



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 6, 2022 – 12:15 PM JST

PDB ID : 7CB3  
Title : Crystal structure of type III polyketide synthase from *Mycobacterium marinum*  
Authors : Pratap, S.; Kant, A.; Saxena, P.; Krishnan, V.  
Deposited on : 2020-06-10  
Resolution : 2.31 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

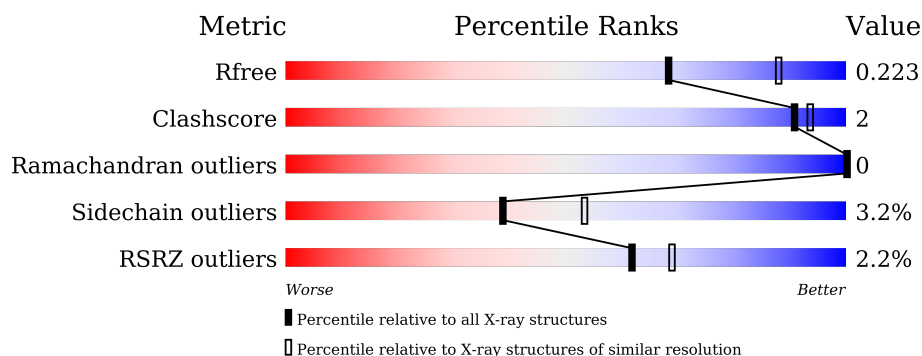
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.31 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	5974 (2.34-2.30)
Clashscore	141614	6604 (2.34-2.30)
Ramachandran outliers	138981	6523 (2.34-2.30)
Sidechain outliers	138945	6523 (2.34-2.30)
RSRZ outliers	127900	5855 (2.34-2.30)

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 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	413	<div> <div>2%</div> <div> <div></div> <div>83%</div> <div>•</div> <div>12%</div> </div> </div>
1	B	413	<div> <div>2%</div> <div> <div></div> <div>84%</div> <div>••</div> <div>12%</div> </div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5403 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 Chalcone/stilbene synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	363	Total	C	N	O	S	0	0	0
			2627	1663	452	497	15			
1	B	363	Total	C	N	O	S	0	0	0
			2636	1667	450	504	15			

There are 40 discrepancies between the modelled and reference sequences:

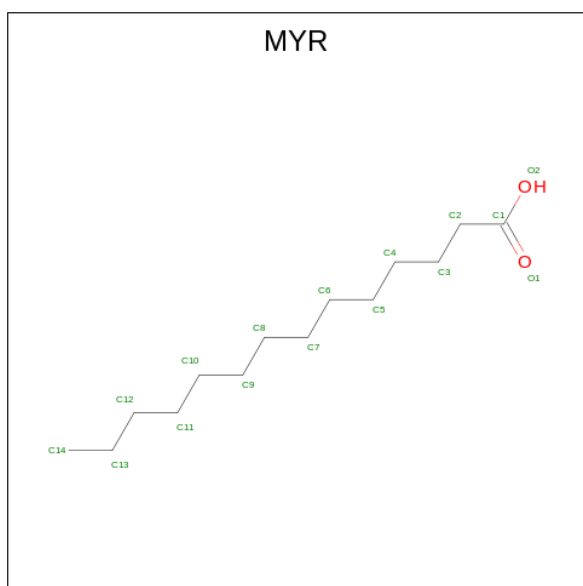
Chain	Residue	Modelled	Actual	Comment	Reference
A	394	GLY	-	expression tag	UNP B2HNE1
A	395	ASN	-	expression tag	UNP B2HNE1
A	396	SER	-	expression tag	UNP B2HNE1
A	397	SER	-	expression tag	UNP B2HNE1
A	398	SER	-	expression tag	UNP B2HNE1
A	399	VAL	-	expression tag	UNP B2HNE1
A	400	ASP	-	expression tag	UNP B2HNE1
A	401	LYS	-	expression tag	UNP B2HNE1
A	402	LEU	-	expression tag	UNP B2HNE1
A	403	ALA	-	expression tag	UNP B2HNE1
A	404	ALA	-	expression tag	UNP B2HNE1
A	405	ALA	-	expression tag	UNP B2HNE1
A	406	LEU	-	expression tag	UNP B2HNE1
A	407	GLU	-	expression tag	UNP B2HNE1
A	408	HIS	-	expression tag	UNP B2HNE1
A	409	HIS	-	expression tag	UNP B2HNE1
A	410	HIS	-	expression tag	UNP B2HNE1
A	411	HIS	-	expression tag	UNP B2HNE1
A	412	HIS	-	expression tag	UNP B2HNE1
A	413	HIS	-	expression tag	UNP B2HNE1
B	394	GLY	-	expression tag	UNP B2HNE1
B	395	ASN	-	expression tag	UNP B2HNE1
B	396	SER	-	expression tag	UNP B2HNE1
B	397	SER	-	expression tag	UNP B2HNE1
B	398	SER	-	expression tag	UNP B2HNE1

