



# Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2022 – 04:05 PM EDT

PDB ID : 7TU4  
Title : Structure of the *L. blandensis* dGTPase del55-58 mutant bound to Mn  
Authors : Sikkema, A.P.; Klemm, B.P.; Horng, J.C.; Hall, T.M.T.  
Deposited on : 2022-02-02  
Resolution : 2.26 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.28.1  
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) : 2.28.1

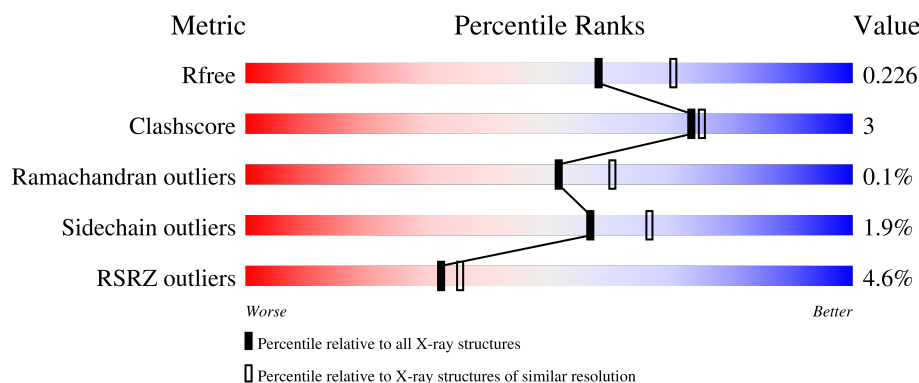
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.26 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	460	<div> <div>2%</div> <div> <div></div> <div>85%</div> <div>8%</div> <div>7%</div> </div> </div>
1	B	460	<div> <div>5%</div> <div> <div></div> <div>86%</div> <div>8%</div> <div>6%</div> </div> </div>
1	C	460	<div> <div>6%</div> <div> <div></div> <div>82%</div> <div>12%</div> <div>7%</div> </div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10330 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 dGTP triphosphohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	430	Total	C	N	O	S	0	0	0
			3388	2150	574	656	8			
1	B	431	Total	C	N	O	S	0	0	0
			3379	2146	571	654	8			
1	C	430	Total	C	N	O	S	0	0	0
			3362	2137	570	647	8			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	expression tag	UNP A3XHN1
A	-20	HIS	-	expression tag	UNP A3XHN1
A	-19	HIS	-	expression tag	UNP A3XHN1
A	-18	HIS	-	expression tag	UNP A3XHN1
A	-17	HIS	-	expression tag	UNP A3XHN1
A	-16	HIS	-	expression tag	UNP A3XHN1
A	-15	HIS	-	expression tag	UNP A3XHN1
A	-14	SER	-	expression tag	UNP A3XHN1
A	-13	SER	-	expression tag	UNP A3XHN1
A	-12	GLY	-	expression tag	UNP A3XHN1
A	-11	VAL	-	expression tag	UNP A3XHN1
A	-10	ASP	-	expression tag	UNP A3XHN1
A	-9	LEU	-	expression tag	UNP A3XHN1
A	-8	GLY	-	expression tag	UNP A3XHN1
A	-7	THR	-	expression tag	UNP A3XHN1
A	-6	GLU	-	expression tag	UNP A3XHN1
A	-5	ASN	-	expression tag	UNP A3XHN1
A	-4	LEU	-	expression tag	UNP A3XHN1
A	-3	TYR	-	expression tag	UNP A3XHN1
A	-2	PHE	-	expression tag	UNP A3XHN1
A	-1	GLN	-	expression tag	UNP A3XHN1
A	0	SER	-	expression tag	UNP A3XHN1
A	1	ASN	-	expression tag	UNP A3XHN1

