



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 12:07 AM GMT

PDB ID : 1ZPU  
Title : Crystal Structure of Fet3p, a Multicopper Oxidase that Functions in Iron Import  
Authors : Taylor, A.B.; Stoj, C.S.; Ziegler, L.; Kosman, D.J.; Hart, P.J.  
Deposited on : 2005-05-17  
Resolution : 2.80 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.

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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 : 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

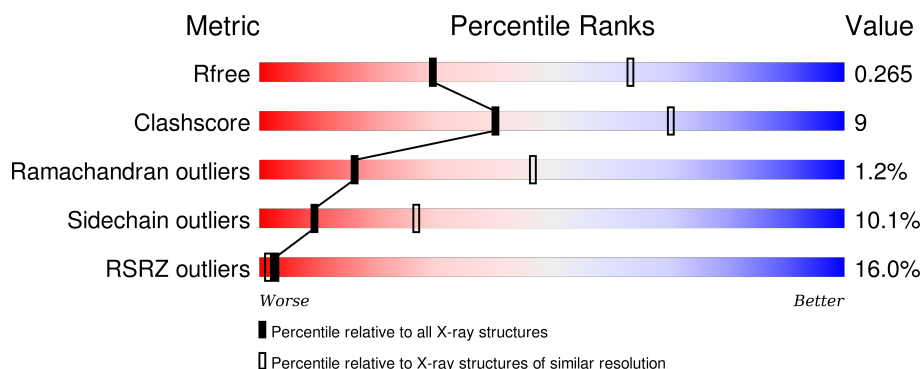
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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	534	<div> <div>8%</div> <div>74%</div> <div>21%</div> <div>• •</div> </div>
1	B	534	<div> <div>7%</div> <div>70%</div> <div>25%</div> <div>• •</div> </div>
1	C	534	<div> <div>12%</div> <div>72%</div> <div>23%</div> <div>• •</div> </div>
1	D	534	<div> <div>13%</div> <div>72%</div> <div>23%</div> <div>• •</div> </div>
1	E	534	<div> <div>23%</div> <div>77%</div> <div>20%</div> <div>• •</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	534	

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	NAG	D	2007	X	-	-	-
2	NAG	E	2007	X	-	-	X
2	NAG	F	2007	-	-	-	X
3	NAG	A	2003	X	-	-	-
3	NAG	B	2003	X	-	-	-
3	NAG	D	2003	X	-	-	-
3	NAG	F	2003	X	-	-	-
4	NAG	C	2008	-	-	-	X
4	NAG	D	2009	X	-	-	-
5	NAG	A	2006	X	-	-	-
5	NAG	A	2012	X	-	-	X
5	NAG	B	2006	X	-	-	-
5	NAG	B	2012	X	-	-	X
5	NAG	B	2014	-	-	-	X
5	NAG	B	2018	X	-	-	-
5	NAG	C	2012	X	-	-	X
5	NAG	C	2018	X	-	-	-
5	NAG	D	2006	X	-	-	-
5	NAG	D	2008	-	-	-	X
5	NAG	D	2012	-	-	-	X
5	NAG	D	2014	-	-	-	X
5	NAG	E	2005	-	-	-	X
5	NAG	E	2008	-	-	-	X
5	NAG	E	2012	X	-	-	-
5	NAG	F	2005	X	-	-	X
5	NAG	F	2006	X	-	-	-
5	NAG	F	2009	X	-	-	-
5	NAG	F	2012	X	-	-	-
6	NAG	A	2010	X	-	-	-
6	NAG	C	2010	X	-	-	-
6	NAG	D	2010	X	-	-	X
6	NAG	E	2010	X	-	-	-
6	NAG	F	2010	X	-	-	X
8	NAG	B	2010	X	-	-	-
9	NAG	C	2003	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
9	NAG	E	2003	X	-	-	-

## 2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 27659 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 Iron transport multicopper oxidase FET3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	529	Total	C	N	O	S	0	0	0
			4254	2701	695	838	20			
1	B	529	Total	C	N	O	S	0	0	0
			4254	2701	695	838	20			
1	C	529	Total	C	N	O	S	0	0	0
			4254	2701	695	838	20			
1	D	529	Total	C	N	O	S	0	0	0
			4254	2701	695	838	20			
1	E	529	Total	C	N	O	S	0	0	0
			4254	2701	695	838	20			
1	F	529	Total	C	N	O	S	0	0	0
			4254	2701	695	838	20			

- Molecule 2 is a polymer of unknown type called SUGAR (5-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	5	Total	C	N	O	0	0
			61	34	2	25		
2	A	5	Total	C	N	O	0	0
			61	34	2	25		
2	B	5	Total	C	N	O	0	0
			61	34	2	25		
2	B	5	Total	C	N	O	0	0
			61	34	2	25		
2	C	5	Total	C	N	O	0	0
			61	34	2	25		
2	C	5	Total	C	N	O	0	0
			61	34	2	25		
2	D	5	Total	C	N	O	0	0
			61	34	2	25		
2	D	5	Total	C	N	O	0	0
			61	34	2	25		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	E	5	Total	C	N	O	0	0
			61	34	2	25		
2	E	5	Total	C	N	O	0	0
			61	34	2	25		
2	F	5	Total	C	N	O	0	0
			61	34	2	25		
2	F	5	Total	C	N	O	0	0
			61	34	2	25		

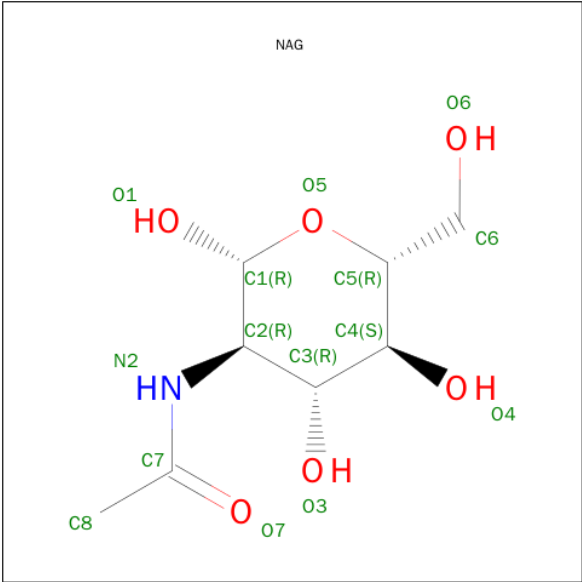
- Molecule 3 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	6	Total	C	N	O	0	0
			72	40	2	30		
3	B	6	Total	C	N	O	0	0
			72	40	2	30		
3	D	6	Total	C	N	O	0	0
			72	40	2	30		
3	F	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 4 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	2	Total	C	N	O	0	0
			28	16	2	10		
4	B	2	Total	C	N	O	0	0
			28	16	2	10		
4	B	2	Total	C	N	O	0	0
			28	16	2	10		
4	C	2	Total	C	N	O	0	0
			28	16	2	10		
4	D	2	Total	C	N	O	0	0
			28	16	2	10		
4	E	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	E	1	Total	C	N	O	0	0
			14	8	1	5		
5	E	1	Total	C	N	O	0	0
			14	8	1	5		
5	E	1	Total	C	N	O	0	0
			14	8	1	5		
5	E	1	Total	C	N	O	0	0
			14	8	1	5		
5	E	1	Total	C	N	O	0	0
			14	8	1	5		
5	F	1	Total	C	N	O	0	0
			14	8	1	5		
5	F	1	Total	C	N	O	0	0
			14	8	1	5		
5	F	1	Total	C	N	O	0	0
			14	8	1	5		
5	F	1	Total	C	N	O	0	0
			14	8	1	5		
5	F	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	3	Total	C	N	O	0	0
			39	22	2	15		
6	A	3	Total	C	N	O	0	0
			39	22	2	15		
6	C	3	Total	C	N	O	0	0
			39	22	2	15		

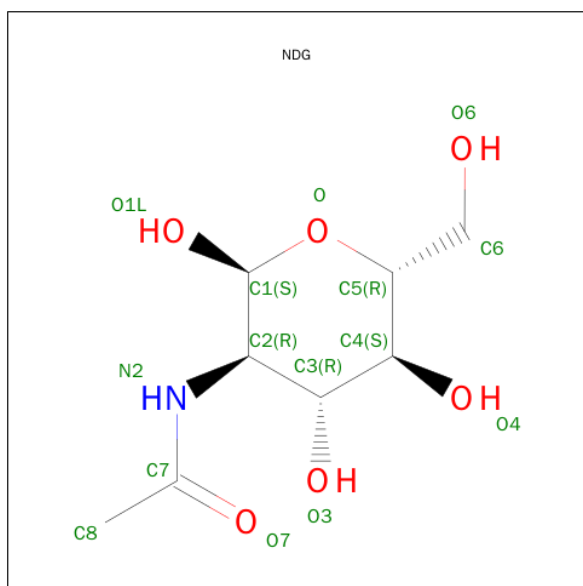
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	C	3	Total	C	N	O	0	0
			39	22	2	15		
6	D	3	Total	C	N	O	0	0
			39	22	2	15		
6	E	3	Total	C	N	O	0	0
			39	22	2	15		
6	F	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 7 is SUGAR (2-(ACETYLAMINO)-2-DEOXY-A-D-GLUCOPYRANOSE) (three-letter code: NDG) (formula:  $C_8H_{15}NO_6$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 8 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	B	4	Total	C	N	O	0	0
			50	28	2	20		

- Molecule 9 is a polymer of unknown type called SUGAR (7-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	C	7	Total	C	N	O	0	0
			83	46	2	35		
9	E	7	Total	C	N	O	0	0
			83	46	2	35		

