



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 08:55 AM GMT

PDB ID : 3GCF  
Title : Terminal oxygenase of carbazole 1,9a-dioxygenase from *Nocardioides aromaticivorans* IC177  
Authors : Inoue, K.; Nojiri, H.  
Deposited on : 2009-02-22  
Resolution : 2.30 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

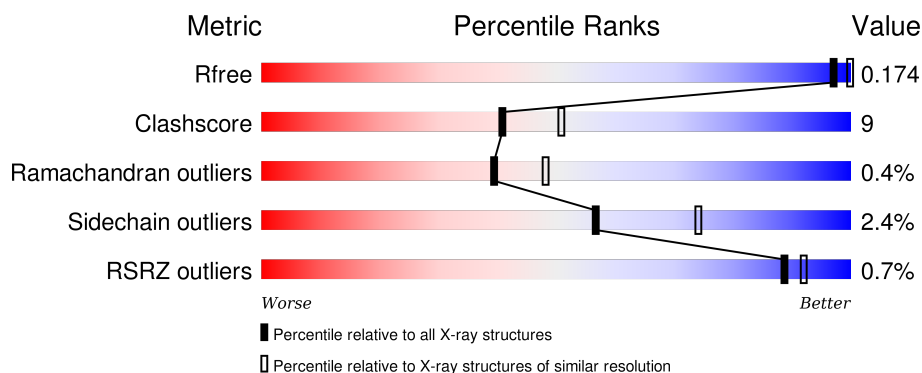
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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.







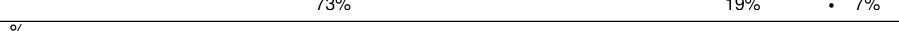



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	394	 76% 16% • 7%
1	B	394	 75% 18% • 7%
1	C	394	 72% 20% • 7%
1	D	394	 75% 19% 7%
1	E	394	 73% 20% • 7%

*Continued on next page...*

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	394	
1	G	394	
1	H	394	
1	I	394	
1	J	394	
1	K	394	
1	L	394	
1	M	394	
1	N	394	
1	O	394	

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
2	FE2	A	501	-	-	-	X
2	FE2	B	501	-	-	-	X
2	FE2	C	501	-	-	-	X
2	FE2	D	501	-	-	-	X
2	FE2	E	501	-	-	-	X
2	FE2	M	501	-	-	-	X
2	FE2	O	501	-	-	-	X
3	FES	B	401	-	-	-	X
3	FES	G	401	-	-	X	-
3	FES	K	401	-	-	X	-
4	CL	A	396	-	-	-	X
4	CL	B	396	-	-	-	X
4	CL	E	397	-	-	-	X
4	CL	H	396	-	-	-	X
4	CL	N	398	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 48063 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 1,9a-dioxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	B	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	C	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	D	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	E	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	F	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	G	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	H	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	I	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	J	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	K	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	L	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	M	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	N	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			
1	O	367	Total	C	N	O	S	0	0	0
			2949	1891	498	551	9			

There are 90 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0
A	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
A	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
A	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
A	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
A	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
B	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0
B	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
B	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
B	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
B	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
B	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
C	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0
C	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
C	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
C	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
C	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
C	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
D	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0
D	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
D	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
D	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
D	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
D	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
E	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0
E	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
E	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
E	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
E	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
E	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
F	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0
F	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
F	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
F	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
F	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
F	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
G	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0
G	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
G	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
G	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
G	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
G	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
H	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
H	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
H	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
H	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
H	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
H	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
I	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0
I	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
I	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
I	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
I	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
I	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
J	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0
J	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
J	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
J	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
J	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
J	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
K	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0
K	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
K	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
K	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
K	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
K	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
L	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0
L	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
L	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
L	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
L	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
L	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
M	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0
M	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
M	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
M	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
M	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
M	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
N	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0
N	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
N	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
N	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
N	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
N	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0
O	389	HIS	-	EXPRESSION TAG	UNP Q2HWI0

*Continued on next page...*

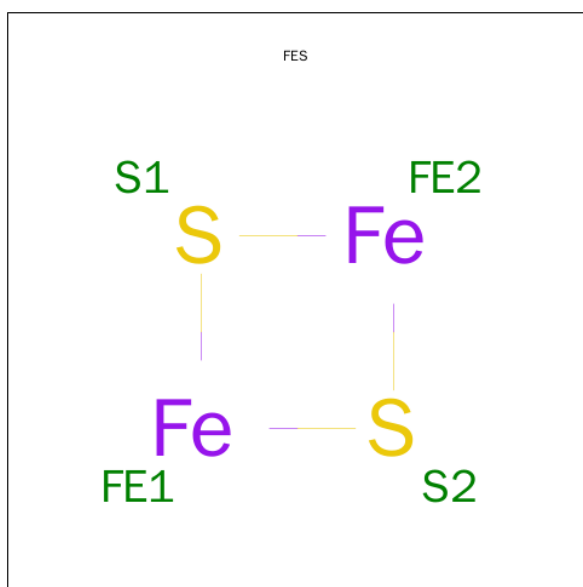
*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
O	390	HIS	-	EXPRESSION TAG	UNP Q2HWI0
O	391	HIS	-	EXPRESSION TAG	UNP Q2HWI0
O	392	HIS	-	EXPRESSION TAG	UNP Q2HWI0
O	393	HIS	-	EXPRESSION TAG	UNP Q2HWI0
O	394	HIS	-	EXPRESSION TAG	UNP Q2HWI0