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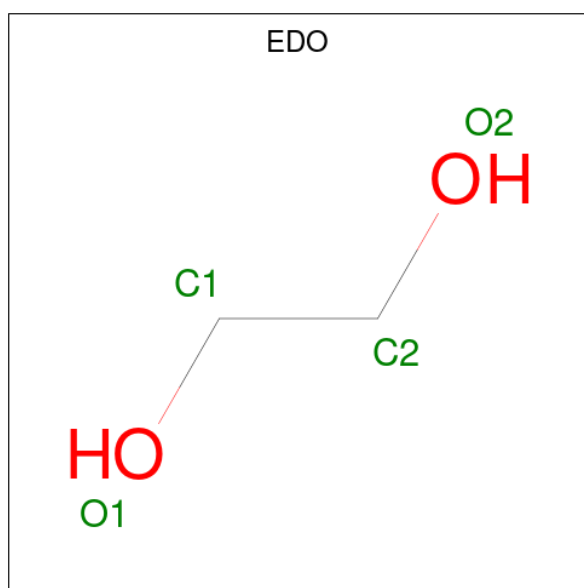
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	ALA	-	expression tag	UNP A3XHN1
A	?	-	PRO	deletion	UNP A3XHN1
A	?	-	LEU	deletion	UNP A3XHN1
A	?	-	SER	deletion	UNP A3XHN1
A	?	-	LYS	deletion	UNP A3XHN1
B	-21	MET	-	expression tag	UNP A3XHN1
B	-20	HIS	-	expression tag	UNP A3XHN1
B	-19	HIS	-	expression tag	UNP A3XHN1
B	-18	HIS	-	expression tag	UNP A3XHN1
B	-17	HIS	-	expression tag	UNP A3XHN1
B	-16	HIS	-	expression tag	UNP A3XHN1
B	-15	HIS	-	expression tag	UNP A3XHN1
B	-14	SER	-	expression tag	UNP A3XHN1
B	-13	SER	-	expression tag	UNP A3XHN1
B	-12	GLY	-	expression tag	UNP A3XHN1
B	-11	VAL	-	expression tag	UNP A3XHN1
B	-10	ASP	-	expression tag	UNP A3XHN1
B	-9	LEU	-	expression tag	UNP A3XHN1
B	-8	GLY	-	expression tag	UNP A3XHN1
B	-7	THR	-	expression tag	UNP A3XHN1
B	-6	GLU	-	expression tag	UNP A3XHN1
B	-5	ASN	-	expression tag	UNP A3XHN1
B	-4	LEU	-	expression tag	UNP A3XHN1
B	-3	TYR	-	expression tag	UNP A3XHN1
B	-2	PHE	-	expression tag	UNP A3XHN1
B	-1	GLN	-	expression tag	UNP A3XHN1
B	0	SER	-	expression tag	UNP A3XHN1
B	1	ASN	-	expression tag	UNP A3XHN1
B	2	ALA	-	expression tag	UNP A3XHN1
B	?	-	PRO	deletion	UNP A3XHN1
B	?	-	LEU	deletion	UNP A3XHN1
B	?	-	SER	deletion	UNP A3XHN1
B	?	-	LYS	deletion	UNP A3XHN1
C	-21	MET	-	expression tag	UNP A3XHN1
C	-20	HIS	-	expression tag	UNP A3XHN1
C	-19	HIS	-	expression tag	UNP A3XHN1
C	-18	HIS	-	expression tag	UNP A3XHN1
C	-17	HIS	-	expression tag	UNP A3XHN1
C	-16	HIS	-	expression tag	UNP A3XHN1
C	-15	HIS	-	expression tag	UNP A3XHN1
C	-14	SER	-	expression tag	UNP A3XHN1
C	-13	SER	-	expression tag	UNP A3XHN1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-12	GLY	-	expression tag	UNP A3XHN1
C	-11	VAL	-	expression tag	UNP A3XHN1
C	-10	ASP	-	expression tag	UNP A3XHN1
C	-9	LEU	-	expression tag	UNP A3XHN1
C	-8	GLY	-	expression tag	UNP A3XHN1
C	-7	THR	-	expression tag	UNP A3XHN1
C	-6	GLU	-	expression tag	UNP A3XHN1
C	-5	ASN	-	expression tag	UNP A3XHN1
C	-4	LEU	-	expression tag	UNP A3XHN1
C	-3	TYR	-	expression tag	UNP A3XHN1
C	-2	PHE	-	expression tag	UNP A3XHN1
C	-1	GLN	-	expression tag	UNP A3XHN1
C	0	SER	-	expression tag	UNP A3XHN1
C	1	ASN	-	expression tag	UNP A3XHN1
C	2	ALA	-	expression tag	UNP A3XHN1
C	?	-	PRO	deletion	UNP A3XHN1
C	?	-	LEU	deletion	UNP A3XHN1
C	?	-	SER	deletion	UNP A3XHN1
C	?	-	LYS	deletion	UNP A3XHN1

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mn	0	0
			1	1		
4	B	1	Total	Mn	0	0
			1	1		
4	C	1	Total	Mn	0	0
			1	1		

