



# Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 07:18 pm BST

PDB ID : 5WGC  
Title : propionyl-DpsC in complex with oxetane-bearing probe  
Authors : Milligan, J.C.; Ellis, B.D.; White, A.R.; Vanderwal, C.D.; Tsai, S.C.  
Deposited on : 2017-07-13  
Resolution : 2.15 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 : 2.11  
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) : 2.11

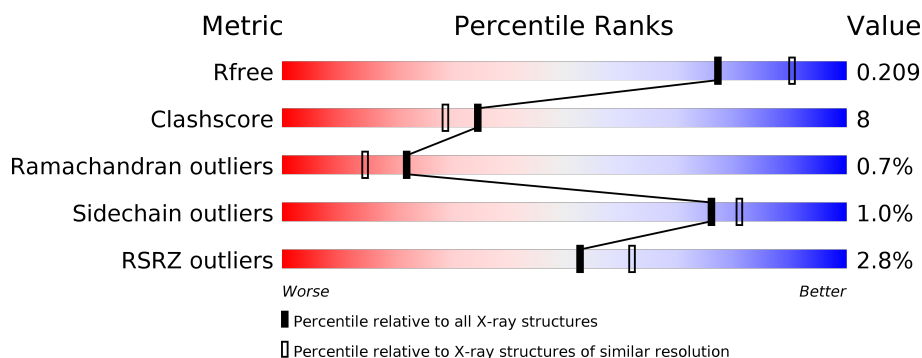
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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 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	356	<div> <div>2%</div> <div> <div></div> <div>86%</div> <div>10%</div> <div>••</div> </div> </div>
1	B	356	<div> <div>3%</div> <div> <div></div> <div>84%</div> <div>11%</div> <div>•••</div> </div> </div>

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	42Y	B	118	-	-	X	-
2	AFY	B	401	-	-	X	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5379 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.

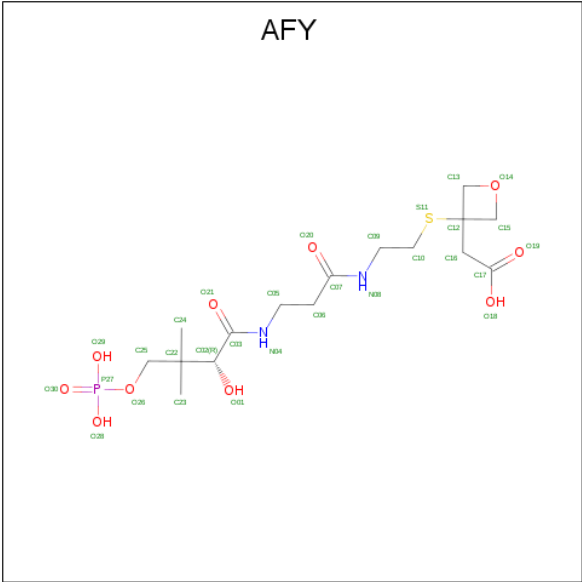
- Molecule 1 is a protein called Daunorubicin-doxorubicin polyketide synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	342	Total	C	N	O	S	0	0	0
			2524	1586	463	463	12			
1	B	343	Total	C	N	O	S	0	0	0
			2549	1600	469	468	12			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q54816
A	-1	SER	-	expression tag	UNP Q54816
A	0	HIS	-	expression tag	UNP Q54816
B	-2	GLY	-	expression tag	UNP Q54816
B	-1	SER	-	expression tag	UNP Q54816
B	0	HIS	-	expression tag	UNP Q54816

- Molecule 2 is (3-{[2-({N-[(2R)-2-hydroxy-3,3-dimethyl-4-(phosphonoxy)butanoyl]-bet a-alanyl}amino)ethyl]sulfanyl}oxetan-3-yl)acetic acid (three-letter code: AFY) (formula: C<sub>16</sub>H<sub>29</sub>N<sub>2</sub>O<sub>10</sub>PS).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			30	16	2	10	1	1		
2	B	1	Total	C	N	O	P	S	0	0
			30	16	2	10	1	1		

