

wwPDB X-ray Structure Validation Summary Report (i)

Mar 27, 2023 – 03:19 PM EDT

PDB ID : 8FUN

Title: Enzymatically Active, Mn/Fe Metallated Form of AibH1H2

Authors: Powell, M.M.; Rittle, J.

Deposited on : 2023-01-17

Resolution : 2.24 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 (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 (1)) were used in the production of this report:

MolProbity : 4.02b-467

Mogul : 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13

EDS : 2.32.2

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)

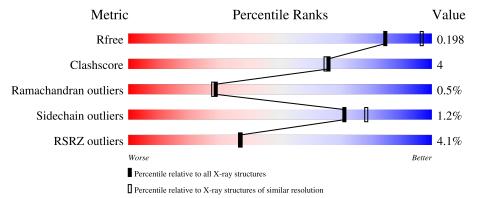
Validation Pipeline (wwPDB-VP) : 2.32.2

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 2.24 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	$(\# ext{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
R_{free}	130704	2391 (2.26-2.22)
Clashscore	141614	2539 (2.26-2.22)
Ramachandran outliers	138981	2489 (2.26-2.22)
Sidechain outliers	138945	2490 (2.26-2.22)
RSRZ outliers	127900	2353 (2.26-2.22)

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	392	84%	• • 11%
1	С	392	82%	7% 10%
2	В	378	84%	11% •••
2	D	378	8%	14% ••



2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 12070 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 Amidohydrolase.

\mathbf{Mol}	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	A	350	Total 2807	C 1791	N 491	O 520	S 5	0	1	0
1	С	351	Total 2817	C 1797	N 495	O 520	S 5	0	2	0

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	-	expression tag	UNP A0A402C2V4
A	-5	GLY	-	expression tag	UNP A0A402C2V4
A	-4	HIS	-	expression tag	UNP A0A402C2V4
A	-3	HIS	-	expression tag	UNP A0A402C2V4
A	-2	HIS	-	expression tag	UNP A0A402C2V4
A	-1	HIS	-	expression tag	UNP A0A402C2V4
A	0	HIS	-	expression tag	UNP A0A402C2V4
A	1	HIS	-	expression tag	UNP A0A402C2V4
A	2	SER	-	expression tag	UNP A0A402C2V4
A	3	GLY	-	expression tag	UNP A0A402C2V4
A	4	GLU	-	expression tag	UNP A0A402C2V4
A	5	ASN	-	expression tag	UNP A0A402C2V4
A	6	LEU	-	expression tag	UNP A0A402C2V4
A	7	TYR	-	expression tag	UNP A0A402C2V4
A	8	PHE	_	expression tag	UNP A0A402C2V4
A	9	GLN	-	expression tag	UNP A0A402C2V4
A	10	SER	-	expression tag	UNP A0A402C2V4
A	11	GLY	_	expression tag	UNP A0A402C2V4
A	12	GLY	-	expression tag	UNP A0A402C2V4
С	-6	MET	_	expression tag	UNP A0A402C2V4
С	-5	GLY	-	expression tag	UNP A0A402C2V4
С	-4	HIS	-	expression tag	UNP A0A402C2V4
С	-3	HIS	-	expression tag	UNP A0A402C2V4
С	-2	HIS	-	expression tag	UNP A0A402C2V4
С	-1	HIS	-	expression tag	UNP A0A402C2V4

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Chain	Residue	Modelled	Actual	Comment	Reference
С	0	HIS	-	expression tag	UNP A0A402C2V4
С	1	HIS	-	expression tag	UNP A0A402C2V4
С	2	SER	-	expression tag	UNP A0A402C2V4
С	3	GLY	-	expression tag	UNP A0A402C2V4
С	4	GLU	-	expression tag	UNP A0A402C2V4
С	5	ASN	-	expression tag	UNP A0A402C2V4
С	6	LEU	-	expression tag	UNP A0A402C2V4
С	7	TYR	-	expression tag	UNP A0A402C2V4
С	8	PHE	-	expression tag	UNP A0A402C2V4
С	9	GLN	-	expression tag	UNP A0A402C2V4
С	10	SER	-	expression tag	UNP A0A402C2V4
С	11	GLY	-	expression tag	UNP A0A402C2V4
C	12	GLY	_	expression tag	UNP A0A402C2V4

• Molecule 2 is a protein called Amidohydrolase.