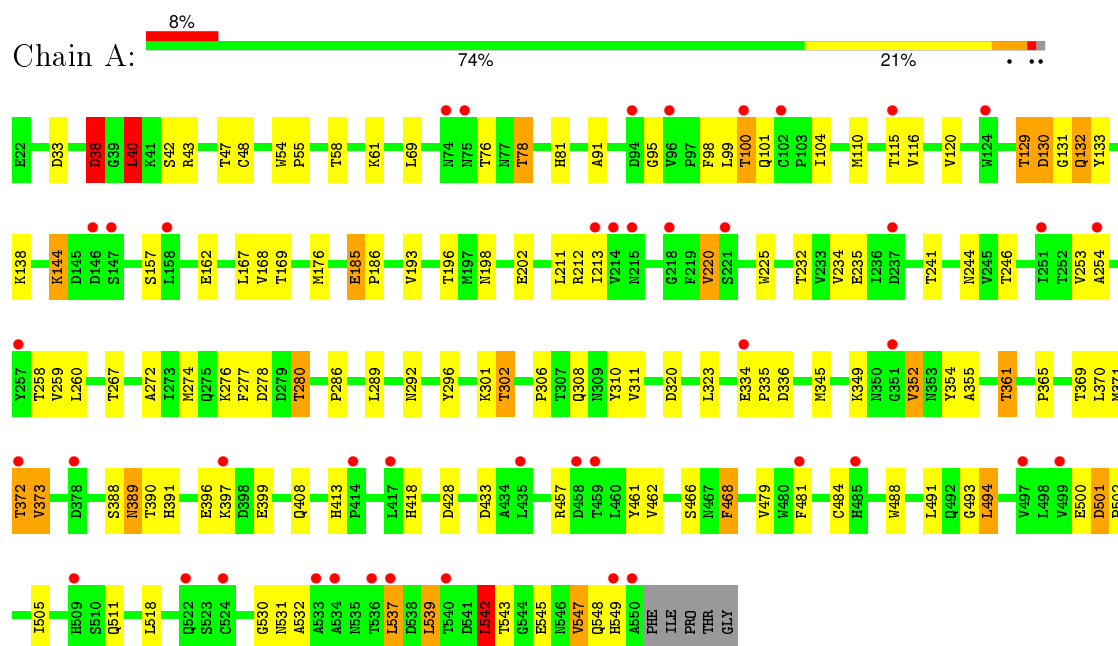
- Molecule 10 is COPPER (I) ION (three-letter code: CU1) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	D	4	Total	Cu	0	0
			4	4		
10	E	4	Total	Cu	0	0
			4	4		
10	B	4	Total	Cu	0	0
			4	4		
10	C	4	Total	Cu	0	0
			4	4		
10	A	4	Total	Cu	0	0
			4	4		
10	F	4	Total	Cu	0	0
			4	4		

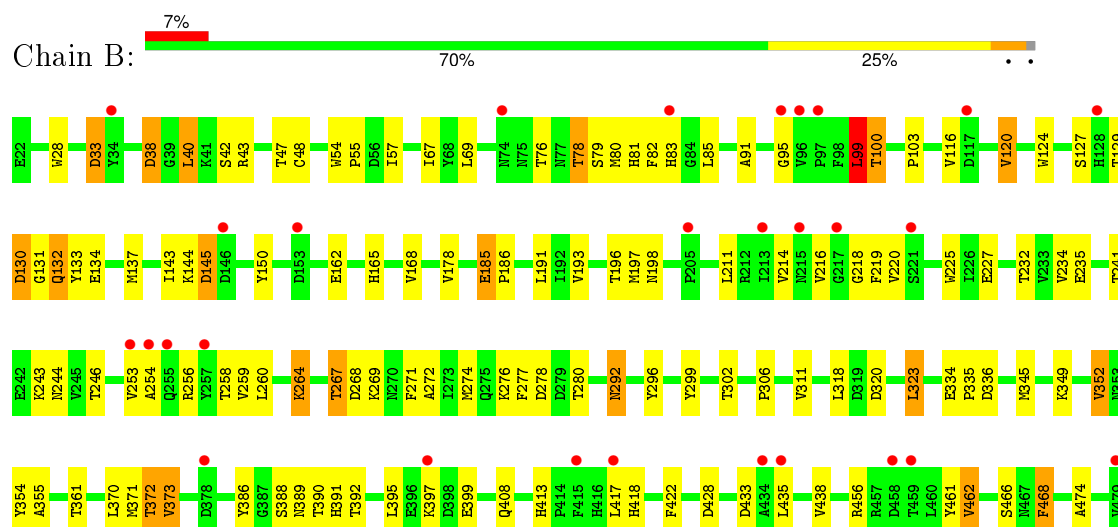
### 3 Residue-property plots

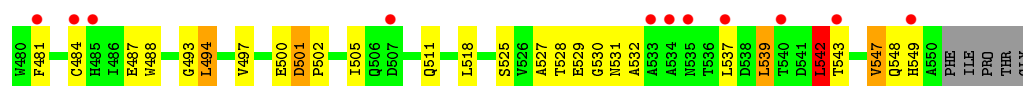
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ( $\text{RSRZ} > 2$ ). 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: Iron transport multicopper oxidase FET3

