

# Full wwPDB X-ray Structure Validation Report (i)

### May 16, 2020 – 05:15 pm BST

PDB ID : 5M6R

Title: Human porphobiling en deaminase in complex with reaction intermediate

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Deposited on : 2016-10-25

Resolution : 2.73 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
A user guide is available at

https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) 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)

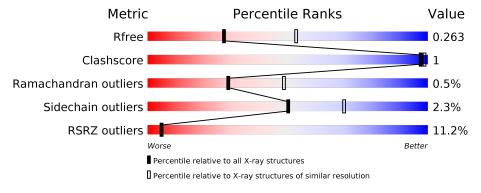
Validation Pipeline (wwPDB-VP) : 2.11

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$ 

The reported resolution of this entry is 2.73 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	Similar resolution		
Metric	$(\# { m Entries})$	$(\#  ext{Entries},  ext{resolution range}( ext{Å}))$		
$R_{free}$	130704	1271 (2.76-2.72)		
Clashscore	141614	1322 (2.76-2.72)		
Ramachandran outliers	138981	1297 (2.76-2.72)		
Sidechain outliers	138945	1298 (2.76-2.72)		
RSRZ outliers	127900	1243 (2.76-2.72)		

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 >=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 <=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		
			10%		
1	A	389	84%	٠	13%
4	D	000	9%		
	В	389	79%		17%



# 2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 10597 atoms, of which 5320 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 Porphobilinogen deaminase.

$\mathbf{Mol}$	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
1	A	340	Total 5307	C 1649	H 2684	N 476	O 488	S 10	0	1	0
1	В	322	Total 5010	C 1560	H 2529	N 450	O 461	S 10	0	0	0

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-27	MET	_	initiating methionine	UNP P08397
A	-26	LYS	-	expression tag	UNP P08397
A	-25	HIS	-	expression tag	UNP P08397
A	-24	HIS	-	expression tag	UNP P08397
A	-23	HIS	-	expression tag	UNP P08397
A	-22	HIS	-	expression tag	UNP P08397
A	-21	HIS	-	expression tag	UNP P08397
A	-20	HIS	-	expression tag	UNP P08397
A	-19	PRO	-	expression tag	UNP P08397
A	-18	MET	-	expression tag	UNP P08397
A	-17	SER	-	expression tag	UNP P08397
A	-16	ASP	-	expression tag	UNP P08397
A	-15	TYR	-	expression tag	UNP P08397
A	-14	ASP	-	expression tag	UNP P08397
A	-13	ILE	-	expression tag	UNP P08397
A	-12	PRO	-	expression tag	UNP P08397
A	-11	THR	-	expression tag	UNP P08397
A	-10	THR	-	expression tag	UNP P08397
A	-9	GLU	-	expression tag	UNP P08397
A	-8	ASN	-	expression tag	UNP P08397
A	-7	LEU	-	expression tag	UNP P08397
A	-6	TYR	-	expression tag	UNP P08397
A	-5	PHE	=	expression tag	UNP P08397
A	-4	GLN	-	expression tag	UNP P08397
A	-3	GLY	-	expression tag	UNP P08397