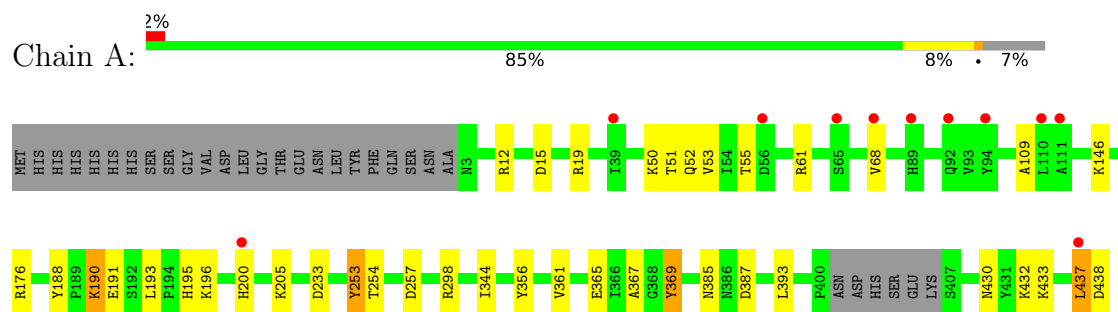
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	62	Total 62	O 62	0	0
5	B	60	Total 60	O 60	0	0
5	C	38	Total 38	O 38	0	0

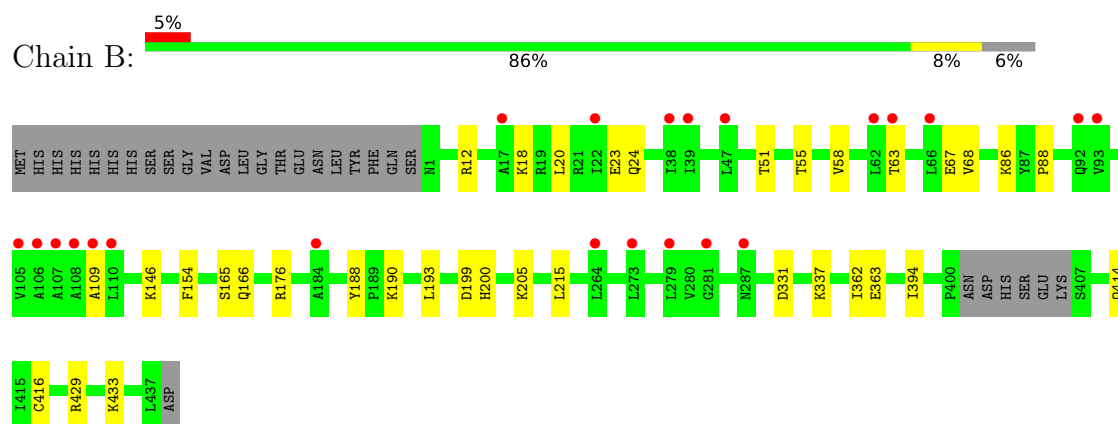
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