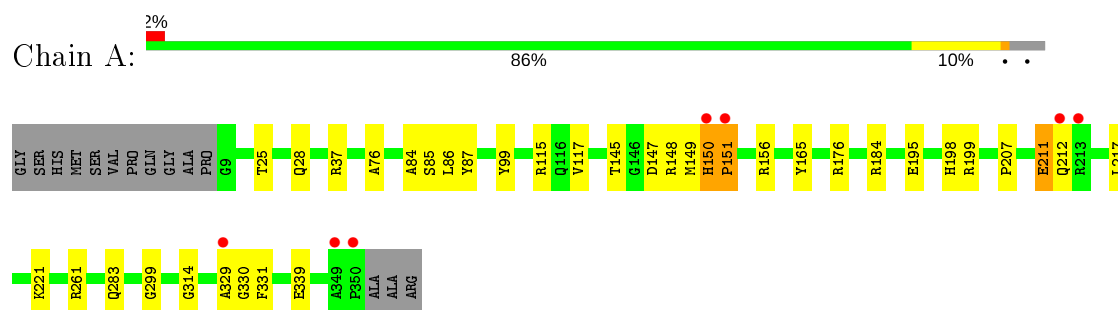
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	109	Total	O	0	0
			109	109		
3	B	137	Total	O	0	0
			137	137		

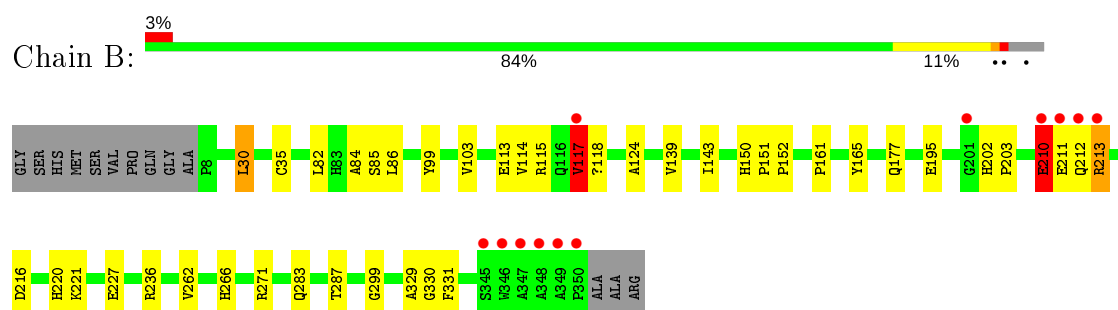
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Daunorubicin-doxorubicin polyketide synthase



- Molecule 1: Daunorubicin-doxorubicin polyketide synthase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	91.62Å 91.62Å 316.06Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	70.91 – 2.15 70.91 – 2.15	Depositor EDS
% Data completeness (in resolution range)	98.5 (70.91-2.15) 98.5 (70.91-2.15)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.52 (at 2.14Å)	Xtriage
Refinement program	PHENIX (1.10pre_2124: ???)	Depositor
R, $R_{free}$	0.178 , 0.213 0.181 , 0.209	Depositor DCC
$R_{free}$ test set	2000 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.1	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 59.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5379	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.60% 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: 42Y, AFY

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	0.49	0/2579	0.67	2/3519 (0.1%)
1	B	0.48	0/2605	0.64	0/3552
All	All	0.49	0/5184	0.65	2/7071 (0.0%)

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	B	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	330	GLY	N-CA-C	5.94	127.95	113.10
1	A	151	PRO	C-N-CD	5.32	139.58	128.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	117	VAL	Mainchain

### 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	2524	0	2444	36	0
1	B	2549	0	2480	51	0
2	A	30	0	0	2	0
2	B	30	0	0	9	0
3	A	109	0	0	8	0
3	B	137	0	0	7	0
All	All	5379	0	4924	83	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 8.

