

Full wwPDB X-ray Structure Validation Report (i)

Aug 1, 2024 – 04:42 pm BST

PDB ID : 8RFR

Title: Low pH (5.5) nitrite-bound MSOX movie series dataset 7 of the copper nitrite

reductase (NirK) from Bradyrhizobium japonicum USDA110 [4.27 MGy] -

NO/water intermediate

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Deposited on : 2023-12-13

Resolution : 1.40 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 (i) 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 (i)) were used in the production of this report:

MolProbity: 4.02b-467

Mogul : 1.8.4, CSD as 541be (2020)

Xtriage (Phenix) : 1.13

EDS : 2.37.1

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac : 5.8.0158

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

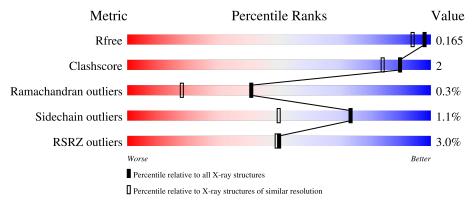
Validation Pipeline (wwPDB-VP) : 2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$

The reported resolution of this entry is 1.40 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
R_{free}	130704	1714 (1.40-1.40)
Clashscore	141614	1812 (1.40-1.40)
Ramachandran outliers	138981	1763 (1.40-1.40)
Sidechain outliers	138945	1762 (1.40-1.40)
RSRZ outliers	127900	1674 (1.40-1.40)

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 <=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	340	92%	6% •				
2	С	2	50%	50%				



2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 3087 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 Copper-containing nitrite reductase.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	334	Total	С	Ν	О	S	0	21	0
1	11	001	2679	1723	452	485	19		21	

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP Q89EJ6
A	310	CYS	GLY	conflict	UNP Q89EJ6

• Molecule 2 is an oligosaccharide called beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose.



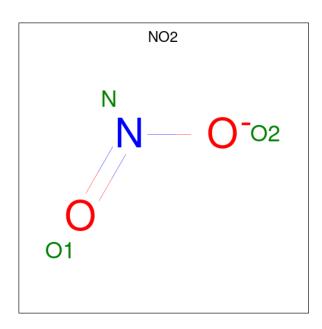
Mol	Chain	Residues	At	oms		ZeroOcc	AltConf	Trace
2	С	2	Total 23	C 12	O 11	0	0	0

• Molecule 3 is COPPER (II) ION (three-letter code: CU) (formula: Cu) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Cu 2 2	0	0

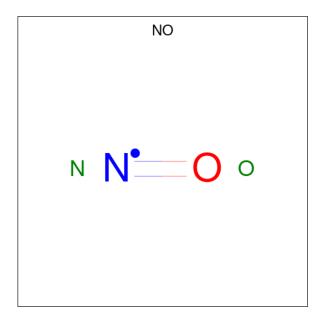
• Molecule 4 is NITRITE ION (three-letter code: NO2) (formula: NO₂).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total N O 3 1 2	0	0
4	A	1	Total N O 3 1 2	0	0

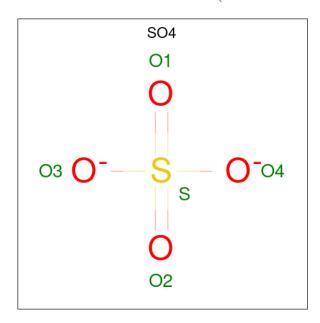
• Molecule 5 is NITRIC OXIDE (three-letter code: NO) (formula: NO) (labeled as "Ligand of Interest" by depositor).



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



 \bullet Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Ator	ns		ZeroOcc	AltConf	
6	A	1	Total	О	S	n	0	
0	Λ	1	5	4	1	U	U	
6	A	1	Total	О	S	0	0	
0	Λ	1	5	4	1	U	U	
6	A	1	Total	О	S	0	0	
0	Λ	1	5	4	1	U		
6	A	1	Total	О	S	0	0	
0	Λ	1	5	4	1	U	U	
6	A	1	Total	О	S	0	0	
0	Λ	1	5	4	1	U	U	
6	A	1	Total	О	S	0	0	
	Λ	1	5	4	1	U	U	

• Molecule 7 is water.

