

wwPDB X-ray Structure Validation Summary Report (i)

Oct 23, 2023 – 02:39 PM JST

PDB ID	:	8J3P
Title	:	Formate dehydrogenase mutant from from Candida dubliniensis M4 complexed
		with NADP+
Authors	:	Ma, W.; Zheng, Y.C.; Geng, Q.; Chen, C.; Xu, J.H.
Deposited on		
Resolution	:	2.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

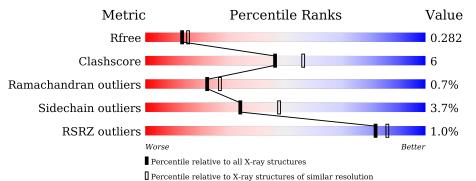
MolProbity Mogul Xtriage (Phenix) EDS	:	4.02b-467 1.8.5 (274361), CSD as541be (2020) 1.13 2.36
buster-report Percentile statistics Refmac	: : :	1.1.7 (2018) 20191225.v01 (using entries in the PDB archive December 25th 2019) 5.8.0158 7.0.044 (Gargrove)
Ideal geometry (DNA, RNA) Validation Pipeline (wwPDB-VP)		Parkinson et al. (1996) 2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.30 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	5042 (2.30-2.30)
Clashscore	141614	5643(2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 >=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		
1	А	382	% 82%	13%	•••
1	В	382	% 8 3%	14%	••
1	С	382	.% 80%	16%	•••



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 9187 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Δ	374	Total	С	Ν	0	S	0	0	0
	A	374	2913	1849	497	554	13	0	0	0
1	D	374	Total	С	Ν	0	S	0	0	0
	D	374	2913	1849	497	554	13	0	0	0
1	C	374	Total	С	Ν	0	S	0	0	0
	U	374	2913	1849	497	554	13	U	0	0

• Molecule 1 is a protein called Formate dehydrogenase.

Chain	Residue	Modelled	Actual	Comment	Reference
А	-7	MET	-	initiating methionine	UNP B9WHT3
А	-6	GLY	-	expression tag	UNP B9WHT3
А	-5	HIS	-	expression tag	UNP B9WHT3
А	-4	HIS	-	expression tag	UNP B9WHT3
A	-3	HIS	-	expression tag	UNP B9WHT3
А	-2	HIS	-	expression tag	UNP B9WHT3
А	-1	HIS	-	expression tag	UNP B9WHT3
А	0	HIS	-	expression tag	UNP B9WHT3
А	16	LEU	HIS	conflict	UNP B9WHT3
А	159	ARG	LYS	conflict	UNP B9WHT3
А	167	ARG	LYS	conflict	UNP B9WHT3
А	197	GLN	ASP	conflict	UNP B9WHT3
А	198	ARG	TYR	conflict	UNP B9WHT3
А	199	ASN	GLN	conflict	UNP B9WHT3
А	371	THR	LYS	conflict	UNP B9WHT3
А	372	SER	ALA	conflict	UNP B9WHT3
В	-7	MET	-	initiating methionine	UNP B9WHT3
В	-6	GLY	-	expression tag	UNP B9WHT3
В	-5	HIS	-	expression tag	UNP B9WHT3
В	-4	HIS	-	expression tag	UNP B9WHT3
В	-3	HIS	-	expression tag	UNP B9WHT3
В	-2	HIS	-	expression tag	UNP B9WHT3
В	-1	HIS	-	expression tag	UNP B9WHT3
				α i	d on nort nago

There are 48 discrepancies between the modelled and reference sequences:

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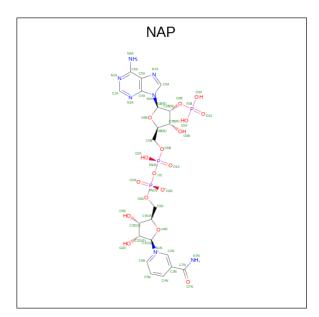


