



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 18, 2021 – 08:13 PM JST

PDB ID : 6LL0
Title : Carbazole-soaked reduced oxygenase in carbazole 1,9a-dioxygenase
Authors : Wang, Y.X.; Suzuki-Minakuchi, C.; Nojiri, H.
Deposited on : 2019-12-21
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 following versions of software and data (see [references ⓘ](#)) 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 : 2.16
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) : 2.16

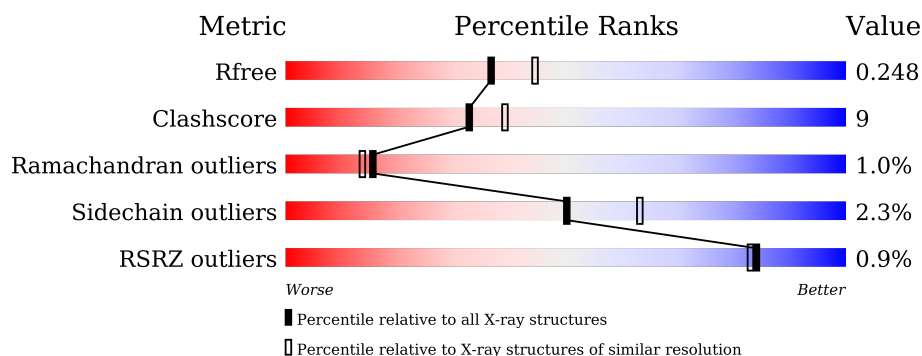
1 Overall quality at a glance

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}	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (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	392	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 17%, green 80%, grey 4%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 80% 17% .. </div> </div>
1	B	392	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 84%, yellow 13%, grey 3%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 84% 13% . </div> </div>
1	C	392	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 14%, green 82%, grey 2%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 82% 14% .. </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	EDO	B	405	-	-	X	-
5	EDO	B	411	-	-	X	-
6	PEG	A	411	-	-	X	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 9950 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 Terminal oxygenase component of carbazole.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	383	Total	C	N	O	S	0	18	0
			3210	2047	548	602	13			
1	B	383	Total	C	N	O	S	0	1	0
			3090	1975	524	578	13			
1	C	383	Total	C	N	O	S	0	7	0
			3133	2002	531	587	13			

There are 24 discrepancies between the modelled and reference sequences:

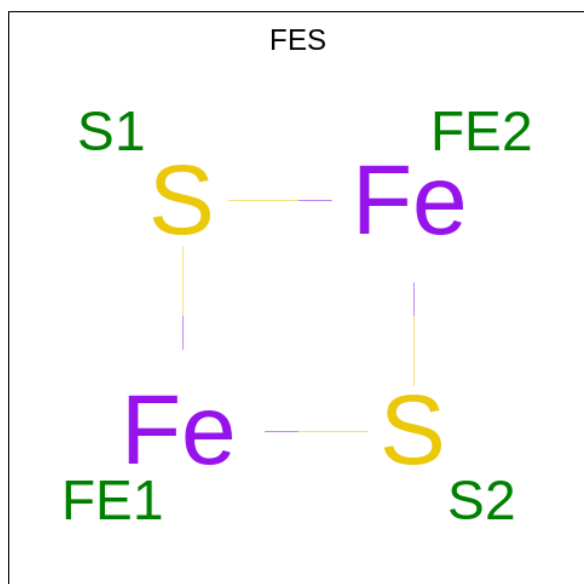
Chain	Residue	Modelled	Actual	Comment	Reference
A	385	LEU	-	expression tag	UNP Q84II6
A	386	GLU	-	expression tag	UNP Q84II6
A	387	HIS	-	expression tag	UNP Q84II6
A	388	HIS	-	expression tag	UNP Q84II6
A	389	HIS	-	expression tag	UNP Q84II6
A	390	HIS	-	expression tag	UNP Q84II6
A	391	HIS	-	expression tag	UNP Q84II6
A	392	HIS	-	expression tag	UNP Q84II6
B	385	LEU	-	expression tag	UNP Q84II6
B	386	GLU	-	expression tag	UNP Q84II6
B	387	HIS	-	expression tag	UNP Q84II6
B	388	HIS	-	expression tag	UNP Q84II6
B	389	HIS	-	expression tag	UNP Q84II6
B	390	HIS	-	expression tag	UNP Q84II6
B	391	HIS	-	expression tag	UNP Q84II6
B	392	HIS	-	expression tag	UNP Q84II6
C	385	LEU	-	expression tag	UNP Q84II6
C	386	GLU	-	expression tag	UNP Q84II6
C	387	HIS	-	expression tag	UNP Q84II6
C	388	HIS	-	expression tag	UNP Q84II6
C	389	HIS	-	expression tag	UNP Q84II6
C	390	HIS	-	expression tag	UNP Q84II6
C	391	HIS	-	expression tag	UNP Q84II6

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	392	HIS	-	expression tag	UNP Q84II6

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe_2S_2).

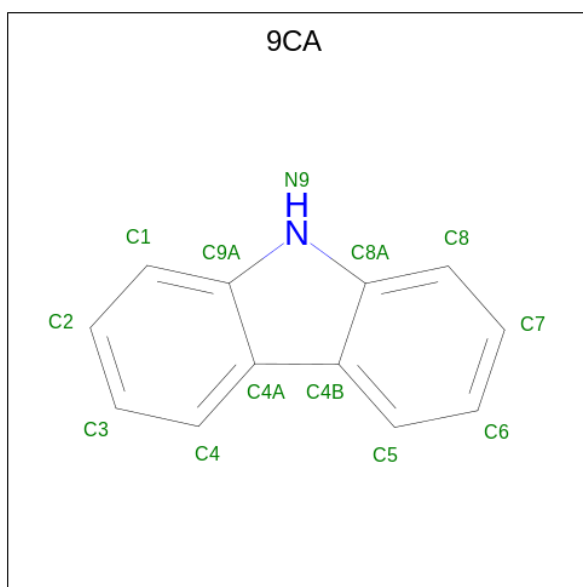


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			4	2	2		
2	B	1	Total	Fe	S	0	0
			4	2	2		
2	C	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 3 is FE (II) ION (three-letter code: FE2) (formula: Fe) (labeled as "Ligand of Interest" by depositor).

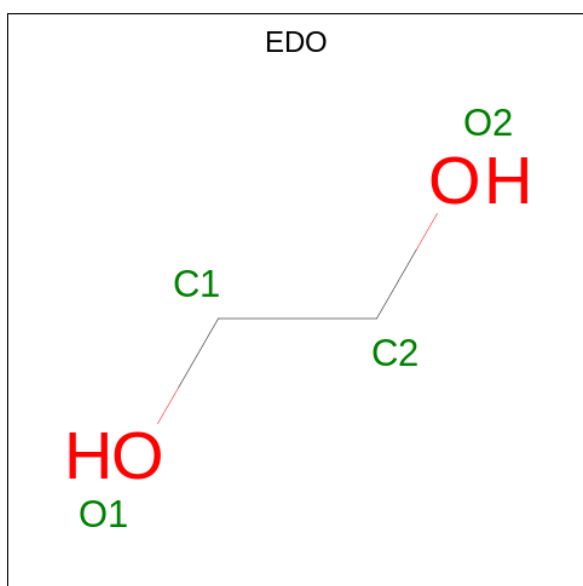
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Fe	0	0
			1	1		
3	A	1	Total	Fe	0	0
			1	1		
3	C	1	Total	Fe	0	0
			1	1		

- Molecule 4 is 9H-CARBAZOLE (three-letter code: 9CA) (formula: $\text{C}_{12}\text{H}_9\text{N}$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	N	0	0
			13	12	1		
4	C	1	Total	C	N	0	0
			13	12	1		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: 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		

Continued on next page...