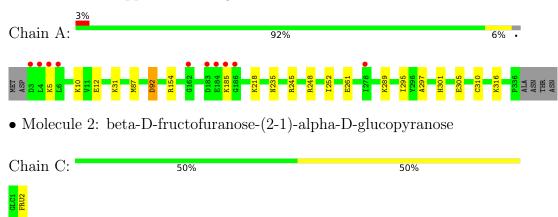
\mathbf{Mol}	Chain	Residues	${f Atoms}$	ZeroOcc	AltConf
7	A	337	Total O 345 345	0	19



3 Residue-property plots (i)

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

• Molecule 1: Copper-containing nitrite reductase





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 63	Depositor
Cell constants	104.36Å 104.36Å 64.51Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 - 1.40	Depositor
rtesolution (A)	45.19 - 1.40	EDS
% Data completeness	99.3 (30.00-1.40)	Depositor
(in resolution range)	99.4 (45.19-1.40)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.41 (at 1.40Å)	Xtriage
Refinement program	REFMAC 5.8.0419	Depositor
D D.	0.145 , 0.165	Depositor
R, R_{free}	0.146 , 0.165	DCC
R_{free} test set	3788 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	17.8	Xtriage
Anisotropy	0.286	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.39 , 50.6	EDS
L-test for twinning ²	$< L >=0.50, < L^2>=0.33$	Xtriage
Estimated twinning fraction	0.034 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	3087	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.14% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

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: SO4, GLC, NO, FRU, NO2, CU

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).

Mal	Chain	Bo	nd lengths	Bo	ond angles
Mol Chain		RMSZ	# Z > 5	RMSZ	# Z > 5
1	A	0.62	$1/2806 \ (0.0\%)$	0.97	7/3802 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(A)	$\operatorname{Ideal}(\text{\AA})$
1	A	261	GLU	CD-OE1	5.94	1.32	1.25

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
1	A	92	ASP	CB-CG-OD1	7.56	125.10	118.30
1	A	261	GLU	GLU CG-CD-OE2		104.86	118.30
1	A	87[A]	MET	CG-SD-CE	-5.82	90.89	100.20
1	A	87[B]	MET	CG-SD-CE	-5.82	90.89	100.20
1	A	245	ARG	NE-CZ-NH2	-5.55	117.53	120.30
1	A	261	GLU	CG-CD-OE1	5.52	129.34	118.30
1	A	154	ARG	CG-CD-NE	-5.49	100.27	111.80

There are no chirality outliers.

All (1) planarity outliers are listed below:



Mol	Chain	Res	Type	Group	
1	A	301	HIS	Peptide	

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	2679	0	2705	10	0
2	С	23	0	21	0	0
3	A	2	0	0	0	0
4	A	6	0	0	0	0
5	A	2	0	0	1	0
6	A	30	0	0	0	0
7	A	345	0	0	3	0
All	All	3087	0	2726	10	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 2.

All (10) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{array}{ll} ext{Interatomic} \ ext{distance } (ext{Å}) \end{array}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$	
1:A:10:LYS:HE2	1:A:12:GLU:OE2	1.91	0.70	
1:A:252:ILE:HD12	1:A:297:ALA:HB3	1.87	0.57	
1:A:10:LYS:CE	1:A:12:GLU:OE2	2.63	0.45	
1:A:218[B]:LYS:HB2	1:A:218[B]:LYS:HE2	1.59	0.44	
1:A:305:GLU:OE2	7:A:689[B]:HOH:O	2.21	0.43	
1:A:235[B]:ASN:OD1	1:A:289[B]:LYS:HB2	2.18	0.42	
1:A:185:LYS:N	7:A:505:HOH:O	2.38	0.42	
1:A:92:ASP:OD2	5:A:405[A]:NO:N	2.52	0.41	
1:A:295[B]:ILE:HG13	7:A:562:HOH:O	2.20	0.41	
1:A:316[A]:LYS:HA	1:A:316[A]:LYS:HD2	1.98	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	354/340 (104%)	348 (98%)	5 (1%)	1 (0%)	41 18	

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	310	CYS

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	290/275 (106%)	287 (99%)	3 (1%)	76 53

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	31	LYS
1	A	248	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

\mathbf{Mol}	Chain	Res	Type
1	A	187	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)

2 monosaccharides 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 Ch	Type	Type Chain	Chain	Chain	Chain	Chain	noin Dog	Res	Pos	Ros	Link	Bo	ond leng	$ ag{ths}$	$ \hspace{.05cm} {f B}$	ond ang	les
		Chain	nes	ites Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2								
	2	GLC	С	1	2	11,11,12	0.60	0	15,15,17	1.16	0							
	2	FRU	С	2	2	11,12,12	1.54	1 (9%)	10,18,18	1.22	1 (10%)							

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
2	GLC	С	1	2	-	0/2/19/22	0/1/1/1
2	FRU	С	2	2	-	0/5/24/24	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(\mathbf{\mathring{A}})$	$\operatorname{Ideal}(\text{\AA})$
2	$^{\mathrm{C}}$	2	FRU	O5-C2	3.92	1.49	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
2	С	2	FRU	O1-C1-C2	-3.00	105.49	111.86



