



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:29 PM UTC

PDB ID : 6EVD / pdb\_00006evd  
Title : Structure of R173A *A. niger* Fdc1 with prFMN in the hydroxylated form  
Authors : Bailey, S.S.; David, L.; Payne, K.A.P.  
Deposited on : 2017-11-01  
Resolution : 1.19 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Buster-report : wwPDB partial adaption of 1.1.7 (2018)  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

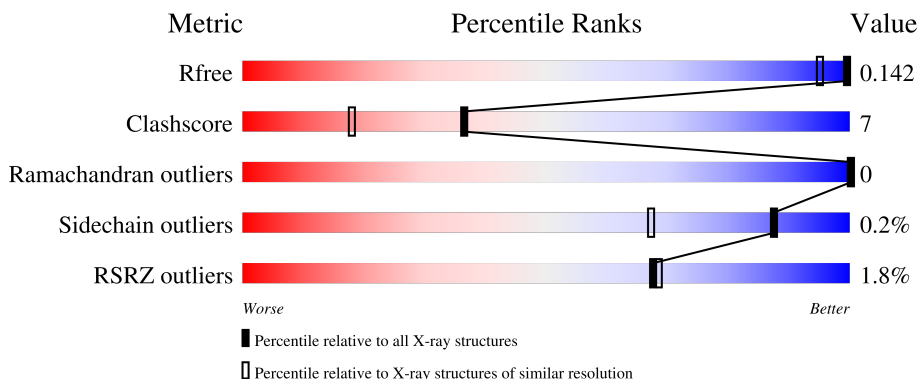
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

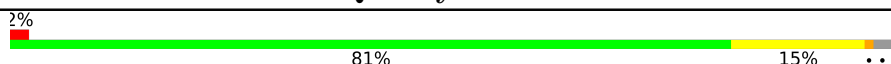
The reported resolution of this entry is 1.19 Å.

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))
$R_{free}$	180053	1216 (1.20-1.20)
Clashscore	190562	1265 (1.20-1.20)
Ramachandran outliers	187476	1226 (1.20-1.20)
Sidechain outliers	187428	1226 (1.20-1.20)
RSRZ outliers	180081	1214 (1.20-1.20)

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	508	

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4888 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 Ferulic acid decarboxylase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	496	4067	2574	693	770	30	0	34	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	173	ALA	ARG	conflict	UNP A2QHE5
A	501	LEU	-	expression tag	UNP A2QHE5
A	502	GLU	-	expression tag	UNP A2QHE5
A	503	HIS	-	expression tag	UNP A2QHE5
A	504	HIS	-	expression tag	UNP A2QHE5
A	505	HIS	-	expression tag	UNP A2QHE5
A	506	HIS	-	expression tag	UNP A2QHE5
A	507	HIS	-	expression tag	UNP A2QHE5
A	508	HIS	-	expression tag	UNP A2QHE5

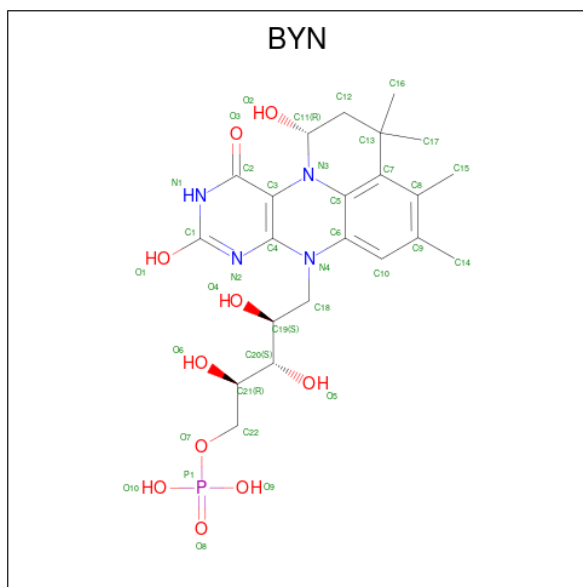
- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mn	0	0
			1	1		