- Molecule 2 is FE (II) ION (three-letter code: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	G	1	Total Fe 1 1	0	0
2	J	1	Total Fe 1 1	0	0
2	D	1	Total Fe 1 1	0	0
2	K	1	Total Fe 1 1	0	0
2	E	1	Total Fe 1 1	0	0
2	H	1	Total Fe 1 1	0	0
2	B	1	Total Fe 1 1	0	0
2	I	1	Total Fe 1 1	0	0
2	C	1	Total Fe 1 1	0	0
2	A	1	Total Fe 1 1	0	0
2	N	1	Total Fe 1 1	0	0
2	O	1	Total Fe 1 1	0	0
2	L	1	Total Fe 1 1	0	0
2	F	1	Total Fe 1 1	0	0
2	M	1	Total Fe 1 1	0	0

- Molecule 3 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe<sub>2</sub>S<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Fe	S	0	0
			4	2	2		
3	B	1	Total	Fe	S	0	0
			4	2	2		
3	C	1	Total	Fe	S	0	0
			4	2	2		
3	D	1	Total	Fe	S	0	0
			4	2	2		
3	E	1	Total	Fe	S	0	0
			4	2	2		
3	F	1	Total	Fe	S	0	0
			4	2	2		
3	G	1	Total	Fe	S	0	0
			4	2	2		
3	H	1	Total	Fe	S	0	0
			4	2	2		
3	I	1	Total	Fe	S	0	0
			4	2	2		
3	J	1	Total	Fe	S	0	0
			4	2	2		
3	K	1	Total	Fe	S	0	0
			4	2	2		
3	L	1	Total	Fe	S	0	0
			4	2	2		
3	M	1	Total	Fe	S	0	0
			4	2	2		
3	N	1	Total	Fe	S	0	0
			4	2	2		

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	O	1	Total 4	Fe 2	S 2	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	1	Total 1	Cl 1	0	0
4	J	2	Total 2	Cl 2	0	0
4	D	2	Total 2	Cl 2	0	0
4	K	2	Total 2	Cl 2	0	0
4	E	3	Total 3	Cl 3	0	0
4	H	2	Total 2	Cl 2	0	0
4	B	3	Total 3	Cl 3	0	0
4	I	1	Total 1	Cl 1	0	0
4	C	2	Total 2	Cl 2	0	0
4	A	3	Total 3	Cl 3	0	0
4	N	4	Total 4	Cl 4	0	0
4	O	2	Total 2	Cl 2	0	0
4	L	2	Total 2	Cl 2	0	0
4	F	2	Total 2	Cl 2	0	0
4	M	3	Total 3	Cl 3	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	276	Total 276	O 276	0	0

*Continued on next page...*

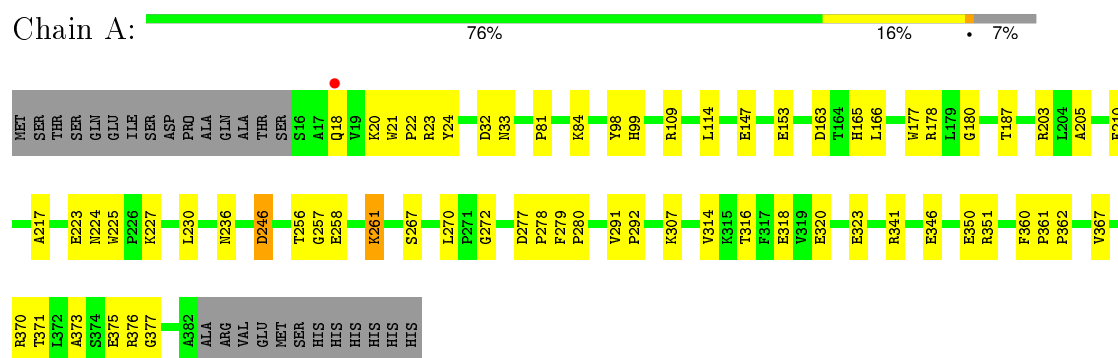
*Continued from previous page...*

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	273	Total 273	O 273	0	0
5	C	272	Total 272	O 272	0	0
5	D	257	Total 257	O 257	0	0
5	E	250	Total 250	O 250	0	0
5	F	252	Total 252	O 252	0	0
5	G	233	Total 233	O 233	0	0
5	H	215	Total 215	O 215	0	0
5	I	222	Total 222	O 222	0	0
5	J	218	Total 218	O 218	0	0
5	K	223	Total 223	O 223	0	0
5	L	231	Total 231	O 231	0	0
5	M	272	Total 272	O 272	0	0
5	N	260	Total 260	O 260	0	0
5	O	265	Total 265	O 265	0	0