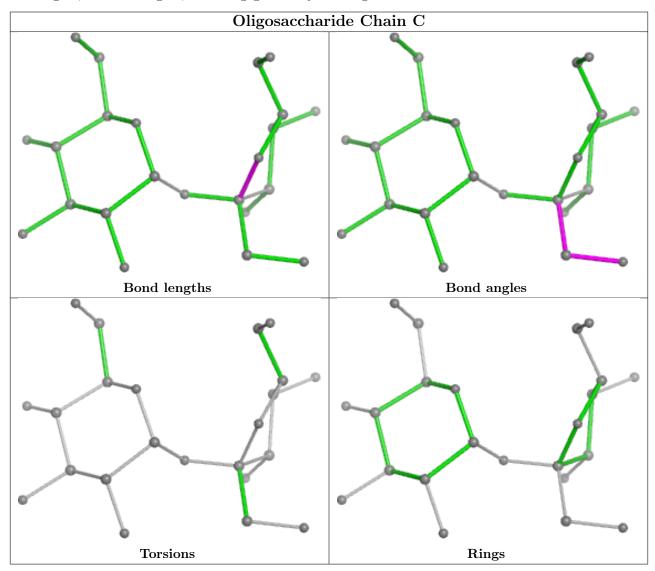
There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry (i)

Of 11 ligands modelled in this entry, 2 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	Tuno	Chain	Res	Link	B	ond leng	gths	В	ond ang	gles
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	SO4	A	406	-	4,4,4	0.38	0	6,6,6	0.20	0
6	SO4	A	408	-	4,4,4	0.43	0	6,6,6	0.11	0
6	SO4	A	410	-	4,4,4	0.46	0	6,6,6	0.13	0
6	SO4	A	407	-	4,4,4	0.54	0	6,6,6	0.20	0
5	NO	A	405[A]	3	0,1,1	-	-	-		
4	NO2	A	403	-	1,2,2	0.75	0	0,1,1	-	-
6	SO4	A	409	-	4,4,4	0.33	0	6,6,6	0.08	0
4	NO2	A	404	-	1,2,2	0.41	0	0,1,1	-	-
6	SO4	A	411	-	4,4,4	0.46	0	6,6,6	0.23	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

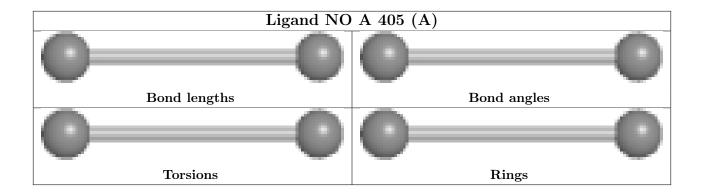
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	405[A]	NO	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ>2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	$\langle { m RSRZ} \rangle$	# RSRZ > 2	$OWAB(A^2)$	Q < 0.9
1	A	334/340 (98%)	-0.13	10 (2%) 50 49	13, 19, 39, 60	13 (3%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	4	LEU	6.1
1	A	5	LYS	4.1
1	A	185	LYS	4.0
1	A	162[A]	GLY	3.8
1	A	186	GLY	3.5
1	A	6	LEU	3.3
1	A	278	ILE	3.1
1	A	3	ASP	2.8
1	A	183	ASP	2.5
1	A	184	GLU	2.5

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

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

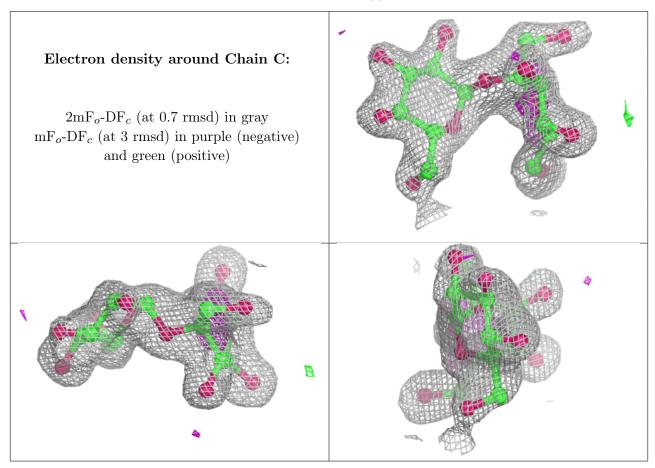
6.3 Carbohydrates (i)

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, 95^{th} 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	$\operatorname{B-factors}({ m \AA}^2)$	Q<0.9
2	FRU	С	2	12/12	0.89	0.19	23,33,43,47	0
2	GLC	С	1	11/12	0.91	0.21	36,47,52,55	0



The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



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, 95^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
4	NO2	A	403	3/3	0.85	0.13	55,55,57,58	0
4	NO2	A	404	3/3	0.87	0.09	33,33,37,38	3
6	SO4	A	411	5/5	0.88	0.26	40,41,45,45	5
6	SO4	A	409	5/5	0.89	0.30	49,57,63,65	5
6	SO4	A	410	5/5	0.95	0.12	36,37,41,43	5
6	SO4	A	408	5/5	0.96	0.23	51,52,56,60	5
6	SO4	A	406	5/5	0.96	0.14	30,36,37,43	5
5	NO	A	405[A]	2/2	0.98	0.12	18,18,18,22	2
6	SO4	A	407	5/5	0.98	0.08	36,38,44,46	5
3	CU	A	401	1/1	0.99	0.06	17,17,17,17	0