- Molecule 3 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	K	0	0
			1	1		

- Molecule 4 is hydroxylated prenyl-FMN (CCD ID: BYN) (formula: C<sub>22</sub>H<sub>31</sub>N<sub>4</sub>O<sub>10</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
4	A	1	37	22	4	10	1	0	0

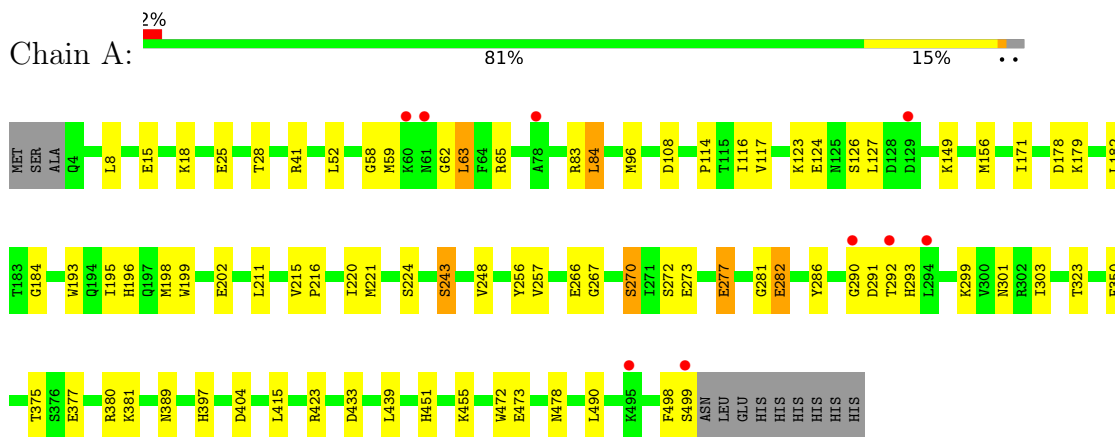
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	782	Total	O	0	0
			782	782		

### 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: Ferulic acid decarboxylase 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.23Å 64.11Å 87.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.85 – 1.19 25.85 – 1.19	Depositor EDS
% Data completeness (in resolution range)	97.2 (25.85-1.19) 97.2 (25.85-1.19)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.44 (at 1.19Å)	Xtrriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.144 , 0.162 0.145 , 0.142	Depositor DCC
$R_{free}$ test set	8454 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	10.3	Xtrriage
Anisotropy	0.273	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 38.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	4888	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	14.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.38% 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: BYN, MN, K

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	1.71	48/4165 (1.2%)	1.35	15/5659 (0.3%)