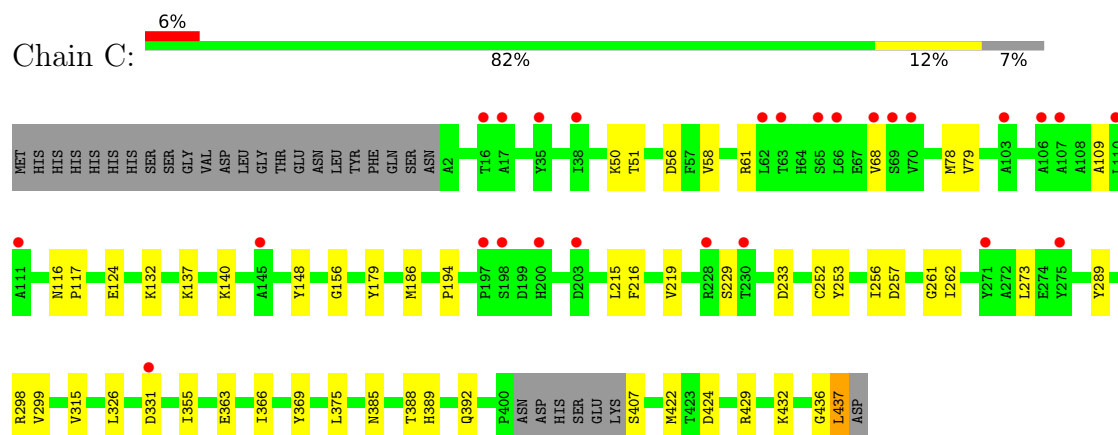
- Molecule 1: dGTP triphosphohydrolase



- Molecule 1: dGTP triphosphohydrolase



- Molecule 1: dGTP triphosphohydrolase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	181.40Å 181.40Å 110.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.56 – 2.26 40.56 – 2.26	Depositor EDS
% Data completeness (in resolution range)	99.2 (40.56-2.26) 94.6 (40.56-2.26)	Depositor EDS
$R_{merge}$	0.16	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.56 (at 2.27Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, $R_{free}$	0.198 , 0.228 0.197 , 0.226	Depositor DCC
$R_{free}$ test set	1405 reflections (1.63%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	50.7	Xtriage
Anisotropy	0.249	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 49.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10330	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MN, EDO, SO4

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	0.48	0/3451	0.63	0/4666
1	B	0.46	0/3441	0.60	0/4654
1	C	0.47	0/3425	0.60	0/4632
All	All	0.47	0/10317	0.61	0/13952

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	3388	0	3301	24	0
1	B	3379	0	3293	21	0
1	C	3362	0	3262	29	0
2	A	4	0	6	0	0
2	B	4	0	6	0	0
3	A	10	0	0	0	0
3	B	10	0	0	0	0
3	C	10	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	1	0	0	0	0
5	A	62	0	0	2	0
5	B	60	0	0	2	0
5	C	38	0	0	1	0
All	All	10330	0	9868	67	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 3.

