



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 30, 2025 – 03:33 am BST

PDB ID : 9ERR / pdb\_00009err  
Title : Hydrogenase-2 Ni-SCO state  
Authors : Carr, S.B.; Li, W.; Wong, K.l.; Ash, P.A.; Vincent, K.A.  
Deposited on : 2024-03-25  
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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : ?? (???), CSD ??CSD?? (????)  
Xtriage (Phenix) : 1.13  
EDS : 3.0  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.003 (Gargrove)  
Density-Fitness : 1.0.11  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.42

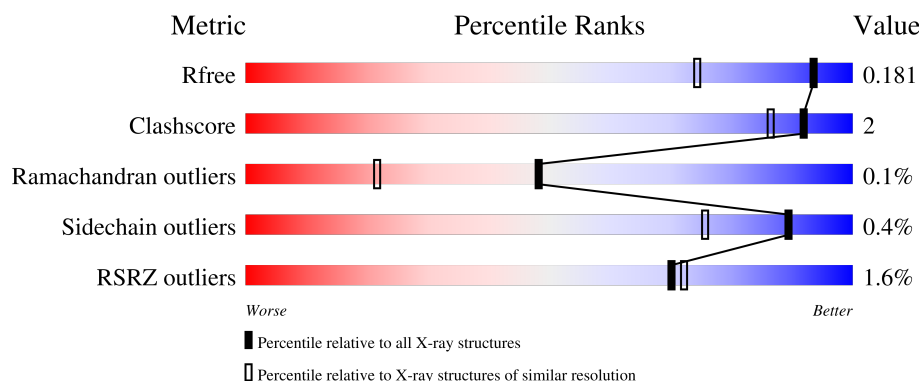
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.40 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	164625	2247 (1.40-1.40)
Clashscore	180529	2446 (1.40-1.40)
Ramachandran outliers	177936	2398 (1.40-1.40)
Sidechain outliers	177891	2397 (1.40-1.40)
RSRZ outliers	164620	2246 (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  $\geq 3$ , 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	S	298	
1	T	298	
2	L	567	
2	M	567	

## 2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 14020 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 Hydrogenase-2 small chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	S	269	Total	C	N	O	S	0	5	0
			2083	1318	368	384	13			
1	T	269	Total	C	N	O	S	0	6	0
			2083	1320	366	384	13			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
S	294	HIS	-	expression tag	UNP P69741
S	295	HIS	-	expression tag	UNP P69741
S	296	HIS	-	expression tag	UNP P69741
S	297	HIS	-	expression tag	UNP P69741
S	298	HIS	-	expression tag	UNP P69741
S	299	HIS	-	expression tag	UNP P69741
T	294	HIS	-	expression tag	UNP P69741
T	295	HIS	-	expression tag	UNP P69741
T	296	HIS	-	expression tag	UNP P69741
T	297	HIS	-	expression tag	UNP P69741
T	298	HIS	-	expression tag	UNP P69741
T	299	HIS	-	expression tag	UNP P69741

- Molecule 2 is a protein called Hydrogenase-2 large chain.

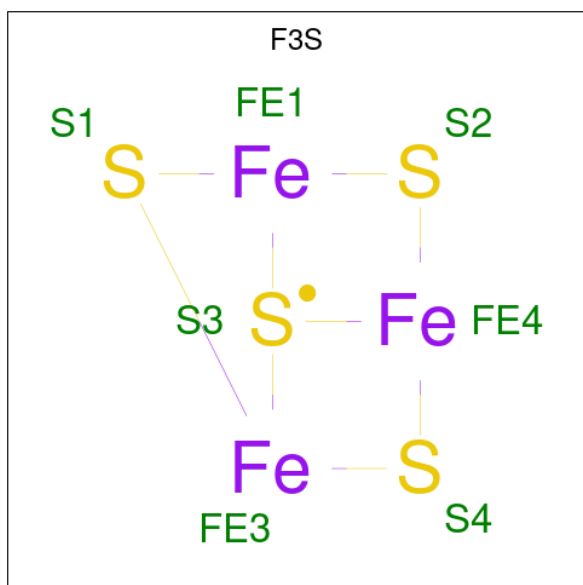
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	551	Total	C	N	O	S	0	6	0
			4325	2751	741	815	18			
2	M	551	Total	C	N	O	S	0	2	0
			4298	2735	739	806	18			

- Molecule 3 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe<sub>4</sub>S<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	S	1	Total	Fe	S	0	0
			8	4	4		
3	S	1	Total	Fe	S	0	0
			8	4	4		
3	T	1	Total	Fe	S	0	0
			8	4	4		
3	T	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 4 is FE3-S4 CLUSTER (CCD ID: F3S) (formula:  $\text{Fe}_3\text{S}_4$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	S	1	Total	Fe	S	0	0
			7	3	4		
4	T	1	Total	Fe	S	0	0
			7	3	4		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	S	1	Total	C	O	0	0
			6	3	3		
5	T	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is CARBONMONOXIDE-(DICYANO) IRON (CCD ID: FCO) (formula:  $C_3FeN_2O$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	L	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		
6	M	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		

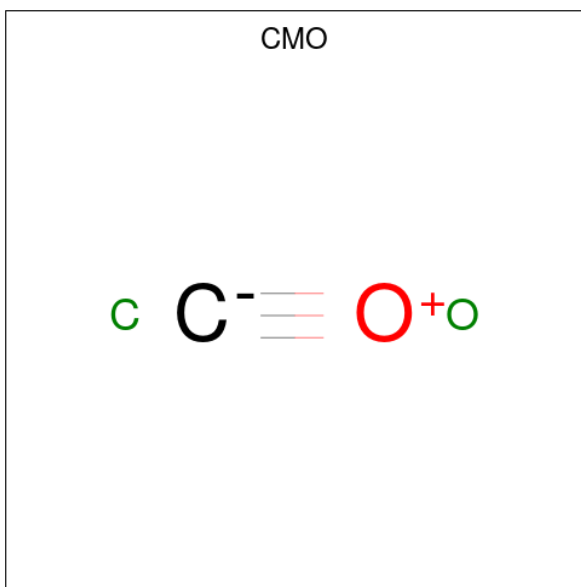
- Molecule 7 is NICKEL (II) ION (CCD ID: NI) (formula: Ni) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	L	1	Total	Ni	0	0
			1	1		
7	M	1	Total	Ni	0	0
			1	1		