All (83) 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:118:42Y:H10	2:B:401:AFY:S11	1.91	1.10
2:A:401:AFY:O20	3:A:501:HOH:O	1.91	0.89
1:A:299:GLY:O	3:A:502:HOH:O	1.91	0.87
1:B:213:ARG:HH12	1:B:220:HIS:CE1	1.98	0.81
1:B:211:GLU:OE1	1:B:213:ARG:N	2.14	0.80
1:A:329:ALA:O	3:A:503:HOH:O	2.05	0.74
1:B:211:GLU:OE1	1:B:213:ARG:HG3	1.86	0.74
1:B:118:42Y:C1	2:B:401:AFY:S11	2.73	0.73
1:B:117:VAL:HG12	1:B:118:42Y:H1	1.53	0.73
1:B:227:GLU:OE2	3:B:501:HOH:O	2.06	0.72
1:B:85:SER:O	1:B:115:ARG:HG2	1.91	0.71
1:B:211:GLU:OE2	1:B:213:ARG:NE	2.25	0.68
1:B:266:HIS:NE2	3:B:505:HOH:O	2.26	0.68
1:B:299:GLY:O	3:B:502:HOH:O	2.12	0.66
1:B:216:ASP:OD1	3:B:503:HOH:O	2.16	0.62
1:B:221:LYS:HZ3	2:B:401:AFY:C06	2.14	0.60
1:A:148:ARG:NH1	1:A:150:HIS:CE1	2.70	0.60
1:B:118:42Y:H10	2:B:401:AFY:C16	2.33	0.59
1:B:117:VAL:HG21	1:B:329:ALA:O	2.04	0.58
1:A:99:TYR:HA	1:B:195:GLU:HG3	1.86	0.56
1:A:86:LEU:HD21	1:A:299:GLY:HA2	1.88	0.56
2:B:401:AFY:C15	2:B:401:AFY:C09	2.84	0.56
1:A:195:GLU:HG3	1:B:99:TYR:HA	1.88	0.56
1:A:25:THR:OG1	1:A:28:GLN:HG3	2.07	0.55
1:A:147:ASP:HB3	1:A:149:MET:CE	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:GLU:OE2	1:B:213:ARG:CZ	2.56	0.54
1:B:152:PRO:HG2	3:B:532:HOH:O	2.08	0.54
1:B:330:GLY:N	1:B:331:PHE:HA	2.24	0.53
1:A:329:ALA:O	1:A:331:PHE:CD2	2.63	0.52
1:B:117:VAL:HG12	1:B:118:42Y:OAC	2.09	0.52
1:A:115:ARG:HB2	1:B:113:GLU:HB2	1.91	0.52
1:B:124:ALA:HB1	1:B:143:ILE:HD12	1.92	0.51
1:A:176:ARG:NH2	3:A:505:HOH:O	2.22	0.51
1:B:117:VAL:HG11	1:B:329:ALA:HB3	1.92	0.51
1:A:221:LYS:NZ	2:A:401:AFY:S11	2.82	0.51
1:B:283:GLN:HG3	3:B:523:HOH:O	2.11	0.51
1:B:210:GLU:OE1	1:B:212:GLN:NE2	2.36	0.51
1:B:211:GLU:OE1	1:B:213:ARG:CG	2.59	0.50
1:A:211:GLU:HG2	3:B:603:HOH:O	2.13	0.49
1:B:118:42Y:C1A	2:B:401:AFY:C16	2.90	0.49
1:B:82:LEU:HB2	1:B:143:ILE:HD13	1.94	0.49
1:A:87:TYR:HB3	1:A:149:MET:HE1	1.94	0.49
1:A:198:HIS:HD2	1:A:217:LEU:HD13	1.78	0.48
1:A:76:ALA:O	3:A:504:HOH:O	2.20	0.48
1:A:99:TYR:CA	1:B:195:GLU:HG3	2.44	0.48
1:A:261:ARG:NH1	3:A:508:HOH:O	2.40	0.48
1:B:82:LEU:HB2	1:B:143:ILE:CD1	2.45	0.47
1:B:84:ALA:HA	1:B:114:VAL:O	2.16	0.46
1:B:117:VAL:HG12	1:B:118:42Y:C1A	2.45	0.46
1:B:30:LEU:HD12	1:B:35:CYS:O	2.16	0.45
1:B:221:LYS:NZ	2:B:401:AFY:C06	2.80	0.45
1:A:148:ARG:NH2	1:A:150:HIS:CD2	2.85	0.45
1:A:156:ARG:HG3	1:A:165:TYR:O	2.16	0.45
1:B:236:ARG:HD3	1:B:236:ARG:HA	1.75	0.44
1:B:139:VAL:HG13	1:B:177:GLN:CG	2.48	0.44
1:B:262:VAL:O	1:B:287:THR:HA	2.17	0.44
1:B:221:LYS:NZ	2:B:401:AFY:C05	2.81	0.44
1:A:195:GLU:HG3	1:B:99:TYR:CA	2.47	0.44
1:B:117:VAL:CG1	1:B:118:42Y:OAC	2.66	0.43
1:A:85:SER:O	1:A:115:ARG:HG2	2.19	0.43
1:A:147:ASP:HB3	1:A:149:MET:HE3	1.99	0.43
1:A:87:TYR:CD2	1:A:149:MET:HE2	2.54	0.43
1:B:99:TYR:CZ	1:B:103:VAL:HG21	2.54	0.43
1:B:139:VAL:HG13	1:B:177:GLN:HG3	2.01	0.43
1:A:217:LEU:O	1:A:221:LYS:HG2	2.19	0.43
1:A:147:ASP:HB3	1:A:149:MET:HE1	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:HIS:CB	1:A:151:PRO:CD	2.97	0.42
1:A:150:HIS:CB	1:A:151:PRO:HD2	2.50	0.42
1:B:213:ARG:HH12	1:B:220:HIS:HE1	1.59	0.42
1:A:283:GLN:NE2	3:A:507:HOH:O	2.40	0.42
1:B:161:PRO:O	1:B:271:ARG:NH2	2.52	0.42
1:B:202:HIS:O	1:B:203:PRO:C	2.58	0.42
1:B:150:HIS:HA	1:B:151:PRO:HD3	1.81	0.41
1:B:221:LYS:HZ1	2:B:401:AFY:C05	2.33	0.41
1:A:150:HIS:HA	1:A:151:PRO:HD3	1.82	0.41
1:B:86:LEU:HD21	1:B:299:GLY:HA2	2.02	0.41
1:A:84:ALA:O	1:A:145:THR:HA	2.20	0.41
1:A:199:ARG:HD2	1:A:207:PRO:HB3	2.03	0.41
1:A:207:PRO:HD3	1:B:99:TYR:CE1	2.56	0.41
1:A:314:GLY:HA2	3:A:531:HOH:O	2.21	0.40
1:B:118:42Y:H8	1:B:165:TYR:OH	2.22	0.40
1:A:148:ARG:NH1	1:A:150:HIS:NE2	2.69	0.40
1:A:184:ARG:HD2	1:A:339:GLU:OE1	2.21	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	339/356 (95%)	323 (95%)	13 (4%)	3 (1%)	17	11
1	B	340/356 (96%)	326 (96%)	12 (4%)	2 (1%)	25	18
All	All	679/712 (95%)	649 (96%)	25 (4%)	5 (1%)	22	15