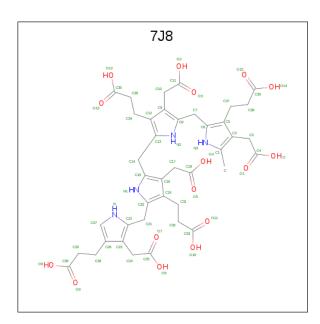


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Chain	Residue	Modelled  Modelled	Actual	Comment	Reference
A	-2	ALA	-	expression tag	UNP P08397
A	-1	MET	-	expression tag	UNP P08397
A	0	ALA	-	- expression tag	
В	-27	MET	-	initiating methionine	UNP P08397
В	-26	LYS	-	expression tag	UNP P08397
В	-25	HIS	-	expression tag	UNP P08397
В	-24	HIS	-	expression tag	UNP P08397
В	-23	HIS	-	expression tag	UNP P08397
В	-22	HIS	-	expression tag	UNP P08397
В	-21	HIS	-	expression tag	UNP P08397
В	-20	HIS	_	expression tag	UNP P08397
В	-19	PRO	_	expression tag	UNP P08397
В	-18	MET	_	expression tag	UNP P08397
В	-17	SER	-	expression tag	UNP P08397
В	-16	ASP	-	expression tag	UNP P08397
В	-15	TYR	-	expression tag	UNP P08397
В	-14	ASP	-	expression tag	UNP P08397
В	-13	ILE	-	expression tag	UNP P08397
В	-12	PRO	-	expression tag	UNP P08397
В	-11	THR	-	expression tag	UNP P08397
В	-10	THR	-	expression tag	UNP P08397
В	-9	GLU	-	expression tag	UNP P08397
В	-8	ASN	-	expression tag	UNP P08397
В	-7	LEU	-	expression tag	UNP P08397
В	-6	TYR	-	expression tag	UNP P08397
В	-5	PHE	-	expression tag	UNP P08397
В	-4	GLN	-	expression tag	UNP P08397
В	-3	GLY	-	expression tag	UNP P08397
В	-2	ALA	-	expression tag	UNP P08397
В	-1	MET	-	expression tag	UNP P08397
В	0	ALA	-	expression tag	UNP P08397

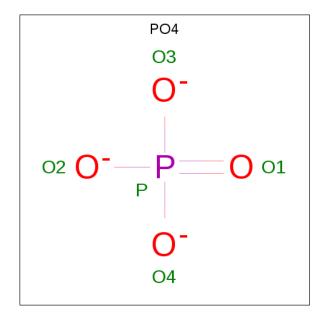
• Molecule 2 is 3-[4-(2-hydroxy-2-oxoethyl)-5-[[4-(2-hydroxy-2-oxoethyl)-5-[[4-(2-hydroxy-2-oxoethyl)-5-[[4-(2-hydroxy-2-oxoethyl)-5-[[4-(2-hydroxy-2-oxoethyl)-5-methyl-1 {H}-pyrrol-2-yl]methyl]-3-(3-hydroxy-3-oxopropyl)-1 {H}-pyrrol-2-yl]methyl]-3-(3-hydroxy-3-oxopropyl)-1 {H}-pyrrol-2-yl]methyl]-1 {H}-pyrrol-3-yl]propanoic acid (three-letter code: 7J8) (formula:  $C_{40}H_{46}N_4O_{16}$ ).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
9	Λ	1	Total	С	H	N	О	0	0
$\begin{array}{ c c c c c c } & Z & A & A & A & A & A & A & A & A & A$	1	105	40	45	4	16	0		
9	D	1	Total	С	Н	N	О	0	0
	Б	1	105	40	45	4	16	0	0

 $\bullet$  Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula:  $\mathrm{O_4P}).$ 



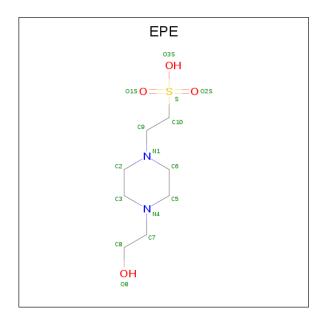
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	
3	A	1	Total O P	0	0	
	9 11	_	5 4 1	Ŭ	Ü	
3	Δ	1	Total O P	0	0	
3   A	Λ	1	$\begin{vmatrix} 5 & 4 & 1 \end{vmatrix}$			



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Mo	Chain	Residues	${f Atoms}$			ZeroOcc	AltConf
3	В	1	Total 5	O 4	P 1	0	0

• Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula:  $C_8H_{18}N_2O_4S$ ).



Mol	Chain	Residues	${f Atoms}$				ZeroOcc	AltConf		
4	A	1	Total 32	_	H 17		O 4	S 1	0	0

• Molecule 5 is water.