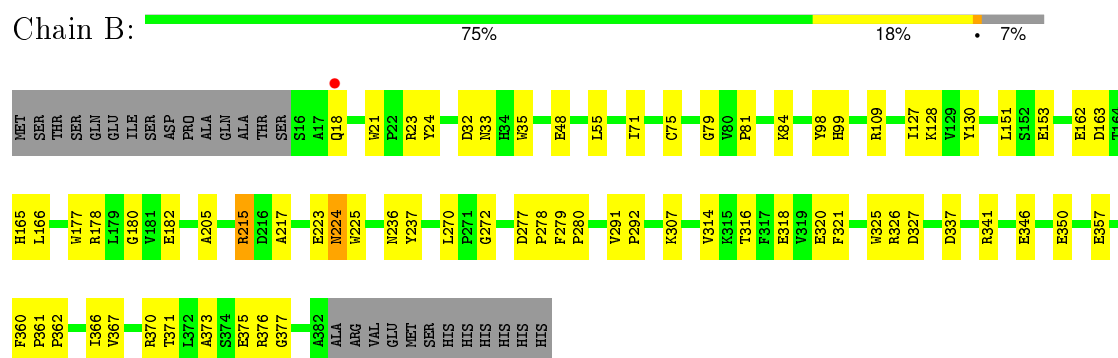
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

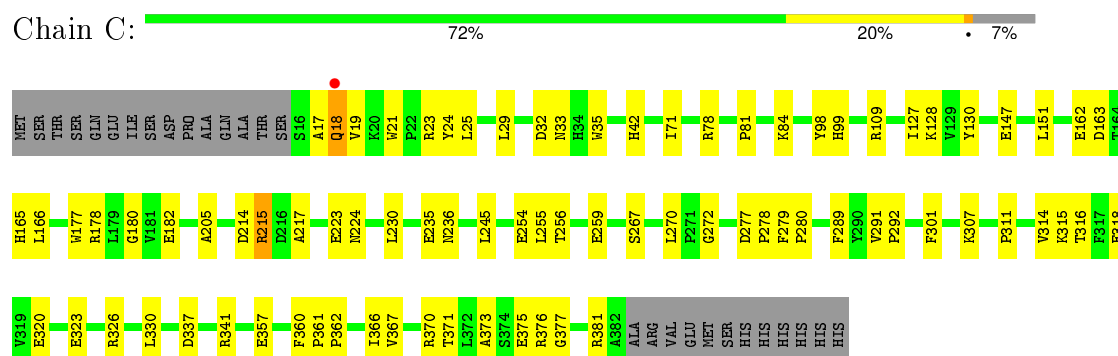
- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase



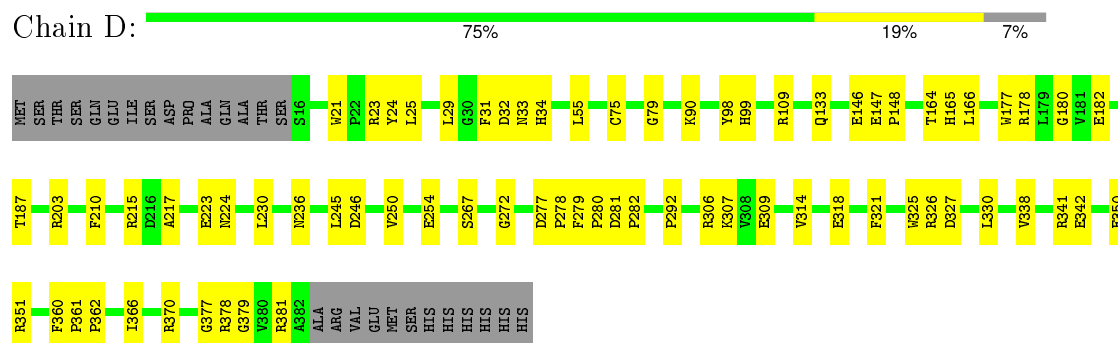
- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase



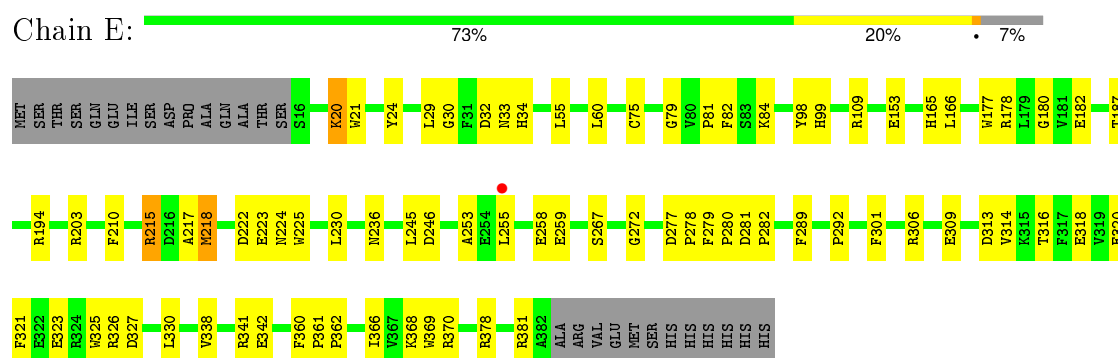
- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase



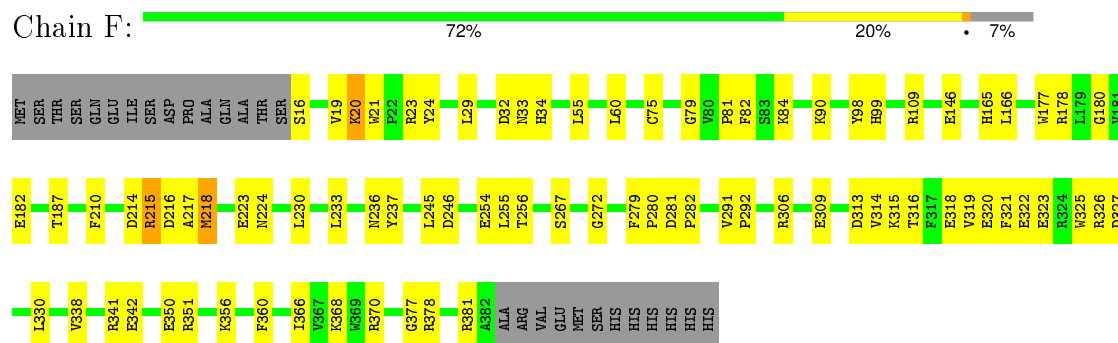
- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase



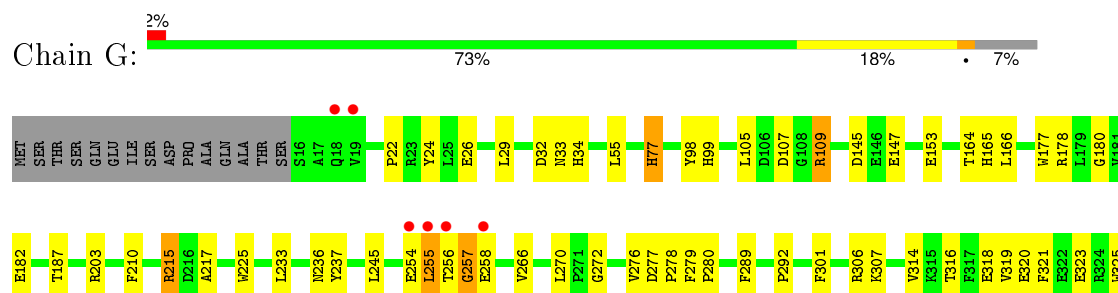
- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase



- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase

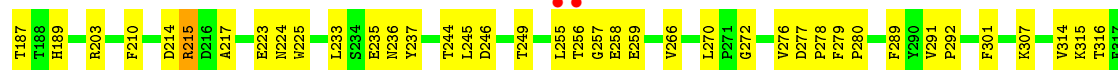
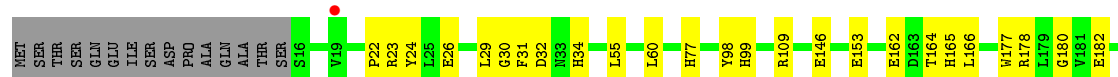


- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase





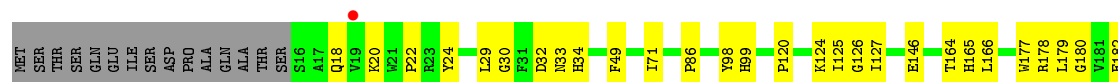
- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase



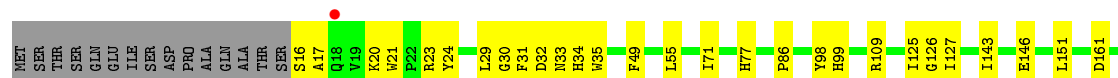
- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase



- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase

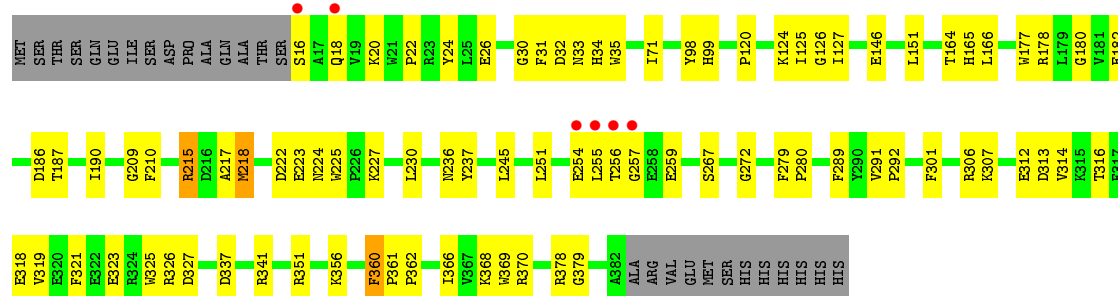


- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase

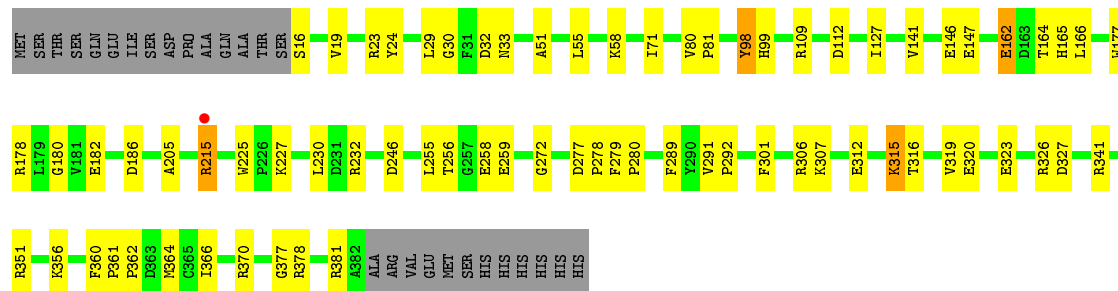




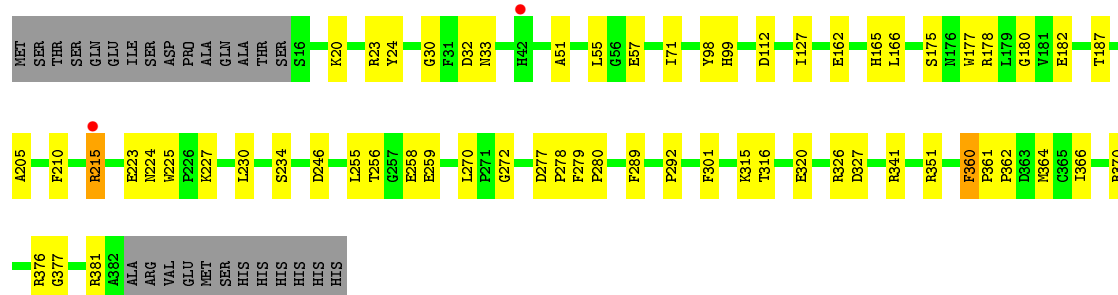
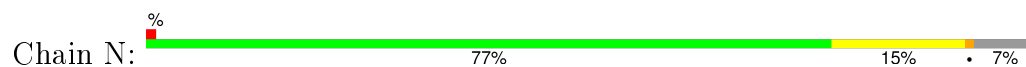
- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase



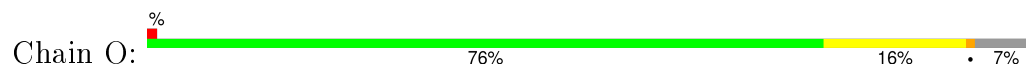
- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase

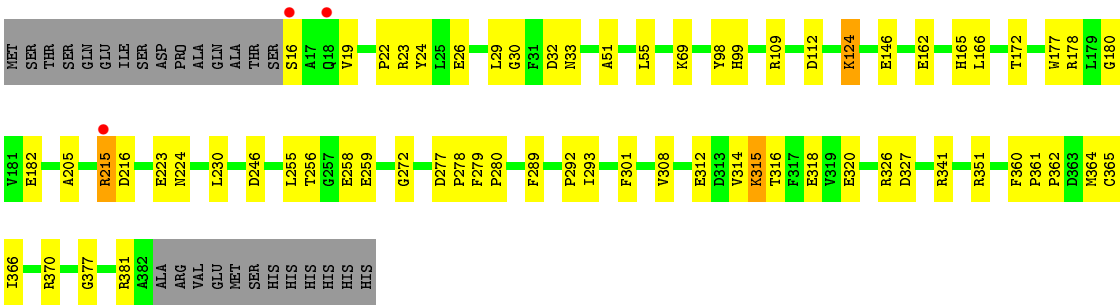


- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase



- Molecule 1: Terminal oxygenase component of carbazole 1,9a-dioxygenase





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	280.00Å 161.66Å 194.66Å 90.00° 118.65° 90.00°	Depositor
Resolution (Å)	48.91 – 2.30 48.91 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.91-2.30) 99.9 (48.91-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.83 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.174 , 0.175 0.173 , 0.174	Depositor DCC
$R_{free}$ test set	16984 reflections (5.32%)	DCC
Wilson B-factor (Å <sup>2</sup> )	20.9	Xtriage
Anisotropy	0.461	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 24.2	EDS
Estimated twinning fraction	0.487 for $1/2^*h+3/2^*k, 1/2^*h-1/2^*k, -1/2^*h-1/2^*k-l$ 0.487 for $1/2^*h-3/2^*k, -1/2^*h-1/2^*k, -1/2^*h+1/2^*k-l$	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	4 of 336543 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	48063	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: FE2, FES, CL

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	0.34	0/3038	0.61	2/4137 (0.0%)
1	B	0.34	0/3038	0.60	1/4137 (0.0%)
1	C	0.34	0/3038	0.60	1/4137 (0.0%)
1	D	0.33	0/3038	0.61	0/4137
1	E	0.33	0/3038	0.61	0/4137
1	F	0.33	0/3038	0.60	0/4137
1	G	0.33	0/3038	0.58	0/4137
1	H	0.33	0/3038	0.57	0/4137
1	I	0.33	0/3038	0.58	0/4137
1	J	0.33	0/3038	0.58	0/4137
1	K	0.33	0/3038	0.58	0/4137
1	L	0.33	0/3038	0.58	0/4137
1	M	0.34	0/3038	0.61	1/4137 (0.0%)
1	N	0.34	0/3038	0.61	1/4137 (0.0%)
1	O	0.34	0/3038	0.61	1/4137 (0.0%)
All	All	0.33	0/45570	0.60	7/62055 (0.0%)

There are no bond length outliers.