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	210	GLU

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Mol	Chain	Res	Type
1	A	212	GLN
1	A	117	VAL
1	A	211	GLU
1	B	117	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	247/262 (94%)	245 (99%)	2 (1%)	81	86
1	B	252/262 (96%)	249 (99%)	3 (1%)	71	76
All	All	499/524 (95%)	494 (99%)	5 (1%)	76	81

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ARG
1	A	150	HIS
1	B	30	LEU
1	B	210	GLU
1	B	213	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	220	HIS

### 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	42Y	B	118	1	8,9,10	1.94	2 (25%)	6,10,12	1.81	2 (33%)
1	42Y	A	118	1	8,9,10	2.01	2 (25%)	6,10,12	1.36	1 (16%)

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	42Y	B	118	1	-	2/7/9/11	-
1	42Y	A	118	1	-	0/7/9/11	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	118	42Y	OG-C1A	4.67	1.47	1.33
1	A	118	42Y	OG-C1A	4.63	1.46	1.33
1	A	118	42Y	OG-CB	-2.85	1.38	1.45
1	B	118	42Y	OG-CB	-2.29	1.39	1.45

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	118	42Y	OG-C1A-C2A	2.66	118.35	111.38
1	B	118	42Y	OAC-C1A-C2A	-2.22	113.21	123.51
1	A	118	42Y	OG-C1A-C2A	2.07	116.82	111.38

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	118	42Y	OAC-C1A-OG-CB
1	B	118	42Y	C2A-C1A-OG-CB

There are no ring outliers.