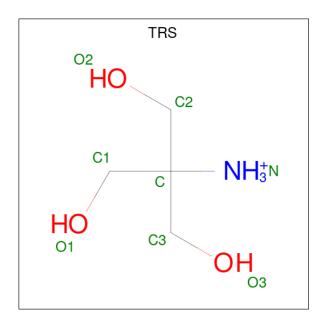
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	D	365	Total	С	N	О	S	0	2	0
2	2 D	303	2883	1836	492	546	9			
2	D	370	Total	С	N	О	S	0	3	0
2	ש	370	2926	1864	504	549	9			

• Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mn 1 1	0	0
3	В	1	Total Mn 1 1	0	0
3	C	1	Total Mn 1 1	0	0
3	D	2	Total Mn 2 2	0	0

• Molecule 4 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	В	1	Total 8	C 4	N 1	O 3	0	0

• Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	В	1	Total Mg 1 1	0	0
5	С	1	Total Mg 1 1	0	0

• Molecule 6 is FE (III) ION (three-letter code: FE) (formula: Fe) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	В	1	Total Fe 1 1	0	0

• Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	164	Total O 164 164	0	0
7	В	161	Total O 161 161	0	0
7	С	156	Total O 156 156	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	D	140	Total O 140 140	0	0

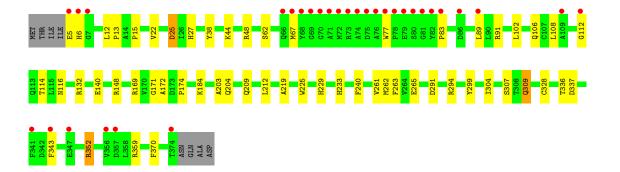


3 Residue-property plots (i)

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: Amidohydrolase Chain A: • Molecule 1: Amidohydrolase Chain C: 82% 10% • Molecule 2: Amidohydrolase Chain B: 11% • Molecule 2: Amidohydrolase Chain D: 83%







4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants	84.14Å 148.92Å 234.09Å	Donositon
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.56 - 2.24	Depositor
Resolution (A)	48.56 - 2.24	EDS
% Data completeness	99.3 (48.56-2.24)	Depositor
(in resolution range)	99.3 (48.56-2.24)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.62 (at 2.24Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
D D.	0.177 , 0.198	Depositor
R, R_{free}	0.177 , 0.198	DCC
R_{free} test set	2000 reflections (2.84%)	wwPDB-VP
Wilson B-factor (Å ²)	25.2	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.36, 37.2	EDS
L-test for twinning ²	$ < L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12070	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 2.73% 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: MN, TRS, FE, MG

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	Bond angles		
Mol Chain	RMSZ	# Z > 5	RMSZ	# Z >5		
1	A	0.53	2/2894~(0.1%)	0.63	$2/3955 \ (0.1\%)$	
1	С	0.56	0/2905	0.64	2/3970 (0.1%)	
2	В	0.67	2/2974~(0.1%)	0.66	2/4063 (0.0%)	
2	D	0.65	0/3021	0.68	2/4125 (0.0%)	
All	All	0.61	4/11794~(0.0%)	0.65	8/16113 (0.0%)	

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
1	A	252	GLU	CD-OE2	-8.83	1.16	1.25
1	A	252	GLU	CD-OE1	-6.03	1.19	1.25
2	В	38	TYR	CE1-CZ	-5.29	1.31	1.38
2	В	228	TYR	CE1-CZ	-5.14	1.31	1.38

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
2	В	323	LEU	CB-CG-CD2	-7.61	98.07	111.00
1	С	350	ARG	NE-CZ-NH2	-5.62	117.49	120.30
2	D	89	LEU	CB-CG-CD2	5.53	120.40	111.00
1	С	83	ASP	CB-CG-OD2	5.47	123.22	118.30
1	A	163	ARG	NE-CZ-NH2	5.42	123.01	120.30