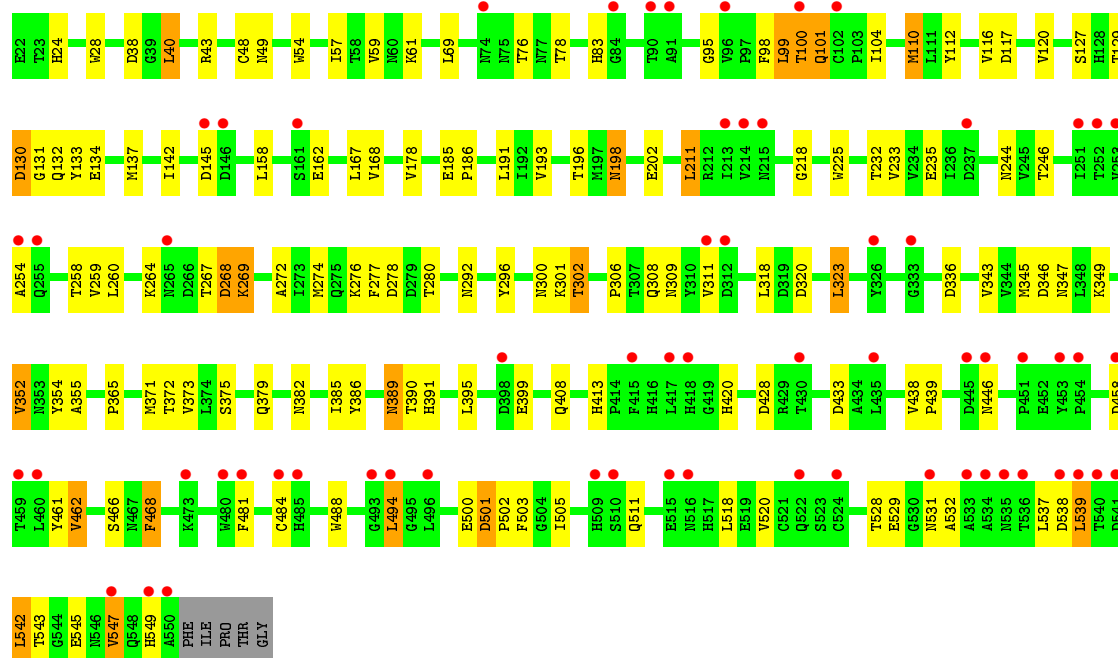


#### • Molecule 1: Iron transport multicopper oxidase FET3

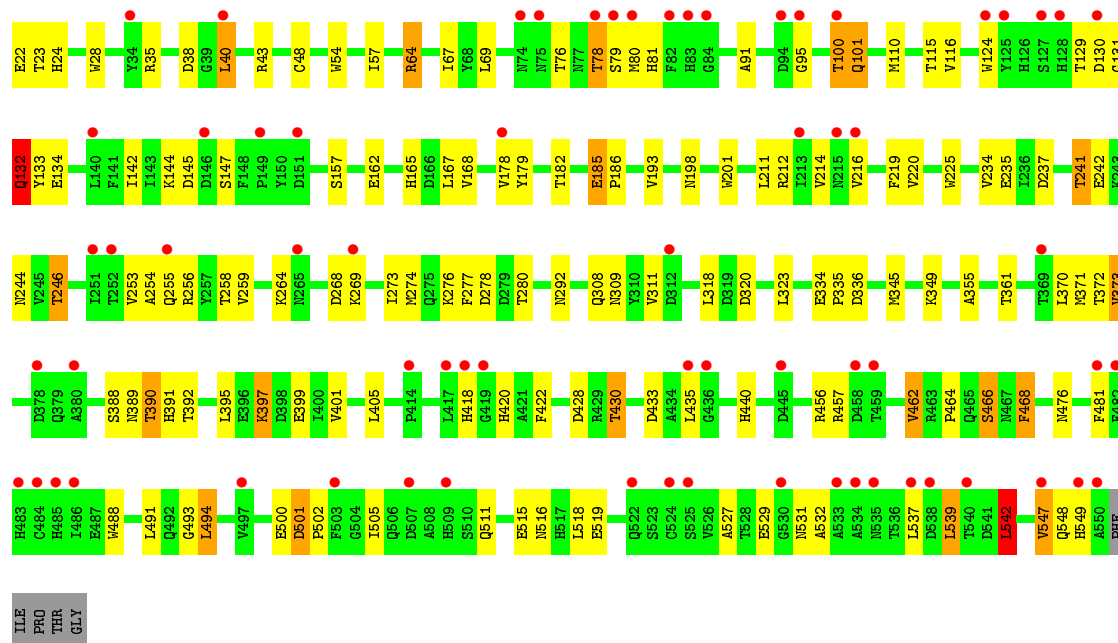




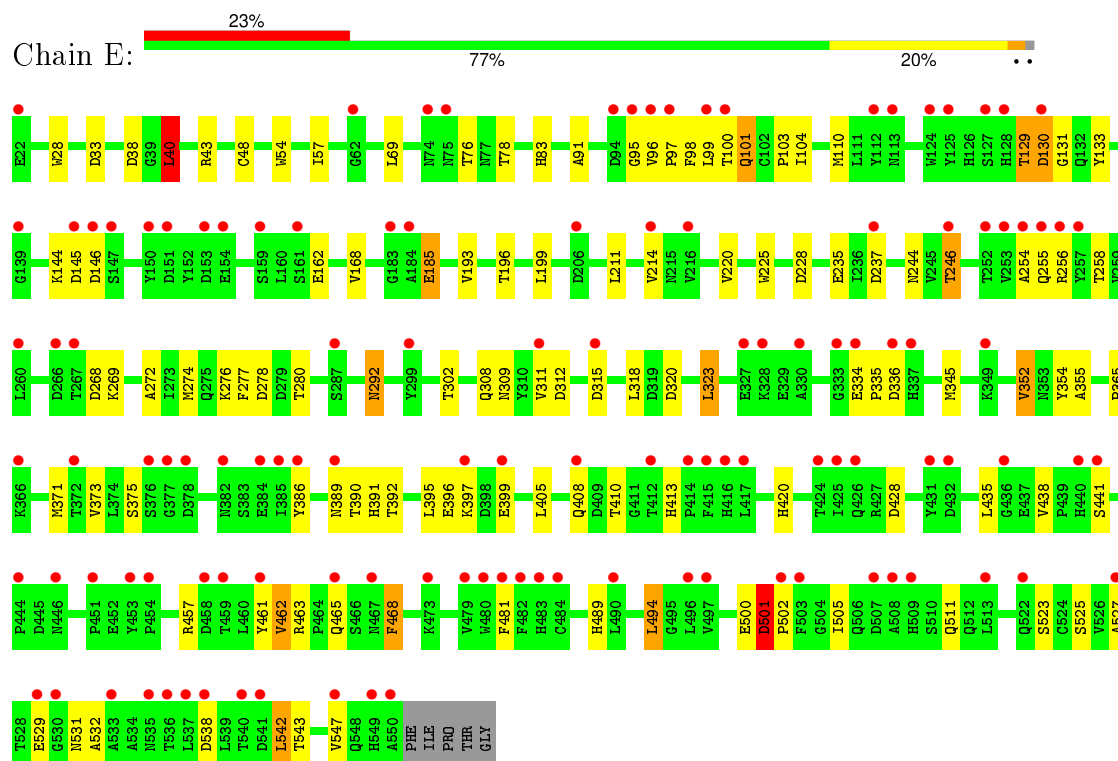
• Molecule 1: Iron transport multicopper oxidase FET3



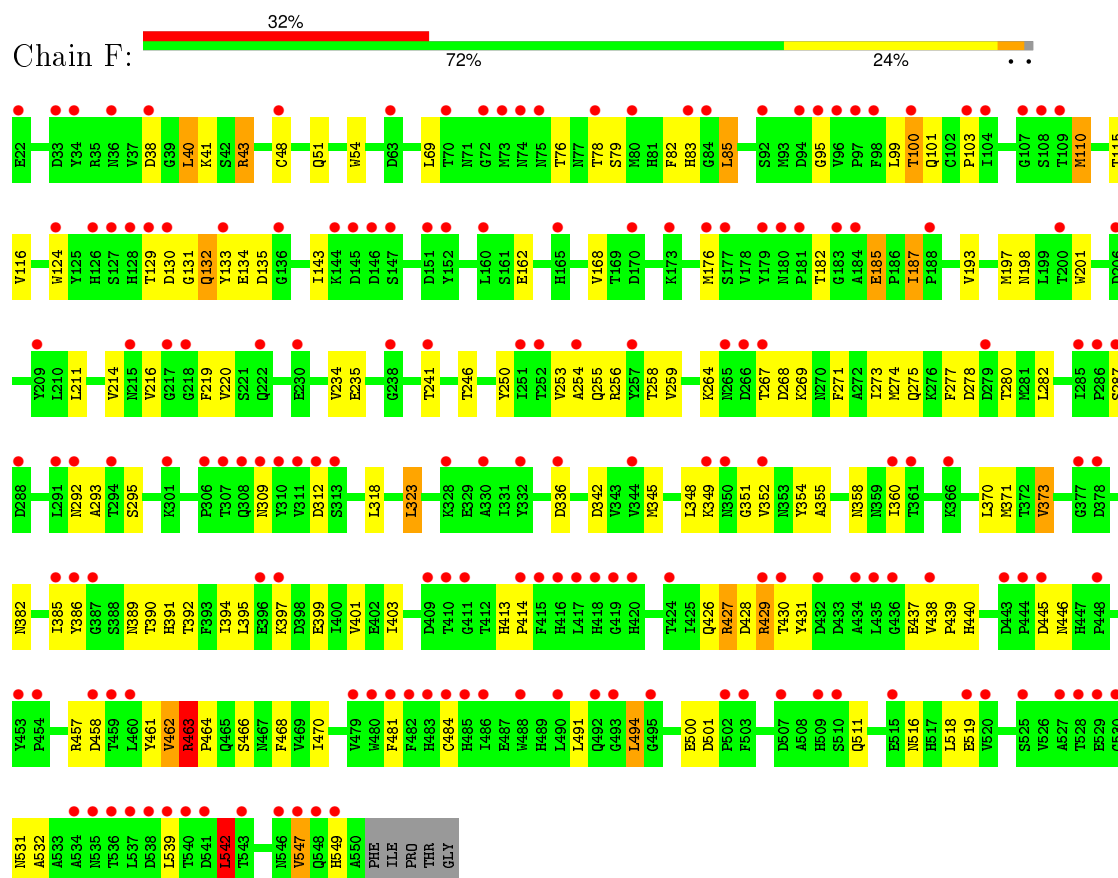
• Molecule 1: Iron transport multicopper oxidase FET3



• Molecule 1: Iron transport multicopper oxidase FET3



• Molecule 1: Iron transport multicopper oxidase FET3



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	168.53Å 168.53Å 174.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.80 36.72 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.2 (50.00-2.80) 98.2 (36.72-2.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.09 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.226 , 0.257 0.235 , 0.265	Depositor DCC
$R_{free}$ test set	6726 reflections (5.28%)	DCC
Wilson B-factor (Å <sup>2</sup> )	62.6	Xtriage
Anisotropy	0.138	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 68.1	EDS
Estimated twinning fraction	0.000 for -h,-k,l 0.010 for h,-h-k,-l 0.000 for -k,-h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 134170 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	27659	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 30.19 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.3641e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CU1, BMA, NAG, NDG, MAN

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	0.50	0/4373	0.66	2/5981 (0.0%)
1	B	0.50	0/4373	0.66	3/5981 (0.1%)
1	C	0.55	5/4373 (0.1%)	0.64	2/5981 (0.0%)
1	D	0.51	3/4373 (0.1%)	0.60	2/5981 (0.0%)
1	E	0.44	1/4373 (0.0%)	0.58	1/5981 (0.0%)
1	F	0.99	20/4373 (0.5%)	0.77	11/5981 (0.2%)
All	All	0.61	29/26238 (0.1%)	0.65	21/35886 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	F	0	1
2	D	1	0
2	E	1	0
3	A	1	0
3	B	1	0
3	D	1	0
3	F	1	0
4	D	1	0
6	A	1	0
6	C	1	0
6	D	1	0
6	E	1	0
6	F	1	0
8	B	1	0
9	C	1	0
9	E	1	0
All	All	15	1

The worst 5 of 29 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	463	ARG	CZ-NH1	40.06	1.85	1.33
1	F	250	TYR	CE2-CZ	15.67	1.58	1.38
1	F	250	TYR	CG-CD2	15.35	1.59	1.39
1	D	515	GLU	CD-OE1	13.48	1.40	1.25
1	F	250	TYR	CE1-CZ	13.27	1.55	1.38

The worst 5 of 21 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	463	ARG	NE-CZ-NH2	-25.91	107.34	120.30
1	F	463	ARG	NE-CZ-NH1	10.71	125.66	120.30
1	F	445	ASP	CB-CG-OD2	-9.31	109.92	118.30
1	F	429	ARG	NE-CZ-NH2	-8.88	115.86	120.30
1	F	463	ARG	NH1-CZ-NH2	-8.13	110.46	119.40

5 of 15 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	2003	NAG	C1
6	A	2010	NAG	C1
3	B	2003	NAG	C1
8	B	2010	NAG	C1
9	C	2003	NAG	C1

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	463	ARG	Sidechain

## 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	4254	0	3960	72	0
1	B	4254	0	3960	97	0
1	C	4254	0	3961	78	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	4254	0	3960	76	0
1	E	4254	0	3962	62	0
1	F	4254	0	3960	78	0
2	A	122	0	104	2	0
2	B	122	0	104	1	0
2	C	122	0	104	1	0
2	D	122	0	104	2	0
2	E	122	0	104	2	0
2	F	122	0	104	1	0
3	A	72	0	61	0	0
3	B	72	0	61	1	0
3	D	72	0	61	0	0
3	F	72	0	61	1	0
4	A	28	0	25	2	0
4	B	56	0	50	0	0
4	C	28	0	25	0	0
4	D	28	0	25	0	0
4	E	28	0	25	0	0
5	A	56	0	52	0	0
5	B	70	0	65	3	0
5	C	70	0	65	2	0
5	D	70	0	64	2	0
5	E	70	0	65	2	0
5	F	84	0	78	4	0
6	A	78	0	68	2	0
6	C	78	0	68	4	0
6	D	39	0	34	0	0
6	E	39	0	34	1	0
6	F	39	0	34	0	0
7	A	14	0	13	0	0
8	B	50	0	43	0	0
9	C	83	0	70	1	0
9	E	83	0	70	2	0
10	A	4	0	0	0	0
10	B	4	0	0	0	0
10	C	4	0	0	0	0
10	D	4	0	0	0	0
10	E	4	0	0	0	0
10	F	4	0	0	0	0
All	All	27659	0	25604	464	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 464 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)
5:F:2006:NAG:C7	5:F:2006:NAG:C8	1.80	1.60
1:F:463:ARG:NH1	1:F:463:ARG:CZ	1.85	1.37
5:F:2005:NAG:C6	5:F:2005:NAG:O6	1.74	1.32
6:C:3013:BMA:C6	6:C:3013:BMA:O6	1.75	1.32
1:F:427:ARG:HD2	1:F:463:ARG:NH1	1.56	1.20

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	527/534 (99%)	497 (94%)	23 (4%)	7 (1%)	15	44
1	B	527/534 (99%)	498 (94%)	23 (4%)	6 (1%)	17	50
1	C	527/534 (99%)	495 (94%)	27 (5%)	5 (1%)	21	55
1	D	527/534 (99%)	497 (94%)	22 (4%)	8 (2%)	13	40
1	E	527/534 (99%)	497 (94%)	24 (5%)	6 (1%)	17	50
1	F	527/534 (99%)	498 (94%)	23 (4%)	6 (1%)	17	50
All	All	3162/3204 (99%)	2982 (94%)	142 (4%)	38 (1%)	16	47