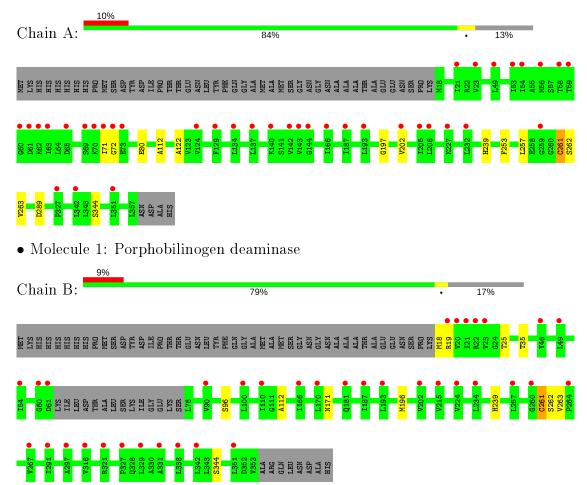
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	15	Total O 15 15	0	0
5	В	8	Total O 8 8	0	0



# 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: Porphobilinogen deaminase





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	68.91Å 81.20Å 79.69Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 93.01° 90.00°	Depositor
Resolution (Å)	56.84 - 2.73	Depositor
Resolution (A)	56.84 - 2.73	EDS
% Data completeness	95.8 (56.84-2.73)	Depositor
(in resolution range)	95.8 (56.84-2.73)	EDS
$R_{merge}$	0.19	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.81 (at 2.73Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
P. P.	0.197 , $0.246$	Depositor
$R, R_{free}$	0.212 , $0.263$	DCC
$R_{free}$ test set	1104 reflections $(4.90\%)$	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	49.4	Xtriage
Anisotropy	0.861	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.40 , 70.2	EDS
L-test for twinning <sup>2</sup>	$ < L > = 0.48, < L^2> = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	10597	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	91.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 7.03% of the height of the origin peak. No significant pseudotranslation is detected.

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



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

# 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: PO4, EPE, 7J8

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		
IVIOI		RMSZ	# Z >5	RMSZ	# Z  > 5	
1	A	0.52	0/2669	0.67	0/3613	
1	В	0.51	0/2523	0.68	0/3417	
All	All	0.51	0/5192	0.68	0/7030	

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	2623	2684	2691	6	0
1	В	2481	2529	2529	3	0
2	A	60	45	0	0	0
2	В	60	45	0	0	0
3	A	10	0	0	0	0
3	В	5	0	0	0	0
4	A	15	17	17	0	0
5	A	15	0	0	0	0
5	В	8	0	0	0	0
All	All	5277	5320	5237	8	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 1.

All (8) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{array}{l}  ext{Interatomic} \  ext{distance} \ ( ext{Å}) \end{array}$	Clash overlap (Å)
1:A:71:ILE:HG22	1:A:72:GLY:H	1.68	0.58
1:B:25:THR:OG1	1:B:35:THR:HG21	2.14	0.48
1:A:71:ILE:HG22	1:A:72:GLY:N	2.31	0.44
1:A:197:GLY:HA2	1:B:196:MET:O	2.19	0.43
1:A:253:PHE:CZ	1:A:257:LEU:HD11	2.54	0.43
1:B:112:ALA:HA	1:B:239:HIS:CE1	2.55	0.42
1:A:112:ALA:HA	1:A:239:HIS:CE1	2.55	0.41
1:A:122:ALA:HB1	1:A:202:VAL:HG13	2.02	0.41

There are no symmetry-related clashes.

### 5.3 Torsion angles (i)

### 5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	$\mathbf{ntiles}$
1	A	339/389 (87%)	320 (94%)	18 (5%)	1 (0%)	41	61
1	В	318/389 (82%)	305 (96%)	11 (4%)	2 (1%)	25	44
All	All	657/778 (84%)	625 (95%)	29 (4%)	3 (0%)	29	48

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	19	ARG
1	A	261	CYS
1	В	261	CYS



### 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	286/324 (88%)	280 (98%)	6 (2%)	53 72		
1	В	270/324 (83%)	263 (97%)	7 (3%)	46 66		
All	All	556/648 (86%)	543 (98%)	13 (2%)	50 70		