- Molecule 8 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	L	2	Total	Mg	0	0
			2	2		
8	M	3	Total	Mg	0	0
			3	3		

- Molecule 9 is CARBON MONOXIDE (CCD ID: CMO) (formula: CO) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	L	1	Total	C	O	0	0
			2	1	1		
9	M	1	Total	C	O	0	0
			2	1	1		

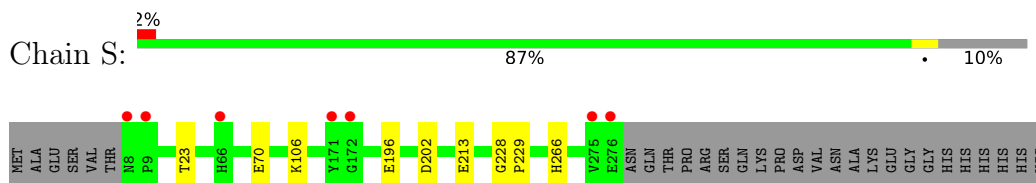
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	S	179	Total	O	0	0
			179	179		
10	L	425	Total	O	0	0
			425	425		
10	T	167	Total	O	0	0
			167	167		
10	M	377	Total	O	0	0
			377	377		

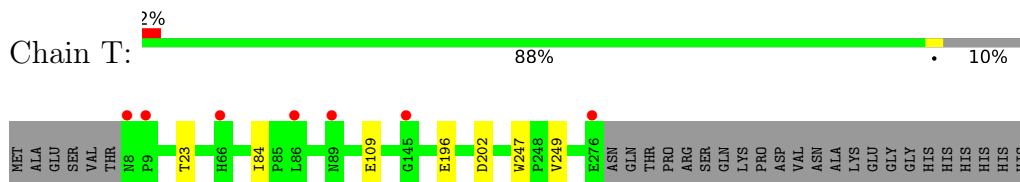
### 3 Residue-property plots [i](#)

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.

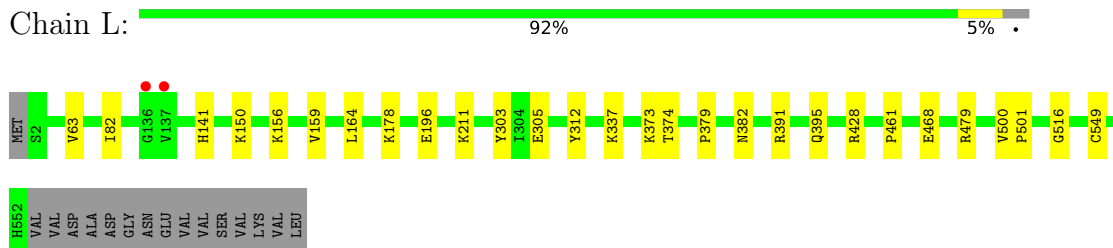
- Molecule 1: Hydrogenase-2 small chain



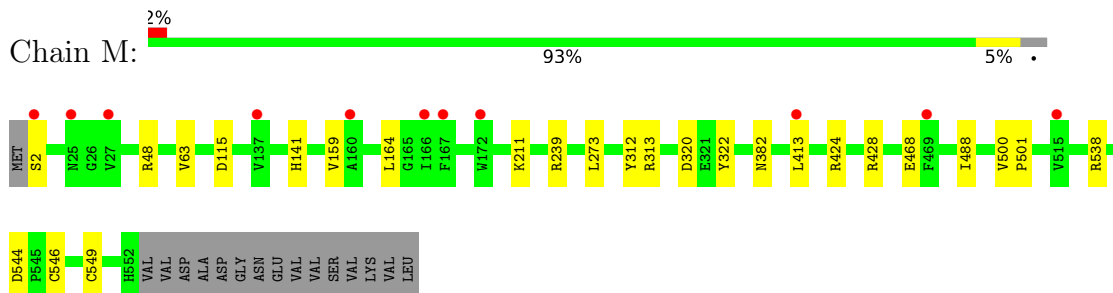
- Molecule 1: Hydrogenase-2 small chain



- Molecule 2: Hydrogenase-2 large chain



- Molecule 2: Hydrogenase-2 large chain





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	100.19Å 99.75Å 168.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	64.50 – 1.40 64.50 – 1.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (64.50-1.40) 99.9 (64.50-1.40)	Depositor EDS
$R_{merge}$	0.26	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.31 (at 1.40Å)	Xtriage
Refinement program	REFMAC 5.8.0419	Depositor
R, $R_{free}$	0.161 , 0.181 0.162 , 0.181	Depositor DCC
$R_{free}$ test set	16433 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	11.7	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 34.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.005 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	14020	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.58% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: F3S, MG, GOL, CMO, NI, SF4, FCO

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	S	0.53	0/2147	0.78	0/2923
1	T	0.49	0/2156	0.79	0/2935
2	L	0.53	1/4444 (0.0%)	0.80	2/6057 (0.0%)
2	M	0.52	0/4411	0.83	7/6013 (0.1%)
All	All	0.52	1/13158 (0.0%)	0.81	9/17928 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	468	GLU	CD-OE1	-6.12	1.19	1.25

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	424	ARG	NE-CZ-NH2	-6.92	116.84	120.30
2	M	382	ASN	N-CA-CB	-6.72	98.50	110.60
2	M	424	ARG	NE-CZ-NH1	6.60	123.60	120.30
2	L	382	ASN	N-CA-CB	-6.21	99.42	110.60
2	M	48	ARG	NE-CZ-NH2	-5.73	117.44	120.30
2	M	428	ARG	NE-CZ-NH2	-5.34	117.63	120.30
2	M	239	ARG	NE-CZ-NH2	-5.30	117.65	120.30
2	M	544	ASP	CB-CG-OD2	-5.10	113.71	118.30
2	L	428	ARG	NE-CZ-NH2	-5.03	117.79	120.30

There are no chirality outliers.

There are no planarity outliers.