Continued from previous page...

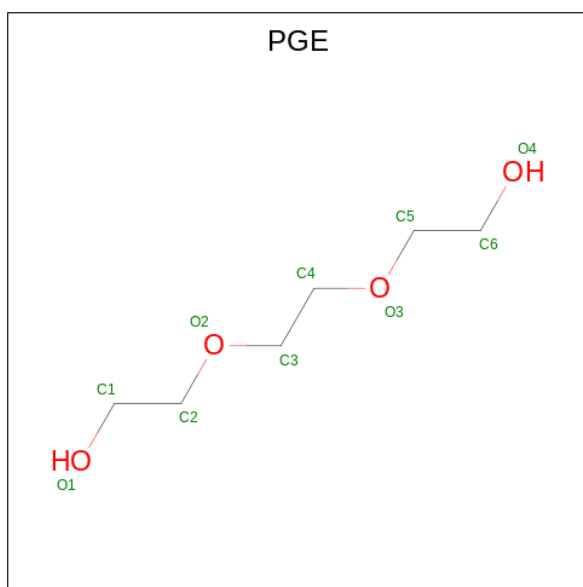
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0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

- Molecule 6 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



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

- Molecule 7 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



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

- Molecule 8 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	1	Total 1	Mg 1	0	0

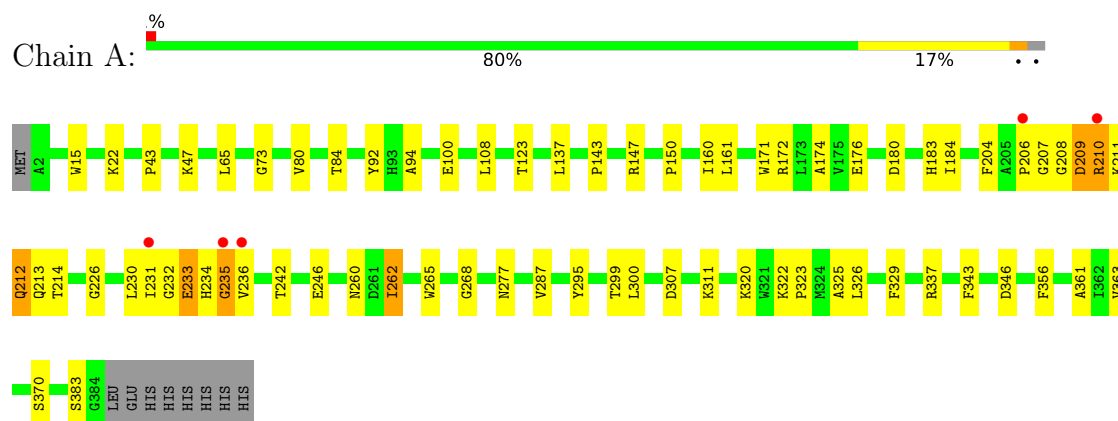
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	124	Total 124	O 124	0	0
9	B	138	Total 138	O 138	0	0
9	C	94	Total 94	O 94	0	0

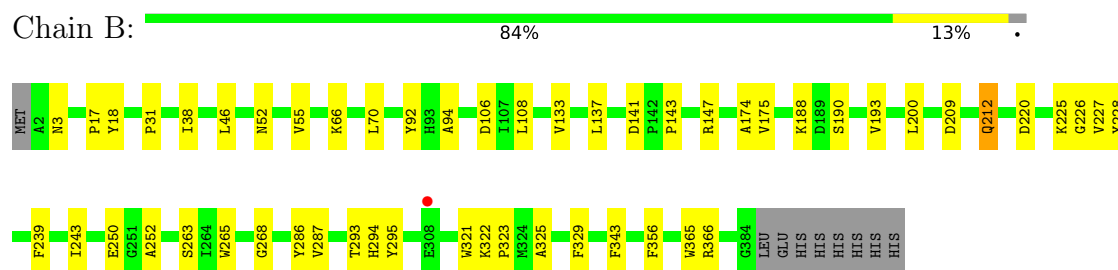
3 Residue-property plots

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.

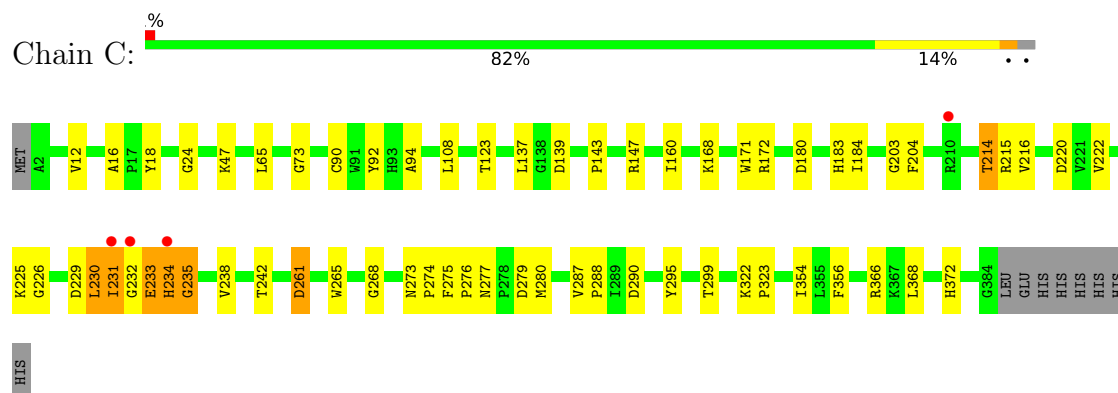
- Molecule 1: Terminal oxygenase component of carbazole



- Molecule 1: Terminal oxygenase component of carbazole



- Molecule 1: Terminal oxygenase component of carbazole



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	92.09Å 92.09Å 243.50Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.44 – 2.20 48.39 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.1 (48.44-2.20) 95.1 (48.39-2.20)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.176 , 0.246 0.182 , 0.248	Depositor DCC
R_{free} test set	2905 reflections (5.19%)	wwPDB-VP
Wilson B-factor (Å ²)	30.4	Xtriage
Anisotropy	0.051	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.075 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9950	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.10% 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: MG, PGE, EDO, FE2, FES, 9CA, PEG

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	0.66	0/3297	0.80	0/4474
1	B	0.66	0/3172	0.81	0/4306
1	C	0.67	0/3219	0.80	0/4370
All	All	0.66	0/9688	0.81	0/13150

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	3210	0	3116	70	0
1	B	3090	0	2997	46	0
1	C	3133	0	3037	48	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
2	C	4	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	13	0	9	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	13	0	9	0	0
5	A	28	0	42	7	0
5	B	44	0	66	21	0
5	C	16	0	24	3	0
6	A	14	0	20	5	0
6	B	7	0	10	0	0
7	A	10	0	14	1	0
8	B	1	0	0	0	0
9	A	124	0	0	2	0
9	B	138	0	0	5	0
9	C	94	0	0	4	0
All	All	9950	0	9344	164	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 9.