1 monomer is involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	118	42Y	9	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

2 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
2	AFY	B	401	-	21,30,30	2.05	4 (19%)	32,43,43	2.41	10 (31%)
2	AFY	A	401	-	21,30,30	1.93	5 (23%)	32,43,43	2.14	11 (34%)

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	AFY	B	401	-	-	12/30/44/44	0/1/1/1
2	AFY	A	401	-	-	16/30/44/44	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	AFY	C03-N04	5.75	1.46	1.33
2	B	401	AFY	C07-N08	5.11	1.45	1.33
2	A	401	AFY	C07-N08	4.96	1.44	1.33
2	A	401	AFY	C03-N04	4.81	1.44	1.33
2	B	401	AFY	O20-C07	-2.55	1.18	1.23
2	A	401	AFY	O21-C03	-2.50	1.18	1.23
2	B	401	AFY	O21-C03	-2.46	1.18	1.23
2	A	401	AFY	O20-C07	-2.34	1.18	1.23
2	A	401	AFY	O14-C15	-2.24	1.42	1.45

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	AFY	C05-C06-C07	8.76	126.94	112.36
2	A	401	AFY	O14-C13-C12	-6.10	89.03	91.28
2	A	401	AFY	C05-C06-C07	4.51	119.86	112.36
2	A	401	AFY	O14-C15-C12	-4.12	89.76	91.28
2	B	401	AFY	O26-P27-O30	4.03	117.78	106.47
2	B	401	AFY	C02-C03-N04	3.51	123.56	116.58
2	A	401	AFY	O26-C25-C22	-3.25	105.33	110.55
2	B	401	AFY	O14-C13-C12	-3.06	90.15	91.28
2	A	401	AFY	C16-C12-C13	-2.92	110.42	125.00
2	B	401	AFY	C16-C12-C15	-2.85	110.76	125.00
2	B	401	AFY	O20-C07-N08	-2.80	117.73	123.01
2	A	401	AFY	O01-C02-C22	2.71	116.63	110.25
2	B	401	AFY	C24-C22-C02	2.62	113.36	108.82
2	B	401	AFY	C05-N04-C03	-2.53	118.08	122.59
2	B	401	AFY	O21-C03-N04	-2.49	117.64	122.99
2	A	401	AFY	C16-C12-C15	-2.41	112.94	125.00
2	A	401	AFY	C06-C05-N04	-2.40	107.05	111.90
2	A	401	AFY	O26-P27-O30	2.29	112.91	106.47
2	A	401	AFY	O28-P27-O26	2.25	112.71	106.73
2	B	401	AFY	C16-C12-C13	-2.19	114.04	125.00
2	A	401	AFY	C15-O14-C13	-2.16	89.17	91.11

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	AFY	C09-C10-S11-C12
2	B	401	AFY	C13-C12-C16-C17
2	B	401	AFY	N04-C05-C06-C07

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Mol	Chain	Res	Type	Atoms
2	B	401	AFY	C22-C25-O26-P27
2	A	401	AFY	C13-C12-C16-C17
2	A	401	AFY	C15-C12-C16-C17
2	A	401	AFY	O01-C02-C22-C23
2	A	401	AFY	C03-C02-C22-C23
2	A	401	AFY	O01-C02-C22-C24
2	A	401	AFY	C03-C02-C22-C24
2	A	401	AFY	O01-C02-C22-C25
2	A	401	AFY	C03-C02-C22-C25
2	A	401	AFY	C25-O26-P27-O28
2	A	401	AFY	C25-O26-P27-O29
2	A	401	AFY	C22-C02-C03-N04
2	A	401	AFY	C22-C02-C03-O21
2	A	401	AFY	O01-C02-C03-N04
2	A	401	AFY	N04-C05-C06-C07
2	A	401	AFY	C06-C07-N08-C09
2	A	401	AFY	O20-C07-N08-C09
2	B	401	AFY	O01-C02-C03-O21
2	B	401	AFY	N08-C09-C10-S11
2	B	401	AFY	C22-C02-C03-N04
2	B	401	AFY	O01-C02-C03-N04
2	B	401	AFY	C23-C22-C25-O26
2	B	401	AFY	C02-C22-C25-O26
2	B	401	AFY	C24-C22-C25-O26
2	B	401	AFY	C22-C02-C03-O21

There are no ring outliers.