## 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	S	2083	0	2005	8	0
1	T	2083	0	2013	7	0
2	L	4325	0	4261	17	0
2	M	4298	0	4240	11	0
3	S	16	0	0	0	0
3	T	16	0	0	0	0
4	S	7	0	0	0	0
4	T	7	0	0	0	0
5	S	6	0	8	0	0
5	T	6	0	8	0	0
6	L	7	0	0	0	0
6	M	7	0	0	0	0
7	L	1	0	0	0	0
7	M	1	0	0	0	0
8	L	2	0	0	0	0
8	M	3	0	0	0	0
9	L	2	0	0	1	0
9	M	2	0	0	1	0
10	L	425	0	0	5	0
10	M	377	0	0	0	0
10	S	179	0	0	4	0
10	T	167	0	0	2	0
All	All	14020	0	12535	39	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 (39) 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:S:213:GLU:OE2	10:S:401:HOH:O	1.79	0.99
1:T:109:GLU:HG2	10:T:556:HOH:O	1.78	0.82
2:L:196[A]:GLU:OE2	10:L:701:HOH:O	2.01	0.77
2:M:273:LEU:HD13	2:M:413:LEU:HD21	1.77	0.67
2:L:374[B]:THR:HG21	2:L:516:GLY:HA2	1.74	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:106:LYS:HD3	10:S:562:HOH:O	1.95	0.66
2:L:305[A]:GLU:CD	2:L:373:LYS:HZ2	2.08	0.58
2:M:313:ARG:HD3	2:M:322:TYR:CE2	2.41	0.55
1:S:70:GLU:HG2	10:S:471:HOH:O	2.08	0.54
2:L:391:ARG:O	2:L:395[A]:GLN:HG3	2.09	0.52
2:L:303:TYR:CE2	2:L:305[B]:GLU:CG	2.95	0.50
1:S:266:HIS:NE2	10:S:402:HOH:O	2.34	0.50
2:L:305[A]:GLU:CD	2:L:373:LYS:NZ	2.65	0.49
2:L:374[B]:THR:CG2	2:L:516:GLY:HA2	2.42	0.49
2:M:159:VAL:HG22	2:M:164:LEU:HD11	1.95	0.48
2:M:546:CYS:CB	9:M:604:CMO:C	2.90	0.48
1:T:109:GLU:CG	10:T:556:HOH:O	2.51	0.47
2:L:379:PRO:HA	10:L:1051:HOH:O	2.15	0.47
2:M:273:LEU:CD1	2:M:413:LEU:HD21	2.43	0.46
2:M:500:VAL:CG1	2:M:501:PRO:HD2	2.45	0.46
2:M:468:GLU:HA	2:M:488:ILE:O	2.17	0.45
2:L:337:LYS:HD2	10:L:1103:HOH:O	2.17	0.44
2:L:479:ARG:CZ	9:L:604:CMO:O	2.65	0.44
1:T:23:THR:HB	2:M:63:VAL:HG23	2.00	0.44
2:M:320:ASP:OD1	2:M:322:TYR:HB3	2.18	0.43
1:S:23:THR:HB	2:L:63:VAL:HG23	2.00	0.43
2:M:115:ASP:HB3	2:M:538:ARG:HG2	2.00	0.43
2:M:500:VAL:HG11	2:M:549:CYS:HB3	2.01	0.42
2:L:159:VAL:HG22	2:L:164:LEU:HD11	2.01	0.42
1:S:202:ASP:OD2	1:T:202:ASP:OD2	2.37	0.42
2:L:82:ILE:HD11	2:L:461:PRO:HB3	2.02	0.42
1:T:84:ILE:HD12	1:T:84:ILE:N	2.35	0.42
1:S:228:GLY:N	1:S:229:PRO:CD	2.83	0.42
1:T:247:TRP:CZ2	1:T:249:VAL:HB	2.55	0.41
2:L:156:LYS:HE2	10:L:849:HOH:O	2.20	0.41
2:L:500:VAL:HG11	2:L:549:CYS:HB3	2.02	0.41
1:S:196:GLU:HA	1:T:196:GLU:HA	2.03	0.41
2:L:178:LYS:HE2	10:L:720:HOH:O	2.20	0.41
2:L:500:VAL:CG1	2:L:501:PRO:HD2	2.51	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	S	272/298 (91%)	262 (96%)	10 (4%)	0	100	100
1	T	273/298 (92%)	263 (96%)	10 (4%)	0	100	100
2	L	555/567 (98%)	532 (96%)	22 (4%)	1 (0%)	44	20
2	M	551/567 (97%)	528 (96%)	22 (4%)	1 (0%)	44	20
All	All	1651/1730 (95%)	1585 (96%)	64 (4%)	2 (0%)	48	22

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	M	211	LYS
2	L	211	LYS

### 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	S	219/239 (92%)	219 (100%)	0	100	100
1	T	220/239 (92%)	220 (100%)	0	100	100
2	L	471/479 (98%)	468 (99%)	3 (1%)	84	66
2	M	467/479 (98%)	464 (99%)	3 (1%)	84	66
All	All	1377/1436 (96%)	1371 (100%)	6 (0%)	89	76

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	L	141	HIS
2	L	150	LYS
2	L	312	TYR
2	M	2	SER
2	M	141	HIS
2	M	312	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 19 ligands modelled in this entry, 7 are monoatomic - leaving 12 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