All (67) 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:C:78:MET:HE2	1:C:315:VAL:HG21	1.63	0.81
1:C:375:LEU:HD13	1:C:422:MET:HE1	1.66	0.78
1:B:68:VAL:HG12	1:B:109:ALA:HB1	1.74	0.69
1:C:68:VAL:HG12	1:C:109:ALA:HB1	1.74	0.69
1:C:50:LYS:HB3	1:C:61:ARG:HD3	1.74	0.68
1:A:51:THR:HG21	1:A:55:THR:HG23	1.75	0.68
1:A:68:VAL:HG12	1:A:109:ALA:HB1	1.80	0.64
1:A:195:HIS:CE1	1:A:196:LYS:HD3	2.34	0.63
1:C:388:THR:O	1:C:392:GLN:HG3	2.01	0.60
1:C:432:LYS:HG2	1:C:437:LEU:HD12	1.87	0.57
1:A:437:LEU:O	1:A:438:ASP:C	2.43	0.57
1:A:253:TYR:O	1:A:257:ASP:HB2	2.05	0.56
1:A:15:ASP:OD2	1:A:19:ARG:HD2	2.09	0.53
1:A:361:VAL:O	1:A:365:GLU:HG3	2.11	0.51
1:B:51:THR:HG21	1:B:55:THR:HG23	1.92	0.50
1:A:298:ARG:HD3	5:A:636:HOH:O	2.10	0.50
1:C:261:GLY:HA3	1:C:355:ILE:HD11	1.94	0.49
1:A:437:LEU:HD13	1:C:437:LEU:HD22	1.95	0.49
1:C:186:MET:HG2	1:C:216:PHE:CE1	2.48	0.49
1:C:140:LYS:HD2	1:C:148:TYR:CE2	2.47	0.49
1:C:331:ASP:OD1	1:C:331:ASP:N	2.41	0.49
1:C:298:ARG:HD2	5:C:627:HOH:O	2.14	0.48
1:B:24:GLN:HG2	5:B:626:HOH:O	2.15	0.47
1:B:205:LYS:NZ	5:B:603:HOH:O	2.48	0.47
1:A:257:ASP:OD2	1:A:356:TYR:OH	2.23	0.47
1:C:262:ILE:HG21	1:C:299:VAL:HG11	1.97	0.47
1:A:146:LYS:HG2	1:A:193:LEU:HB2	1.96	0.47
1:A:191:GLU:OE2	1:A:233:ASP:HA	2.16	0.46
1:C:289:TYR:CE1	1:C:298:ARG:HG2	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:LYS:HB2	1:A:190:LYS:HE2	1.53	0.46
1:B:63:THR:O	1:B:67:GLU:HG3	2.16	0.46
1:A:254:THR:HG21	1:A:344:ILE:HG23	1.98	0.46
1:B:20:LEU:HB2	1:B:23:GLU:HG3	1.98	0.45
1:B:166:GLN:HE22	1:B:414:GLN:HE22	1.62	0.45
1:C:132:LYS:O	1:C:137:LYS:HG3	2.17	0.45
1:C:215:LEU:HD12	1:C:215:LEU:O	2.16	0.45
1:B:188:TYR:CZ	1:B:205:LYS:HD3	2.51	0.45
1:B:190:LYS:HB2	1:B:190:LYS:HE2	1.85	0.44
1:A:369:TYR:HE1	1:C:389:HIS:HD2	1.64	0.44
1:B:146:LYS:HG2	1:B:193:LEU:HB2	1.98	0.44
1:A:367:ALA:HB1	1:C:436:GLY:HA2	1.99	0.44
1:B:86:LYS:O	1:B:88:PRO:HD3	2.18	0.44
1:A:430:ASN:ND2	5:A:601:HOH:O	2.38	0.44
1:B:18:LYS:HE3	1:B:18:LYS:HB3	1.92	0.43
1:B:331:ASP:OD1	1:B:331:ASP:N	2.44	0.43
1:C:117:PRO:HG3	1:C:124:GLU:HG3	2.00	0.43
1:B:394:ILE:HD13	1:C:366:ILE:HD13	2.01	0.43
1:B:12:ARG:HD3	1:B:176:ARG:O	2.18	0.43
1:C:179:TYR:CD2	1:C:219:VAL:HA	2.55	0.42
1:C:252:CYS:O	1:C:256:ILE:HB	2.20	0.42
1:A:393:LEU:HD22	1:B:362:ILE:HG23	2.02	0.42
1:B:433:LYS:HD2	1:C:363:GLU:HB3	2.01	0.42
1:C:194:PRO:HD3	1:C:233:ASP:HB3	2.02	0.42
1:C:253:TYR:O	1:C:257:ASP:HB2	2.19	0.41
1:A:432:LYS:HG2	1:A:437:LEU:HD21	2.01	0.41
1:A:188:TYR:CE2	1:A:205:LYS:HD2	2.55	0.41
1:C:79:VAL:HG23	1:C:315:VAL:HG22	2.01	0.41
1:A:433:LYS:HG3	1:B:363:GLU:HG2	2.01	0.41
1:A:12:ARG:HD3	1:A:176:ARG:O	2.20	0.41
1:B:429:ARG:O	1:B:433:LYS:HG2	2.21	0.41
1:C:116:ASN:OD1	1:C:156:GLY:HA3	2.21	0.41
1:C:273:LEU:HD23	1:C:273:LEU:HA	1.89	0.41
1:B:165:SER:HB2	1:B:215:LEU:HD21	2.02	0.40
1:A:50:LYS:HB3	1:A:61:ARG:HD3	2.03	0.40
1:B:154:PHE:CE1	1:B:416:CYS:HB3	2.56	0.40
1:C:56:ASP:C	1:C:58:VAL:H	2.25	0.40
1:A:387:ASP:OD1	1:A:387:ASP:N	2.43	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	426/460 (93%)	419 (98%)	6 (1%)	1 (0%)	47	55
1	B	427/460 (93%)	416 (97%)	11 (3%)	0	100	100
1	C	426/460 (93%)	416 (98%)	10 (2%)	0	100	100
All	All	1279/1380 (93%)	1251 (98%)	27 (2%)	1 (0%)	51	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	53	VAL

### 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	356/396 (90%)	349 (98%)	7 (2%)	55	64
1	B	353/396 (89%)	349 (99%)	4 (1%)	73	82
1	C	348/396 (88%)	339 (97%)	9 (3%)	46	55
All	All	1057/1188 (89%)	1037 (98%)	20 (2%)	57	66

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	GLN
1	A	190	LYS

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Mol	Chain	Res	Type
1	A	200	HIS
1	A	253	TYR
1	A	369	TYR
1	A	385	ASN
1	A	437	LEU
1	B	58	VAL
1	B	199	ASP
1	B	200	HIS
1	B	337	LYS
1	C	51	THR
1	C	229	SER
1	C	326	LEU
1	C	369	TYR
1	C	385	ASN
1	C	407	SER
1	C	424	ASP
1	C	429	ARG
1	C	437	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	ASN
1	A	92	GLN
1	A	293	GLN
1	B	221	GLN
1	B	332	GLN
1	B	357	ASN
1	B	414	GLN
1	C	392	GLN
1	C	414	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 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 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	SO4	C	501	-	4,4,4	0.21	0	6,6,6	0.52	0
2	EDO	B	501	-	3,3,3	0.12	0	2,2,2	0.13	0
3	SO4	C	502	-	4,4,4	0.19	0	6,6,6	0.23	0
3	SO4	B	503	-	4,4,4	0.45	0	6,6,6	0.05	0
3	SO4	A	503	-	4,4,4	0.45	0	6,6,6	0.06	0
3	SO4	B	502	-	4,4,4	0.50	0	6,6,6	0.06	0
3	SO4	A	502	-	4,4,4	0.58	0	6,6,6	0.08	0
2	EDO	A	501	-	3,3,3	0.25	0	2,2,2	0.25	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
2	EDO	A	501	-	-	1/1/1/1	-
2	EDO	B	501	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	A	501	EDO	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