All (164) 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:210[A]:ARG:HG3	1:A:211[A]:LYS:HE3	1.30	1.12
1:A:232[A]:GLY:O	1:A:234[A]:HIS:N	1.88	1.04
1:B:188:LYS:HA	5:B:411:EDO:H12	1.40	1.01
1:A:210[A]:ARG:CG	1:A:211[A]:LYS:HE3	1.92	0.99
1:A:212[A]:GLN:NE2	1:A:212[A]:GLN:O	1.99	0.94
1:A:210[A]:ARG:O	1:A:213[A]:GLN:N	2.01	0.94
1:A:212[B]:GLN:OE1	1:A:234[B]:HIS:NE2	2.02	0.92
1:B:52:ASN:HD21	5:B:405:EDO:H21	1.39	0.87
1:B:52:ASN:HD21	5:B:405:EDO:C2	1.91	0.83
1:A:210[A]:ARG:O	1:A:212[A]:GLN:N	2.12	0.82
1:B:66:LYS:NZ	9:B:501:HOH:O	2.12	0.81
1:B:252:ALA:HB2	5:B:411:EDO:H21	1.61	0.81
1:A:207[B]:GLY:O	1:A:209[B]:ASP:N	2.13	0.80
1:A:212[B]:GLN:O	1:A:234[B]:HIS:CD2	2.35	0.79
1:A:210[A]:ARG:HG3	1:A:211[A]:LYS:CE	2.11	0.79
1:C:232[B]:GLY:O	1:C:233[B]:GLU:HG3	1.83	0.78
1:B:188:LYS:CA	5:B:411:EDO:H12	2.15	0.75
1:C:230[B]:LEU:O	1:C:231[B]:ILE:HG22	1.86	0.74
1:A:206[A]:PRO:HB3	1:A:231[A]:ILE:HG13	1.68	0.74
1:C:214:THR:CG2	9:C:503:HOH:O	2.36	0.73
1:C:234[B]:HIS:O	1:C:235:GLY:O	2.08	0.71
1:A:232[A]:GLY:C	1:A:234[A]:HIS:H	1.95	0.69