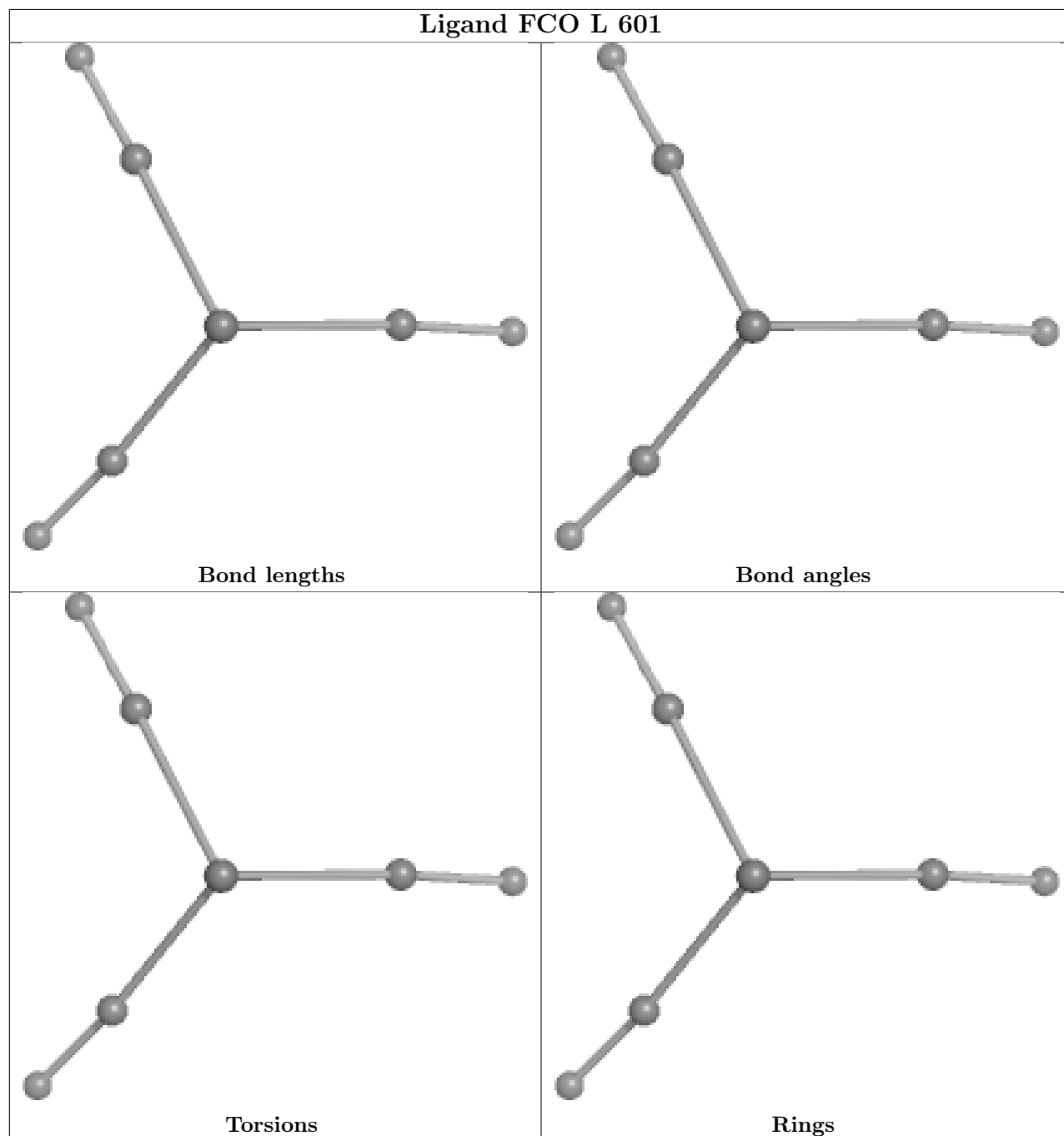
There are no torsion outliers.

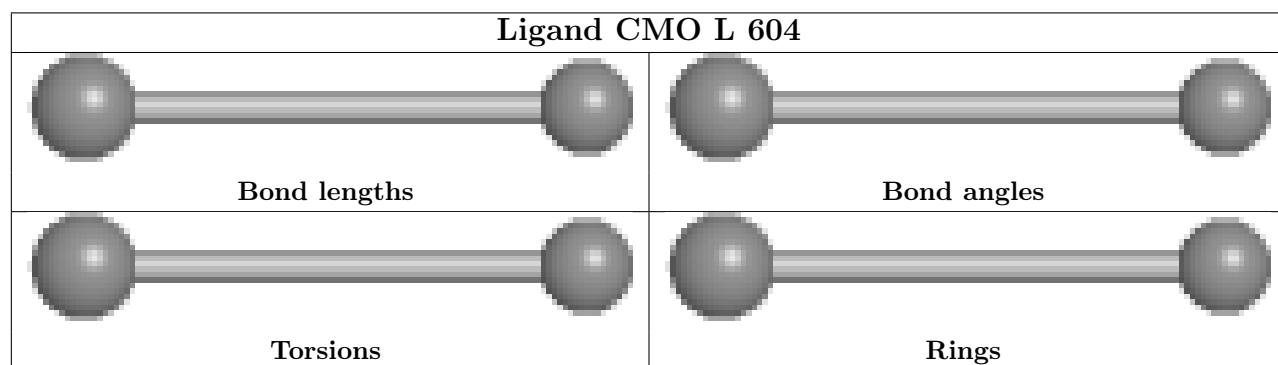
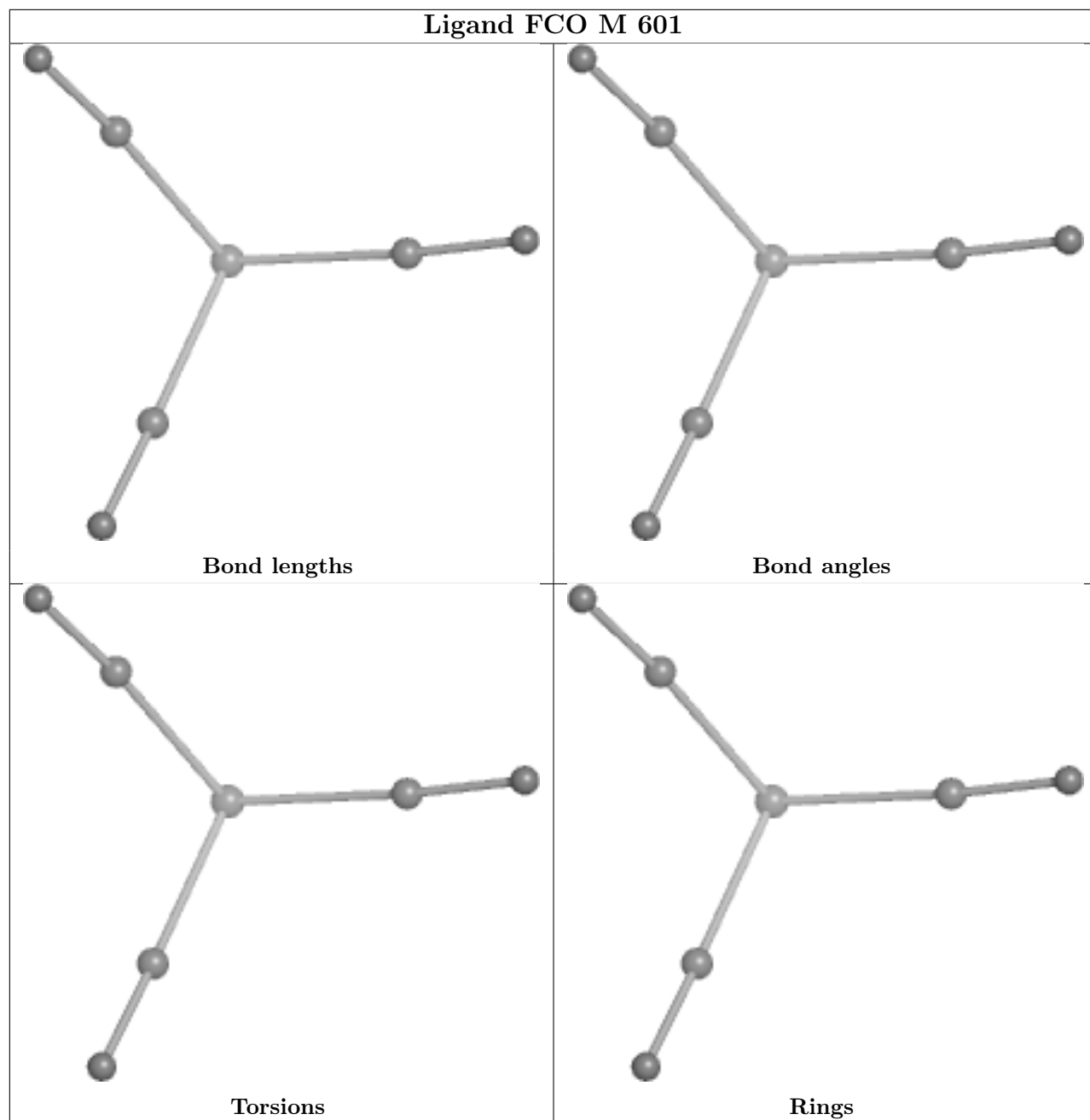
There are no ring outliers.