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	257	GLY	N-CA-C	6.70	129.84	113.10
1	A	205	ALA	N-CA-C	-5.85	95.20	111.00
1	C	205	ALA	N-CA-C	-5.81	95.31	111.00
1	B	205	ALA	N-CA-C	-5.71	95.59	111.00
1	M	205	ALA	N-CA-C	-5.16	97.08	111.00

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	A	2949	0	2805	39	0
1	B	2949	0	2805	44	0
1	C	2949	0	2805	52	0
1	D	2949	0	2805	47	0
1	E	2949	0	2805	53	0
1	F	2949	0	2805	57	0
1	G	2949	0	2805	48	0
1	H	2949	0	2805	55	0
1	I	2949	0	2805	51	0
1	J	2949	0	2805	56	0
1	K	2949	0	2805	68	0
1	L	2949	0	2805	61	0
1	M	2949	0	2805	51	0
1	N	2949	0	2805	38	0
1	O	2949	0	2805	41	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
2	M	1	0	0	0	0
2	N	1	0	0	0	0
2	O	1	0	0	0	0
3	A	4	0	0	1	0
3	B	4	0	0	1	0
3	C	4	0	0	1	0
3	D	4	0	0	1	0
3	E	4	0	0	1	0
3	F	4	0	0	1	0
3	G	4	0	0	2	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	4	0	0	1	0
3	I	4	0	0	1	0
3	J	4	0	0	1	0
3	K	4	0	0	2	0
3	L	4	0	0	1	0
3	M	4	0	0	1	0
3	N	4	0	0	1	0
3	O	4	0	0	1	0
4	A	3	0	0	0	0
4	B	3	0	0	0	0
4	C	2	0	0	0	0
4	D	2	0	0	0	0
4	E	3	0	0	0	0
4	F	2	0	0	0	0
4	G	1	0	0	0	0
4	H	2	0	0	0	0
4	I	1	0	0	0	0
4	J	2	0	0	0	0
4	K	2	0	0	0	0
4	L	2	0	0	0	0
4	M	3	0	0	0	0
4	N	4	0	0	0	0
4	O	2	0	0	0	0
5	A	276	0	0	1	0
5	B	273	0	0	2	0
5	C	272	0	0	3	0
5	D	257	0	0	4	0
5	E	250	0	0	3	0
5	F	252	0	0	2	0
5	G	233	0	0	2	0
5	H	215	0	0	3	0
5	I	222	0	0	3	0
5	J	218	0	0	3	0
5	K	223	0	0	5	0
5	L	231	0	0	3	0
5	M	272	0	0	6	0
5	N	260	0	0	2	0
5	O	265	0	0	5	0
All	All	48063	0	42075	747	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.

The worst 5 of 747 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:K:215:ARG:HD3	1:K:216:ASP:H	1.21	1.00
1:D:164:THR:HG22	1:D:307:LYS:HD3	1.52	0.91
1:C:215:ARG:H	1:C:215:ARG:HE	1.16	0.88
1:L:215:ARG:HA	1:L:215:ARG:HE	1.39	0.87
1:F:254:GLU:HG3	1:F:255:LEU:H	1.41	0.86

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	365/394 (93%)	347 (95%)	17 (5%)	1 (0%)	46	57
1	B	365/394 (93%)	348 (95%)	15 (4%)	2 (0%)	34	41
1	C	365/394 (93%)	347 (95%)	16 (4%)	2 (0%)	34	41
1	D	365/394 (93%)	349 (96%)	14 (4%)	2 (0%)	34	41
1	E	365/394 (93%)	348 (95%)	16 (4%)	1 (0%)	46	57
1	F	365/394 (93%)	348 (95%)	16 (4%)	1 (0%)	46	57
1	G	365/394 (93%)	344 (94%)	16 (4%)	5 (1%)	14	13
1	H	365/394 (93%)	345 (94%)	17 (5%)	3 (1%)	24	27
1	I	365/394 (93%)	348 (95%)	16 (4%)	1 (0%)	46	57
1	J	365/394 (93%)	345 (94%)	19 (5%)	1 (0%)	46	57
1	K	365/394 (93%)	343 (94%)	21 (6%)	1 (0%)	46	57
1	L	365/394 (93%)	342 (94%)	22 (6%)	1 (0%)	46	57
1	M	365/394 (93%)	350 (96%)	14 (4%)	1 (0%)	46	57
1	N	365/394 (93%)	346 (95%)	18 (5%)	1 (0%)	46	57

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	O	365/394 (93%)	349 (96%)	15 (4%)	1 (0%)	46	57
All	All	5475/5910 (93%)	5199 (95%)	252 (5%)	24 (0%)	39	48