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	2807	0	2705	10	0
1	С	2817	0	2710	15	0
2	В	2883	0	2756	27	0
2	D	2926	0	2806	42	0
3	A	1	0	0	0	0
3	В	1	0	0	0	0
3	С	1	0	0	0	0
3	D	2	0	0	0	0
4	В	8	0	12	1	0
5	В	1	0	0	0	0
5	С	1	0	0	0	0
6	В	1	0	0	0	0
7	A	164	0	0	2	0
7	В	161	0	0	1	0
7	С	156	0	0	0	0
7	D	140	0	0	4	0
All	All	12070	0	10989	89	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 4.

The worst 5 of 89 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{aligned} & ext{Interatomic} \ & ext{distance (Å)} \end{aligned}$	Clash overlap (Å)	
2:D:352:ARG:HH11	2:D:352:ARG:HG3	1.10	1.07	
2:B:113:GLN:CB	2:B:116:ASN:HD21	1.84	0.91	
2:D:6:HIS:ND1	2:D:6:HIS:O	2.06	0.88	
1:A:270:PHE:HB3	1:A:308:VAL:HG11	1.56	0.87	
2:D:352:ARG:HG3	2:D:352:ARG:NH1	1.88	0.85	

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	Percen	tiles
1	A	349/392 (89%)	345 (99%)	3 (1%)	1 (0%)	41	44
1	С	351/392 (90%)	346 (99%)	4 (1%)	1 (0%)	41	44
2	В	365/378~(97%)	342 (94%)	21 (6%)	2 (0%)	29	28
2	D	371/378 (98%)	345 (93%)	23 (6%)	3 (1%)	19	16
All	All	1436/1540 (93%)	1378 (96%)	51 (4%)	7 (0%)	29	28

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	В	328	CYS
2	D	112	GLY
2	D	328	CYS
1	С	270	PHE
1	A	270	PHE

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	Percer	ntiles
1	A	300/332~(90%)	299 (100%)	1 (0%)	92	95
1	С	300/332 (90%)	296 (99%)	4 (1%)	69	76
2	В	307/318 (96%)	301 (98%)	6 (2%)	55	62
2	D	310/318 (98%)	306 (99%)	4 (1%)	69	76
All	All	1217/1300 (94%)	1202 (99%)	15 (1%)	71	78



5 of 15 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	С	82	SER
2	D	309	GLN
1	С	127	ARG
2	D	352	ARG
2	D	108	LEU

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	58	GLN
2	В	31	GLN
2	D	106	GLN
2	D	116	ASN
2	D	309	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 9 ligands modelled in this entry, 8 are monoatomic - leaving 1 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	\mathbf{B}_{0}	ond leng	gths	В	ond ang	gles
Moi Type Cr	Chain Res	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2		
4	TRS	В	501	-	7,7,7	0.45	0	9,9,9	0.66	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
4	TRS	В	501	-	-	1/9/9/9	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	В	501	TRS	N-C-C1-O1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	В	501	TRS	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, 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	<rsrz></rsrz>	$\# \mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q<0.9
1	A	350/392 (89%)	-0.20	10 (2%) 51 52	16, 23, 38, 68	0
1	С	351/392 (89%)	-0.20	5 (1%) 75 76	17, 25, 38, 67	0
2	В	365/378~(96%)	-0.14	13 (3%) 42 42	15, 23, 44, 72	0
2	D	370/378 (97%)	0.14	31 (8%) 11 11	16, 26, 67, 98	0
All	All	1436/1540 (93%)	-0.10	59 (4%) 37 37	15, 24, 50, 98	0

The worst 5 of 59 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	77	TRP	7.8
2	D	69	GLY	6.0
2	D	80	SER	5.3
2	D	68	TYR	5.3
2	D	78	PRO	5.3

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 monosaccharides in this entry.