5 of 38 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	130	ASP
1	B	531	ASN
1	C	130	ASP
1	D	130	ASP
1	E	130	ASP

### 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	477/481 (99%)	430 (90%)	47 (10%)	10	28
1	B	477/481 (99%)	425 (89%)	52 (11%)	8	23
1	C	477/481 (99%)	431 (90%)	46 (10%)	10	29
1	D	477/481 (99%)	425 (89%)	52 (11%)	8	23
1	E	477/481 (99%)	432 (91%)	45 (9%)	11	31
1	F	477/481 (99%)	429 (90%)	48 (10%)	9	27
All	All	2862/2886 (99%)	2572 (90%)	290 (10%)	9	27

5 of 290 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	389	ASN
1	D	185	GLU
1	F	292	ASN
1	C	446	ASN
1	D	64	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 15 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	517	HIS
1	C	51	GLN
1	E	382	ASN
1	B	324	GLN
1	E	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 ⓘ

135 carbohydrates are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	NAG	A	2001	1,2	14,14,15	0.69	0	15,19,21	1.32	2 (13%)
2	NAG	A	2002	2	14,14,15	0.52	0	15,19,21	0.99	1 (6%)
3	NAG	A	2003	1,3	14,14,15	0.75	0	15,19,21	1.95	4 (26%)
3	NAG	A	2004	3	14,14,15	0.61	0	15,19,21	0.96	0
4	NAG	A	2005	1,4	14,14,15	0.61	0	15,19,21	1.86	4 (26%)
2	NAG	A	2007	1,2	14,14,15	0.69	0	15,19,21	1.74	3 (20%)
6	NAG	A	2009	1,6	14,14,15	0.49	0	15,19,21	1.67	2 (13%)
6	NAG	A	2010	1,6	14,14,15	0.48	0	15,19,21	1.92	3 (20%)
6	NAG	A	2011	6	14,14,15	0.52	0	15,19,21	1.04	1 (6%)
2	NAG	A	2013	2	14,14,15	0.60	0	15,19,21	2.16	3 (20%)
4	NAG	A	2015	4	14,14,15	1.06	2 (14%)	15,19,21	1.33	1 (6%)
6	NAG	A	2016	6	14,14,15	0.45	0	15,19,21	0.88	0
2	BMA	A	3001	2	11,11,12	0.80	0	14,15,17	1.94	2 (14%)
6	BMA	A	3002	6	11,11,12	0.62	0	14,15,17	2.23	2 (14%)
2	BMA	A	3003	2	11,11,12	0.71	0	14,15,17	1.56	2 (14%)
3	BMA	A	3004	3	11,11,12	0.76	0	14,15,17	2.07	3 (21%)
2	MAN	A	3005	2	11,11,12	0.71	0	14,15,17	1.85	3 (21%)
2	MAN	A	3006	2	11,11,12	0.68	0	14,15,17	1.89	1 (7%)
2	MAN	A	3007	2	11,11,12	0.58	0	14,15,17	1.42	1 (7%)
2	MAN	A	3008	2	11,11,12	0.70	0	14,15,17	2.27	2 (14%)
3	MAN	A	3009	3	11,11,12	0.67	0	14,15,17	1.75	3 (21%)
3	MAN	A	3010	3	11,11,12	0.58	0	14,15,17	1.13	1 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	MAN	A	3011	3	11,11,12	0.69	0	14,15,17	1.17	1 (7%)
6	BMA	A	3013	6	11,11,12	0.91	0	14,15,17	1.24	1 (7%)
2	NAG	B	2001	1,2	14,14,15	0.69	0	15,19,21	0.97	0
2	NAG	B	2002	2	14,14,15	0.59	0	15,19,21	1.16	1 (6%)
3	NAG	B	2003	1,3	14,14,15	0.67	0	15,19,21	1.27	2 (13%)
3	NAG	B	2004	3	14,14,15	0.55	0	15,19,21	1.04	1 (6%)
2	NAG	B	2007	1,2	14,14,15	0.54	0	15,19,21	1.84	3 (20%)
4	NAG	B	2008	1,4	14,14,15	0.54	0	15,19,21	0.86	0
4	NAG	B	2009	1,4	14,14,15	0.45	0	15,19,21	1.87	4 (26%)
8	NAG	B	2010	1,8	14,14,15	0.72	0	15,19,21	1.18	1 (6%)
8	NAG	B	2011	8	14,14,15	0.61	0	15,19,21	1.26	2 (13%)
2	NAG	B	2013	2	14,14,15	0.71	0	15,19,21	1.24	2 (13%)
4	NAG	B	2016	4	14,14,15	0.46	0	15,19,21	1.26	1 (6%)
4	NAG	B	2017	4	14,14,15	0.48	0	15,19,21	1.02	2 (13%)
2	BMA	B	3001	2	11,11,12	0.74	0	14,15,17	1.03	1 (7%)
8	BMA	B	3002	8	11,11,12	0.53	0	14,15,17	1.06	1 (7%)
2	BMA	B	3003	2	11,11,12	0.72	0	14,15,17	1.73	5 (35%)
3	BMA	B	3004	3	11,11,12	0.77	0	14,15,17	1.60	2 (14%)
2	MAN	B	3005	2	11,11,12	0.59	0	14,15,17	2.19	5 (35%)
2	MAN	B	3006	2	11,11,12	1.70	2 (18%)	14,15,17	2.39	3 (21%)
2	MAN	B	3007	2	11,11,12	0.68	0	14,15,17	2.86	5 (35%)
2	MAN	B	3008	2	11,11,12	0.43	0	14,15,17	1.44	3 (21%)
3	MAN	B	3009	3	11,11,12	0.55	0	14,15,17	1.18	2 (14%)
3	MAN	B	3010	3	11,11,12	0.59	0	14,15,17	2.19	2 (14%)
3	MAN	B	3011	3	11,11,12	0.54	0	14,15,17	1.35	2 (14%)
8	MAN	B	3014	8	11,11,12	0.53	0	14,15,17	2.28	4 (28%)
2	NAG	C	2001	1,2	14,14,15	0.71	0	15,19,21	1.17	3 (20%)
2	NAG	C	2002	2	14,14,15	0.61	0	15,19,21	1.46	3 (20%)
9	NAG	C	2003	1,9	14,14,15	0.88	1 (7%)	15,19,21	1.75	6 (40%)
9	NAG	C	2004	9	14,14,15	0.43	0	15,19,21	0.82	0
2	NAG	C	2007	1,2	14,14,15	0.46	0	15,19,21	2.15	2 (13%)
4	NAG	C	2008	1,4	14,14,15	0.50	0	15,19,21	1.18	1 (6%)
6	NAG	C	2009	1,6	14,14,15	1.97	3 (21%)	15,19,21	0.88	0
6	NAG	C	2010	1,6	14,14,15	0.47	0	15,19,21	1.48	4 (26%)
6	NAG	C	2011	6	14,14,15	0.41	0	15,19,21	1.15	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAG	C	2013	2	14,14,15	0.72	0	15,19,21	1.00	2 (13%)
6	NAG	C	2016	6	14,14,15	1.73	3 (21%)	15,19,21	2.37	4 (26%)
4	NAG	C	2017	4	14,14,15	1.19	2 (14%)	15,19,21	1.49	2 (13%)
2	BMA	C	3001	2	11,11,12	0.54	0	14,15,17	2.03	4 (28%)
6	BMA	C	3002	6	11,11,12	0.65	0	14,15,17	0.95	0
2	BMA	C	3003	2	11,11,12	0.68	0	14,15,17	1.40	2 (14%)
9	BMA	C	3004	9	11,11,12	0.57	0	14,15,17	2.04	3 (21%)
2	MAN	C	3005	2	11,11,12	0.75	0	14,15,17	1.67	2 (14%)
2	MAN	C	3006	2	11,11,12	0.59	0	14,15,17	1.69	2 (14%)
2	MAN	C	3007	2	11,11,12	0.55	0	14,15,17	1.53	3 (21%)
2	MAN	C	3008	2	11,11,12	0.62	0	14,15,17	1.43	3 (21%)
9	MAN	C	3009	9	11,11,12	0.76	0	14,15,17	1.95	3 (21%)
9	MAN	C	3010	9	11,11,12	0.58	0	14,15,17	1.37	1 (7%)
9	MAN	C	3011	9	11,11,12	0.69	0	14,15,17	0.83	0
9	MAN	C	3012	9	11,11,12	0.56	0	14,15,17	2.24	3 (21%)
6	BMA	C	3013	6	11,11,12	2.42	1 (9%)	14,15,17	1.92	4 (28%)
2	NAG	D	2001	1,2	14,14,15	0.52	0	15,19,21	0.78	0
2	NAG	D	2002	2	14,14,15	0.57	0	15,19,21	1.28	2 (13%)
3	NAG	D	2003	1,3	14,14,15	0.60	0	15,19,21	1.74	2 (13%)
3	NAG	D	2004	3	14,14,15	0.67	0	15,19,21	0.89	1 (6%)
2	NAG	D	2007	1,2	14,14,15	0.40	0	15,19,21	1.29	1 (6%)
4	NAG	D	2009	1,4	14,14,15	1.39	2 (14%)	15,19,21	1.41	1 (6%)
6	NAG	D	2010	1,6	14,14,15	0.51	0	15,19,21	1.04	0
6	NAG	D	2011	6	14,14,15	0.52	0	15,19,21	0.74	0
2	NAG	D	2013	2	14,14,15	0.56	0	15,19,21	0.84	1 (6%)
4	NAG	D	2016	4	14,14,15	0.49	0	15,19,21	0.72	0
2	BMA	D	3001	2	11,11,12	0.56	0	14,15,17	1.46	3 (21%)
6	BMA	D	3002	6	11,11,12	0.53	0	14,15,17	1.01	2 (14%)
2	BMA	D	3003	2	11,11,12	0.65	0	14,15,17	1.30	2 (14%)
3	BMA	D	3004	3	11,11,12	0.72	0	14,15,17	1.97	4 (28%)
2	MAN	D	3005	2	11,11,12	0.52	0	14,15,17	1.79	3 (21%)
2	MAN	D	3006	2	11,11,12	0.60	0	14,15,17	1.59	2 (14%)
2	MAN	D	3007	2	11,11,12	0.56	0	14,15,17	1.36	1 (7%)
2	MAN	D	3008	2	11,11,12	0.62	0	14,15,17	1.23	2 (14%)
3	MAN	D	3009	3	11,11,12	1.05	1 (9%)	14,15,17	2.16	3 (21%)
3	MAN	D	3010	3	11,11,12	0.58	0	14,15,17	1.18	1 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	MAN	D	3011	3	11,11,12	0.61	0	14,15,17	1.00	1 (7%)
2	NAG	E	2001	1,2	14,14,15	0.58	0	15,19,21	1.01	0
2	NAG	E	2002	2	14,14,15	0.53	0	15,19,21	0.74	0
9	NAG	E	2003	1,9	14,14,15	0.55	0	15,19,21	0.90	0
9	NAG	E	2004	9	14,14,15	0.48	0	15,19,21	1.30	2 (13%)
2	NAG	E	2007	1,2	14,14,15	0.53	0	15,19,21	1.30	3 (20%)
4	NAG	E	2009	1,4	14,14,15	0.52	0	15,19,21	1.06	1 (6%)
6	NAG	E	2010	1,6	14,14,15	0.58	0	15,19,21	1.75	4 (26%)
6	NAG	E	2011	6	14,14,15	0.59	0	15,19,21	1.05	2 (13%)
2	NAG	E	2013	2	14,14,15	0.49	0	15,19,21	0.91	1 (6%)
4	NAG	E	2016	4	14,14,15	0.56	0	15,19,21	1.01	1 (6%)
2	BMA	E	3001	2	11,11,12	0.77	0	14,15,17	2.23	4 (28%)
6	BMA	E	3002	6	11,11,12	0.87	0	14,15,17	2.03	4 (28%)
2	BMA	E	3003	2	11,11,12	0.61	0	14,15,17	1.13	1 (7%)