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 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.







Ligand CMO M 604			
 Bond lengths		 Bond angles	
 Torsions		 Rings	

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	S	269/298 (90%)	-0.17	7 (2%) 57 58	5, 13, 27, 47	5 (1%)
1	T	269/298 (90%)	-0.02	7 (2%) 57 58	7, 14, 29, 49	6 (2%)
2	L	551/567 (97%)	-0.39	2 (0%) 89 90	6, 11, 23, 33	6 (1%)
2	M	551/567 (97%)	-0.17	11 (1%) 64 66	7, 13, 27, 43	2 (0%)
All	All	1640/1730 (94%)	-0.22	27 (1%) 70 72	5, 13, 26, 49	19 (1%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	M	2	SER	4.5
1	T	9	PRO	4.1
1	S	9	PRO	3.4
2	M	27	VAL	3.3
1	T	8	ASN	3.1
2	M	137	VAL	3.1
2	M	166	ILE	3.0
2	L	136	GLY	3.0
2	L	137	VAL	2.8
2	M	413	LEU	2.7
1	S	8	ASN	2.6
2	M	167	PHE	2.6
1	S	276	GLU	2.6
1	T	145	GLY	2.5
1	T	86	LEU	2.5
1	S	171	TYR	2.4
2	M	160	ALA	2.4
1	S	66	HIS	2.4
1	S	275	VAL	2.3
2	M	25	ASN	2.2
1	T	66	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	T	89	ASN	2.2
1	T	276	GLU	2.2
1	S	172	GLY	2.2
2	M	515	VAL	2.1
2	M	469	PHE	2.1
2	M	172	TRP	2.0

## 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<sup>th</sup> 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.