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:65:LEU:HD23	1:C:123:THR:HG22	1.74	0.69
1:C:230[B]:LEU:O	1:C:231[B]:ILE:CB	2.45	0.65
1:A:212[B]:GLN:O	1:A:234[B]:HIS:HD2	1.80	0.64
1:A:206[A]:PRO:HA	1:A:231[A]:ILE:HG12	1.80	0.64
5:A:410:EDO:H22	6:A:411:PEG:H11	1.80	0.64
1:C:230[B]:LEU:O	1:C:231[B]:ILE:CG2	2.46	0.63
1:A:370:SER:HB3	7:A:413:PGE:H2	1.79	0.63
1:C:230[B]:LEU:O	1:C:231[B]:ILE:HB	1.98	0.62
1:A:210[A]:ARG:C	1:A:212[A]:GLN:H	2.02	0.62
1:B:38:ILE:HD13	1:B:55:VAL:HG12	1.80	0.62
1:B:188:LYS:HA	5:B:411:EDO:C1	2.24	0.61
1:C:277:ASN:ND2	1:C:279[A]:ASP:OD1	2.33	0.61
1:C:229[B]:ASP:O	1:C:230[B]:LEU:O	2.19	0.61
1:A:307:ASP:O	1:A:311:LYS:HG2	2.01	0.60
1:B:52:ASN:ND2	5:B:413:EDO:H22	2.17	0.60
1:A:232[A]:GLY:C	1:A:234[A]:HIS:N	2.48	0.60
1:A:211[B]:LYS:HD3	1:A:212[B]:GLN:HG2	1.83	0.60
1:C:233[B]:GLU:O	1:C:234[B]:HIS:HB2	2.01	0.59
1:B:209:ASP:H	1:B:212:GLN:HE21	1.51	0.59
1:C:204:PHE:CE2	1:C:231[B]:ILE:HG13	2.38	0.59
1:C:180:ASP:HB3	1:C:183:HIS:HB3	1.85	0.57
1:C:216:VAL:HG13	1:C:368:LEU:HD23	1.85	0.57
1:B:252:ALA:CB	5:B:411:EDO:H21	2.32	0.57
1:B:94:ALA:HB1	1:B:108:LEU:HB2	1.87	0.57
1:A:214[A]:THR:HG21	1:A:361:ALA:HA	1.87	0.56
1:A:210[A]:ARG:HG2	1:A:211[A]:LYS:HE3	1.84	0.56
1:A:161:LEU:HA	5:A:404:EDO:H12	1.88	0.55
1:B:209:ASP:O	1:B:212:GLN:HG3	2.05	0.55
1:C:171:TRP:CE2	1:C:172:ARG:HG3	2.42	0.54
1:B:190:SER:HB2	5:B:408:EDO:H21	1.90	0.54
1:C:290:ASP:HB2	5:C:406:EDO:C1	2.38	0.54
1:C:214:THR:HG22	9:C:503:HOH:O	2.06	0.53
1:C:290:ASP:HB2	5:C:406:EDO:H11	1.90	0.53
1:A:84:THR:HG21	5:A:409:EDO:H12	1.91	0.53
1:A:204:PHE:CD1	1:A:231[B]:ILE:HG23	2.43	0.53
1:B:209:ASP:OD2	1:B:212:GLN:NE2	2.41	0.53
1:B:188:LYS:CB	5:B:411:EDO:H12	2.38	0.53
1:A:322:LYS:HB3	1:A:323:PRO:HD3	1.89	0.53
1:B:239:PHE:O	1:B:250:GLU:HA	2.09	0.53
1:B:252:ALA:HB2	5:B:411:EDO:C2	2.35	0.52
1:A:94:ALA:HB1	1:A:108:LEU:HB2	1.91	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210[A]:ARG:C	1:A:212[A]:GLN:N	2.63	0.51
1:A:210[A]:ARG:HG2	1:A:211[A]:LYS:H	1.74	0.51
1:C:232[A]:GLY:O	1:C:234[A]:HIS:N	2.43	0.51
1:B:174:ALA:HB1	1:B:286:TYR:CD1	2.46	0.51
1:A:171:TRP:CE2	1:A:172:ARG:HG3	2.46	0.50
1:A:277:ASN:HD21	6:A:411:PEG:C3	2.24	0.50
1:C:139:ASP:OD2	9:C:501:HOH:O	2.19	0.50
1:A:211[B]:LYS:HD3	1:A:211[B]:LYS:C	2.32	0.50
1:B:252:ALA:N	5:B:411:EDO:H21	2.27	0.50
1:A:361:ALA:HB3	9:A:507:HOH:O	2.11	0.49
1:A:234[B]:HIS:ND1	1:A:234[B]:HIS:N	2.61	0.48
1:A:206[A]:PRO:CA	1:A:231[A]:ILE:HG12	2.43	0.48
1:A:242:THR:HA	1:A:246:GLU:O	2.13	0.48
1:B:52:ASN:HD21	5:B:405:EDO:H22	1.76	0.48
1:A:226:GLY:HA3	1:A:265:TRP:CE3	2.48	0.48
1:B:17:PRO:HD3	5:B:410:EDO:H22	1.96	0.48
1:C:24:GLY:HA3	1:C:288:PRO:HG2	1.94	0.48
1:C:143:PRO:HG3	1:C:147:ARG:NH1	2.29	0.48
1:B:188:LYS:O	5:B:411:EDO:H22	2.13	0.48
1:A:232[B]:GLY:HA3	1:A:260:ASN:ND2	2.29	0.47
1:A:277:ASN:HD21	6:A:411:PEG:H31	1.78	0.47
1:A:287:VAL:HB	1:A:295:TYR:HB2	1.96	0.47
1:A:232[A]:GLY:C	1:A:235[A]:GLY:H	2.18	0.47
1:B:108:LEU:HD12	5:B:406:EDO:H22	1.96	0.47
1:B:287:VAL:HB	1:B:295:TYR:HB2	1.96	0.47
1:A:325:ALA:O	1:A:329:PHE:HB3	2.15	0.47
1:B:322:LYS:HB3	1:B:323:PRO:HD3	1.97	0.47
1:B:143:PRO:HG3	1:B:147:ARG:CZ	2.46	0.46
1:A:73:GLY:HA3	1:B:343:PHE:CG	2.50	0.46
1:A:232[A]:GLY:HA2	1:A:235[A]:GLY:CA	2.45	0.46
1:A:232[B]:GLY:HA3	1:A:260:ASN:HD22	1.80	0.46
1:B:52:ASN:ND2	5:B:405:EDO:H21	2.18	0.46
1:A:65:LEU:HD23	1:A:123:THR:HG22	1.98	0.46
1:B:190:SER:OG	1:B:193:VAL:HG23	2.15	0.46
1:A:180:ASP:HB3	1:A:183:HIS:HB3	1.96	0.46
1:B:225:LYS:HE2	9:B:609:HOH:O	2.16	0.46
1:B:3:ASN:OD1	9:B:502:HOH:O	2.20	0.46
1:A:143:PRO:HG3	1:A:147:ARG:CZ	2.46	0.46
1:A:212[B]:GLN:C	1:A:234[B]:HIS:HD2	2.19	0.46
1:A:231[B]:ILE:HD12	1:A:262:ILE:HB	1.97	0.46
1:A:235[A]:GLY:C	1:A:236[A]:VAL:HG13	2.36	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:406:EDO:H22	1:C:242:THR:O	2.16	0.46
1:C:261:ASP:HA	9:C:525:HOH:O	2.15	0.46
1:B:227:VAL:O	1:B:263:SER:HA	2.16	0.45
1:C:215:ARG:HB2	1:C:230[A]:LEU:HD11	1.98	0.45
1:C:273:ASN:HA	1:C:274:PRO:HA	1.82	0.45
1:A:230[A]:LEU:HD22	1:A:233[A]:GLU:HG3	1.99	0.45
1:B:106:ASP:OD1	5:B:406:EDO:H11	2.16	0.45
1:B:287:VAL:O	1:B:294:HIS:HA	2.17	0.45
1:A:22:LYS:HG3	5:A:408:EDO:H22	1.99	0.45
1:C:94:ALA:HB1	1:C:108:LEU:HB2	1.99	0.45
1:C:12:VAL:HG23	1:C:16:ALA:HA	2.00	0.44
1:A:212[B]:GLN:OE1	1:A:234[B]:HIS:CE1	2.69	0.44
1:A:320:LYS:HD3	5:A:410:EDO:O1	2.18	0.44
1:A:15:TRP:CZ3	1:A:363:VAL:HG13	2.53	0.44
1:C:275:PHE:CG	1:C:276:PRO:HA	2.53	0.44
1:A:233[B]:GLU:HG2	1:A:233[B]:GLU:H	1.51	0.43
1:A:160:ILE:HG23	1:A:299:THR:HB	2.00	0.43
1:A:43:PRO:HG2	1:A:80:VAL:HG11	2.01	0.43
5:B:412:EDO:C2	9:B:541:HOH:O	2.66	0.43
1:C:230[B]:LEU:HD22	1:C:234[B]:HIS:CB	2.48	0.43
5:A:409:EDO:H12	1:B:243:ILE:HG22	2.01	0.43
1:B:193:VAL:HG21	1:B:200:LEU:HD22	2.01	0.43
1:A:277:ASN:ND2	6:A:411:PEG:H31	2.34	0.42
1:B:226:GLY:HA3	1:B:265:TRP:CE3	2.54	0.42
1:C:230[B]:LEU:CD1	1:C:231[B]:ILE:N	2.82	0.42
1:C:226:GLY:HA3	1:C:265:TRP:CE3	2.54	0.42
1:C:160:ILE:HG23	1:C:299:THR:HB	2.02	0.42
1:A:174:ALA:HB2	1:A:337:ARG:HD3	2.00	0.42
1:C:184:ILE:HD11	1:C:203:GLY:HA2	2.02	0.42
1:C:232[A]:GLY:C	1:C:234[A]:HIS:N	2.72	0.42
1:B:70:LEU:HB3	1:C:354:ILE:HD12	2.01	0.42
1:C:287:VAL:HB	1:C:295:TYR:HB2	2.00	0.42
5:A:408:EDO:H21	9:A:539:HOH:O	2.20	0.41
1:A:176:GLU:O	1:A:180:ASP:HB2	2.20	0.41
1:C:18:TYR:CE2	1:C:366:ARG:HG2	2.56	0.41
1:A:343:PHE:CG	1:C:73:GLY:HA3	2.55	0.41
1:B:228:TYR:CD1	1:B:263:SER:HB3	2.54	0.41
1:B:321:TRP:O	1:B:325:ALA:HB3	2.20	0.41
1:C:220:ASP:OD1	1:C:222:VAL:HG22	2.20	0.41
1:C:322:LYS:HB3	1:C:323:PRO:HD3	2.03	0.41
1:C:225:LYS:HG3	1:C:372:HIS:HB3	2.01	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:ALA:CB	1:A:108:LEU:HB2	2.50	0.41
1:A:150:PRO:HB3	1:A:265:TRP:CD1	2.55	0.41
1:A:346:ASP:OD1	1:A:346:ASP:N	2.53	0.41
1:B:108:LEU:O	1:C:238:VAL:HG11	2.21	0.41
1:C:47:LYS:NZ	5:C:404:EDO:H21	2.35	0.41
1:A:211[A]:LYS:HD3	1:A:211[A]:LYS:HA	1.96	0.41
5:B:412:EDO:H21	9:B:541:HOH:O	2.20	0.41
1:A:204:PHE:CE1	1:A:231[B]:ILE:HG23	2.55	0.41
1:A:47:LYS:NZ	6:A:412:PEG:H22	2.36	0.41
1:B:18:TYR:CE2	1:B:366:ARG:HG2	2.56	0.41
1:C:204:PHE:CE2	1:C:231[B]:ILE:CG1	3.03	0.41
1:C:230[B]:LEU:HD13	1:C:231[B]:ILE:N	2.35	0.41
1:A:204:PHE:CG	1:A:231[B]:ILE:HG23	2.55	0.41
1:A:300:LEU:HD21	1:A:326:LEU:HD21	2.03	0.40
1:B:175:VAL:HG11	1:B:365:TRP:CE2	2.56	0.40
1:B:325:ALA:O	1:B:329:PHE:HB3	2.22	0.40
1:B:31:PRO:HA	1:B:133:VAL:O	2.21	0.40
1:C:90:CYS:O	1:C:94:ALA:HA	2.21	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	399/392 (102%)	353 (88%)	38 (10%)	8 (2%)	7	4
1	B	382/392 (97%)	358 (94%)	23 (6%)	1 (0%)	41	46
1	C	388/392 (99%)	356 (92%)	22 (6%)	10 (3%)	5	3
All	All	1169/1176 (99%)	1067 (91%)	83 (7%)	19 (2%)	15	7

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	208[A]	GLY
1	A	208[B]	GLY
1	C	230[A]	LEU
1	C	230[B]	LEU
1	C	231[A]	ILE
1	C	231[B]	ILE
1	C	234[A]	HIS
1	C	234[B]	HIS
1	C	235	GLY
1	A	184	ILE
1	A	268	GLY
1	C	268	GLY
1	A	233[A]	GLU
1	A	233[B]	GLU
1	B	268	GLY
1	C	233[A]	GLU
1	C	233[B]	GLU
1	A	235[A]	GLY
1	A	235[B]	GLY