9	BMA	E	3004	9	11,11,12	0.61	0	14,15,17	1.62	4 (28%)
2	MAN	E	3005	2	11,11,12	0.54	0	14,15,17	1.48	2 (14%)
2	MAN	E	3006	2	11,11,12	0.56	0	14,15,17	1.52	2 (14%)
2	MAN	E	3007	2	11,11,12	0.58	0	14,15,17	0.66	0
2	MAN	E	3008	2	11,11,12	0.53	0	14,15,17	1.49	1 (7%)
9	MAN	E	3009	9	11,11,12	0.53	0	14,15,17	1.66	1 (7%)
9	MAN	E	3010	9	11,11,12	0.62	0	14,15,17	1.60	3 (21%)
9	MAN	E	3011	9	11,11,12	0.54	0	14,15,17	1.01	1 (7%)
9	MAN	E	3012	9	11,11,12	0.69	0	14,15,17	1.64	2 (14%)
2	NAG	F	2001	1,2	14,14,15	0.59	0	15,19,21	0.95	0
2	NAG	F	2002	2	14,14,15	0.58	0	15,19,21	1.21	1 (6%)
3	NAG	F	2003	1,3	14,14,15	0.61	0	15,19,21	1.07	2 (13%)
3	NAG	F	2004	3	14,14,15	0.71	0	15,19,21	1.47	2 (13%)
2	NAG	F	2007	1,2	14,14,15	0.50	0	15,19,21	1.33	1 (6%)
6	NAG	F	2010	1,6	14,14,15	0.49	0	15,19,21	1.11	1 (6%)
6	NAG	F	2011	6	14,14,15	0.56	0	15,19,21	1.39	3 (20%)
2	NAG	F	2013	2	14,14,15	0.53	0	15,19,21	1.09	2 (13%)
2	BMA	F	3001	2	11,11,12	0.63	0	14,15,17	1.50	3 (21%)
6	BMA	F	3002	6	11,11,12	1.81	2 (18%)	14,15,17	1.10	0
2	BMA	F	3003	2	11,11,12	0.69	0	14,15,17	1.36	1 (7%)
3	BMA	F	3004	3	11,11,12	0.78	0	14,15,17	1.03	2 (14%)
2	MAN	F	3005	2	11,11,12	0.65	0	14,15,17	0.92	1 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	MAN	F	3006	2	11,11,12	1.54	1 (9%)	14,15,17	1.90	2 (14%)
2	MAN	F	3007	2	11,11,12	0.62	0	14,15,17	1.84	2 (14%)
2	MAN	F	3008	2	11,11,12	0.47	0	14,15,17	1.69	1 (7%)
3	MAN	F	3009	3	11,11,12	0.58	0	14,15,17	1.80	2 (14%)
3	MAN	F	3010	3	11,11,12	1.16	1 (9%)	14,15,17	1.50	2 (14%)
3	MAN	F	3011	3	11,11,12	0.93	1 (9%)	14,15,17	1.02	1 (7%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	A	2001	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	2002	2	-	0/6/23/26	0/1/1/1
3	NAG	A	2003	1,3	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	A	2004	3	-	0/6/23/26	0/1/1/1
4	NAG	A	2005	1,4	-	0/6/23/26	0/1/1/1
2	NAG	A	2007	1,2	-	0/6/23/26	0/1/1/1
6	NAG	A	2009	1,6	-	2/6/23/26	0/1/1/1
6	NAG	A	2010	1,6	1/1/5/7	0/6/23/26	0/1/1/1
6	NAG	A	2011	6	-	0/6/23/26	0/1/1/1
2	NAG	A	2013	2	-	0/6/23/26	0/1/1/1
4	NAG	A	2015	4	-	0/6/23/26	0/1/1/1
6	NAG	A	2016	6	-	0/6/23/26	0/1/1/1
2	BMA	A	3001	2	-	0/2/19/22	0/1/1/1
6	BMA	A	3002	6	-	0/2/19/22	0/1/1/1
2	BMA	A	3003	2	-	0/2/19/22	0/1/1/1
3	BMA	A	3004	3	-	0/2/19/22	0/1/1/1
2	MAN	A	3005	2	-	0/2/19/22	0/1/1/1
2	MAN	A	3006	2	-	0/2/19/22	0/1/1/1
2	MAN	A	3007	2	-	0/2/19/22	0/1/1/1
2	MAN	A	3008	2	-	0/2/19/22	1/1/1/1
3	MAN	A	3009	3	-	0/2/19/22	0/1/1/1
3	MAN	A	3010	3	-	0/2/19/22	0/1/1/1
3	MAN	A	3011	3	-	0/2/19/22	0/1/1/1
6	BMA	A	3013	6	-	0/2/19/22	0/1/1/1
2	NAG	B	2001	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	2002	2	-	0/6/23/26	0/1/1/1
3	NAG	B	2003	1,3	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	B	2004	3	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	2007	1,2	-	0/6/23/26	0/1/1/1
4	NAG	B	2008	1,4	-	0/6/23/26	0/1/1/1
4	NAG	B	2009	1,4	-	0/6/23/26	0/1/1/1
8	NAG	B	2010	1,8	1/1/5/7	0/6/23/26	0/1/1/1
8	NAG	B	2011	8	-	0/6/23/26	0/1/1/1
2	NAG	B	2013	2	-	0/6/23/26	0/1/1/1
4	NAG	B	2016	4	-	0/6/23/26	0/1/1/1
4	NAG	B	2017	4	-	0/6/23/26	0/1/1/1
2	BMA	B	3001	2	-	0/2/19/22	0/1/1/1
8	BMA	B	3002	8	-	0/2/19/22	0/1/1/1
2	BMA	B	3003	2	-	0/2/19/22	0/1/1/1
3	BMA	B	3004	3	-	0/2/19/22	0/1/1/1
2	MAN	B	3005	2	-	0/2/19/22	0/1/1/1
2	MAN	B	3006	2	-	0/2/19/22	0/1/1/1
2	MAN	B	3007	2	-	0/2/19/22	0/1/1/1
2	MAN	B	3008	2	-	0/2/19/22	1/1/1/1
3	MAN	B	3009	3	-	0/2/19/22	0/1/1/1
3	MAN	B	3010	3	-	0/2/19/22	0/1/1/1
3	MAN	B	3011	3	-	0/2/19/22	0/1/1/1
8	MAN	B	3014	8	-	0/2/19/22	0/1/1/1
2	NAG	C	2001	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2002	2	-	0/6/23/26	0/1/1/1
9	NAG	C	2003	1,9	1/1/5/7	0/6/23/26	0/1/1/1
9	NAG	C	2004	9	-	0/6/23/26	0/1/1/1
2	NAG	C	2007	1,2	-	0/6/23/26	0/1/1/1
4	NAG	C	2008	1,4	-	0/6/23/26	0/1/1/1
6	NAG	C	2009	1,6	-	0/6/23/26	0/1/1/1
6	NAG	C	2010	1,6	1/1/5/7	0/6/23/26	0/1/1/1
6	NAG	C	2011	6	-	0/6/23/26	0/1/1/1
2	NAG	C	2013	2	-	0/6/23/26	0/1/1/1
6	NAG	C	2016	6	-	1/6/23/26	0/1/1/1
4	NAG	C	2017	4	-	0/6/23/26	0/1/1/1
2	BMA	C	3001	2	-	0/2/19/22	0/1/1/1
6	BMA	C	3002	6	-	0/2/19/22	0/1/1/1
2	BMA	C	3003	2	-	0/2/19/22	0/1/1/1
9	BMA	C	3004	9	-	0/2/19/22	0/1/1/1
2	MAN	C	3005	2	-	0/2/19/22	0/1/1/1
2	MAN	C	3006	2	-	0/2/19/22	0/1/1/1
2	MAN	C	3007	2	-	0/2/19/22	0/1/1/1
2	MAN	C	3008	2	-	0/2/19/22	0/1/1/1
9	MAN	C	3009	9	-	0/2/19/22	0/1/1/1
9	MAN	C	3010	9	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	MAN	C	3011	9	-	0/2/19/22	0/1/1/1
9	MAN	C	3012	9	-	0/2/19/22	0/1/1/1
6	BMA	C	3013	6	-	0/2/19/22	0/1/1/1
2	NAG	D	2001	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	2002	2	-	0/6/23/26	0/1/1/1
3	NAG	D	2003	1,3	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	D	2004	3	-	0/6/23/26	0/1/1/1
2	NAG	D	2007	1,2	1/1/5/7	0/6/23/26	0/1/1/1
4	NAG	D	2009	1,4	1/1/5/7	0/6/23/26	0/1/1/1
6	NAG	D	2010	1,6	1/1/5/7	0/6/23/26	0/1/1/1
6	NAG	D	2011	6	-	0/6/23/26	0/1/1/1
2	NAG	D	2013	2	-	0/6/23/26	0/1/1/1
4	NAG	D	2016	4	-	0/6/23/26	0/1/1/1
2	BMA	D	3001	2	-	0/2/19/22	0/1/1/1
6	BMA	D	3002	6	-	0/2/19/22	0/1/1/1
2	BMA	D	3003	2	-	0/2/19/22	0/1/1/1
3	BMA	D	3004	3	-	0/2/19/22	0/1/1/1
2	MAN	D	3005	2	-	0/2/19/22	0/1/1/1
2	MAN	D	3006	2	-	0/2/19/22	0/1/1/1
2	MAN	D	3007	2	-	0/2/19/22	0/1/1/1
2	MAN	D	3008	2	-	0/2/19/22	0/1/1/1
3	MAN	D	3009	3	-	0/2/19/22	0/1/1/1
3	MAN	D	3010	3	-	0/2/19/22	0/1/1/1
3	MAN	D	3011	3	-	0/2/19/22	0/1/1/1
2	NAG	E	2001	1,2	-	0/6/23/26	0/1/1/1
2	NAG	E	2002	2	-	0/6/23/26	0/1/1/1
9	NAG	E	2003	1,9	1/1/5/7	0/6/23/26	0/1/1/1
9	NAG	E	2004	9	-	0/6/23/26	0/1/1/1
2	NAG	E	2007	1,2	1/1/5/7	0/6/23/26	0/1/1/1
4	NAG	E	2009	1,4	-	1/6/23/26	0/1/1/1
6	NAG	E	2010	1,6	1/1/5/7	0/6/23/26	0/1/1/1
6	NAG	E	2011	6	-	0/6/23/26	0/1/1/1
2	NAG	E	2013	2	-	0/6/23/26	0/1/1/1
4	NAG	E	2016	4	-	0/6/23/26	0/1/1/1
2	BMA	E	3001	2	-	0/2/19/22	0/1/1/1
6	BMA	E	3002	6	-	0/2/19/22	0/1/1/1
2	BMA	E	3003	2	-	0/2/19/22	0/1/1/1
9	BMA	E	3004	9	-	0/2/19/22	0/1/1/1
2	MAN	E	3005	2	-	0/2/19/22	1/1/1/1
2	MAN	E	3006	2	-	0/2/19/22	0/1/1/1
2	MAN	E	3007	2	-	0/2/19/22	0/1/1/1
2	MAN	E	3008	2	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	MAN	E	3009	9	-	0/2/19/22	0/1/1/1
9	MAN	E	3010	9	-	0/2/19/22	0/1/1/1
9	MAN	E	3011	9	-	0/2/19/22	0/1/1/1
9	MAN	E	3012	9	-	0/2/19/22	0/1/1/1
2	NAG	F	2001	1,2	-	0/6/23/26	0/1/1/1
2	NAG	F	2002	2	-	0/6/23/26	0/1/1/1
3	NAG	F	2003	1,3	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	F	2004	3	-	0/6/23/26	0/1/1/1
2	NAG	F	2007	1,2	-	2/6/23/26	0/1/1/1
6	NAG	F	2010	1,6	1/1/5/7	0/6/23/26	0/1/1/1
6	NAG	F	2011	6	-	0/6/23/26	0/1/1/1
2	NAG	F	2013	2	-	0/6/23/26	0/1/1/1
2	BMA	F	3001	2	-	0/2/19/22	0/1/1/1
6	BMA	F	3002	6	-	0/2/19/22	0/1/1/1
2	BMA	F	3003	2	-	0/2/19/22	0/1/1/1
3	BMA	F	3004	3	-	0/2/19/22	0/1/1/1
2	MAN	F	3005	2	-	0/2/19/22	0/1/1/1
2	MAN	F	3006	2	-	0/2/19/22	1/1/1/1
2	MAN	F	3007	2	-	0/2/19/22	0/1/1/1
2	MAN	F	3008	2	-	0/2/19/22	1/1/1/1
3	MAN	F	3009	3	-	0/2/19/22	0/1/1/1
3	MAN	F	3010	3	-	0/2/19/22	0/1/1/1
3	MAN	F	3011	3	-	0/2/19/22	0/1/1/1