Chain	Residue	Modelled	Actual	Comment	Reference
В	0	HIS	-	expression tag	UNP B9WHT3
В	16	LEU	HIS	conflict	UNP B9WHT3
В	159	ARG	LYS	conflict	UNP B9WHT3
В	167	ARG	LYS	conflict	UNP B9WHT3
В	197	GLN	ASP	conflict	UNP B9WHT3
В	198	ARG	TYR	conflict	UNP B9WHT3
В	199	ASN	GLN	conflict	UNP B9WHT3
В	371	THR	LYS	conflict	UNP B9WHT3
В	372	SER	ALA	conflict	UNP B9WHT3
С	-7	MET	-	initiating methionine	UNP B9WHT3
С	-6	GLY	-	expression tag	UNP B9WHT3
С	-5	HIS	-	expression tag	UNP B9WHT3
С	-4	HIS	-	expression tag	UNP B9WHT3
С	-3	HIS	-	expression tag	UNP B9WHT3
С	-2	HIS	-	expression tag	UNP B9WHT3
С	-1	HIS	-	expression tag	UNP B9WHT3
С	0	HIS	-	expression tag	UNP B9WHT3
С	16	LEU	HIS	conflict	UNP B9WHT3
С	159	ARG	LYS	conflict	UNP B9WHT3
С	167	ARG	LYS	conflict	UNP B9WHT3
С	197	GLN	ASP	conflict	UNP B9WHT3
С	198	ARG	TYR	conflict	UNP B9WHT3
С	199	ASN	GLN	conflict	UNP B9WHT3
С	371	THR	LYS	conflict	UNP B9WHT3
С	372	SER	ALA	conflict	UNP B9WHT3

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• Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (threeletter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf		
2	Δ	1	Total	С	Ν	Ο	Р	0	0
2	Λ	T	48	21	7	17	3	0	0
9	В	1	Total	С	Ν	Ο	Р	0	0
	D	1	48	21	7	17	3	0	0
9	С	1	Total	С	Ν	Ο	Р	0	0
	U	1	48	21	7	17	3	0	U

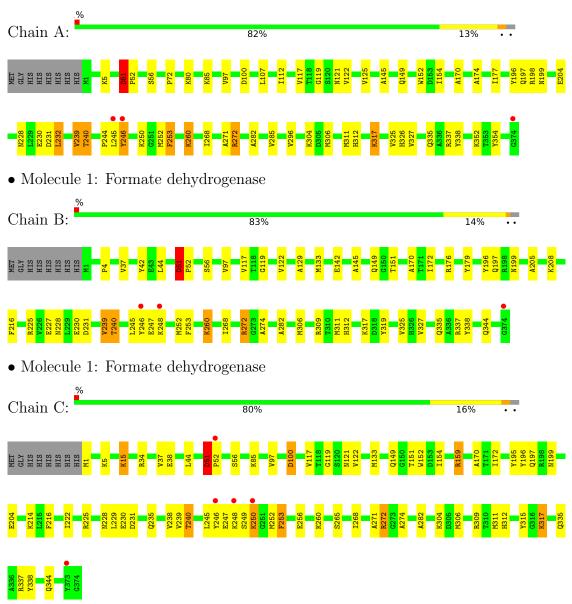
• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	104	Total O 104 104	0	0
3	В	97	Total O 97 97	0	0
3	С	103	Total O 103 103	0	0



3 Residue-property plots (i)

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



• Molecule 1: Formate dehydrogenase



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants	65.22Å 113.02Å 401.29Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.44 - 2.30	Depositor
Resolution (A)	49.24 - 2.30	EDS
% Data completeness	99.9(33.44-2.30)	Depositor
(in resolution range)	$99.8 \ (49.24 - 2.30)$	EDS
R _{merge}	0.07	Depositor
R_{sum}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$6.53 (at 2.29 \text{\AA})$	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
P. P.	0.230 , 0.283	Depositor
R, R_{free}	0.229 , 0.282	DCC
R_{free} test set	3403 reflections $(5.12%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	37.5	Xtriage
Anisotropy	0.238	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35 , 24.8	EDS
L-test for $twinning^2$	$< L > = 0.49, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.487 for 1/2 *h-1/2 *k,-3/2 *h-1/2 *k,-1	Xtriage
	0.487 for $1/2$ *h+ $1/2$ *k, $3/2$ *h- $1/2$ *k,-l	Attrage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9187	wwPDB-VP
Average B, all atoms $(Å^2)$	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.38% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities 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: NAP

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	Chain Bond lengths		Bond angles		
10101	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.29	0/2974	0.55	2/4032~(0.0%)	
1	В	0.25	0/2974	0.52	1/4032~(0.0%)	
1	С	0.28	0/2974	0.57	6/4032~(0.1%)	
All	All	0.27	0/8922	0.55	9/12096~(0.1%)	

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	А	0	1
1	В	0	1
1	С	0	1
All	All	0	3

There are no bond length outliers.