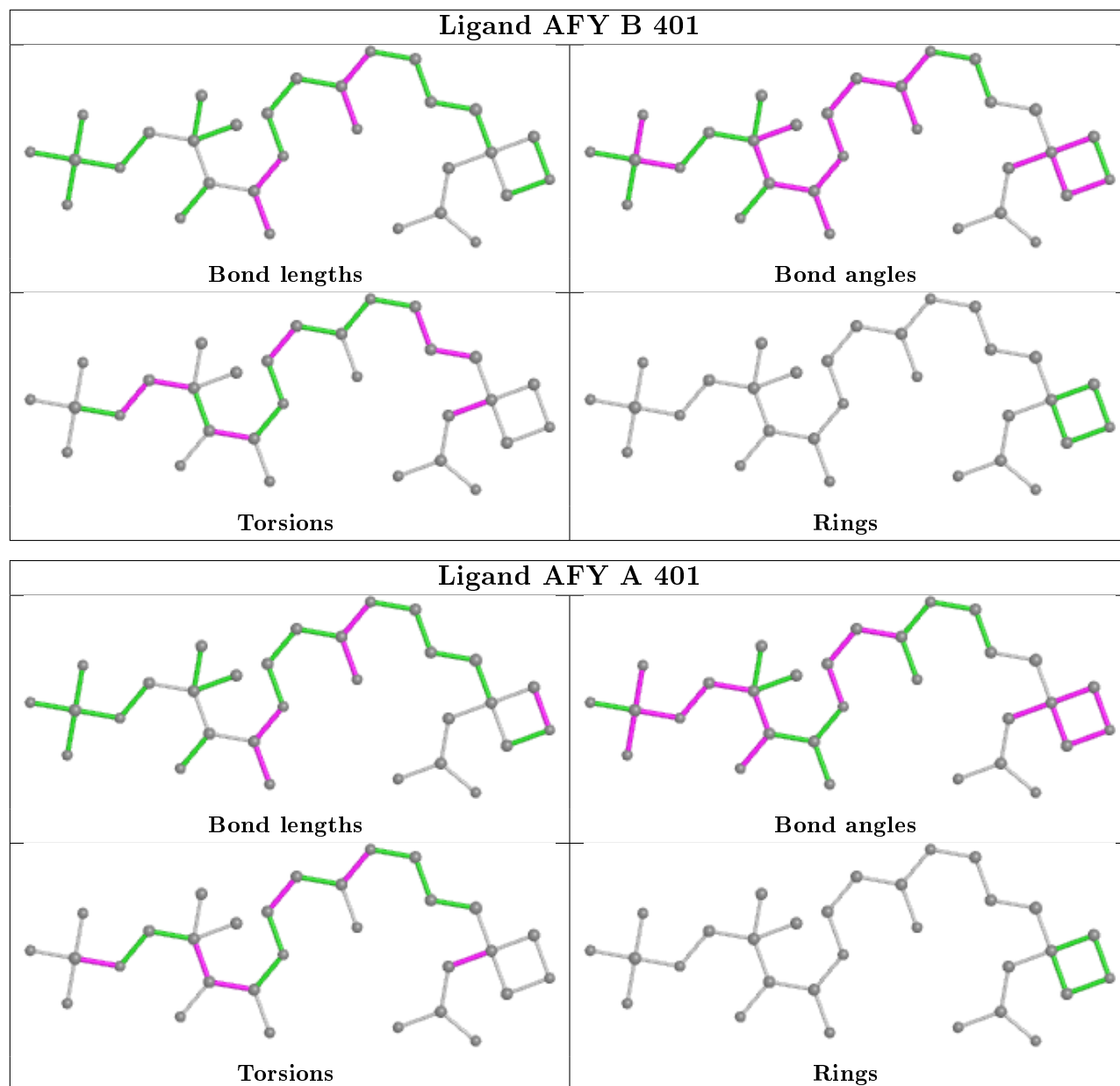
2 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	AFY	9	0
2	A	401	AFY	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and



any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



## 5.7 Other polymers ⓘ

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	341/356 (95%)	0.06	7 (2%) 63 71	23, 37, 63, 107	0
1	B	342/356 (96%)	0.09	12 (3%) 44 52	24, 35, 58, 99	0
All	All	683/712 (95%)	0.07	19 (2%) 53 62	23, 36, 63, 107	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	350	PRO	5.9
1	B	212	GLN	4.9
1	A	350	PRO	4.2
1	A	151	PRO	4.0
1	A	213	ARG	3.9
1	B	213	ARG	3.9
1	A	349	ALA	3.8
1	B	349	ALA	3.8
1	B	345	SER	3.0
1	B	210	GLU	2.9
1	A	150	HIS	2.9
1	B	347	ALA	2.7
1	A	212	GLN	2.5
1	A	329	ALA	2.5
1	B	211	GLU	2.4
1	B	117	VAL	2.4
1	B	201	GLY	2.3
1	B	346	TRP	2.3
1	B	348	ALA	2.2