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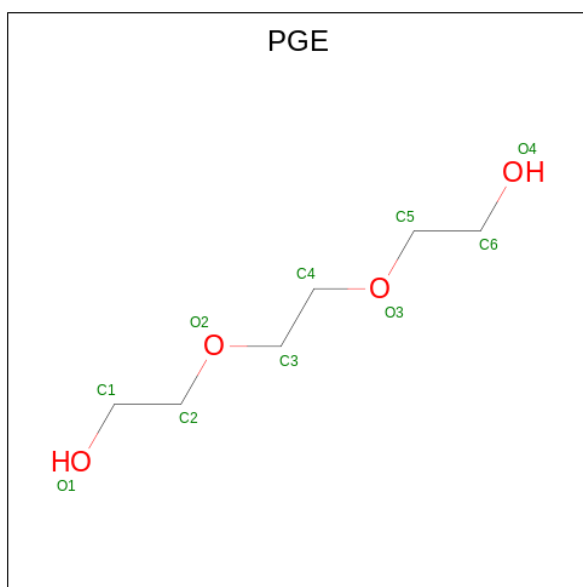
Chain	Residue	Modelled	Actual	Comment	Reference
B	399	VAL	-	expression tag	UNP B2HNE1
B	400	ASP	-	expression tag	UNP B2HNE1
B	401	LYS	-	expression tag	UNP B2HNE1
B	402	LEU	-	expression tag	UNP B2HNE1
B	403	ALA	-	expression tag	UNP B2HNE1
B	404	ALA	-	expression tag	UNP B2HNE1
B	405	ALA	-	expression tag	UNP B2HNE1
B	406	LEU	-	expression tag	UNP B2HNE1
B	407	GLU	-	expression tag	UNP B2HNE1
B	408	HIS	-	expression tag	UNP B2HNE1
B	409	HIS	-	expression tag	UNP B2HNE1
B	410	HIS	-	expression tag	UNP B2HNE1
B	411	HIS	-	expression tag	UNP B2HNE1
B	412	HIS	-	expression tag	UNP B2HNE1
B	413	HIS	-	expression tag	UNP B2HNE1

- Molecule 2 is MYRISTIC ACID (three-letter code: MYR) (formula:  $C_{14}H_{28}O_2$ ) (labeled as "Ligand of Interest" by depositor).



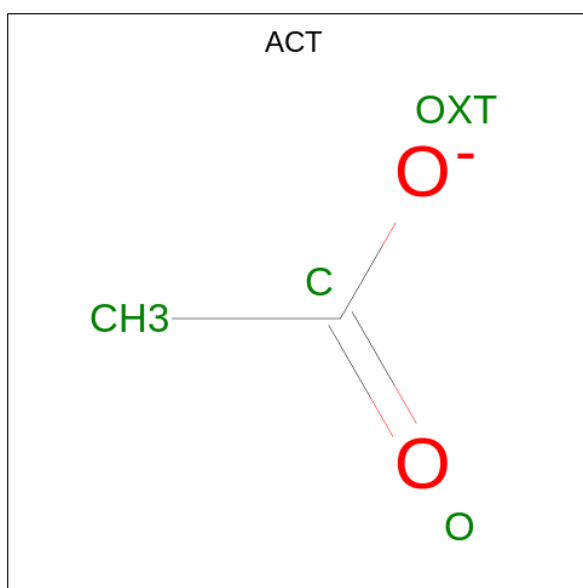
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			16	14	2		
2	B	1	Total	C	O	0	0
			16	14	2		

- Molecule 3 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula:  $C_6H_{14}O_4$ ).