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	344/339 (102%)	332 (96%)	12 (4%)	36	46
1	B	331/339 (98%)	323 (98%)	8 (2%)	49	62
1	C	336/339 (99%)	330 (98%)	6 (2%)	59	72
All	All	1011/1017 (99%)	985 (97%)	26 (3%)	50	58

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	92	TYR
1	A	100	GLU
1	A	137	LEU
1	A	209[A]	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	209[B]	ASP
1	A	210[A]	ARG
1	A	210[B]	ARG
1	A	212[A]	GLN
1	A	212[B]	GLN
1	A	262	ILE
1	A	356	PHE
1	A	383	SER
1	B	46	LEU
1	B	92	TYR
1	B	137	LEU
1	B	141	ASP
1	B	212	GLN
1	B	220	ASP
1	B	293	THR
1	B	356	PHE
1	C	92	TYR
1	C	137	LEU
1	C	168	LYS
1	C	214	THR
1	C	261	ASP
1	C	356	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	260	ASN
1	B	52	ASN
1	B	212	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 35 ligands modelled in this entry, 4 are monoatomic - leaving 31 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	0.14	0	2,2,2	0.31	0
6	PEG	B	415	-	6,6,6	0.23	0	5,5,5	0.14	0
4	9CA	A	403	-	15,15,15	1.36	3 (20%)	21,21,21	1.69	5 (23%)
5	EDO	A	410	-	3,3,3	0.09	0	2,2,2	0.26	0
5	EDO	B	405	-	3,3,3	0.17	0	2,2,2	0.30	0
5	EDO	B	413	-	3,3,3	0.06	0	2,2,2	0.28	0
7	PGE	A	413	-	9,9,9	0.21	0	8,8,8	0.23	0
6	PEG	A	411	-	6,6,6	0.16	0	5,5,5	0.11	0
5	EDO	C	406	-	3,3,3	0.16	0	2,2,2	0.51	0
2	FES	C	401	1	0,4,4	0.00	-	-		
2	FES	B	401	1	0,4,4	0.00	-	-		
2	FES	A	401	1	0,4,4	0.00	-	-		
5	EDO	C	407	-	3,3,3	0.10	0	2,2,2	0.08	0
5	EDO	A	404	-	3,3,3	0.08	0	2,2,2	0.30	0
5	EDO	C	404	-	3,3,3	0.32	0	2,2,2	0.04	0
5	EDO	A	409	-	3,3,3	0.10	0	2,2,2	0.32	0
5	EDO	B	414	-	3,3,3	0.10	0	2,2,2	0.27	0
5	EDO	B	408	-	3,3,3	0.13	0	2,2,2	0.48	0
4	9CA	C	403	-	15,15,15	1.41	3 (20%)	21,21,21	1.70	6 (28%)
5	EDO	A	406	-	3,3,3	0.19	0	2,2,2	0.13	0
5	EDO	A	407	-	3,3,3	0.11	0	2,2,2	0.10	0
5	EDO	B	412	-	3,3,3	0.14	0	2,2,2	0.23	0
5	EDO	B	410	-	3,3,3	0.16	0	2,2,2	0.57	0
5	EDO	B	406	-	3,3,3	0.04	0	2,2,2	0.36	0
5	EDO	B	411	-	3,3,3	0.49	0	2,2,2	0.71	0
5	EDO	A	408	-	3,3,3	0.08	0	2,2,2	0.19	0
5	EDO	B	404	-	3,3,3	0.11	0	2,2,2	0.22	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	C	405	-	3,3,3	0.19	0	2,2,2	0.18	0
5	EDO	B	409	-	3,3,3	0.10	0	2,2,2	0.37	0
5	EDO	B	407	-	3,3,3	0.27	0	2,2,2	0.20	0
6	PEG	A	412	-	6,6,6	0.15	0	5,5,5	0.19	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	-	-	1/1/1/1	-
6	PEG	B	415	-	-	3/4/4/4	-
4	9CA	A	403	-	-	-	0/3/3/3
5	EDO	A	410	-	-	1/1/1/1	-
5	EDO	B	405	-	-	0/1/1/1	-
5	EDO	B	413	-	-	0/1/1/1	-
7	PGE	A	413	-	-	5/7/7/7	-
4	9CA	C	403	-	-	-	0/3/3/3
5	EDO	C	406	-	-	0/1/1/1	-
2	FES	C	401	1	-	-	0/1/1/1
2	FES	B	401	1	-	-	0/1/1/1
2	FES	A	401	1	-	-	0/1/1/1
5	EDO	C	407	-	-	0/1/1/1	-
5	EDO	A	404	-	-	1/1/1/1	-
5	EDO	C	404	-	-	1/1/1/1	-
5	EDO	A	409	-	-	1/1/1/1	-
5	EDO	B	414	-	-	1/1/1/1	-
5	EDO	B	408	-	-	1/1/1/1	-
6	PEG	A	411	-	-	0/4/4/4	-
5	EDO	A	406	-	-	0/1/1/1	-
5	EDO	A	407	-	-	1/1/1/1	-
5	EDO	B	412	-	-	1/1/1/1	-
5	EDO	B	410	-	-	0/1/1/1	-
5	EDO	B	406	-	-	1/1/1/1	-
5	EDO	B	411	-	-	0/1/1/1	-
5	EDO	A	408	-	-	1/1/1/1	-
5	EDO	B	404	-	-	1/1/1/1	-
5	EDO	C	405	-	-	0/1/1/1	-
5	EDO	B	409	-	-	1/1/1/1	-
5	EDO	B	407	-	-	0/1/1/1	-
6	PEG	A	412	-	-	1/4/4/4	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	403	9CA	C4B-C8A	2.63	1.49	1.42
4	C	403	9CA	C7-C8	2.44	1.42	1.36
4	A	403	9CA	C4A-C9A	2.41	1.49	1.42
4	A	403	9CA	C4B-C8A	2.36	1.49	1.42
4	C	403	9CA	C4A-C9A	2.15	1.48	1.42
4	A	403	9CA	C7-C8	2.02	1.41	1.36

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	403	9CA	C4-C4A-C9A	3.99	123.46	118.17
4	A	403	9CA	C5-C4B-C8A	3.85	123.27	118.17
4	A	403	9CA	C4-C4A-C9A	3.09	122.27	118.17
4	C	403	9CA	C5-C4B-C8A	3.01	122.16	118.17
4	A	403	9CA	C8A-N9-C9A	2.65	112.53	107.09
4	C	403	9CA	C8A-N9-C9A	2.52	112.26	107.09
4	C	403	9CA	C8-C8A-C4B	-2.48	116.23	120.76
4	A	403	9CA	C6-C5-C4B	-2.34	116.93	120.86
4	C	403	9CA	C1-C9A-C4A	-2.31	116.54	120.76
4	A	403	9CA	C1-C9A-C4A	-2.23	116.69	120.76
4	C	403	9CA	C3-C4-C4A	-2.14	117.26	120.86

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	415	PEG	O2-C3-C4-O4
7	A	413	PGE	O1-C1-C2-O2
7	A	413	PGE	C1-C2-O2-C3
7	A	413	PGE	O3-C5-C6-O4
5	A	407	EDO	O1-C1-C2-O2
5	B	412	EDO	O1-C1-C2-O2
6	A	412	PEG	O2-C3-C4-O4
5	A	404	EDO	O1-C1-C2-O2
5	B	414	EDO	O1-C1-C2-O2
5	B	406	EDO	O1-C1-C2-O2
5	A	409	EDO	O1-C1-C2-O2
5	B	404	EDO	O1-C1-C2-O2
6	B	415	PEG	C1-C2-O2-C3
5	A	405	EDO	O1-C1-C2-O2
5	C	404	EDO	O1-C1-C2-O2

Continued on next page...