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	C	2003	NAG	O5-C1	-2.16	1.40	1.43
4	A	2015	NAG	C8-C7	2.03	1.54	1.50
4	A	2015	NAG	C1-C2	2.40	1.55	1.52
6	C	2016	NAG	C7-N2	2.45	1.43	1.34
6	F	3002	BMA	C4-C3	2.47	1.58	1.52

The worst 5 of 258 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	2016	NAG	C2-N2-C7	-6.72	114.40	123.04
2	D	3005	MAN	C1-C2-C3	-5.11	103.50	109.54
2	B	3005	MAN	C1-C2-C3	-4.85	103.81	109.54
2	A	3005	MAN	C1-C2-C3	-4.51	104.20	109.54
9	C	3009	MAN	O5-C1-C2	-4.34	103.82	110.86

5 of 15 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	2003	NAG	C1
8	B	2010	NAG	C1
3	D	2003	NAG	C1
3	B	2003	NAG	C1
6	E	2010	NAG	C1

5 of 6 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	2009	NAG	O7-C7-N2-C2
6	C	2016	NAG	O7-C7-N2-C2
6	A	2009	NAG	O7-C7-N2-C2
2	F	2007	NAG	O7-C7-N2-C2
6	A	2009	NAG	C8-C7-N2-C2

All (5) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	3006	MAN	C1-C2-C3-C4-C5-O5
2	A	3008	MAN	C1-C2-C3-C4-C5-O5
2	E	3005	MAN	C1-C2-C3-C4-C5-O5
2	F	3008	MAN	C1-C2-C3-C4-C5-O5
2	B	3008	MAN	C1-C2-C3-C4-C5-O5

21 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	2005	NAG	2	0
6	A	2010	NAG	2	0
2	A	2013	NAG	1	0
2	A	3001	BMA	1	0
2	A	3007	MAN	1	0
3	B	2003	NAG	1	0
2	B	2007	NAG	1	0
2	C	2002	NAG	1	0
9	C	2003	NAG	1	0
6	C	2010	NAG	1	0
6	C	2016	NAG	1	0
6	C	3013	BMA	2	0
2	D	2002	NAG	1	0
2	D	3008	MAN	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	2001	NAG	1	0
9	E	2003	NAG	1	0
9	E	2004	NAG	2	0
2	E	2007	NAG	1	0
6	E	2010	NAG	1	0
3	F	2003	NAG	1	0
2	F	2007	NAG	1	0

## 5.6 Ligand geometry [i](#)

Of 55 ligands modelled in this entry, 24 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	NAG	A	2006	1	14,14,15	0.85	0	15,19,21	0.89	0
5	NAG	A	2008	1	14,14,15	0.46	0	15,19,21	0.75	0
5	NAG	A	2012	1	14,14,15	0.53	0	15,19,21	1.12	1 (6%)
5	NAG	A	2014	1	14,14,15	0.61	0	15,19,21	1.05	1 (6%)
7	NDG	A	2018	1	14,14,15	0.67	0	15,19,21	1.55	2 (13%)
5	NAG	B	2005	1	14,14,15	0.56	0	15,19,21	0.78	0
5	NAG	B	2006	1	14,14,15	0.66	0	15,19,21	0.67	0
5	NAG	B	2012	1	14,14,15	0.52	0	15,19,21	1.23	1 (6%)
5	NAG	B	2014	1	14,14,15	0.45	0	15,19,21	1.83	1 (6%)
5	NAG	B	2018	1	14,14,15	0.76	1 (7%)	15,19,21	1.22	1 (6%)
5	NAG	C	2005	1	14,14,15	0.52	0	15,19,21	1.74	4 (26%)
5	NAG	C	2006	1	14,14,15	0.45	0	15,19,21	1.44	1 (6%)
5	NAG	C	2012	1	14,14,15	0.51	0	15,19,21	1.70	2 (13%)
5	NAG	C	2014	1	14,14,15	0.53	0	15,19,21	1.90	2 (13%)
5	NAG	C	2018	1	14,14,15	5.02	3 (21%)	15,19,21	2.16	4 (26%)
5	NAG	D	2005	1	14,14,15	0.51	0	15,19,21	1.00	0
5	NAG	D	2006	1	14,14,15	3.78	3 (21%)	15,19,21	1.98	3 (20%)
5	NAG	D	2008	1	14,14,15	0.47	0	15,19,21	0.91	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	NAG	D	2012	1	14,14,15	0.65	0	15,19,21	0.82	0
5	NAG	D	2014	1	14,14,15	0.51	0	15,19,21	0.85	0
5	NAG	E	2005	1	14,14,15	0.54	0	15,19,21	0.85	0
5	NAG	E	2006	1	14,14,15	0.64	0	15,19,21	1.74	3 (20%)
5	NAG	E	2008	1	14,14,15	0.50	0	15,19,21	0.83	0
5	NAG	E	2012	1	14,14,15	3.46	2 (14%)	15,19,21	1.52	3 (20%)
5	NAG	E	2014	1	14,14,15	0.46	0	15,19,21	1.93	2 (13%)
5	NAG	F	2005	1	14,14,15	4.21	4 (28%)	15,19,21	2.01	3 (20%)
5	NAG	F	2006	1	14,14,15	4.19	4 (28%)	15,19,21	1.86	3 (20%)
5	NAG	F	2008	1	14,14,15	0.53	0	15,19,21	1.23	1 (6%)
5	NAG	F	2009	1	14,14,15	0.85	0	15,19,21	1.43	2 (13%)
5	NAG	F	2012	1	14,14,15	0.47	0	15,19,21	1.57	3 (20%)
5	NAG	F	2014	1	14,14,15	0.44	0	15,19,21	1.21	1 (6%)