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	GLU
1	A	261	CYS
1	A	262	SER
1	A	263	VAL
1	A	289	ASP
1	A	344	SER
1	В	18	MET
1	В	96	SER
1	В	171	ASN
1	В	261	CYS
1	В	262	SER
1	В	263	VAL
1	В	344	SER

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

#### 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 carbohydrates in this entry.

## 5.6 Ligand geometry (i)

6 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	Tuna	Chain	ain Res Link Bugg   Bond lengths				ths	Bond angles		
10101	Type	Chain	nes	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	7J8	A	401	1	32,63,63	1.08	2 (6%)	30,89,89	1.32	4 (13%)
3	PO4	A	403	-	4,4,4	2.73	2 (50%)	6,6,6	1.08	0
4	EPE	A	404	-	15,15,15	1.59	1 (6%)	18,20,20	0.65	0
3	PO4	A	402	-	4,4,4	2.55	1 (25%)	6,6,6	0.51	0
2	7J8	В	401	1	32,63,63	1.23	3 (9%)	30,89,89	0.87	2 (6%)
3	PO4	В	402	-	4,4,4	1.83	0	6,6,6	0.46	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	7J8	A	401	1	-	1/26/48/48	0/4/4/4
2	7J8	В	401	1	-	1/26/48/48	0/4/4/4
4	EPE	A	404	-	-	0/9/19/19	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(\mathbf{\mathring{A}})$	$Ideal( ext{\AA})$
4	A	404	EPE	C10-S	-6.01	1.69	1.77
3	A	402	PO4	P-O1	4.23	1.60	1.50
3	A	403	PO4	P-O1	4.18	1.60	1.50
2	В	401	7J8	C16-C19	2.97	1.46	1.39
2	A	401	7J8	C9-C12	2.48	1.45	1.39



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Mol	Chain	Res	Type	${f Atoms}$	${f Z}$	${ m Observed}({ m \AA})$	$\operatorname{Ideal}( ext{\AA})$
2	В	401	7J8	C2-C5	2.41	1.44	1.39
2	В	401	7J8	C9-C12	2.37	1.44	1.39
2	A	401	7J8	C26-C23	2.25	1.44	1.40
3	A	403	PO4	P-O3	2.17	1.61	1.54

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
2	A	401	7J8	C34-C12-C13	3.58	129.81	127.30
2	A	401	7J8	C17-C16-C15	3.16	129.19	126.41
2	В	401	7J8	C34-C12-C13	2.65	129.16	127.30
2	A	401	7J8	C31-C19-C20	-2.65	125.44	127.30
2	В	401	7J8	C24-C23-C22	-2.20	124.47	126.41
2	A	401	7J8	C37-C5-C6	2.12	128.78	127.30

There are no chirality outliers.

All (2) torsion outliers are listed below:

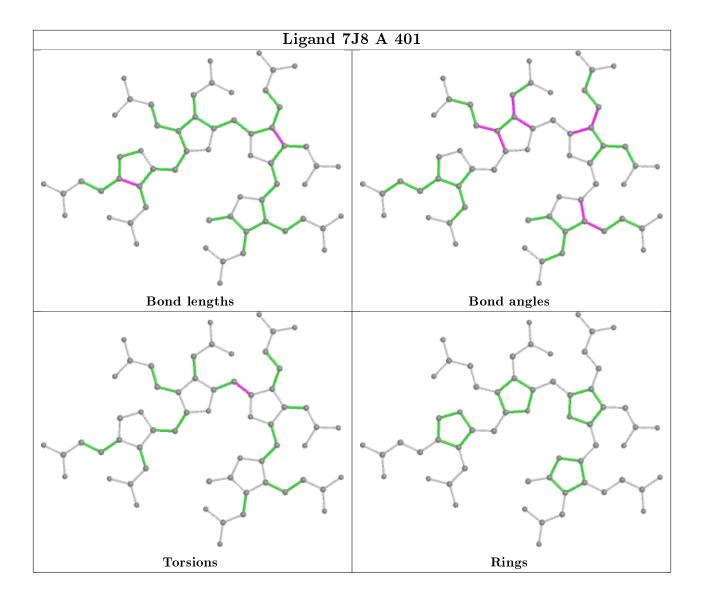
Mol	Chain	Res	Type	Atoms
2	A	401	7J8	C12-C13-C14-C15
2	В	401	7J8	C12-C13-C14-C15

There are no ring outliers.