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

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula:  $C_2H_3O_2$ ).



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

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	37	Total	O	0	0
			37	37		
5	B	39	Total	O	0	0
			39	39		



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	131.41Å 131.41Å 187.19Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	72.29 – 2.31 97.24 – 2.31	Depositor EDS
% Data completeness (in resolution range)	53.0 (72.29-2.31) 53.1 (97.24-2.31)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.12 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, $R_{free}$	0.198 , 0.217 0.202 , 0.223	Depositor DCC
$R_{free}$ test set	1552 reflections (5.50%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	49.4	Xtriage
Anisotropy	0.091	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 37.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.024 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5403	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MYR, MCS, ACT, PGE

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	0.58	0/2660	0.76	0/3634
1	B	0.56	0/2669	0.76	0/3647
All	All	0.57	0/5329	0.76	0/7281

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

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	2627	0	2540	11	0
1	B	2636	0	2554	9	0
2	A	16	0	27	0	0
2	B	16	0	27	0	0
3	A	10	0	14	0	0
3	B	10	0	14	0	0
4	A	4	0	3	0	0
4	B	8	0	6	0	0
5	A	37	0	0	0	0
5	B	39	0	0	0	0
All	All	5403	0	5185	18	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

All (18) 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:B:371:THR:HB	1:B:389:ASP:OD1	1.91	0.71
1:B:220:ILE:HG13	1:B:224:PHE:HE2	1.62	0.64
1:B:175:MCS:CAG	1:B:346:ASN:HD21	2.13	0.61
1:A:175:MCS:CAG	1:A:346:ASN:HD21	2.14	0.60
1:B:175:MCS:HAG1	1:B:346:ASN:HD21	1.67	0.58
1:A:175:MCS:HAG1	1:A:346:ASN:HD21	1.69	0.58
1:A:128:LEU:HD22	1:A:129:PRO:HD2	1.85	0.57
1:A:183:ARG:NH2	1:A:384:GLU:OE1	2.36	0.54
1:A:128:LEU:CD2	1:A:129:PRO:HD2	2.39	0.52
1:B:133:ASP:N	1:B:133:ASP:OD1	2.43	0.51
1:A:133:ASP:N	1:A:133:ASP:OD1	2.43	0.50
1:A:148:ALA:HB3	1:B:266:LEU:HB2	1.93	0.50
1:A:266:LEU:HB2	1:B:148:ALA:HB3	1.98	0.46
1:A:297:LEU:O	1:A:301:GLY:N	2.52	0.43
1:A:125:LEU:O	1:A:128:LEU:HB2	2.19	0.43
1:B:297:LEU:O	1:B:301:GLY:N	2.52	0.42
1:A:128:LEU:HD23	1:A:128:LEU:HA	1.93	0.42
1:B:176:ALA:HB1	1:B:384:GLU:HB2	2.03	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	360/413 (87%)	349 (97%)	11 (3%)	0	100	100
1	B	360/413 (87%)	350 (97%)	10 (3%)	0	100	100
All	All	720/826 (87%)	699 (97%)	21 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	263/334 (79%)	254 (97%)	9 (3%)	37	51
1	B	268/334 (80%)	260 (97%)	8 (3%)	41	56
All	All	531/668 (80%)	514 (97%)	17 (3%)	39	53

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	133	ASP
1	A	143	SER
1	A	220	ILE
1	A	250	ARG
1	A	251	SER
1	A	252	SER
1	A	274	THR
1	A	370	SER
1	A	371	THR
1	B	133	ASP
1	B	143	SER
1	B	220	ILE
1	B	250	ARG
1	B	251	SER
1	B	274	THR
1	B	370	SER
1	B	371	THR