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	NAG	A	2006	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	A	2008	1	-	0/6/23/26	0/1/1/1
5	NAG	A	2012	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	A	2014	1	-	0/6/23/26	0/1/1/1
7	NDG	A	2018	1	-	0/6/23/26	0/1/1/1
5	NAG	B	2005	1	-	0/6/23/26	0/1/1/1
5	NAG	B	2006	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	B	2012	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	B	2014	1	-	0/6/23/26	0/1/1/1
5	NAG	B	2018	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	C	2005	1	-	0/6/23/26	0/1/1/1
5	NAG	C	2006	1	-	0/6/23/26	0/1/1/1
5	NAG	C	2012	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	C	2014	1	-	0/6/23/26	0/1/1/1
5	NAG	C	2018	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	D	2005	1	-	0/6/23/26	0/1/1/1
5	NAG	D	2006	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	D	2008	1	-	0/6/23/26	0/1/1/1
5	NAG	D	2012	1	-	0/6/23/26	0/1/1/1
5	NAG	D	2014	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	E	2005	1	-	0/6/23/26	0/1/1/1
5	NAG	E	2006	1	-	0/6/23/26	0/1/1/1
5	NAG	E	2008	1	-	0/6/23/26	0/1/1/1
5	NAG	E	2012	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	E	2014	1	-	0/6/23/26	0/1/1/1
5	NAG	F	2005	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	F	2006	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	F	2008	1	-	0/6/23/26	0/1/1/1
5	NAG	F	2009	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	F	2012	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	F	2014	1	-	0/6/23/26	0/1/1/1

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	2006	NAG	O7-C7	-3.29	1.15	1.23
5	F	2006	NAG	C1-C2	2.10	1.55	1.52
5	B	2018	NAG	C1-C2	2.16	1.55	1.52
5	D	2006	NAG	C1-C2	2.37	1.55	1.52
5	F	2006	NAG	C7-N2	2.51	1.44	1.34

The worst 5 of 44 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	2006	NAG	C8-C7-N2	-4.60	107.31	116.11
5	F	2006	NAG	C2-N2-C7	-4.51	117.24	123.04
5	C	2018	NAG	O7-C7-N2	-4.40	112.90	121.86
5	F	2006	NAG	O7-C7-N2	-3.95	113.82	121.86
5	D	2006	NAG	C2-N2-C7	-3.58	118.44	123.04

5 of 13 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	A	2012	NAG	C1
5	B	2012	NAG	C1
5	F	2012	NAG	C1
5	B	2018	NAG	C1
5	B	2006	NAG	C1

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	D	2014	NAG	C8-C7-N2-C2
5	D	2014	NAG	O7-C7-N2-C2

There are no ring outliers.

9 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	2005	NAG	3	0
5	C	2005	NAG	2	0
5	D	2005	NAG	1	0
5	D	2008	NAG	1	0
5	E	2005	NAG	1	0
5	E	2006	NAG	1	0
5	F	2005	NAG	2	0
5	F	2006	NAG	1	0
5	F	2012	NAG	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.



## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<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	529/534 (99%)	0.71	44 (8%)	14 7	67, 69, 71, 74	0
1	B	529/534 (99%)	0.70	39 (7%)	17 9	67, 69, 71, 74	0
1	C	529/534 (99%)	0.80	64 (12%)	6 3	67, 69, 71, 74	0
1	D	529/534 (99%)	0.86	67 (12%)	5 2	67, 69, 71, 73	0
1	E	529/534 (99%)	1.31	121 (22%)	1 1	68, 69, 71, 73	0
1	F	529/534 (99%)	1.63	172 (32%)	1 0	68, 69, 71, 73	0
All	All	3174/3204 (99%)	1.00	507 (15%)	3 1	67, 69, 71, 74	0

The worst 5 of 507 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	549	HIS	10.2
1	F	266	ASP	8.7
1	F	443	ASP	8.1
1	E	536	THR	7.2
1	E	549	HIS	7.2

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

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

### 6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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( $\text{\AA}^2$ )	Q<0.9
6	NAG	D	2010	14/15	0.85	0.36	3.54	72,73,75,76	0
4	NAG	C	2008	14/15	0.81	0.40	3.22	75,77,78,80	0
2	NAG	E	2007	14/15	0.73	0.47	2.15	72,73,74,74	0
2	NAG	C	2007	14/15	0.75	0.37	1.28	69,73,76,76	0
6	NAG	E	2010	14/15	0.88	0.36	1.26	63,66,69,69	0
6	NAG	F	2010	14/15	0.73	0.44	1.09	75,76,77,77	0
4	NAG	B	2008	14/15	0.88	0.28	1.07	76,79,81,86	0
4	NAG	A	2005	14/15	0.93	0.24	1.04	71,76,80,87	0
2	NAG	F	2007	14/15	0.79	0.41	0.74	71,72,75,76	0
3	NAG	F	2003	14/15	0.84	0.27	0.18	78,81,83,83	0
2	NAG	A	2007	14/15	0.78	0.27	0.13	74,76,80,83	0
3	MAN	F	3011	11/12	0.73	0.32	0.12	75,76,77,77	0
2	NAG	B	2007	14/15	0.86	0.23	0.10	74,77,81,84	0
9	MAN	E	3011	11/12	0.92	0.22	0.00	71,72,73,73	0
6	NAG	C	2010	14/15	0.92	0.22	-0.35	62,64,67,68	0
2	NAG	E	2002	14/15	0.85	0.25	-0.53	75,77,79,79	0
9	NAG	C	2003	14/15	0.92	0.20	-0.71	69,70,76,79	0
2	NAG	D	2007	14/15	0.89	0.22	-0.72	73,75,78,78	0
3	NAG	D	2003	14/15	0.90	0.18	-0.74	64,68,70,71	0
3	MAN	D	3011	11/12	0.93	0.19	-0.84	63,64,65,65	0
8	NAG	B	2010	14/15	0.89	0.22	-1.11	68,69,71,72	0
2	NAG	F	2001	14/15	0.87	0.18	-1.21	70,71,72,72	0
2	NAG	C	2002	14/15	0.90	0.15	-1.38	71,74,78,82	0
3	NAG	B	2003	14/15	0.94	0.18	-1.48	70,74,77,77	0
2	NAG	A	2001	14/15	0.96	0.17	-1.48	63,65,68,68	0
3	NAG	A	2003	14/15	0.93	0.17	-1.51	67,72,78,80	0
2	NAG	D	2001	14/15	0.95	0.14	-1.57	68,69,69,71	0
6	NAG	A	2010	14/15	0.94	0.19	-1.72	64,67,69,70	0
2	NAG	F	2002	14/15	0.92	0.17	-1.79	71,71,74,75	0
2	NAG	A	2002	14/15	0.95	0.15	-1.79	70,72,73,77	0
2	NAG	D	2002	14/15	0.93	0.14	-1.83	70,72,74,74	0
2	NAG	B	2002	14/15	0.92	0.16	-1.85	72,74,76,79	0
2	NAG	B	2001	14/15	0.96	0.17	-1.93	67,68,70,71	0
9	NAG	E	2003	14/15	0.92	0.18	-2.11	76,77,79,79	0
3	MAN	B	3011	11/12	0.95	0.14	-2.24	68,70,71,73	0
2	NAG	C	2001	14/15	0.95	0.13	-2.77	67,69,71,72	0
2	NAG	E	2001	14/15	0.95	0.10	-3.53	70,71,73,75	0
9	MAN	C	3011	11/12	0.97	0.11	-3.96	64,65,66,67	0
3	MAN	A	3011	11/12	0.94	0.13	-4.29	67,72,73,75	0
6	BMA	F	3002	11/12	0.64	0.67	-	81,82,82,83	0
2	BMA	B	3003	11/12	0.84	0.31	-	71,72,73,74	0
9	BMA	C	3004	11/12	0.93	0.21	-	77,81,85,88	0
9	NAG	E	2004	14/15	0.86	0.28	-	80,82,82,83	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NAG	F	2004	14/15	0.72	0.43	-	82,83,85,85	0
2	NAG	F	2013	14/15	0.77	0.45	-	68,70,70,71	0
4	NAG	B	2016	14/15	0.86	0.33	-	74,76,77,78	0
2	NAG	E	2013	14/15	0.86	0.43	-	73,74,74,74	0
6	BMA	C	3013	11/12	0.87	0.50	-	47,49,50,50	0
2	MAN	F	3008	11/12	0.71	0.34	-	82,84,84,85	0
9	NAG	C	2004	14/15	0.95	0.17	-	72,75,77,79	0
2	MAN	B	3005	11/12	0.62	0.28	-	70,73,74,75	0
3	MAN	D	3010	11/12	0.85	0.31	-	67,68,69,69	0
2	BMA	F	3003	11/12	0.68	0.46	-	66,67,68,68	0
2	BMA	C	3003	11/12	0.90	0.29	-	81,83,86,88	0
2	MAN	D	3008	11/12	0.74	0.29	-	77,78,78,78	0
2	BMA	E	3001	11/12	0.82	0.26	-	81,83,86,87	0
2	MAN	D	3007	11/12	0.66	0.47	-	78,78,79,79	0
6	BMA	A	3002	11/12	0.61	0.56	-	78,80,81,81	0
3	BMA	D	3004	11/12	0.91	0.34	-	70,72,73,73	0
2	MAN	C	3006	11/12	0.65	0.51	-	89,90,91,91	0
4	NAG	B	2009	14/15	0.87	0.24	-	70,71,73,74	0
2	MAN	D	3005	11/12	0.70	0.27	-	74,76,76,77	0
2	MAN	A	3007	11/12	0.79	0.26	-	92,93,94,95	0
2	MAN	C	3007	11/12	0.72	0.46	-	96,97,98,99	0
4	NAG	C	2017	14/15	0.69	0.64	-	82,83,84,85	0
9	MAN	C	3009	11/12	0.71	0.29	-	92,95,96,97	0
2	BMA	B	3001	11/12	0.84	0.18	-	82,85,89,90	0
2	NAG	A	2013	14/15	0.90	0.25	-	71,73,75,75	0
2	MAN	C	3008	11/12	0.75	0.30	-	95,95,96,96	0
4	NAG	B	2017	14/15	0.85	0.55	-	89,90,92,93	0
6	NAG	A	2009	14/15	0.90	0.22	-	64,69,71,72	0
4	NAG	D	2009	14/15	0.81	0.32	-	68,69,70,72	0
2	MAN	A	3008	11/12	0.56	0.36	-	90,91,94,94	0
4	NAG	E	2009	14/15	0.83	0.37	-	65,67,67,67	0
2	BMA	E	3003	11/12	0.72	0.28	-	75,76,77,78	0
3	BMA	B	3004	11/12	0.86	0.25	-	79,85,88,92	0
2	NAG	C	2013	14/15	0.83	0.32	-	76,78,80,80	0
2	MAN	E	3006	11/12	0.59	0.56	-	78,79,79,79	0
3	MAN	A	3009	11/12	0.73	0.31	-	98,100,100,100	0
2	BMA	F	3001	11/12	0.86	0.29	-	76,77,79,82	0
6	BMA	D	3002	11/12	0.57	0.66	-	80,80,80,81	0
3	MAN	B	3010	11/12	0.88	0.23	-	73,77,78,79	0
2	MAN	B	3007	11/12	0.79	0.34	-	91,92,93,94	0
3	MAN	D	3009	11/12	0.66	0.36	-	73,75,76,76	0
2	MAN	B	3008	11/12	0.83	0.19	-	93,94,95,95	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MAN	F	3005	11/12	0.60	0.47	-	63,65,65,65	0
4	NAG	D	2016	14/15	0.73	0.56	-	73,74,75,76	0
9	BMA	E	3004	11/12	0.77	0.34	-	82,86,89,92	0
6	NAG	A	2011	14/15	0.92	0.36	-	71,72,73,75	0
2	BMA	C	3001	11/12	0.82	0.39	-	87,90,93,94	0
6	BMA	E	3002	11/12	0.77	0.36	-	65,66,66,66	0
2	MAN	F	3007	11/12	0.79	0.41	-	77,78,79,79	0
9	MAN	C	3010	11/12	0.95	0.20	-	72,76,80,83	0
2	NAG	D	2013	14/15	0.84	0.41	-	71,74,75,77	0
4	NAG	A	2015	14/15	0.58	0.58	-	92,95,96,96	0
3	BMA	F	3004	11/12	0.57	0.66	-	80,82,82,83	0
6	NAG	C	2009	14/15	0.91	0.40	-	58,62,65,66	0
3	BMA	A	3004	11/12	0.94	0.24	-	80,86,89,93	0
9	MAN	E	3010	11/12	0.89	0.22	-	76,80,82,84	0
6	NAG	D	2011	14/15	0.89	0.33	-	75,76,76,78	0
8	BMA	B	3002	11/12	0.53	0.53	-	82,84,85,87	0
3	NAG	B	2004	14/15	0.93	0.24	-	78,79,80,82	0
2	MAN	C	3005	11/12	0.64	0.29	-	82,84,86,86	0
2	BMA	A	3003	11/12	0.89	0.27	-	73,74,75,77	0
3	NAG	A	2004	14/15	0.96	0.23	-	79,82,86,86	0
6	NAG	E	2011	14/15	0.84	0.35	-	65,65,66,66	0
2	MAN	B	3006	11/12	0.78	0.42	-	74,75,75,75	0
3	MAN	F	3010	11/12	0.69	0.36	-	76,78,79,79	0
6	NAG	F	2011	14/15	0.71	0.57	-	77,78,79,81	0
2	BMA	A	3001	11/12	0.84	0.17	-	80,84,88,89	0
3	NAG	D	2004	14/15	0.89	0.21	-	67,68,69,71	0
2	NAG	B	2013	14/15	0.83	0.29	-	68,71,71,72	0
2	MAN	E	3005	11/12	0.65	0.35	-	74,76,76,76	0
2	MAN	E	3008	11/12	0.67	0.41	-	88,89,89,90	0
2	MAN	D	3006	11/12	0.60	0.34	-	81,82,83,83	0
3	MAN	B	3009	11/12	0.78	0.42	-	96,100,100,100	0
6	NAG	C	2016	14/15	0.80	0.43	-	52,53,54,55	0
3	MAN	F	3009	11/12	0.54	0.50	-	80,80,81,82	0
6	NAG	A	2016	14/15	0.81	0.30	-	73,73,74,76	0
2	MAN	E	3007	11/12	0.69	0.45	-	87,88,89,89	0
9	MAN	E	3009	11/12	0.29	0.58	-	95,97,97,97	0
8	NAG	B	2011	14/15	0.90	0.32	-	73,75,76,79	0
2	MAN	F	3006	11/12	0.71	0.54	-	67,69,69,69	0
9	MAN	E	3012	11/12	0.64	0.46	-	83,85,86,87	0
6	BMA	A	3013	11/12	0.76	0.53	-	73,74,75,75	0
8	MAN	B	3014	11/12	0.66	0.54	-	88,88,89,89	0
2	MAN	A	3006	11/12	0.56	0.50	-	77,78,79,79	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	BMA	D	3003	11/12	0.67	0.42	-	78,78,80,81	0
4	NAG	E	2016	14/15	0.87	0.36	-	64,65,66,66	0
3	MAN	A	3010	11/12	0.93	0.23	-	74,77,78,79	0
2	BMA	D	3001	11/12	0.87	0.22	-	75,76,77,77	0
9	MAN	C	3012	11/12	0.78	0.37	-	86,87,88,90	0
6	NAG	C	2011	14/15	0.91	0.28	-	68,71,75,76	0
2	MAN	A	3005	11/12	0.64	0.30	-	74,76,77,78	0
6	BMA	C	3002	11/12	0.67	0.51	-	78,80,81,82	0