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	A	0	1

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	65	ARG	NE-CZ	9.63	1.43	1.33
1	A	380	ARG	CA-CB	8.77	1.67	1.53
1	A	301	ASN	CA-C	-7.83	1.42	1.52
1	A	116	ILE	C-O	7.65	1.32	1.24
1	A	156	MET	CA-CB	7.35	1.63	1.53
1	A	62	GLY	CA-C	6.72	1.60	1.51
1	A	83	ARG	CA-C	-6.67	1.44	1.52
1	A	198	MET	C-O	6.65	1.31	1.24
1	A	116	ILE	C-N	-6.54	1.26	1.33
1	A	65	ARG	CZ-NH1	-6.53	1.23	1.32
1	A	380	ARG	CA-C	-6.45	1.44	1.52
1	A	126	SER	C-O	6.40	1.31	1.23
1	A	323	THR	C-O	6.30	1.32	1.24
1	A	8	LEU	CA-C	-6.22	1.44	1.52
1	A	389	ASN	CG-ND2	-6.12	1.20	1.33
1	A	178	ASP	CA-C	-6.08	1.45	1.53
1	A	52	LEU	C-O	6.04	1.31	1.24
1	A	256	TYR	C-N	-6.04	1.26	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	84[A]	LEU	N-CA	5.87	1.53	1.46
1	A	84[B]	LEU	N-CA	5.87	1.53	1.46
1	A	292	THR	CA-C	-5.85	1.45	1.52
1	A	127	LEU	CA-C	-5.84	1.45	1.52
1	A	193	TRP	NE1-CE2	-5.83	1.31	1.37
1	A	272	SER	CB-OG	-5.77	1.30	1.42
1	A	8	LEU	CB-CG	-5.74	1.42	1.53
1	A	199	TRP	C-O	5.69	1.30	1.24
1	A	211	LEU	CA-C	-5.55	1.46	1.52
1	A	277	GLU	CD-OE2	5.54	1.35	1.25
1	A	123	LYS	C-O	5.48	1.31	1.24
1	A	451	HIS	CA-CB	5.45	1.61	1.53
1	A	286	TYR	CA-C	-5.43	1.46	1.52
1	A	65	ARG	CD-NE	-5.42	1.38	1.46
1	A	114	PRO	C-N	-5.41	1.26	1.33
1	A	293	HIS	C-O	-5.36	1.17	1.23
1	A	182	LEU	CA-C	-5.33	1.46	1.52
1	A	28	THR	C-O	5.31	1.31	1.23
1	A	270	SER	CA-C	-5.31	1.46	1.52
1	A	281	GLY	CA-C	-5.25	1.44	1.51
1	A	303	ILE	CA-C	-5.20	1.46	1.52
1	A	41	ARG	CA-C	-5.18	1.46	1.52
1	A	472	TRP	CA-C	-5.15	1.46	1.52
1	A	124	GLU	CA-C	-5.15	1.45	1.52
1	A	195	ILE	C-O	5.14	1.29	1.24
1	A	404	ASP	CA-C	-5.13	1.45	1.52
1	A	116	ILE	CA-C	-5.09	1.47	1.52
1	A	243	SER	CB-OG	5.07	1.52	1.42
1	A	473	GLU	C-N	-5.07	1.26	1.33
1	A	273	GLU	CA-CB	-5.03	1.45	1.53

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	83	ARG	CB-CA-C	6.81	122.42	110.85
1	A	380	ARG	CG-CD-NE	6.23	125.70	112.00
1	A	156	MET	N-CA-CB	-5.91	101.27	110.55
1	A	490	LEU	O-C-N	5.81	128.28	122.12
1	A	65	ARG	CD-NE-CZ	5.78	132.49	124.40
1	A	52	LEU	CA-C-O	-5.64	114.43	120.46
1	A	423	ARG	NE-CZ-NH2	-5.52	114.23	119.20
1	A	257	VAL	O-C-N	-5.25	117.04	121.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	277	GLU	CG-CD-OE2	-5.22	106.38	118.40
1	A	116	ILE	CA-C-N	5.21	129.90	122.24
1	A	116	ILE	C-N-CA	5.21	129.90	122.24
1	A	266	GLU	CA-C-O	-5.17	115.11	120.70
1	A	63	LEU	O-C-N	-5.14	116.65	123.13
1	A	375	THR	CA-CB-OG1	5.11	117.27	109.60
1	A	404	ASP	CB-CA-C	-5.04	99.88	110.17

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	243	SER	Mainchain

## 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	4067	0	3946	57	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	37	0	0	1	0
5	A	782	0	0	33	8
All	All	4888	0	3946	58	8

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 7.