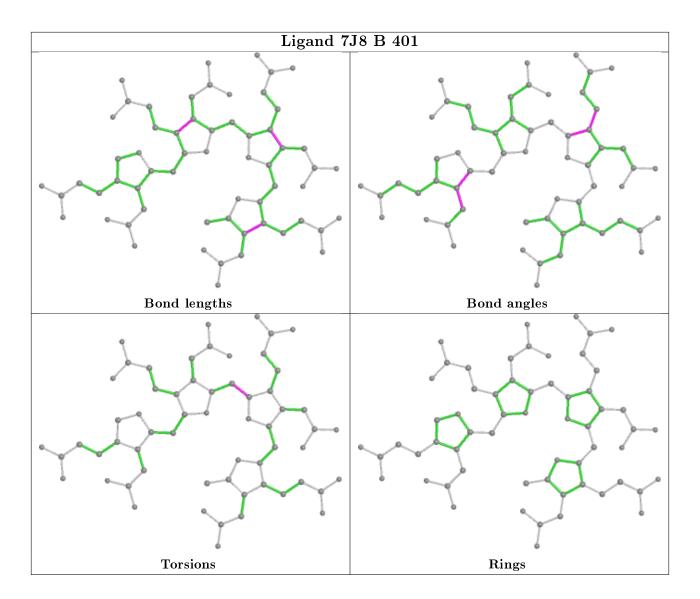
No monomer is involved in short contacts.

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 then 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)

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,  $95^{th}$  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 $>$	$\#\mathrm{RSRZ}{>}2$	$OWAB(\AA^2)$	Q < 0.9
1	A	340/389 (87%)	0.84	38 (11%) 5 5	53, 83, 145, 177	0
1	В	322/389 (82%)	0.83	36 (11%) 5 5	62, 86, 133, 157	0
All	All	662/778 (85%)	0.83	74 (11%) 5 5	53, 85, 138, 177	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	61	ASP	7.9
1	A	60	GLY	6.0
1	В	351	LEU	5.4
1	A	49	LEU	4.2
1	A	70	LYS	4.1
1	A	59	THR	4.0
1	A	71	ILE	3.9
1	A	62	LYS	3.6
1	A	143	VAL	3.6
1	В	46	TYR	3.5
1	A	259	GLY	3.3
1	A	129	PHE	3.3
1	В	100	LEU	3.3
1	A	124	VAL	3.3
1	A	58	THR	3.1
1	В	260	GLY	3.1
1	В	342	LEU	3.1
1	A	23	VAL	3.1
1	В	316	VAL	3.1
1	В	110	ILE	3.1
1	В	61	ASP	3.0
1	В	54	ILE	3.0
1	A	63	ILE	2.9
1	A	342	LEU	2.9



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$oxed{Continued\ from\ previous\ page}} oxed{oxed{Mol\ Chain\ Res\ Type\ RSRZ}}$								
			Type					
1	В	49	LEU	2.8				
1	A	65	ASP	2.8				
1	A	166	ILE	2.8				
1	A	187	ILE	2.8				
1	В	21	ILE	2.8				
1	В	257	LEU	2.7				
1	A	54	ILE	2.7				
1	В	327	PRO	2.7				
1	A	205	ILE	2.6				
1	В	23	VAL	2.6				
1	В	187	ILE	2.6				
1	A	206	LEU	2.6				
1	A	227	LYS	2.6				
1	A	69	SER	2.5				
1	A	56	MET	2.5				
1	В	193	LEU	2.5				
1	В	166	ILE	2.5				
1	В	202	VAL	2.4				
1	В	19	ARG	2.4				
1	В	264	PRO	2.4				
1	В	291	ILE	2.4				
1	В	329	LEU	2.4				
1	A	134	LEU	2.4				
1	A	137	LEU	2.4				
1	A	53	ILE	2.4				
1	В	338	LEU	2.4				
1	A	73	GLU	2.4				
1	A	351	LEU	2.4				
1	В	297	ALA	2.4				
1	В	224	VAL	2.4				
1	A	202	VAL	2.3				
1	В	90	VAL	2.3				
1	A	72	GLY	2.3				
1	A	144	GLY	2.3				
1	A	193	LEU	2.2				
1	В	321	ARG	2.2				
1	В	267	VAL	2.2				
1	A	140	LYS	2.2				
1	В	22	ARG	2.2				
1	A	232	LEU	2.1				
1	В	60	GLY	2.1				
1	В	181	GLN	2.1				