## 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	430/460 (93%)	0.15	11 (2%) 56 59	42, 62, 91, 133	0
1	B	431/460 (93%)	0.31	22 (5%) 28 30	46, 65, 92, 148	0
1	C	430/460 (93%)	0.33	26 (6%) 21 23	50, 71, 99, 123	0
All	All	1291/1380 (93%)	0.26	59 (4%) 32 35	42, 66, 94, 148	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	16	THR	6.1
1	A	437	LEU	4.8
1	C	110	LEU	4.3
1	C	17	ALA	3.6
1	C	66	LEU	3.5
1	B	279	LEU	3.4
1	C	230	THR	3.4
1	C	68	VAL	3.3
1	B	273	LEU	3.2
1	B	93	VAL	3.0
1	C	70	VAL	3.0
1	B	92	GLN	2.9
1	B	109	ALA	2.8
1	B	287	ASN	2.8
1	B	39	ILE	2.8
1	C	145	ALA	2.8
1	A	39	ILE	2.8
1	C	228	ARG	2.7
1	C	271	TYR	2.7
1	A	200	HIS	2.7
1	C	69	SER	2.7
1	C	203	ASP	2.6
1	C	198	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	17	ALA	2.6
1	C	35	TYR	2.6
1	A	94	TYR	2.6
1	B	107	ALA	2.5
1	C	111	ALA	2.5
1	B	22	ILE	2.5
1	A	92	GLN	2.4
1	A	68	VAL	2.4
1	C	275	TYR	2.3
1	C	62	LEU	2.3
1	C	103	ALA	2.3
1	B	281	GLY	2.3
1	C	197	PRO	2.3
1	B	184	ALA	2.3
1	B	110	LEU	2.3
1	B	38	ILE	2.3
1	C	38	ILE	2.3
1	B	63	THR	2.2
1	A	110	LEU	2.2
1	B	105	VAL	2.2
1	B	62	LEU	2.2
1	B	264	LEU	2.2
1	A	56	ASP	2.2
1	A	89	HIS	2.2
1	C	106	ALA	2.2
1	C	107	ALA	2.2
1	B	66	LEU	2.2
1	A	111	ALA	2.1
1	B	47	LEU	2.1
1	C	200	HIS	2.1
1	B	106	ALA	2.1
1	B	108	ALA	2.1
1	C	331	ASP	2.0
1	A	65	SER	2.0
1	C	63	THR	2.0
1	C	65	SER	2.0

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

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [i](#)