5 of 24 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	18	GLN
1	G	272	GLY
1	H	77	HIS
1	H	272	GLY
1	I	272	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	312/336 (93%)	306 (98%)	6 (2%)	65	81
1	B	312/336 (93%)	308 (99%)	4 (1%)	76	87
1	C	312/336 (93%)	307 (98%)	5 (2%)	70	84
1	D	312/336 (93%)	305 (98%)	7 (2%)	60	77
1	E	312/336 (93%)	304 (97%)	8 (3%)	54	71
1	F	312/336 (93%)	304 (97%)	8 (3%)	54	71
1	G	312/336 (93%)	304 (97%)	8 (3%)	54	71
1	H	312/336 (93%)	305 (98%)	7 (2%)	60	77
1	I	312/336 (93%)	305 (98%)	7 (2%)	60	77
1	J	312/336 (93%)	305 (98%)	7 (2%)	60	77
1	K	312/336 (93%)	304 (97%)	8 (3%)	54	71
1	L	312/336 (93%)	305 (98%)	7 (2%)	60	77
1	M	312/336 (93%)	302 (97%)	10 (3%)	46	62
1	N	312/336 (93%)	303 (97%)	9 (3%)	50	66
1	O	312/336 (93%)	301 (96%)	11 (4%)	43	58

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	4680/5040 (93%)	4568 (98%)	112 (2%)	57 74

5 of 112 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	245	LEU
1	J	218	MET
1	O	124	LYS
1	H	360	PHE
1	I	162	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 41 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	358	GLN
1	F	358	GLN
1	N	358	GLN
1	F	18	GLN
1	F	195	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 64 ligands modelled in this entry, 49 are monoatomic - leaving 15 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	$\# Z  > 2$
3	FES	A	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	B	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	C	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	D	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	E	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	F	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	G	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	H	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	I	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	J	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	K	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	L	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	M	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	N	401	1	0,4,4	0.00	-	0,4,4	0.00	-
3	FES	O	401	1	0,4,4	0.00	-	0,4,4	0.00	-

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
3	FES	A	401	1	-	0/0/4/4	0/1/1/1
3	FES	B	401	1	-	0/0/4/4	0/1/1/1
3	FES	C	401	1	-	0/0/4/4	0/1/1/1
3	FES	D	401	1	-	0/0/4/4	0/1/1/1
3	FES	E	401	1	-	0/0/4/4	0/1/1/1
3	FES	F	401	1	-	0/0/4/4	0/1/1/1
3	FES	G	401	1	-	0/0/4/4	0/1/1/1
3	FES	H	401	1	-	0/0/4/4	0/1/1/1
3	FES	I	401	1	-	0/0/4/4	0/1/1/1
3	FES	J	401	1	-	0/0/4/4	0/1/1/1
3	FES	K	401	1	-	0/0/4/4	0/1/1/1
3	FES	L	401	1	-	0/0/4/4	0/1/1/1
3	FES	M	401	1	-	0/0/4/4	0/1/1/1
3	FES	N	401	1	-	0/0/4/4	0/1/1/1
3	FES	O	401	1	-	0/0/4/4	0/1/1/1

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.

15 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	FES	1	0
3	B	401	FES	1	0
3	C	401	FES	1	0
3	D	401	FES	1	0
3	E	401	FES	1	0
3	F	401	FES	1	0
3	G	401	FES	2	0
3	H	401	FES	1	0
3	I	401	FES	1	0
3	J	401	FES	1	0
3	K	401	FES	2	0
3	L	401	FES	1	0
3	M	401	FES	1	0
3	N	401	FES	1	0
3	O	401	FES	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

### 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	A	367/394 (93%)	-0.57	1 (0%) 94 96	6, 14, 32, 65	0
1	B	367/394 (93%)	-0.58	1 (0%) 94 96	6, 14, 33, 65	0
1	C	367/394 (93%)	-0.57	1 (0%) 94 96	6, 14, 32, 64	0
1	D	367/394 (93%)	-0.56	0 100 100	8, 17, 39, 51	0
1	E	367/394 (93%)	-0.52	1 (0%) 94 96	9, 17, 39, 62	0
1	F	367/394 (93%)	-0.53	0 100 100	9, 17, 39, 55	0
1	G	367/394 (93%)	-0.51	6 (1%) 74 80	15, 22, 44, 72	0
1	H	367/394 (93%)	-0.48	3 (0%) 87 90	15, 22, 45, 70	0
1	I	367/394 (93%)	-0.50	4 (1%) 82 86	15, 22, 45, 69	0
1	J	367/394 (93%)	-0.51	4 (1%) 82 86	14, 22, 43, 64	0
1	K	367/394 (93%)	-0.46	5 (1%) 78 83	14, 23, 44, 65	0
1	L	367/394 (93%)	-0.47	6 (1%) 74 80	15, 23, 44, 63	0
1	M	367/394 (93%)	-0.50	1 (0%) 94 96	8, 17, 37, 65	0
1	N	367/394 (93%)	-0.49	2 (0%) 91 94	8, 16, 36, 65	0
1	O	367/394 (93%)	-0.50	3 (0%) 87 90	8, 17, 36, 65	0
All	All	5505/5910 (93%)	-0.52	38 (0%) 89 92	6, 19, 39, 72	0

The worst 5 of 38 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	254	GLU	6.8
1	J	254	GLU	5.1
1	E	255	LEU	4.5
1	H	256	THR	4.3
1	L	254	GLU	4.0