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
1	В	51	ASP	C-N-CD	-11.07	96.25	120.60
1	С	85	LYS	CD-CE-NZ	-9.54	89.77	111.70
1	С	214	LYS	CA-CB-CG	7.25	129.35	113.40
1	А	85	LYS	CD-CE-NZ	-6.92	95.78	111.70
1	С	250	LYS	CA-CB-CG	-6.12	99.93	113.40

There are no chirality outliers.

All (3) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	А	51	ASP	Peptide
1	В	51	ASP	Peptide
1	С	51	ASP	Peptide

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	А	2913	0	2907	39	0
1	В	2913	0	2907	34	0
1	С	2913	0	2907	44	0
2	А	48	0	25	6	0
2	В	48	0	25	2	0
2	С	48	0	25	1	0
3	А	104	0	0	3	0
3	В	97	0	0	3	0
3	С	103	0	0	6	0
All	All	9187	0	8796	112	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 6.

The worst 5 of 112 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:51:ASP:O	1:B:56:SER:HB3	1.78	0.84	
1:A:228:ASN:HD22	1:A:231:ASP:H	1.24	0.83	
1:A:317:LYS:HD3	1:A:317:LYS:H	1.50	0.76	
1:B:240:THR:HG23	1:B:268:ILE:HB	1.72	0.71	
1:B:337:ARG:NH2	3:B:502:HOH:O	2.25	0.69	

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	ntiles
1	А	372/382~(97%)	355~(95%)	15 (4%)	2~(0%)	29	35
1	В	372/382~(97%)	354 (95%)	15 (4%)	3(1%)	19	23
1	С	372/382~(97%)	353~(95%)	16 (4%)	3 (1%)	19	23
All	All	1116/1146~(97%)	1062 (95%)	46 (4%)	8 (1%)	22	26

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	\mathbf{Res}	Type
1	А	52	PRO
1	В	51	ASP
1	В	52	PRO
1	С	52	PRO
1	В	247	GLU

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	А	309/316~(98%)	295~(96%)	14 (4%)	27 39	
1	В	309/316~(98%)	302~(98%)	7 (2%)	50 67	
1	С	309/316~(98%)	296~(96%)	13 (4%)	30 42	
All	All	927/948~(98%)	893~(96%)	34~(4%)	34 48	

5 of 34 residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	С	256	GLU
1	С	265	SER
1	С	306	MET
1	А	306	MET
1	А	272	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	В	340	ASN
1	В	344	GLN
1	С	344	GLN
1	С	121	ASN
1	В	228	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

3 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	ol Type Chain Res L		Link	Bond lengths			Bond angles			
IVIOI	Type	Unam	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
2	NAP	С	401	-	$45,\!52,\!52$	0.81	1 (2%)	$56,\!80,\!80$	1.16	4 (7%)
2	NAP	В	401	-	45,52,52	0.81	2 (4%)	56,80,80	1.13	4 (7%)
2	NAP	А	401	-	45,52,52	0.84	1 (2%)	56,80,80	1.24	4 (7%)

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	NAP	С	401	-	-	7/31/67/67	0/5/5/5
2	NAP	В	401	-	-	6/31/67/67	0/5/5/5
2	NAP	А	401	-	-	7/31/67/67	0/5/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	Observed(Å)	Ideal(Å)
2	А	401	NAP	C5A-C4A	2.52	1.47	1.40
2	С	401	NAP	C5A-C4A	2.43	1.47	1.40
2	В	401	NAP	C5A-C4A	2.43	1.47	1.40
2	В	401	NAP	O4D-C1D	2.04	1.43	1.41

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$\mathbf{Ideal}(^{o})$
2	А	401	NAP	C3D-C2D-C1D	3.78	106.67	100.98
2	С	401	NAP	C3D-C2D-C1D	3.42	106.13	100.98
2	В	401	NAP	PN-O3-PA	-3.40	121.17	132.83
2	В	401	NAP	N3A-C2A-N1A	-3.22	123.64	128.68
2	С	401	NAP	N3A-C2A-N1A	-3.22	123.65	128.68

There are no chirality outliers.