Continued from previous page...

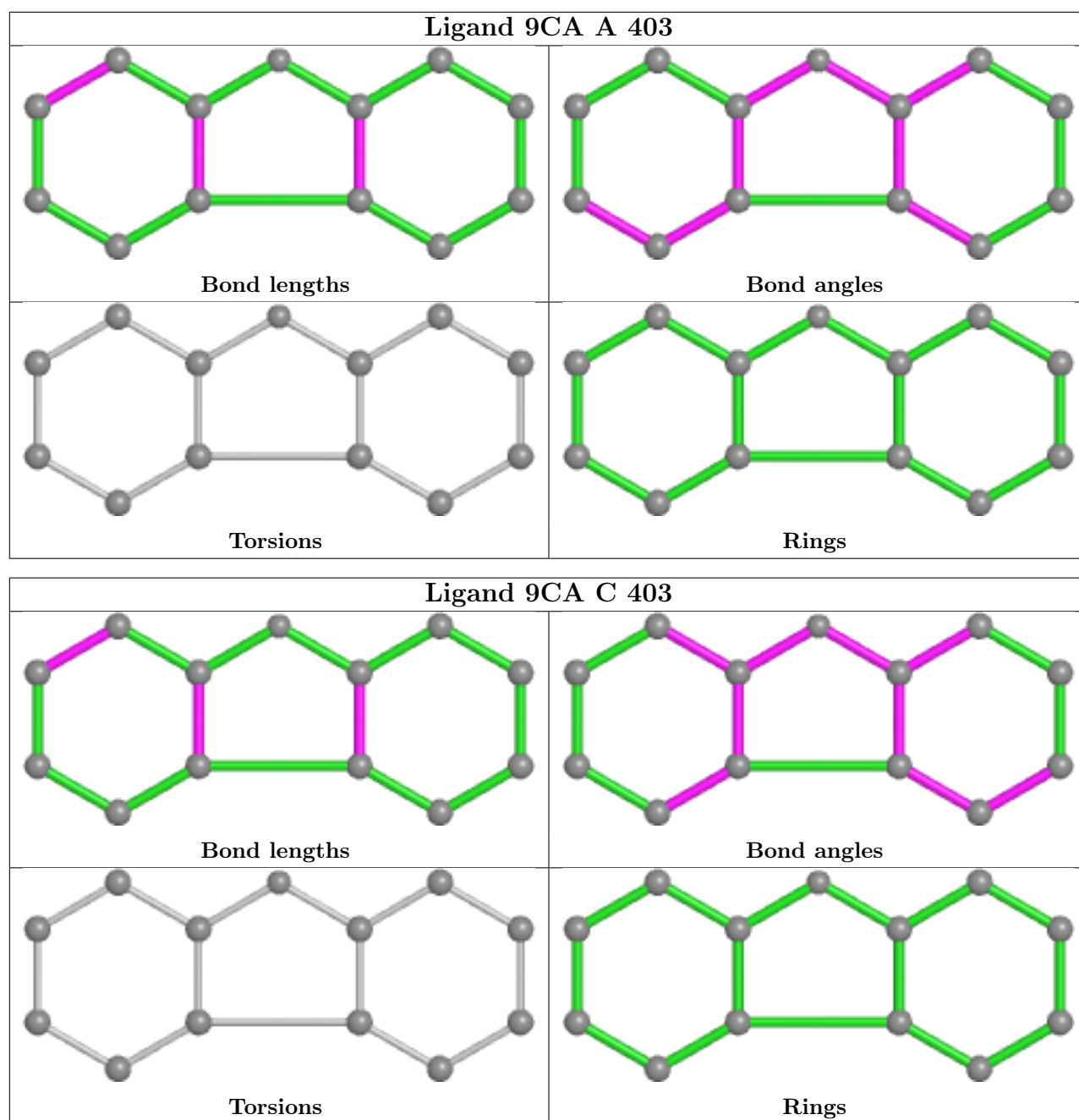
Mol	Chain	Res	Type	Atoms
5	B	408	EDO	O1-C1-C2-O2
5	B	409	EDO	O1-C1-C2-O2
7	A	413	PGE	O2-C3-C4-O3
6	B	415	PEG	C4-C3-O2-C2
7	A	413	PGE	C4-C3-O2-C2
5	A	410	EDO	O1-C1-C2-O2
5	A	408	EDO	O1-C1-C2-O2

There are no ring outliers.

16 monomers are involved in 36 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	410	EDO	2	0
5	B	405	EDO	4	0
5	B	413	EDO	1	0
7	A	413	PGE	1	0
6	A	411	PEG	4	0
5	C	406	EDO	2	0
5	A	404	EDO	1	0
5	C	404	EDO	1	0
5	A	409	EDO	2	0
5	B	408	EDO	1	0
5	B	412	EDO	2	0
5	B	410	EDO	1	0
5	B	406	EDO	3	0
5	B	411	EDO	9	0
5	A	408	EDO	2	0
6	A	412	PEG	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 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.



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	383/392 (97%)	-0.47	5 (1%) 77 75	18, 34, 61, 81	0
1	B	383/392 (97%)	-0.57	1 (0%) 94 93	17, 30, 53, 80	0
1	C	383/392 (97%)	-0.46	4 (1%) 82 81	21, 36, 61, 80	0
All	All	1149/1176 (97%)	-0.50	10 (0%) 84 83	17, 33, 60, 81	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	236[A]	VAL	4.1
1	A	206[A]	PRO	3.1
1	C	234[A]	HIS	3.1
1	A	231[A]	ILE	2.8
1	C	231[A]	ILE	2.6
1	A	235[A]	GLY	2.6
1	C	210	ARG	2.6
1	C	232[A]	GLY	2.6
1	B	308	GLU	2.1
1	A	210[A]	ARG	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

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.