## 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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	CL	E	397	1/1	0.91	0.21	17.53	57,57,57,57	0
4	CL	N	398	1/1	0.89	0.23	11.63	56,56,56,56	0
4	CL	A	396	1/1	0.93	0.14	5.45	55,55,55,55	0
2	FE2	O	501	1/1	1.00	0.14	5.37	10,10,10,10	0
2	FE2	D	501	1/1	1.00	0.12	4.47	11,11,11,11	0
2	FE2	C	501	1/1	1.00	0.12	4.44	7,7,7,7	0
4	CL	B	396	1/1	0.94	0.14	3.62	48,48,48,48	0
2	FE2	E	501	1/1	1.00	0.12	3.21	10,10,10,10	0
2	FE2	M	501	1/1	0.99	0.12	2.96	11,11,11,11	0
2	FE2	B	501	1/1	1.00	0.12	2.75	8,8,8,8	0
4	CL	H	396	1/1	0.92	0.14	2.42	59,59,59,59	0
2	FE2	A	501	1/1	1.00	0.12	2.21	8,8,8,8	0
3	FES	B	401	4/4	1.00	0.11	2.06	10,10,10,11	0
2	FE2	L	501	1/1	1.00	0.11	1.82	14,14,14,14	0
4	CL	D	396	1/1	0.98	0.22	1.48	37,37,37,37	0
3	FES	A	401	4/4	0.99	0.12	1.47	9,10,10,12	0
4	CL	O	395	1/1	0.98	0.20	1.38	36,36,36,36	0
3	FES	E	401	4/4	1.00	0.11	1.38	11,12,12,12	0
4	CL	M	395	1/1	0.98	0.20	1.23	36,36,36,36	0
2	FE2	K	501	1/1	1.00	0.11	1.11	14,14,14,14	0
4	CL	E	396	1/1	0.97	0.22	1.06	36,36,36,36	0
4	CL	N	395	1/1	0.98	0.18	1.00	38,38,38,38	0
3	FES	D	401	4/4	1.00	0.11	0.91	11,12,12,13	0
3	FES	F	401	4/4	1.00	0.10	0.91	10,12,13,13	0
2	FE2	N	501	1/1	1.00	0.11	0.80	10,10,10,10	0

*Continued on next page...*

*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	FES	C	401	4/4	1.00	0.10	0.45	9,10,10,11	0
4	CL	F	396	1/1	0.97	0.18	0.36	39,39,39,39	0
3	FES	O	401	4/4	0.99	0.10	0.07	10,11,13,14	0
3	FES	N	401	4/4	0.99	0.10	0.01	10,10,11,14	0
2	FE2	J	501	1/1	1.00	0.09	-0.02	14,14,14,14	0
4	CL	K	395	1/1	0.98	0.15	-0.04	44,44,44,44	0
4	CL	J	396	1/1	0.87	0.16	-0.07	46,46,46,46	0
2	FE2	F	501	1/1	0.99	0.10	-0.10	10,10,10,10	0
3	FES	J	401	4/4	1.00	0.08	-0.48	17,17,18,19	0
2	FE2	I	501	1/1	0.99	0.09	-0.58	15,15,15,15	0
3	FES	K	401	4/4	1.00	0.08	-0.68	17,18,18,18	0
2	FE2	G	501	1/1	1.00	0.08	-0.71	16,16,16,16	0
4	CL	L	396	1/1	0.94	0.14	-0.79	46,46,46,46	0
2	FE2	H	501	1/1	1.00	0.08	-0.83	16,16,16,16	0
3	FES	M	401	4/4	1.00	0.09	-0.96	10,13,13,14	0
3	FES	L	401	4/4	1.00	0.08	-1.21	18,18,18,19	0
3	FES	I	401	4/4	1.00	0.07	-1.35	16,16,18,19	0
3	FES	H	401	4/4	1.00	0.07	-1.38	15,16,18,19	0
3	FES	G	401	4/4	0.99	0.06	-1.82	16,16,18,20	0
4	CL	N	396	1/1	0.97	0.20	-	34,34,34,34	0
4	CL	I	395	1/1	0.98	0.21	-	37,37,37,37	0
4	CL	L	395	1/1	0.97	0.18	-	35,35,35,35	0
4	CL	N	397	1/1	0.91	0.13	-	61,61,61,61	0
4	CL	H	395	1/1	0.99	0.20	-	39,39,39,39	0
4	CL	C	396	1/1	0.97	0.17	-	35,35,35,35	0
4	CL	B	397	1/1	0.96	0.16	-	34,34,34,34	0
4	CL	O	396	1/1	0.98	0.20	-	31,31,31,31	0
4	CL	C	395	1/1	0.93	0.18	-	47,47,47,47	0
4	CL	A	395	1/1	0.95	0.16	-	48,48,48,48	0
4	CL	D	395	1/1	0.97	0.24	-	33,33,33,33	0
4	CL	M	397	1/1	0.88	0.10	-	63,63,63,63	0
4	CL	A	397	1/1	0.97	0.14	-	33,33,33,33	0
4	CL	E	395	1/1	0.98	0.23	-	33,33,33,33	0
4	CL	M	396	1/1	0.99	0.17	-	32,32,32,32	0
4	CL	F	395	1/1	0.97	0.23	-	33,33,33,33	0
4	CL	B	395	1/1	0.96	0.14	-	48,48,48,48	0
4	CL	K	396	1/1	0.98	0.15	-	36,36,36,36	0
4	CL	G	395	1/1	0.98	0.19	-	38,38,38,38	0
4	CL	J	395	1/1	0.99	0.19	-	35,35,35,35	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.