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

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\# Z  > 2$
1	MCS	B	175	1	6,11,12	3.03	2 (33%)	4,13,15	6.17	4 (100%)
1	MCS	A	175	1	6,11,12	3.00	3 (50%)	4,13,15	5.62	3 (75%)

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
1	MCS	B	175	1	-	3/6/11/13	-
1	MCS	A	175	1	-	3/6/11/13	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	175	MCS	CAK-SAI	-6.49	1.60	1.76
1	A	175	MCS	CAK-SAI	-5.73	1.62	1.76
1	A	175	MCS	CB-SAI	-3.41	1.73	1.81
1	B	175	MCS	CB-SAI	-3.08	1.74	1.81
1	A	175	MCS	OAC-CAK	2.54	1.25	1.21

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	175	MCS	CB-SAI-CAK	10.70	115.80	100.84
1	A	175	MCS	CB-SAI-CAK	10.00	114.83	100.84
1	B	175	MCS	OAC-CAK-SAI	-5.22	115.84	122.61
1	A	175	MCS	OAC-CAK-SAI	-4.31	117.02	122.61
1	B	175	MCS	OAC-CAK-CAG	2.44	127.57	123.35
1	A	175	MCS	OAC-CAK-CAG	2.39	127.50	123.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	175	MCS	CAG-CAK-SAI	2.16	116.39	113.69

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	175	MCS	CA-CB-SAI-CAK
1	A	175	MCS	OAC-CAK-SAI-CB
1	A	175	MCS	CAG-CAK-SAI-CB
1	B	175	MCS	CA-CB-SAI-CAK
1	B	175	MCS	OAC-CAK-SAI-CB
1	B	175	MCS	CAG-CAK-SAI-CB

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	175	MCS	2	0
1	A	175	MCS	2	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

7 ligands 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	$\# Z  > 2$
4	ACT	B	503	-	1,3,3	1.30	0	0,3,3	-	-
2	MYR	A	501	-	12,15,15	0.25	0	11,15,15	0.55	0
4	ACT	B	504	-	1,3,3	1.69	0	0,3,3	-	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	MYR	B	501	-	12,15,15	0.14	0	11,15,15	0.84	0
4	ACT	A	503	-	1,3,3	1.36	0	0,3,3	-	-
3	PGE	B	502	-	9,9,9	0.52	0	8,8,8	0.44	0
3	PGE	A	502	-	9,9,9	0.55	0	8,8,8	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
3	PGE	B	502	-	-	4/7/7/7	-
2	MYR	B	501	-	-	2/11/13/13	-
3	PGE	A	502	-	-	3/7/7/7	-
2	MYR	A	501	-	-	4/11/13/13	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

