3,3,3	0.06	0	2,2,2	0.08	0
3	PEG	B	404	-	6,6,6	0.15	0	5,5,5	0.07	0
4	PGE	A	303	-	9,9,9	0.18	0	8,8,8	0.09	0
3	PEG	A	302	-	6,6,6	0.14	0	5,5,5	0.08	0
2	SO4	B	402	-	4,4,4	0.34	0	6,6,6	0.07	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	B	403	-	-	0/1/1/1	-
5	EDO	A	305	-	-	1/1/1/1	-
3	PEG	B	404	-	-	3/4/4/4	-
4	PGE	A	303	-	-	3/7/7/7	-
3	PEG	A	302	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	404	PEG	O2-C3-C4-O4
4	A	303	PGE	O3-C5-C6-O4
3	A	302	PEG	O1-C1-C2-O2
3	B	404	PEG	O1-C1-C2-O2
4	A	303	PGE	O2-C3-C4-O3
3	A	302	PEG	C4-C3-O2-C2
3	B	404	PEG	C1-C2-O2-C3
5	A	305	EDO	O1-C1-C2-O2
4	A	303	PGE	C1-C2-O2-C3

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	403	EDO	1	0
3	B	404	PEG	1	0
4	A	303	PGE	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	275/275 (100%)	-0.76	0 100 100	31, 54, 79, 101	2 (0%)
1	B	275/275 (100%)	-0.78	0 100 100	32, 54, 81, 102	2 (0%)
All	All	550/550 (100%)	-0.77	0 100 100	31, 54, 81, 102	4 (0%)

There are no RSRZ outliers to report.

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(Å ²)	Q<0.9
3	PEG	A	302	7/7	0.95	0.10	70,74,85,85	0
5	EDO	B	403	4/4	0.95	0.09	63,68,72,75	0
4	PGE	A	303	10/10	0.96	0.13	67,82,93,95	0
3	PEG	B	404	7/7	0.96	0.09	71,76,82,85	0
2	SO4	A	304	5/5	0.98	0.06	64,80,84,95	0
5	EDO	A	305	4/4	0.98	0.10	59,62,62,65	0
2	SO4	B	401	5/5	0.98	0.06	65,83,91,98	0

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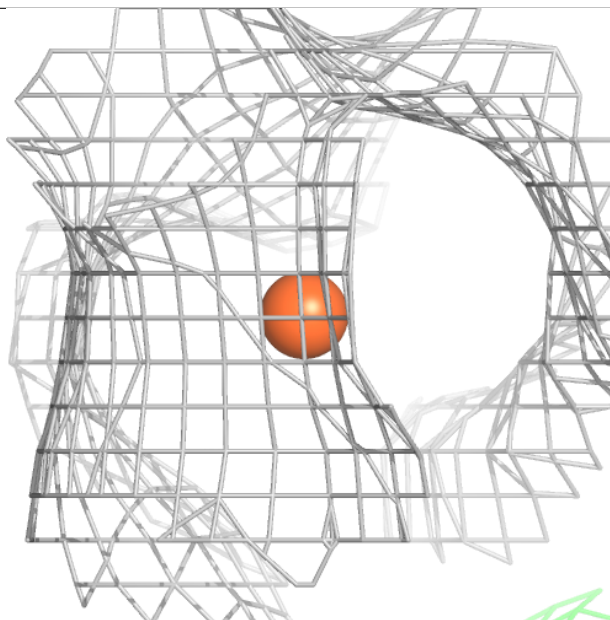
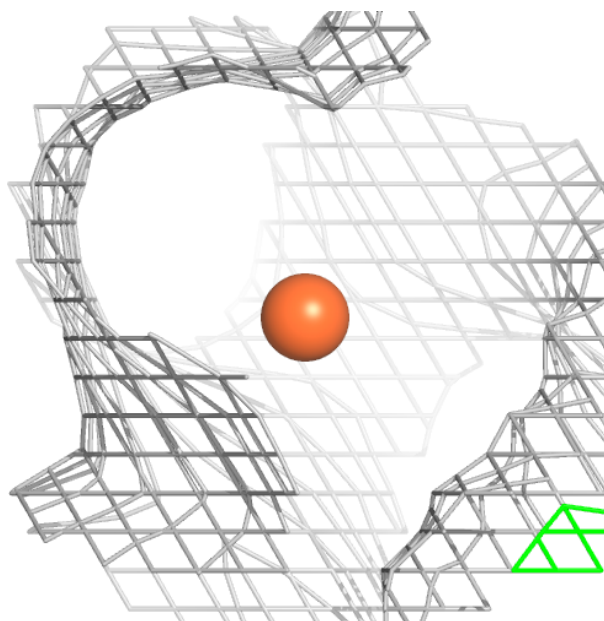
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	B	402	5/5	0.99	0.04	76,82,83,88	0
2	SO4	A	301	5/5	0.99	0.04	70,75,77,85	0
6	FE	A	306	1/1	1.00	0.01	51,51,51,51	0
6	FE	B	405	1/1	1.00	0.01	49,49,49,49	0
7	CO	A	307	1/1	1.00	0.01	48,48,48,48	0
7	CO	B	406	1/1	1.00	0.01	47,47,47,47	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.