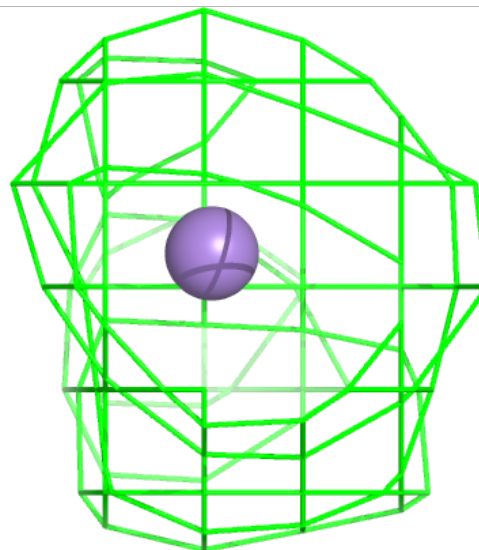
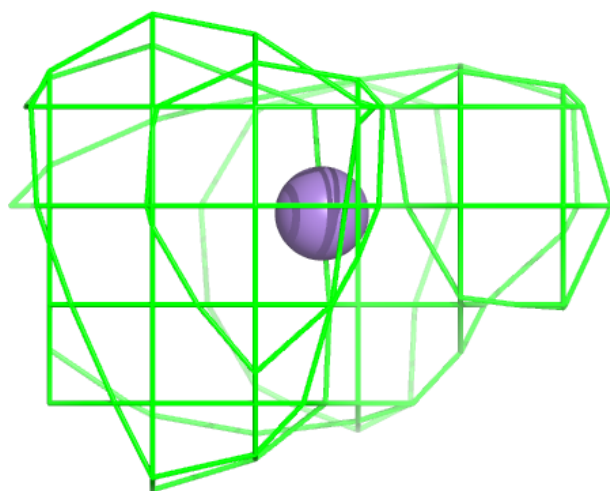
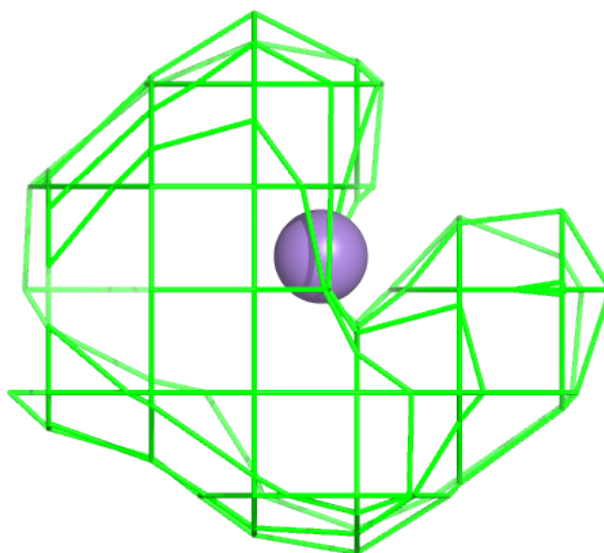
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	EDO	A	501	4/4	0.80	0.22	63,64,67,70	0
3	SO4	A	503	5/5	0.83	0.24	102,103,112,112	5
2	EDO	B	501	4/4	0.88	0.20	72,76,79,81	0
3	SO4	C	502	5/5	0.93	0.23	97,99,107,108	0
3	SO4	B	503	5/5	0.94	0.21	88,92,99,102	5
3	SO4	C	501	5/5	0.95	0.10	74,75,78,81	0
3	SO4	B	502	5/5	0.95	0.09	74,79,83,85	0
3	SO4	A	502	5/5	0.98	0.09	64,66,69,69	0
4	MN	B	504	1/1	0.99	0.17	52,52,52,52	0
4	MN	C	503	1/1	0.99	0.23	63,63,63,63	0
4	MN	A	504	1/1	1.00	0.14	44,44,44,44	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

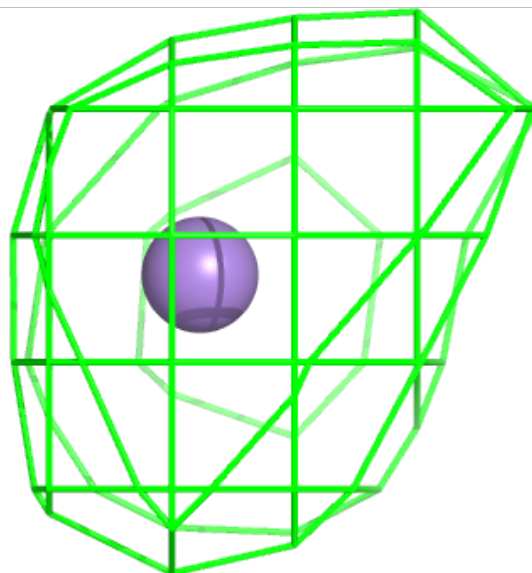
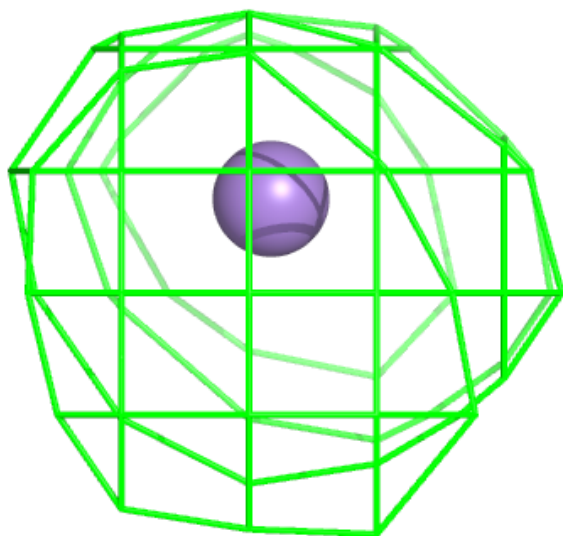
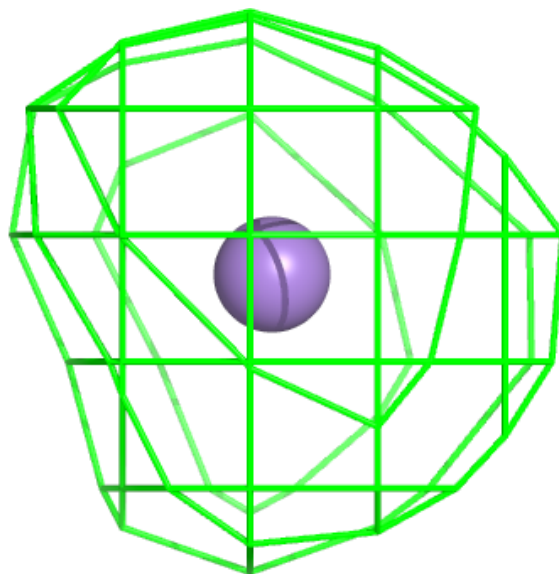
**Electron density around MN B 504:**