## 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. 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( $\text{\AA}^2$ )	Q<0.9
5	NAG	E	2005	14/15	0.69	0.53	5.12	70,71,71,71	0
5	NAG	F	2005	14/15	0.74	0.52	4.77	67,70,71,72	0
5	NAG	B	2014	14/15	0.79	0.36	4.56	67,70,71,71	0
5	NAG	B	2012	14/15	0.90	0.28	3.88	74,75,79,79	0
5	NAG	D	2014	14/15	0.75	0.43	3.82	65,66,67,67	0
5	NAG	E	2008	14/15	0.68	0.43	3.39	72,72,73,73	0
5	NAG	C	2012	14/15	0.84	0.35	2.86	75,77,79,80	0
5	NAG	A	2012	14/15	0.83	0.36	2.31	76,79,80,81	0
5	NAG	D	2012	14/15	0.85	0.33	2.07	73,74,76,77	0
5	NAG	D	2008	14/15	0.83	0.38	2.01	70,71,73,73	0
5	NAG	C	2005	14/15	0.86	0.31	1.84	64,68,70,70	0
5	NAG	A	2014	14/15	0.87	0.25	1.53	67,67,71,72	0
5	NAG	D	2005	14/15	0.86	0.27	1.27	71,72,74,75	0
5	NAG	F	2008	14/15	0.84	0.35	0.99	71,71,72,72	0
5	NAG	F	2014	14/15	0.89	0.31	0.77	59,62,63,64	0
5	NAG	C	2014	14/15	0.89	0.23	0.11	53,55,57,58	0
5	NAG	E	2014	14/15	0.91	0.26	0.02	52,56,57,58	0
5	NAG	A	2008	14/15	0.90	0.22	-0.06	68,69,70,70	0
5	NAG	B	2005	14/15	0.91	0.20	-0.78	64,67,69,70	0
10	CU1	E	1001	1/1	0.92	0.23	-1.10	88,88,88,88	0
10	CU1	E	1004	1/1	0.85	0.17	-2.13	100,100,100,100	0
10	CU1	F	1001	1/1	0.85	0.20	-2.34	90,90,90,90	0
10	CU1	E	1002	1/1	0.95	0.09	-2.79	93,93,93,93	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
10	CU1	A	1001	1/1	0.98	0.12	-2.85	82,82,82,82	0
10	CU1	E	1003	1/1	0.91	0.15	-2.99	99,99,99,99	0
10	CU1	C	1001	1/1	0.97	0.07	-3.12	85,85,85,85	0
10	CU1	B	1002	1/1	0.98	0.17	-3.62	88,88,88,88	0
10	CU1	D	1001	1/1	0.93	0.09	-3.73	86,86,86,86	0
10	CU1	C	1002	1/1	0.98	0.09	-3.80	85,85,85,85	0
10	CU1	C	1004	1/1	0.93	0.11	-3.81	100,100,100,100	0
10	CU1	A	1002	1/1	0.98	0.16	-3.92	87,87,87,87	0
10	CU1	A	1003	1/1	0.99	0.18	-4.20	89,89,89,89	0
10	CU1	F	1003	1/1	0.95	0.09	-4.27	91,91,91,91	0
10	CU1	B	1001	1/1	0.97	0.10	-4.41	87,87,87,87	0
10	CU1	F	1004	1/1	0.95	0.09	-4.53	100,100,100,100	0
10	CU1	F	1002	1/1	0.94	0.07	-4.92	96,96,96,96	0
10	CU1	D	1002	1/1	0.99	0.07	-5.01	87,87,87,87	0
10	CU1	B	1003	1/1	0.98	0.16	-5.05	87,87,87,87	0
10	CU1	D	1004	1/1	0.95	0.07	-5.63	100,100,100,100	0
10	CU1	C	1003	1/1	0.97	0.10	-6.12	83,83,83,83	0
10	CU1	D	1003	1/1	0.96	0.10	-7.02	89,89,89,89	0
10	CU1	A	1004	1/1	0.90	0.12	-9.75	100,100,100,100	0
10	CU1	B	1004	1/1	0.98	0.10	-10.23	100,100,100,100	0
5	NAG	E	2012	14/15	0.54	0.53	-	73,74,77,77	0
7	NDG	A	2018	14/15	0.63	0.44	-	78,80,82,82	0
5	NAG	B	2018	14/15	0.84	0.40	-	80,83,83,83	0
5	NAG	C	2018	14/15	0.69	0.52	-	74,76,77,78	0
5	NAG	D	2006	14/15	0.81	0.33	-	65,68,70,71	0
5	NAG	F	2006	14/15	0.69	0.51	-	67,68,70,70	0
5	NAG	C	2006	14/15	0.84	0.33	-	74,76,77,78	0
5	NAG	B	2006	14/15	0.84	0.27	-	74,76,78,79	0
5	NAG	F	2009	14/15	0.84	0.28	-	66,69,69,69	0
5	NAG	F	2012	14/15	0.69	0.57	-	72,73,75,75	0
5	NAG	A	2006	14/15	0.75	0.30	-	71,73,74,74	0
5	NAG	E	2006	14/15	0.72	0.43	-	73,74,75,76	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.