6.4 Ligands (i)

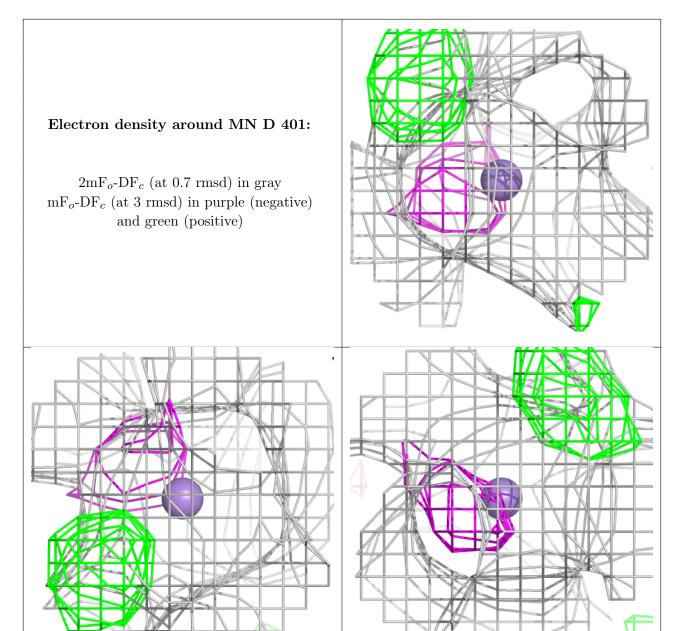
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.



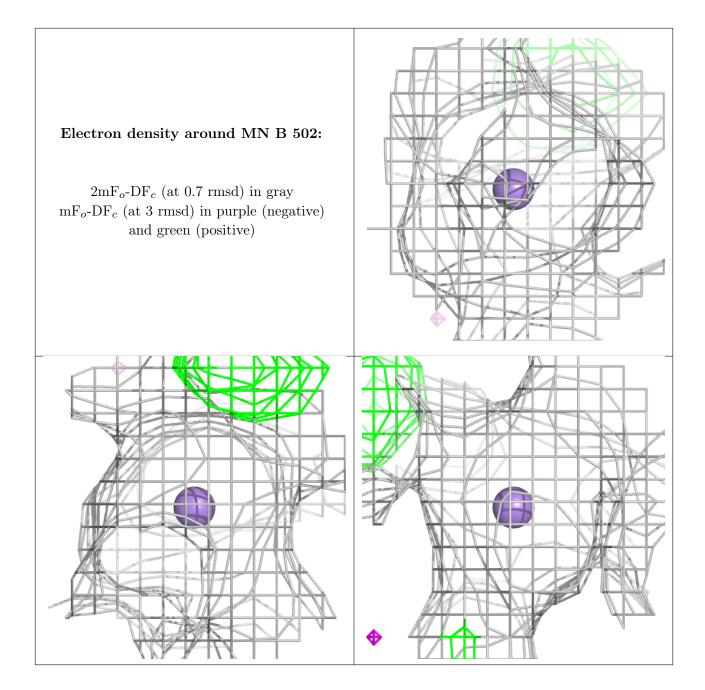
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
4	TRS	В	501	8/8	0.84	0.32	35,42,48,56	0
5	MG	С	402	1/1	0.87	0.16	35,35,35,35	0
3	MN	D	401	1/1	0.93	0.06	35,35,35,35	0
5	MG	В	503	1/1	0.94	0.08	36,36,36,36	1
3	MN	В	502	1/1	0.94	0.04	36,36,36,36	0
3	MN	D	402	1/1	0.96	0.06	46,46,46,46	1
3	MN	A	401	1/1	0.99	0.06	24,24,24,24	0
6	FE	В	504	1/1	0.99	0.04	33,33,33,33	0
3	MN	С	401	1/1	1.00	0.06	25,25,25,25	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.