All (58) 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:A:415[A]:LEU:HD12	5:A:1314:HOH:O	1.23	1.28
1:A:381:LYS:HG3	5:A:715:HOH:O	1.63	0.98
1:A:290[A]:GLY:HA2	5:A:703:HOH:O	1.64	0.96
1:A:290[A]:GLY:CA	5:A:703:HOH:O	2.18	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59[A]:MET:SD	5:A:1075:HOH:O	2.33	0.86
1:A:478:ASN:HB2	5:A:1188:HOH:O	1.77	0.84
1:A:59[A]:MET:CE	5:A:1075:HOH:O	2.24	0.84
1:A:499[B]:SER:N	5:A:701:HOH:O	2.10	0.83
1:A:415[A]:LEU:CD1	5:A:1314:HOH:O	1.94	0.81
1:A:478:ASN:HB2	5:A:1071:HOH:O	1.78	0.81
1:A:149[B]:LYS:HE2	5:A:705:HOH:O	1.85	0.76
1:A:290[A]:GLY:C	5:A:703:HOH:O	2.30	0.74
1:A:498:PHE:C	1:A:499[B]:SER:N	2.47	0.72
1:A:499[A]:SER:CA	5:A:701:HOH:O	2.38	0.71
1:A:433:ASP:HB2	5:A:1240:HOH:O	1.90	0.70
1:A:202:GLU:OE1	5:A:702:HOH:O	2.11	0.69
1:A:179:LYS:CD	5:A:1358:HOH:O	2.41	0.67
1:A:196[B]:HIS:HE1	1:A:270:SER:O	1.79	0.65
1:A:381:LYS:HE2	5:A:1276:HOH:O	1.97	0.64
1:A:381:LYS:HE3	5:A:715:HOH:O	1.98	0.63
1:A:59[A]:MET:SD	5:A:709:HOH:O	2.56	0.62
1:A:290[A]:GLY:O	1:A:291:ASP:C	2.42	0.62
1:A:498:PHE:C	1:A:499[B]:SER:CA	2.73	0.62
1:A:377[B]:GLU:HG3	5:A:715:HOH:O	2.00	0.61
1:A:290[A]:GLY:O	5:A:703:HOH:O	2.16	0.59
1:A:59[A]:MET:HG2	5:A:709:HOH:O	2.04	0.58
1:A:25:GLU:OE1	5:A:704:HOH:O	2.18	0.57
1:A:96[B]:MET:HG2	1:A:350:PHE:CE1	2.41	0.55
1:A:478:ASN:CB	5:A:1071:HOH:O	2.44	0.54
1:A:149[A]:LYS:NZ	5:A:705:HOH:O	2.36	0.52
1:A:196[A]:HIS:HE1	1:A:270:SER:O	1.92	0.52
1:A:15[B]:GLU:OE1	1:A:18[B]:LYS:HE3	2.10	0.51
1:A:84[B]:LEU:HD22	1:A:96[B]:MET:HE1	1.92	0.51
1:A:59[A]:MET:CG	5:A:709:HOH:O	2.59	0.49
1:A:117:VAL:HG23	1:A:248[B]:VAL:HG22	1.96	0.48
1:A:184:GLY:HA2	1:A:277:GLU:HG3	1.95	0.47
1:A:108[A]:ASP:OD1	1:A:108[A]:ASP:C	2.58	0.46
1:A:59[A]:MET:HE3	5:A:1075:HOH:O	2.07	0.46
4:A:603:BYN:C17	4:A:603:BYN:C15	2.93	0.46
1:A:381:LYS:CG	5:A:715:HOH:O	2.42	0.45
1:A:171:ILE:HD12	1:A:224[A]:SER:OG	2.17	0.45
1:A:196[A]:HIS:CD2	5:A:1043:HOH:O	2.71	0.44
1:A:290[A]:GLY:C	1:A:291:ASP:O	2.61	0.44
1:A:290[A]:GLY:O	1:A:291:ASP:O	2.35	0.44
1:A:478:ASN:CB	5:A:1188:HOH:O	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:GLY:O	1:A:59[B]:MET:C	2.56	0.44
1:A:196[A]:HIS:HD2	5:A:1043:HOH:O	2.01	0.43
1:A:282[A]:GLU:OE2	1:A:439[A]:LEU:HG	2.18	0.43
1:A:171:ILE:HD11	1:A:220:ILE:HG13	2.00	0.42
1:A:59[B]:MET:HG3	1:A:63:LEU:C	2.44	0.42
1:A:377[B]:GLU:CG	5:A:715:HOH:O	2.63	0.41
1:A:397:HIS:CD2	1:A:455:LYS:HE2	2.56	0.41
1:A:18[B]:LYS:HE2	1:A:18[B]:LYS:HB2	1.11	0.41
1:A:221[B]:MET:HA	1:A:221[B]:MET:HE2	2.02	0.41
1:A:499[B]:SER:CA	5:A:701:HOH:O	2.62	0.41
1:A:96[A]:MET:HG2	1:A:350:PHE:CE1	2.56	0.41
1:A:267:GLY:HA3	1:A:299:LYS:O	2.21	0.40
1:A:215:VAL:HB	1:A:216:PRO:HD2	2.03	0.40