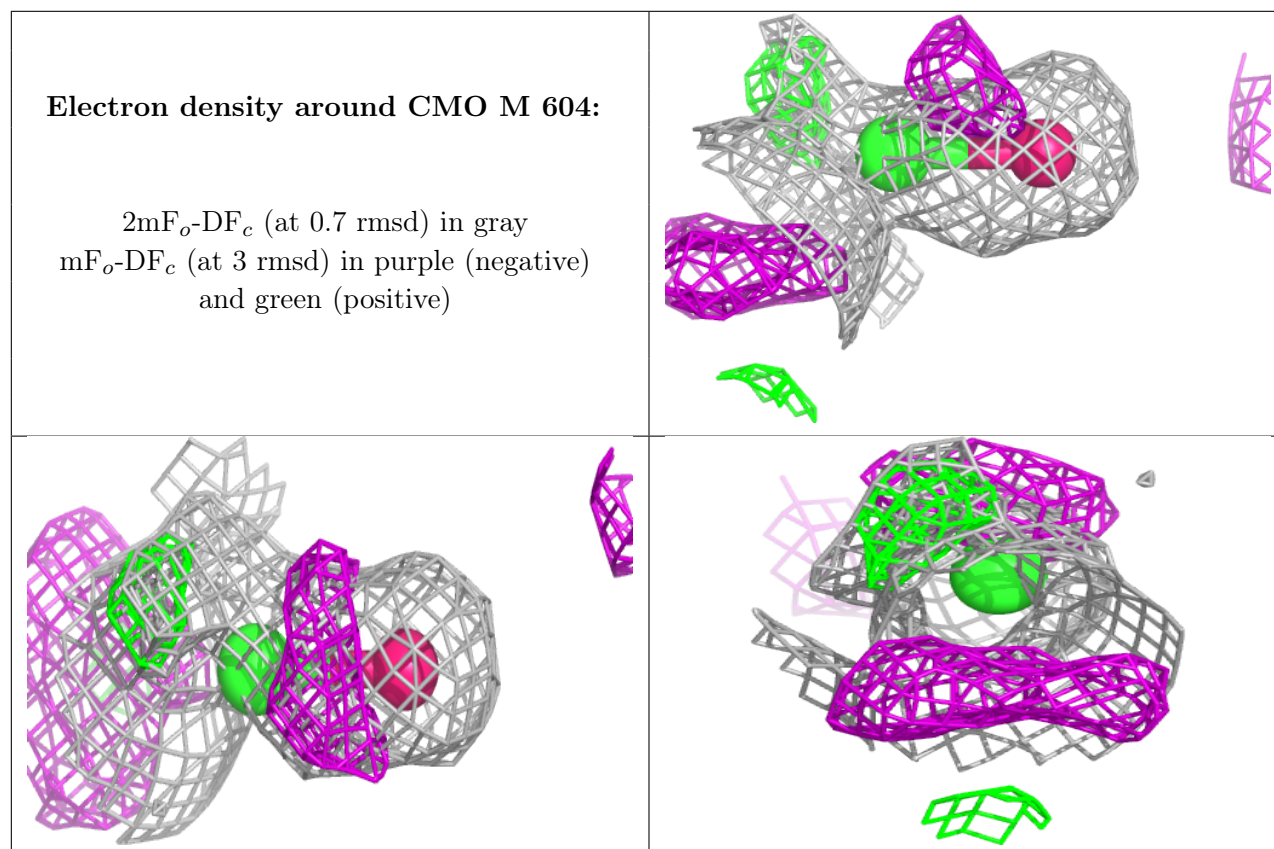
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
8	MG	M	606	1/1	0.87	0.20	43,43,43,43	0
9	CMO	M	604	2/2	0.87	0.23	20,20,20,24	0
8	MG	L	605	1/1	0.91	0.11	30,30,30,30	0
9	CMO	L	604	2/2	0.93	0.22	18,18,18,20	0
5	GOL	T	304	6/6	0.95	0.09	17,22,22,24	0
8	MG	M	605	1/1	0.96	0.07	18,18,18,18	0
5	GOL	S	304	6/6	0.96	0.07	18,20,22,22	0
3	SF4	S	303	8/8	1.00	0.02	8,9,9,9	0
6	FCO	L	601	7/7	1.00	0.03	7,8,8,9	0
6	FCO	M	601	7/7	1.00	0.03	8,8,10,11	0
7	NI	L	602	1/1	1.00	0.02	10,10,10,10	0
7	NI	M	602	1/1	1.00	0.03	11,11,11,11	0
8	MG	L	603	1/1	1.00	0.05	4,4,4,4	0
3	SF4	T	301	8/8	1.00	0.02	10,11,11,11	0
8	MG	M	603	1/1	1.00	0.06	6,6,6,6	0
3	SF4	T	303	8/8	1.00	0.02	9,10,11,11	0
4	F3S	S	302	7/7	1.00	0.02	9,9,9,10	0

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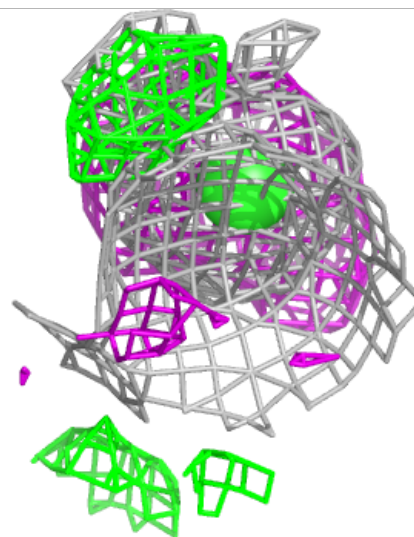
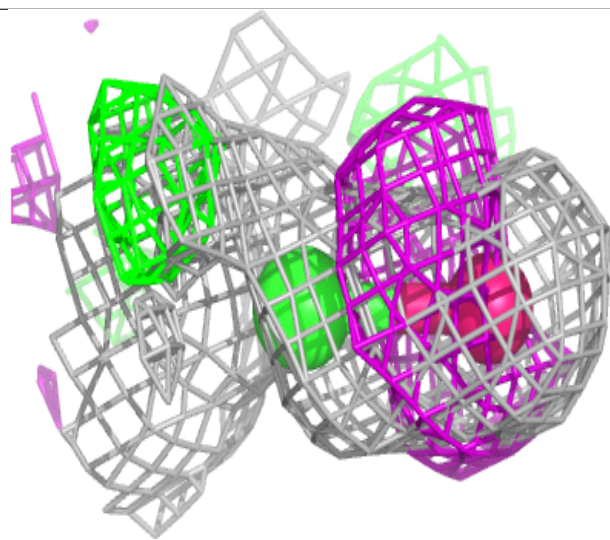
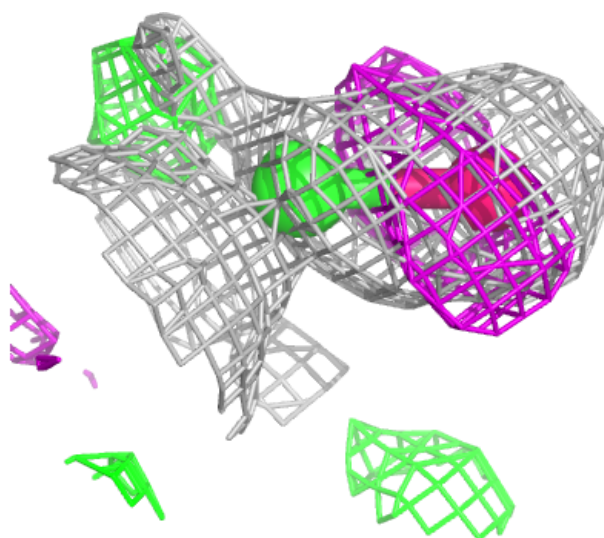
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	F3S	T	302	7/7	1.00	0.02	9,9,10,10	0
3	SF4	S	301	8/8	1.00	0.02	10,11,11,11	0

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.