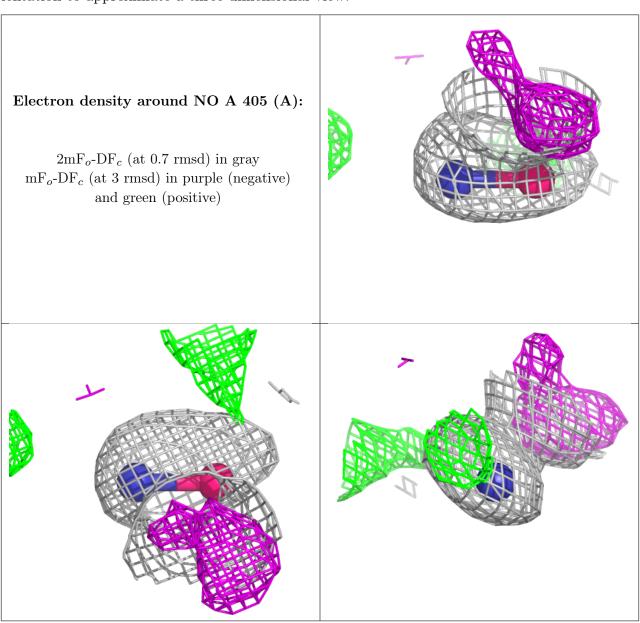
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	${f B-factors}({f \AA}^2)$	Q < 0.9
3	CU	A	402	1/1	1.00	0.05	15,15,15,15	0

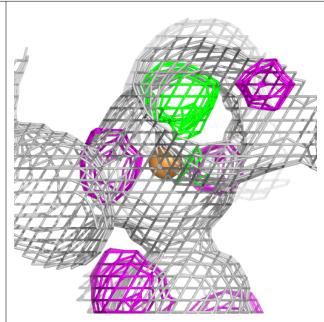
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

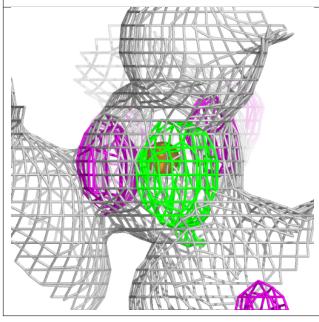


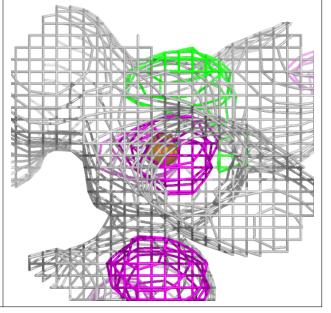


Electron density around CU A 401:

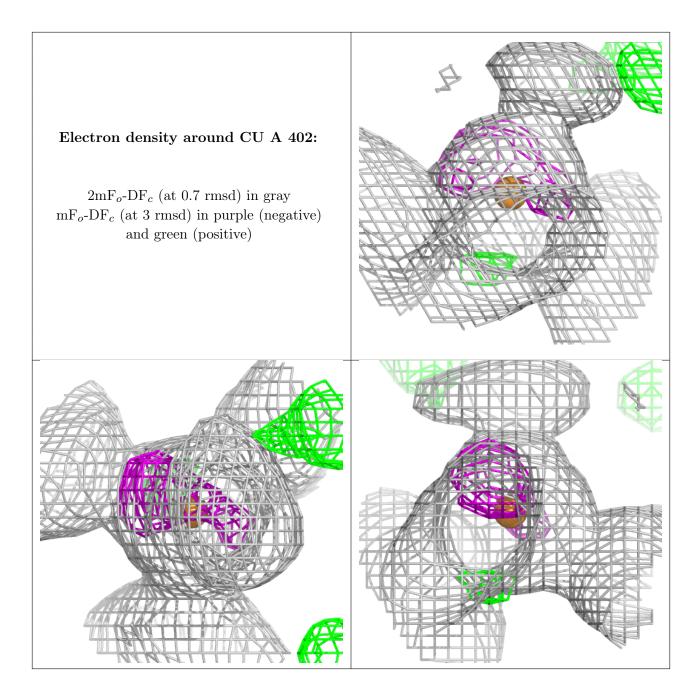
 $2 {
m mF}_o {
m -DF}_c$ (at 0.7 rmsd) in gray ${
m mF}_o {
m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)











6.5 Other polymers (i)

There are no such residues in this entry.