All (8) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:729:HOH:O	5:A:729:HOH:O[2_555]	0.73	1.47
5:A:961:HOH:O	5:A:961:HOH:O[2_655]	0.95	1.25
5:A:1359:HOH:O	5:A:1437:HOH:O[2_555]	1.22	0.98
5:A:800:HOH:O	5:A:953:HOH:O[2_655]	1.66	0.54
5:A:1151:HOH:O	5:A:1211:HOH:O[2_655]	1.74	0.46
5:A:1154:HOH:O	5:A:1365:HOH:O[2_655]	1.98	0.22
5:A:898:HOH:O	5:A:1171:HOH:O[3_545]	2.00	0.20
5:A:1020:HOH:O	5:A:1212:HOH:O[3_555]	2.19	0.01

## 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	527/508 (104%)	517 (98%)	10 (2%)	0	<a href="#">100</a> <a href="#">100</a>

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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	438/429 (102%)	436 (100%)	2 (0%)	81 59

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	282[A]	GLU
1	A	282[B]	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	ASN
1	A	160	GLN
1	A	197	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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
4	BYN	A	603	3,2	36,40,40	1.36	4 (11%)	47,64,64	1.56	5 (10%)

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
4	BYN	A	603	3,2	-	2/18/33/33	0/4/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	603	BYN	C3-N3	3.96	1.44	1.37
4	A	603	BYN	O3-C2	3.45	1.30	1.23
4	A	603	BYN	C12-C11	-3.15	1.49	1.52
4	A	603	BYN	C3-C2	-2.56	1.37	1.43

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	603	BYN	C13-C7-C5	-6.29	114.61	121.48
4	A	603	BYN	C3-C2-N1	3.51	117.12	110.94
4	A	603	BYN	O2-C11-C12	2.73	117.04	110.83
4	A	603	BYN	C17-C13-C7	2.59	119.74	112.06
4	A	603	BYN	C1-N2-C4	2.56	118.34	113.47

There are no chirality outliers.

All (2) torsion outliers are listed below:

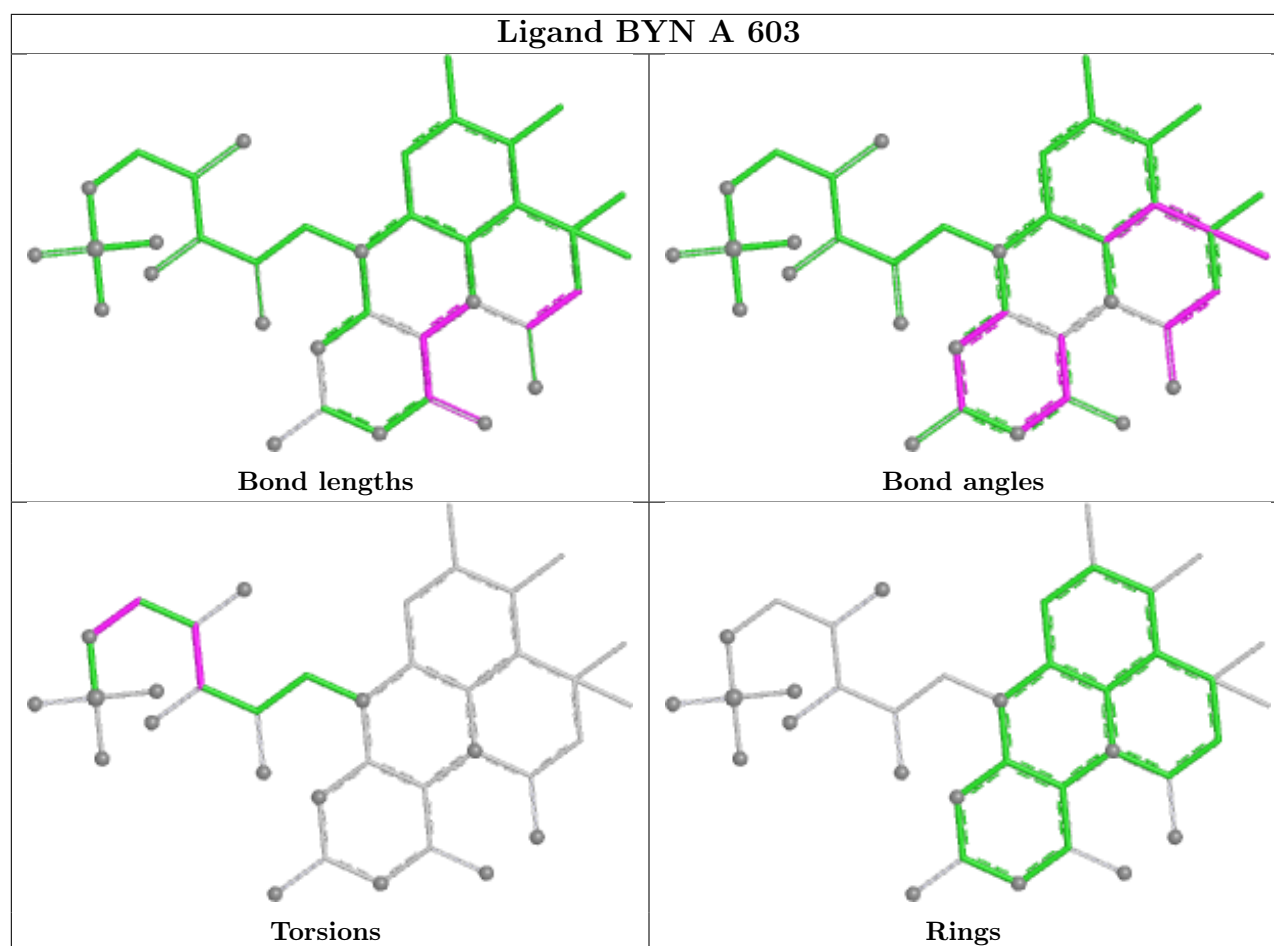
Mol	Chain	Res	Type	Atoms
4	A	603	BYN	C19-C20-C21-C22
4	A	603	BYN	C21-C22-O7-P1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	603	BYN	1	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 [\(i\)](#)

There are no such residues in this entry.

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

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	498:PHE	C	499[B]:SER	N	2.47

## 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	496/508 (97%)	-0.02	9 (1%) 67 68	4, 10, 19, 32	34 (6%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	292	THR	4.4
1	A	129[A]	ASP	3.9
1	A	61	ASN	3.0
1	A	294	LEU	2.5
1	A	60	LYS	2.5
1	A	499[A]	SER	2.4
1	A	78	ALA	2.3
1	A	290[A]	GLY	2.2
1	A	495	LYS	2.0