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Mol	Chain	Res	Type	RSRZ
1	A	142	VAL	2.1
1	В	215	VAL	2.1
1	В	170	LEU	2.0
1	В	331	ALA	2.0
1	В	234	LEU	2.0
1	A	327	PRO	2.0
1	В	20	VAL	2.0
1	A	21	ILE	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 carbohydrates in this entry.

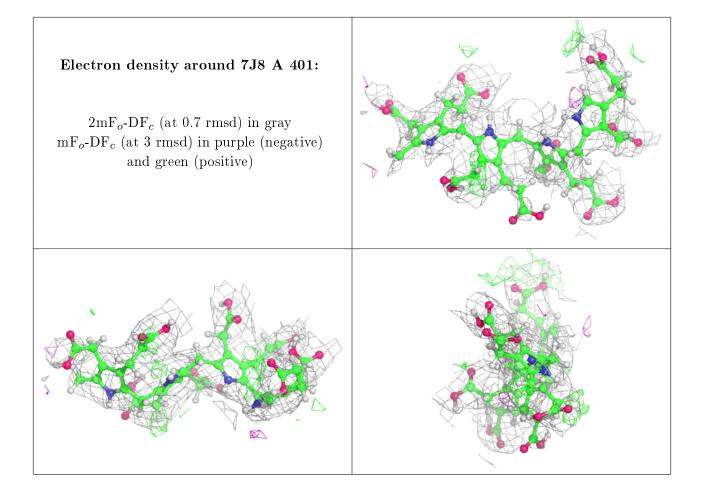
## 6.4 Ligands (i)

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,  $95^{th}$  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.

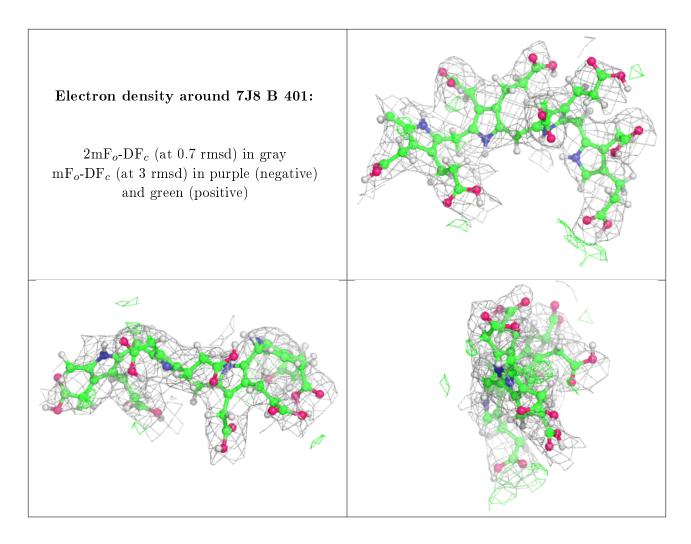
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\operatorname{B-factors}(  ext{\AA}^2 )$	Q<0.9
4	EPE	A	404	15/15	0.56	0.28	133,159,176,179	0
2	7J8	A	401	60/60	0.87	0.21	65,87,109,116	0
2	7J8	В	401	60/60	0.89	0.22	68,97,114,133	0
3	PO4	A	403	5/5	0.92	0.23	94,97,99,102	0
3	PO4	В	402	5/5	0.94	0.23	96,99,102,104	0
3	PO4	A	402	5/5	0.95	0.17	$99,\!103,\!105,\!106$	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.