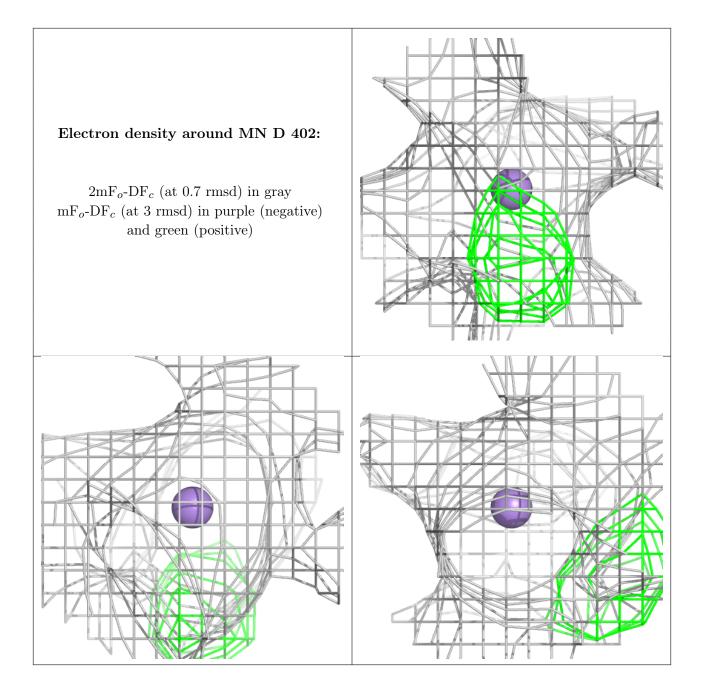








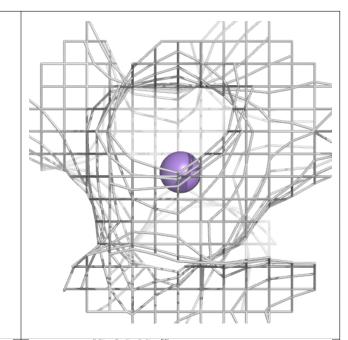


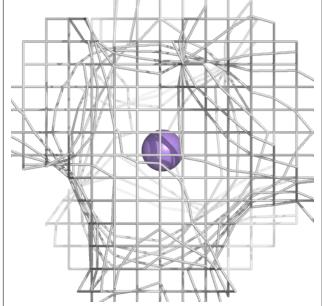


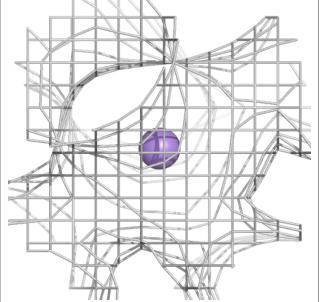


Electron density around MN A 401:

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

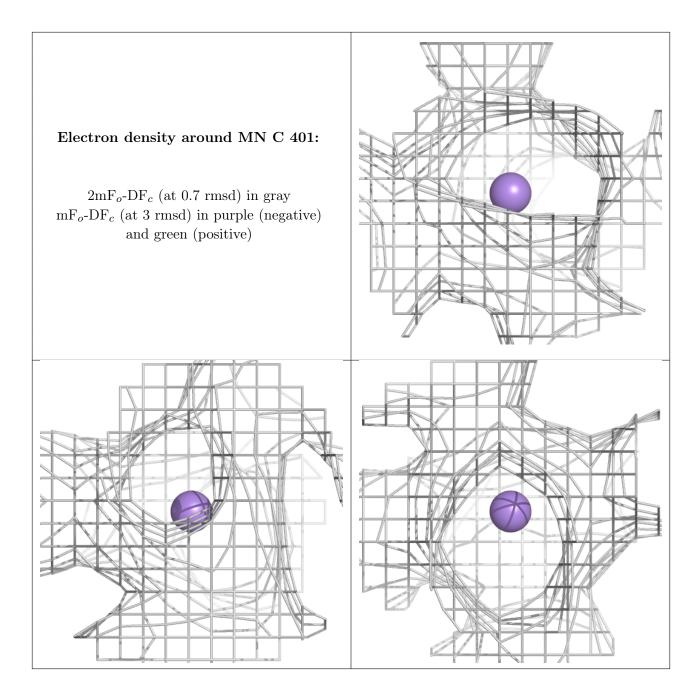






Electron density around FE B 504: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray $\mathrm{mF}_{o}\text{-}\mathrm{DF}_{c}$ (at 3 rmsd) in purple (negative) and green (positive)





6.5 Other polymers (i)

There are no such residues in this entry.