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

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

### 6.3 Carbohydrates [i](#)

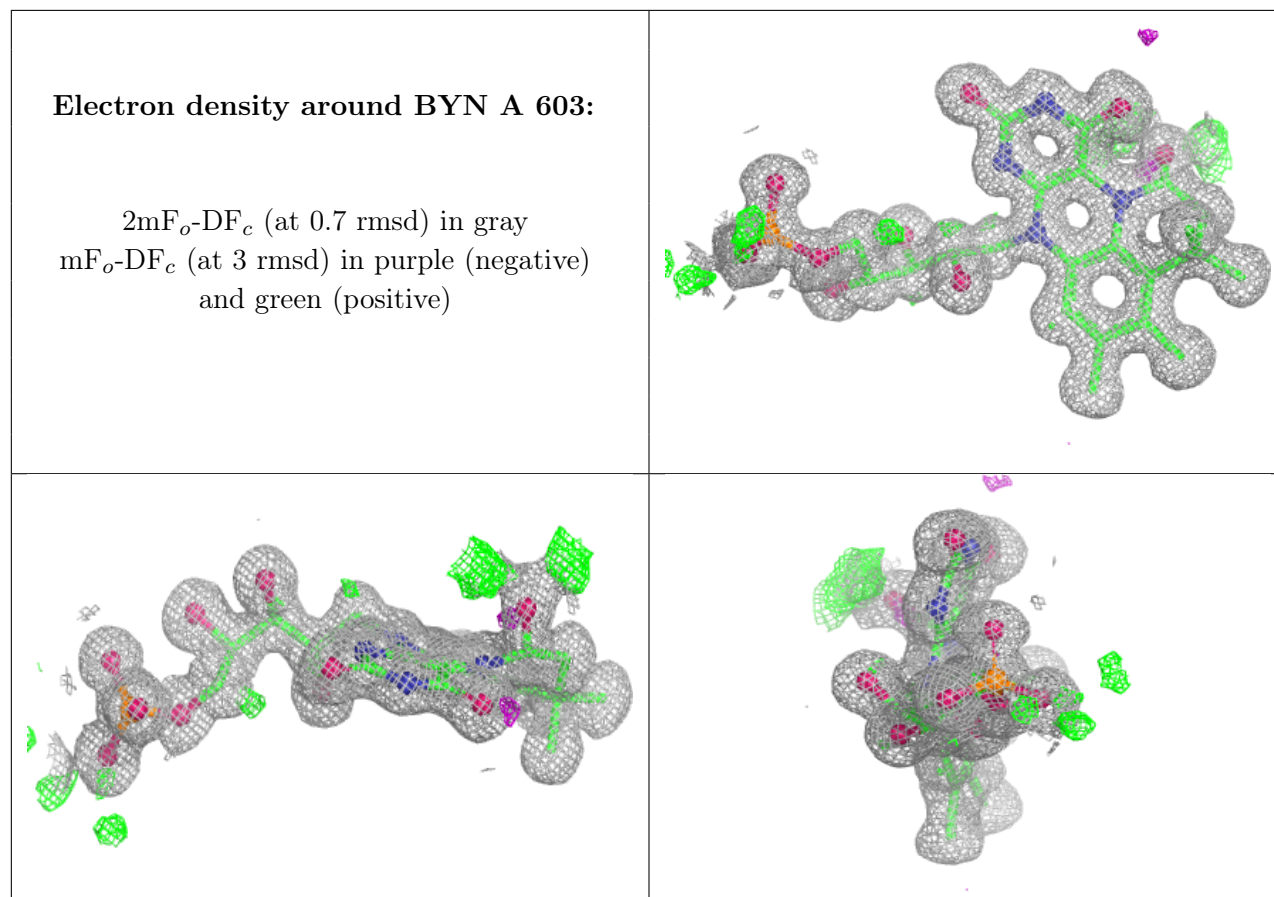
There are no oligosaccharides in this entry.

### 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
4	BYN	A	603	37/37	0.99	0.04	6,7,10,26	0
3	K	A	602	1/1	1.00	0.01	6,6,6,6	0
2	MN	A	601	1/1	1.00	0.01	6,6,6,6	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.



## 6.5 Other polymers [i](#)

There are no such residues in this entry.