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



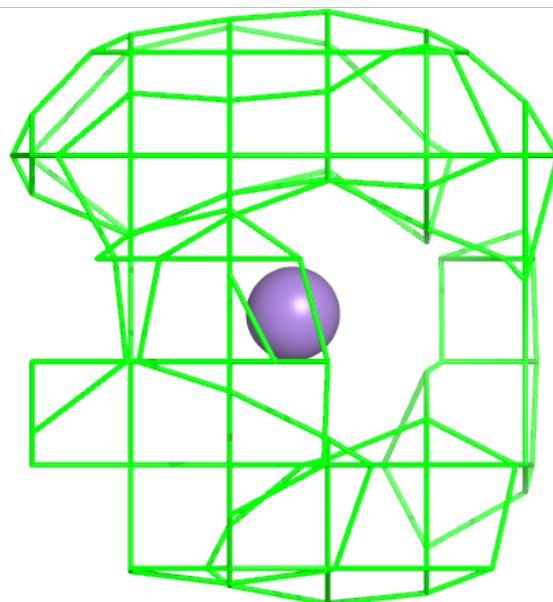
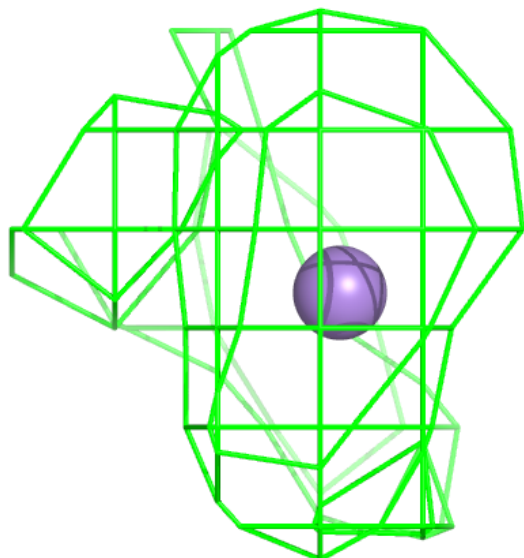
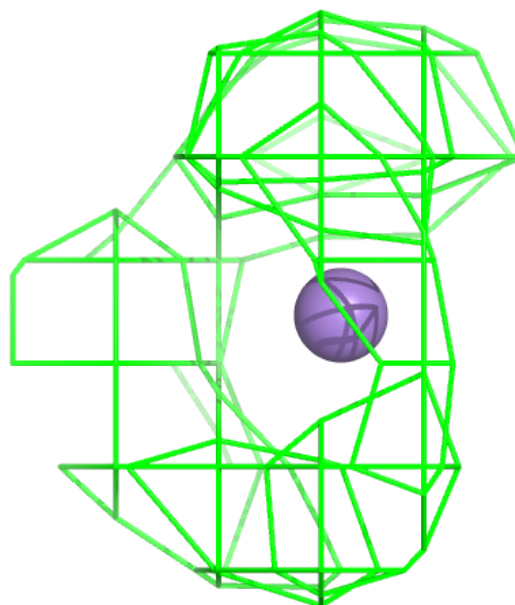
**Electron density around MN C 503:**

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



**Electron density around MN A 504:**

$2mF_o-DF_c$  (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.