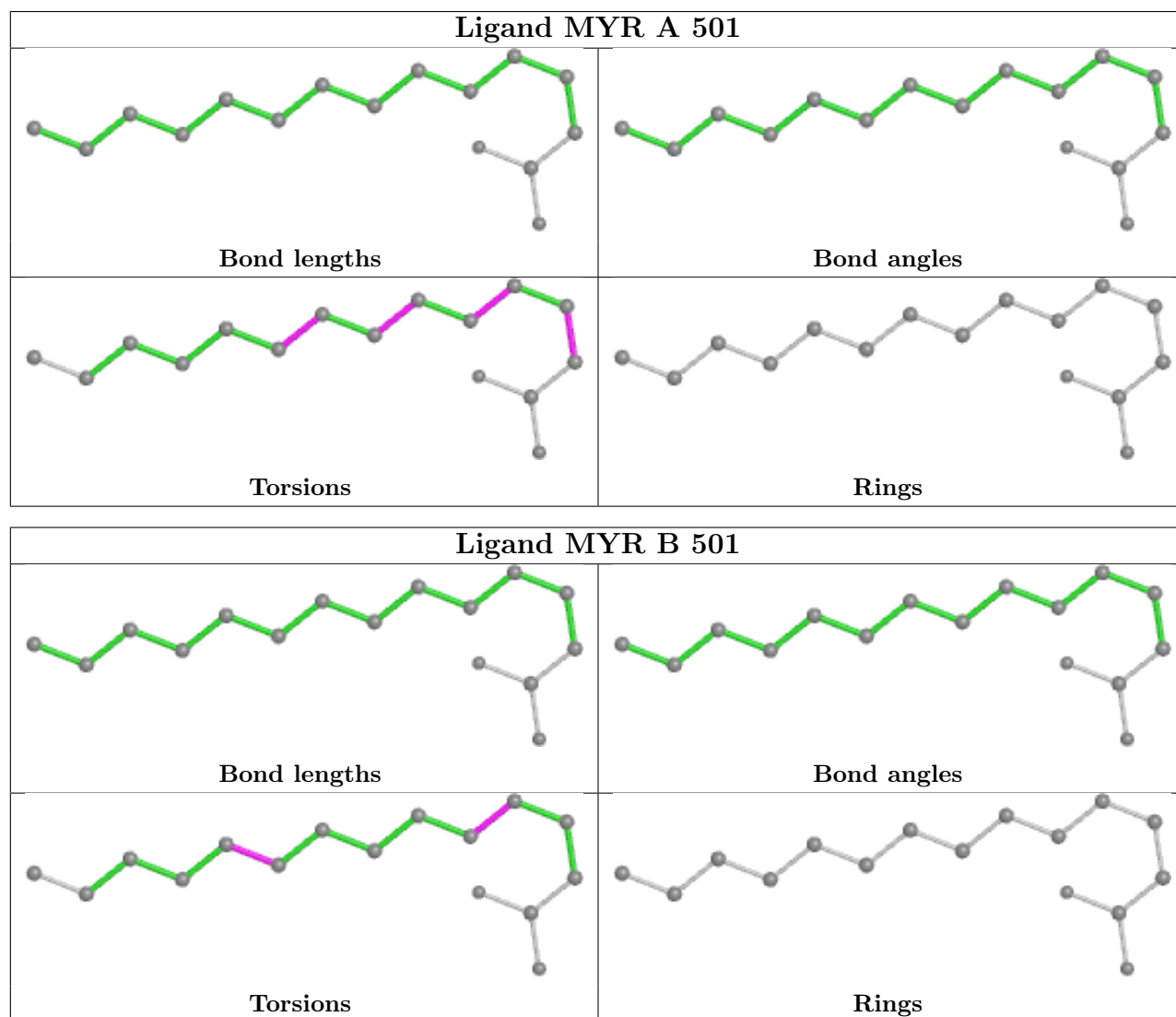
Mol	Chain	Res	Type	Atoms
2	A	501	MYR	C1-C2-C3-C4
3	A	502	PGE	O2-C3-C4-O3
2	A	501	MYR	C7-C8-C9-C10
3	A	502	PGE	O1-C1-C2-O2
3	B	502	PGE	O1-C1-C2-O2
2	A	501	MYR	C5-C6-C7-C8
3	A	502	PGE	O3-C5-C6-O4
3	B	502	PGE	O3-C5-C6-O4
2	A	501	MYR	C3-C4-C5-C6
3	B	502	PGE	C3-C4-O3-C5
2	B	501	MYR	C3-C4-C5-C6
2	B	501	MYR	C11-C10-C9-C8
3	B	502	PGE	O2-C3-C4-O3

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	362/413 (87%)	0.30	9 (2%) 57 64	28, 52, 108, 135	0
1	B	362/413 (87%)	0.28	7 (1%) 66 73	28, 53, 109, 134	0
All	All	724/826 (87%)	0.29	16 (2%) 62 69	28, 52, 109, 135	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	366	ASP	6.5
1	B	60	PHE	4.7
1	A	130	TYR	4.0
1	B	211	PHE	3.4
1	A	211	PHE	3.3
1	B	366	ASP	3.0
1	B	241	PRO	2.5
1	A	88	LEU	2.5
1	B	63	PRO	2.5
1	A	241	PRO	2.4
1	B	88	LEU	2.3
1	A	239	GLN	2.3
1	A	302	LEU	2.3
1	A	238	VAL	2.1
1	A	95	PHE	2.1
1	B	239	GLN	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [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
1	MCS	B	175	12/13	0.96	0.17	29,37,51,52	0
1	MCS	A	175	12/13	0.97	0.17	29,36,49,51	0