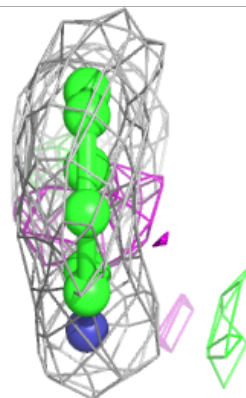
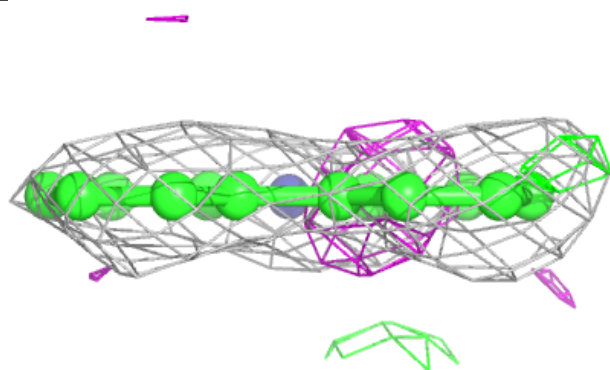
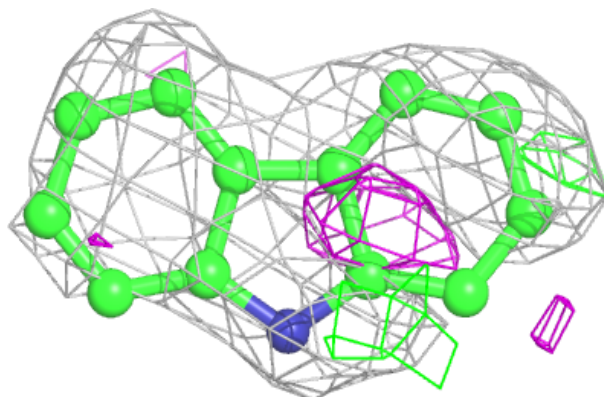
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	9CA	C	403	13/13	0.69	0.29	47,57,60,62	0
5	EDO	A	410	4/4	0.77	0.19	53,55,55,55	0
5	EDO	C	404	4/4	0.81	0.14	47,49,50,54	0
4	9CA	A	403	13/13	0.82	0.21	43,50,53,54	0
6	PEG	A	411	7/7	0.82	0.16	50,54,56,57	0
5	EDO	A	405	4/4	0.83	0.30	43,45,48,48	0
5	EDO	C	406	4/4	0.85	0.19	45,49,49,49	0
6	PEG	B	415	7/7	0.86	0.17	57,58,60,61	0
7	PGE	A	413	10/10	0.86	0.21	53,62,65,67	0
5	EDO	A	406	4/4	0.87	0.15	50,51,52,54	0
5	EDO	A	407	4/4	0.87	0.12	56,57,60,62	0
5	EDO	B	409	4/4	0.87	0.21	48,52,52,60	0
5	EDO	A	408	4/4	0.88	0.15	42,43,44,45	0
5	EDO	B	408	4/4	0.88	0.16	36,39,39,40	0
5	EDO	C	405	4/4	0.89	0.09	58,58,59,61	0
5	EDO	B	405	4/4	0.89	0.11	44,47,47,48	0
5	EDO	B	404	4/4	0.90	0.13	39,40,41,43	0
5	EDO	B	412	4/4	0.90	0.13	47,51,55,59	0
5	EDO	B	414	4/4	0.90	0.12	60,63,63,64	0
5	EDO	A	409	4/4	0.91	0.20	43,45,46,48	0
5	EDO	B	406	4/4	0.91	0.28	51,52,53,60	0
5	EDO	C	407	4/4	0.92	0.10	46,46,46,50	0
5	EDO	B	411	4/4	0.93	0.22	29,33,37,40	0
6	PEG	A	412	7/7	0.93	0.13	43,44,49,55	0
5	EDO	A	404	4/4	0.94	0.10	35,35,37,39	0
5	EDO	B	407	4/4	0.94	0.10	49,49,50,53	0
8	MG	B	403	1/1	0.94	0.06	34,34,34,34	0
5	EDO	B	413	4/4	0.95	0.19	45,47,47,50	0
5	EDO	B	410	4/4	0.95	0.24	33,40,42,44	0
3	FE2	C	402	1/1	0.98	0.07	42,42,42,42	0
2	FES	B	401	4/4	0.99	0.08	24,27,27,28	0
2	FES	A	401	4/4	0.99	0.10	19,20,21,23	0
2	FES	C	401	4/4	0.99	0.06	31,33,33,37	0
3	FE2	A	402	1/1	0.99	0.08	50,50,50,50	0
3	FE2	B	402	1/1	0.99	0.12	23,23,23,23	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.

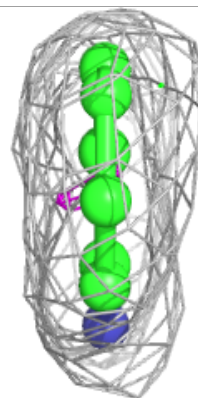
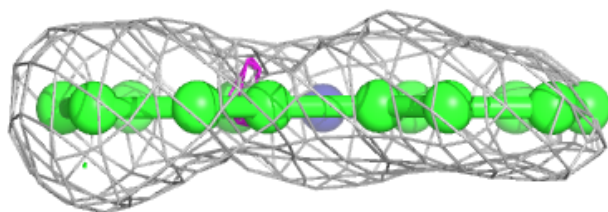
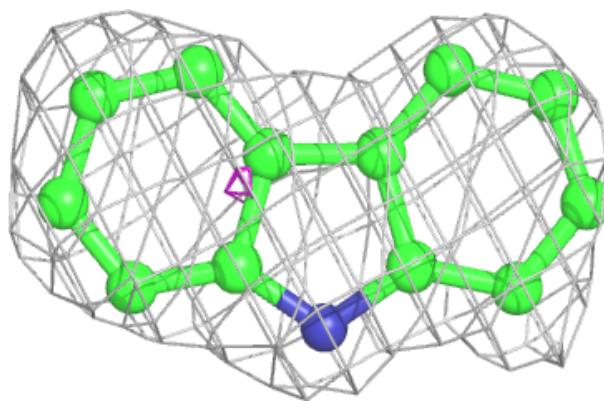
Electron density around 9CA C 403:

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



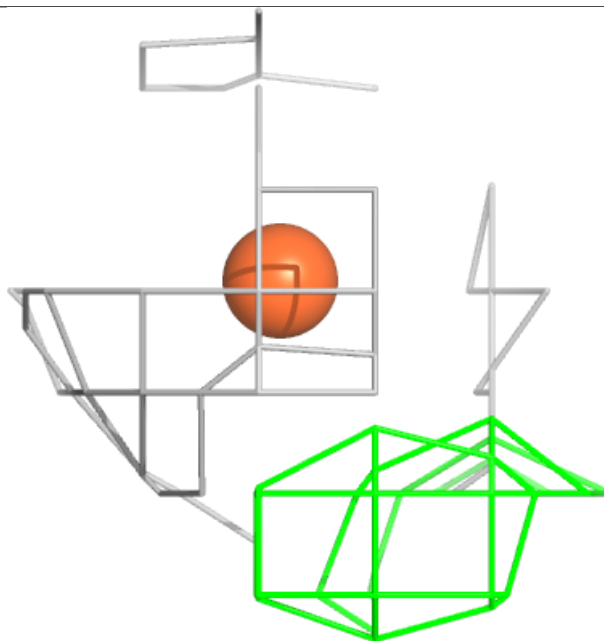
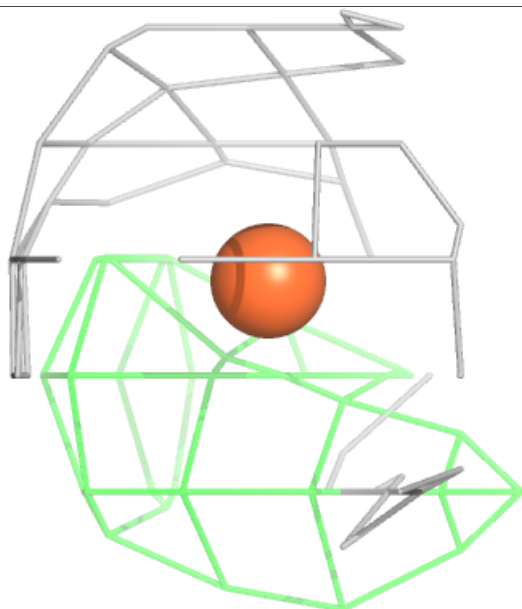
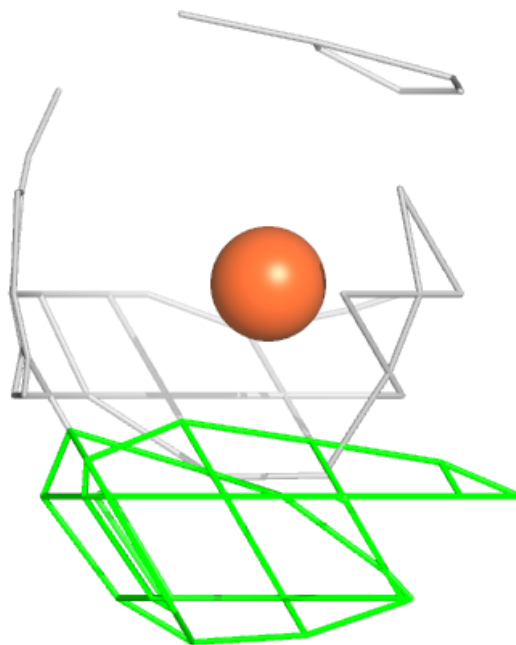
Electron density around 9CA A 403:

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



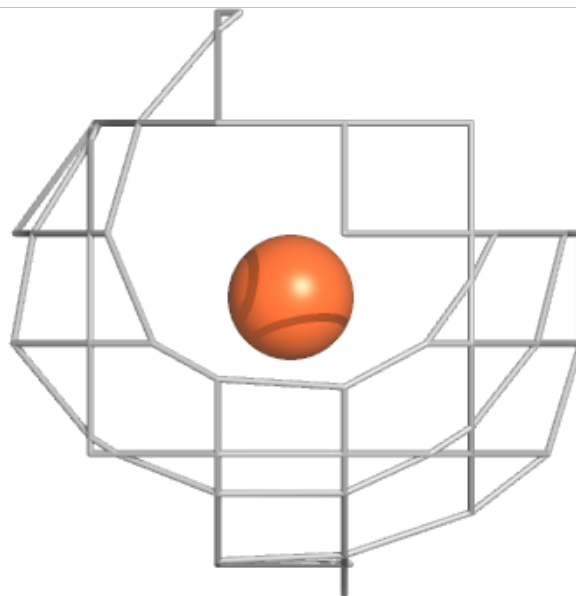
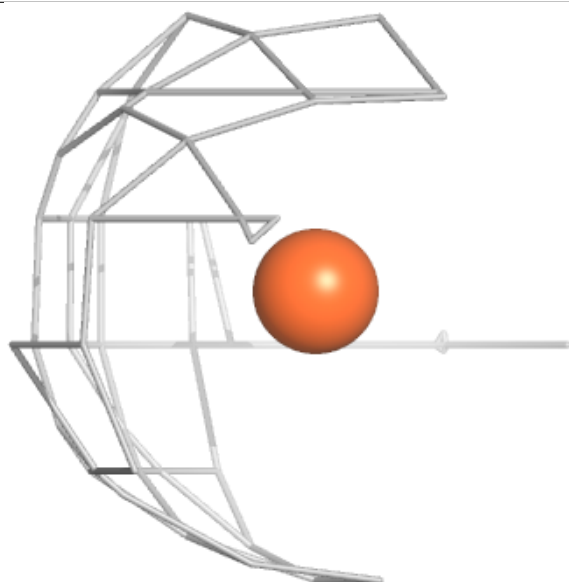
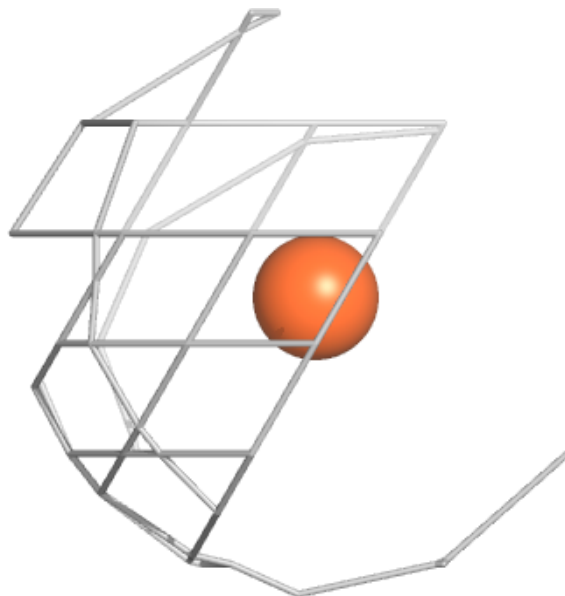
Electron density around FE2 C 402:

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



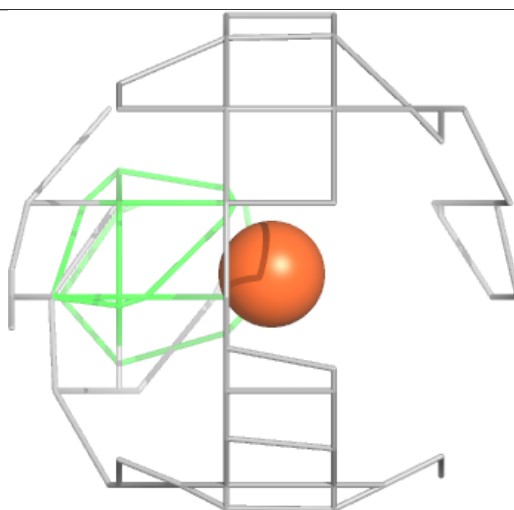
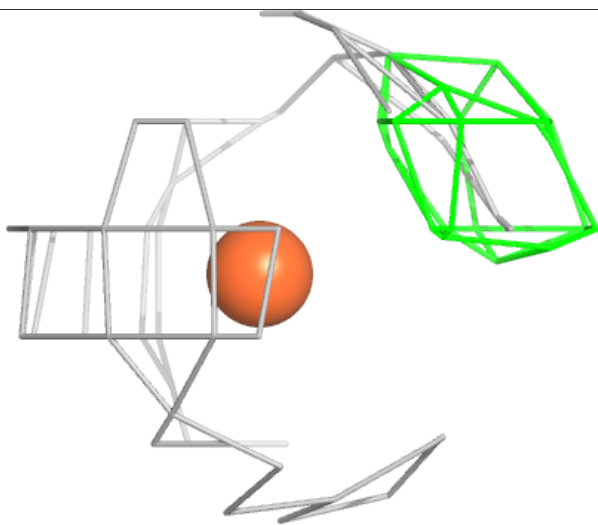
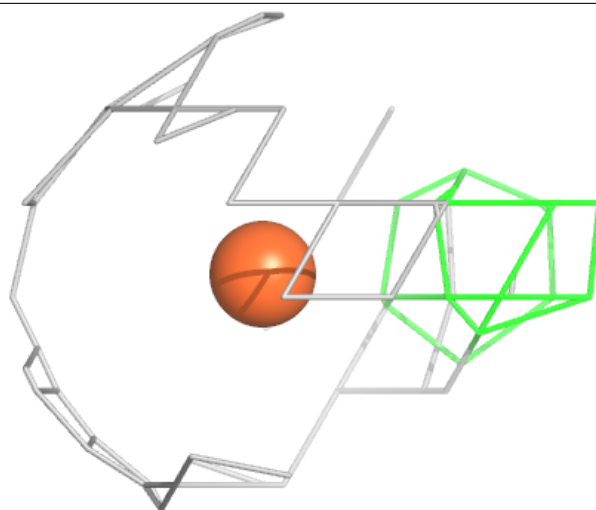
Electron density around FE2 A 402:

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



Electron density around FE2 B 402:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.