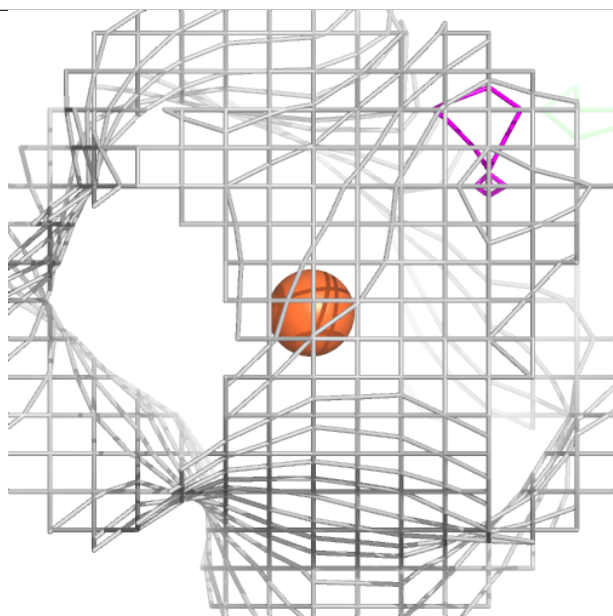
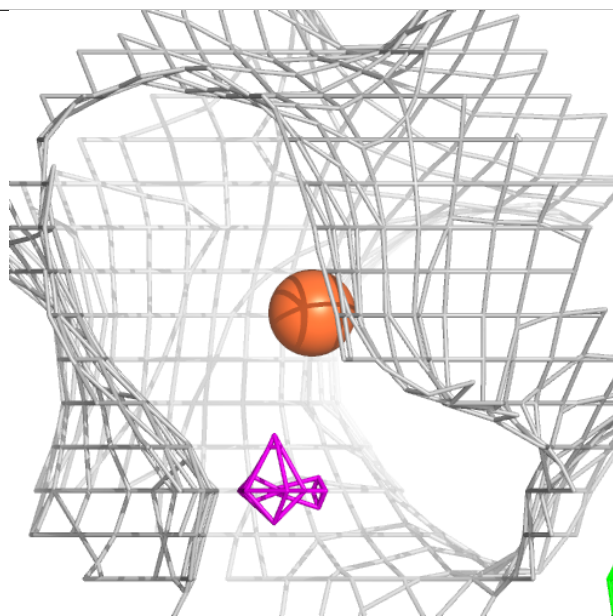
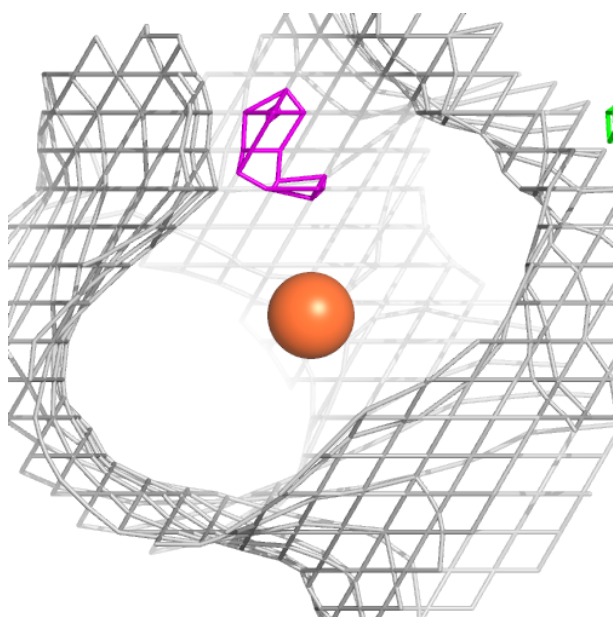
Electron density around FE A 306:

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



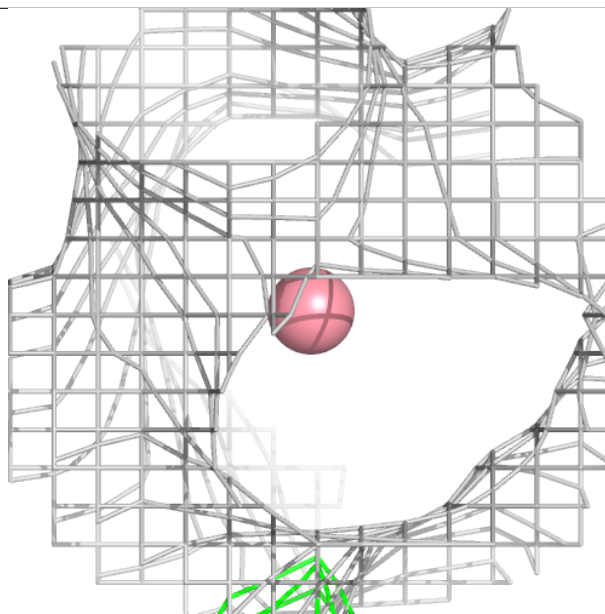
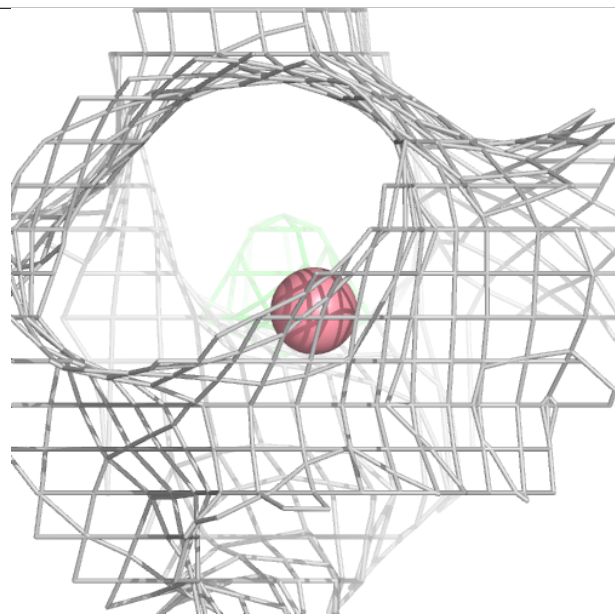
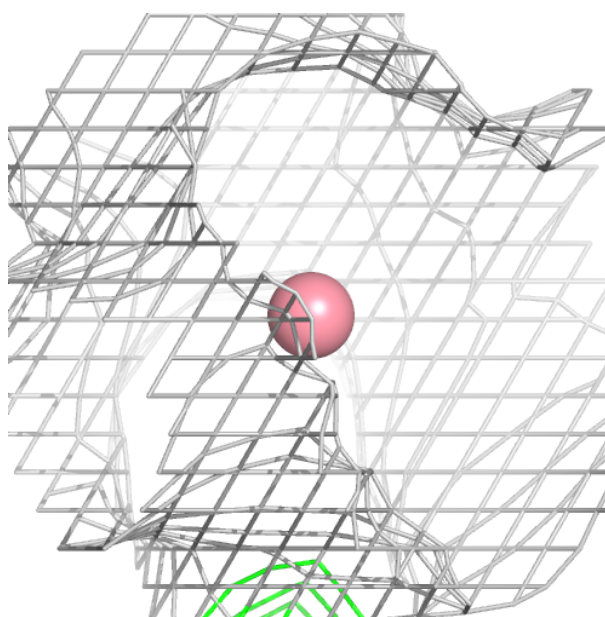
Electron density around FE B 405:

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



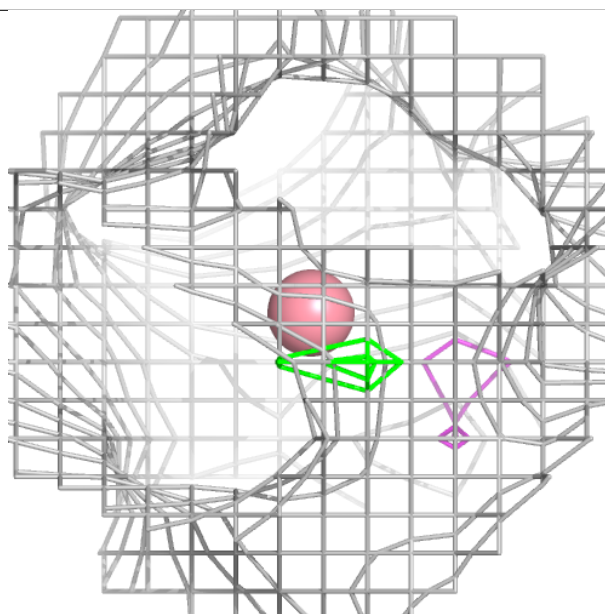
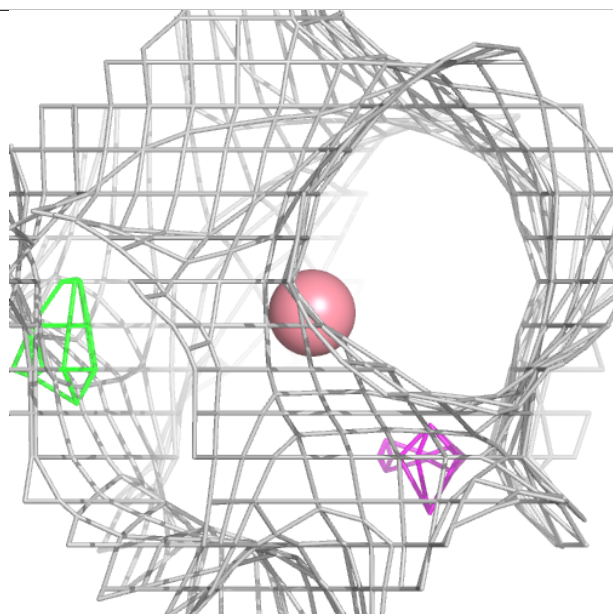
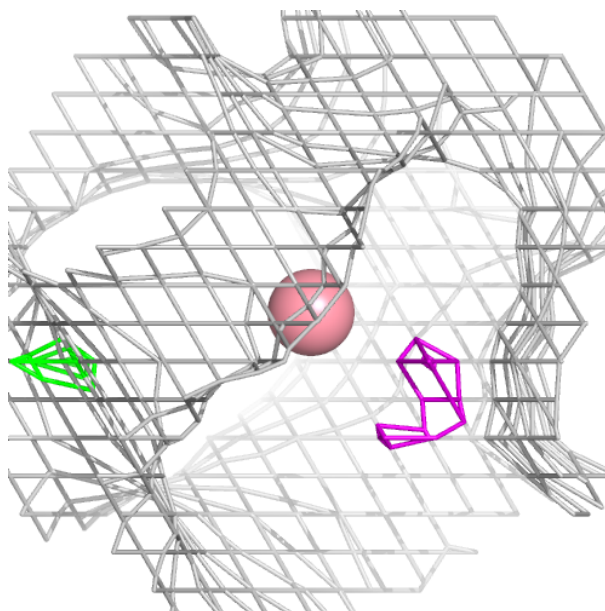
Electron density around CO A 307:

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



Electron density around CO B 406:

$2mF_o-DF_c$ (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.