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [i](#)

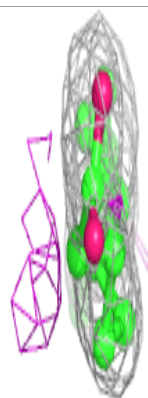
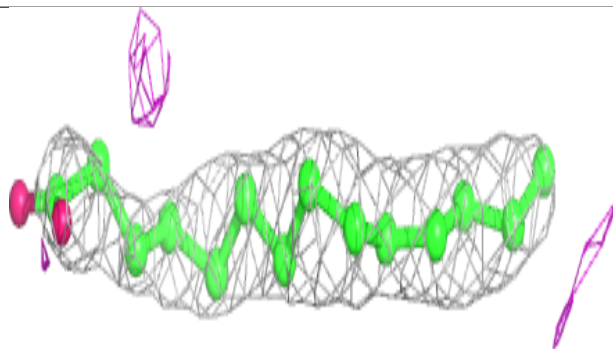
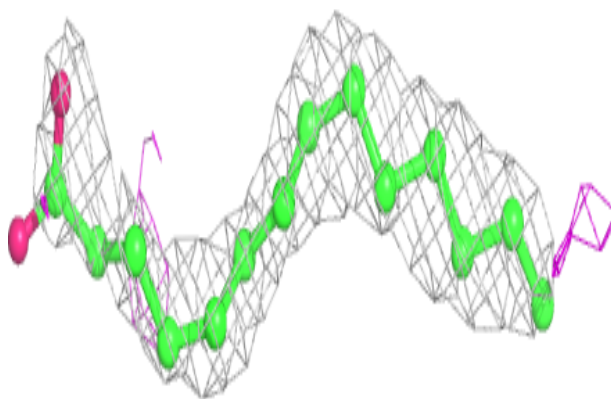
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
3	PGE	A	502	10/10	0.71	0.18	69,88,90,92	0
3	PGE	B	502	10/10	0.77	0.23	70,82,87,88	0
4	ACT	A	503	4/4	0.84	0.18	62,63,68,70	0
4	ACT	B	504	4/4	0.87	0.14	66,67,71,74	0
2	MYR	B	501	16/16	0.90	0.38	56,68,98,100	0
4	ACT	B	503	4/4	0.91	0.14	52,53,56,58	0
2	MYR	A	501	16/16	0.92	0.23	48,61,79,80	0

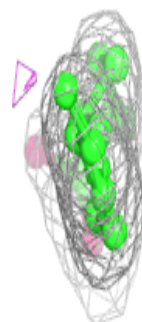
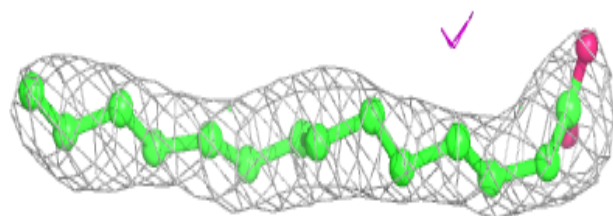
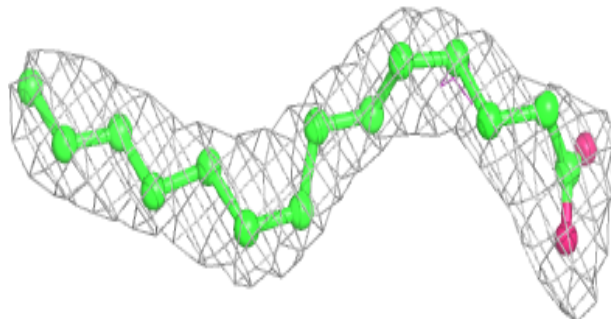
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around MYR B 501:**

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

**Electron density around MYR A 501:**

$2mF_o-DF_c$  (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.