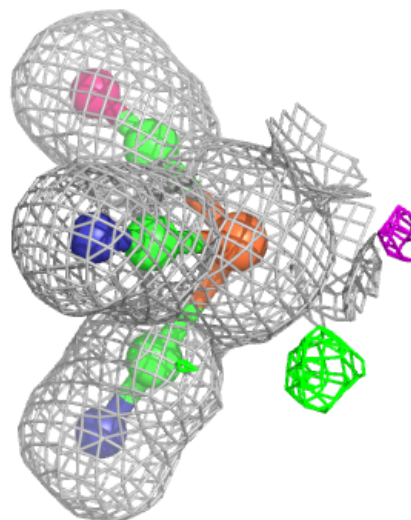
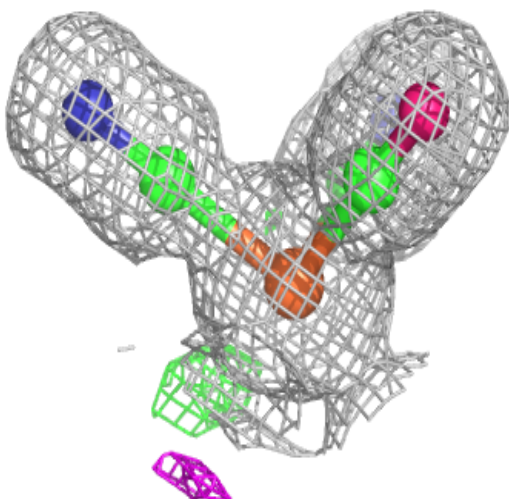
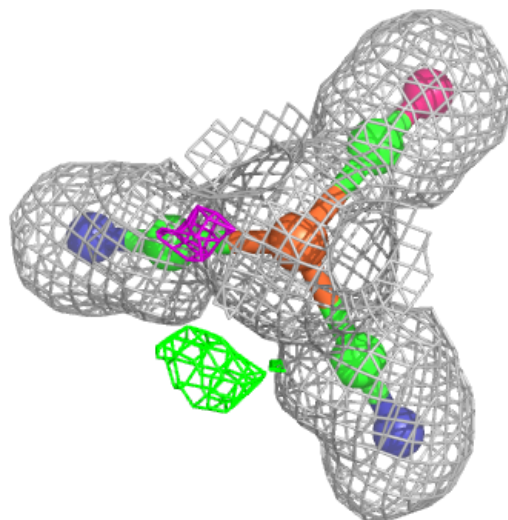
**Electron density around CMO L 604:**

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



**Electron density around FCO L 601:**

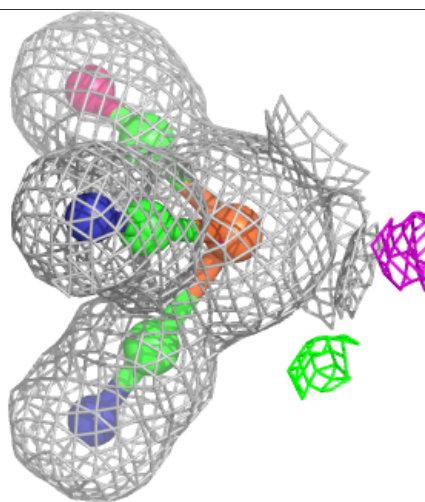
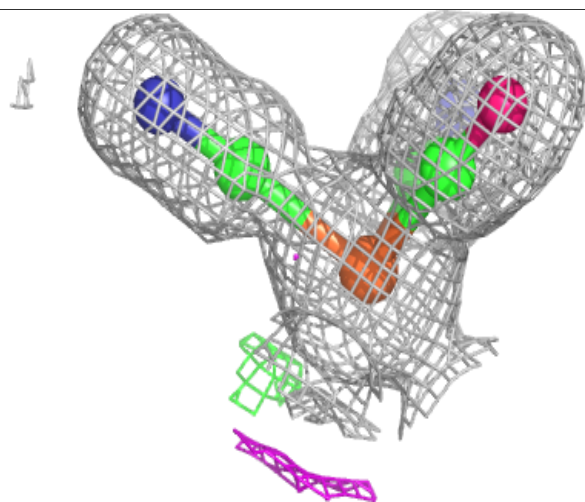
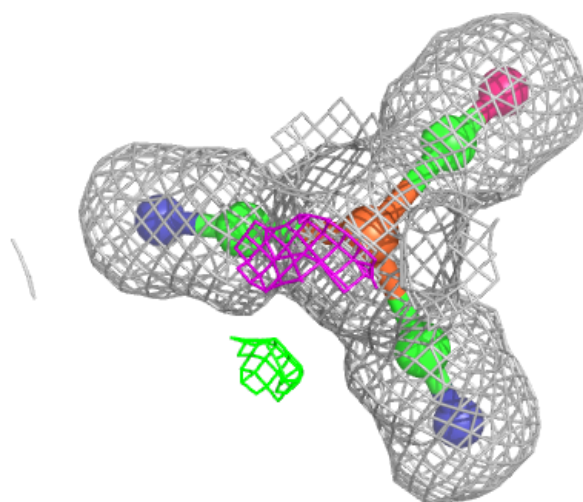
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