### 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(Å <sup>2</sup> )	Q<0.9
1	42Y	B	118	10/11	0.78	0.24	29,31,34,35	0
1	42Y	A	118	10/11	0.82	0.23	24,30,36,37	0

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

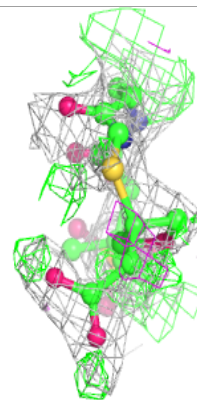
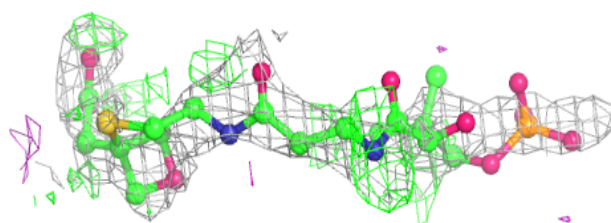
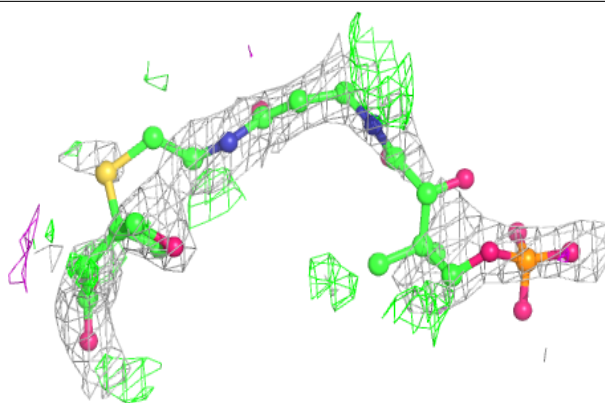
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	AFY	B	401	30/30	0.62	0.40	80,92,120,126	16
2	AFY	A	401	30/30	0.68	0.35	79,89,128,132	12

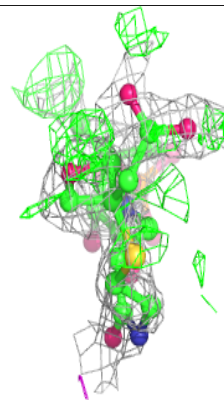
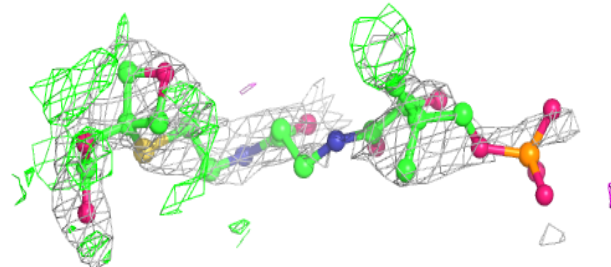
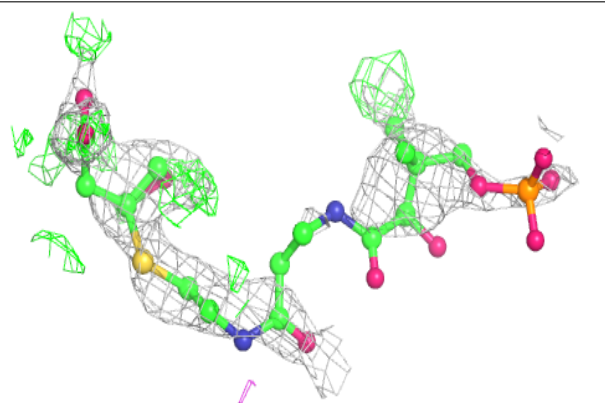
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around AFY B 401:**

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

**Electron density around AFY A 401:**

$2mF_o-DF_c$  (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.