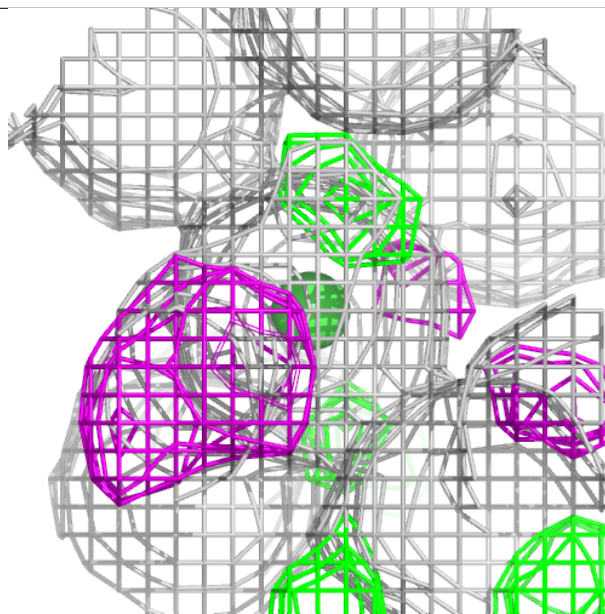
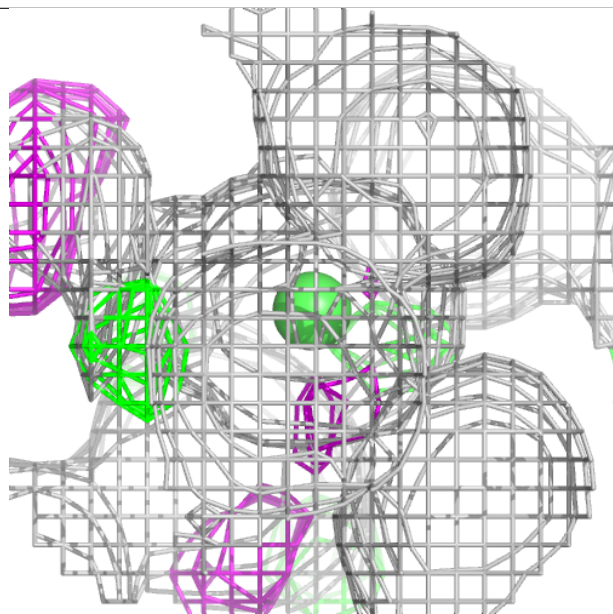
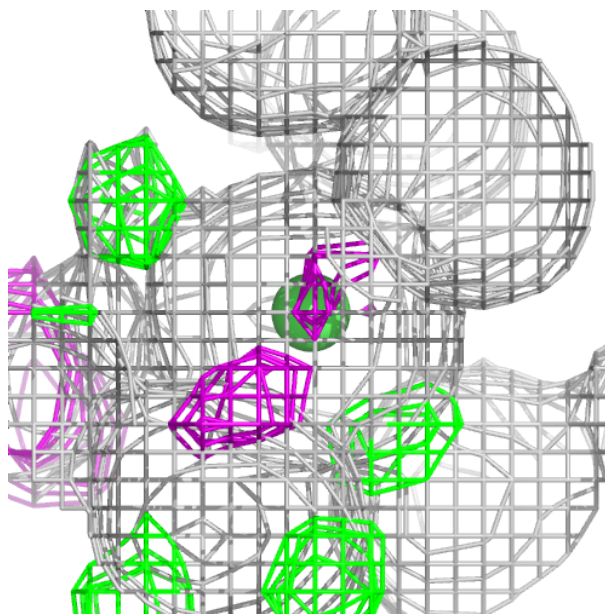
**Electron density around FCO M 601:**

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



**Electron density around NI L 602:**

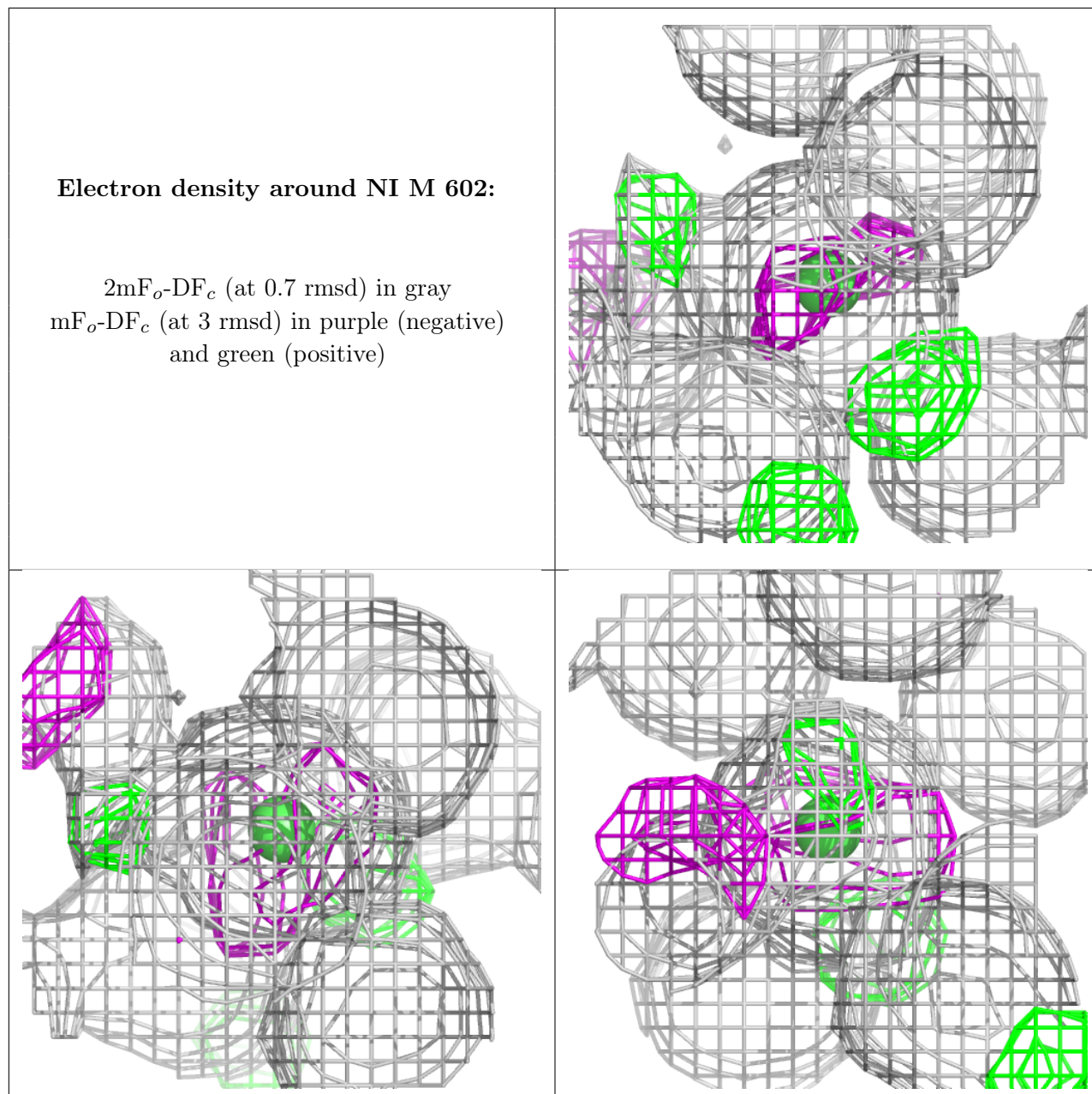
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around NI M 602:**

$2mF_o-DF_c$  (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.