5 of 20 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	А	401	NAP	C5D-O5D-PN-O3
2	А	401	NAP	C5D-O5D-PN-O2N
2	А	401	NAP	O4D-C4D-C5D-O5D
2	А	401	NAP	O4D-C1D-N1N-C6N
2	В	401	NAP	C5D-O5D-PN-O3

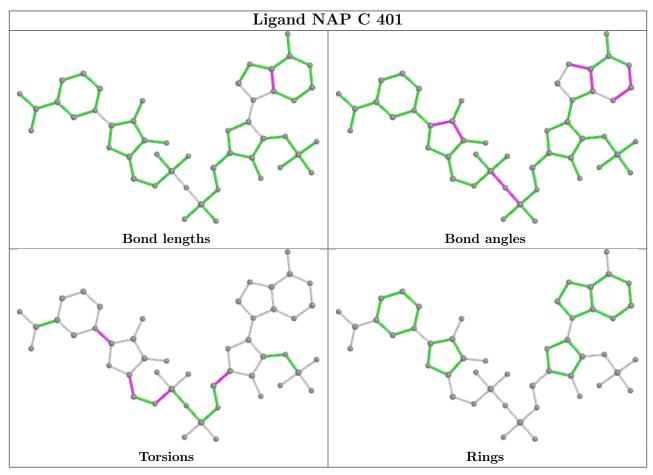


There are no ring outliers.

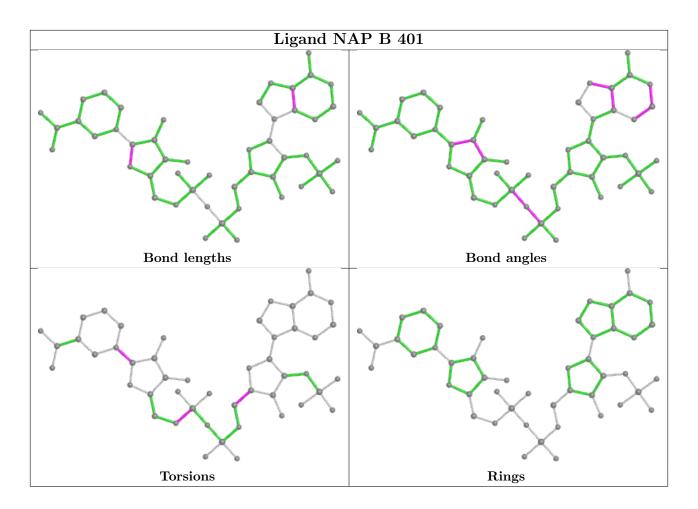
3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	С	401	NAP	1	0
2	В	401	NAP	2	0
2	А	401	NAP	6	0

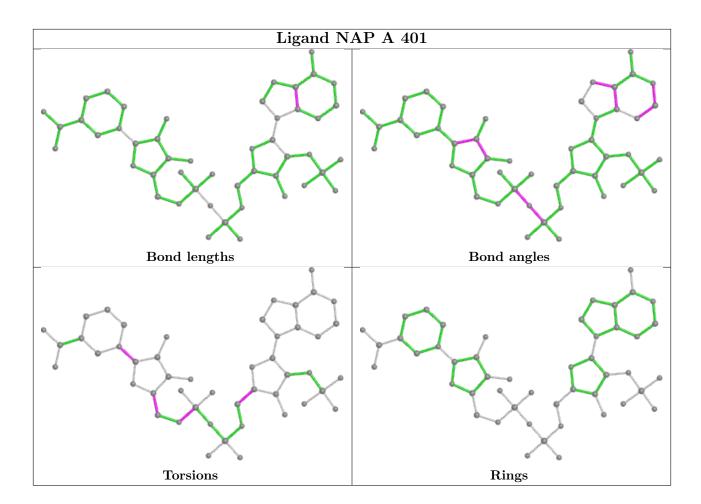
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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 sufficient 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 $>$	# RSRZ > 2	$OWAB(Å^2)$	Q<0.9
1	А	374/382~(97%)	0.18	3 (0%) 86 89	26, 38, 59, 85	0
1	В	374/382~(97%)	0.22	3 (0%) 86 89	25, 37, 59, 86	0
1	С	374/382~(97%)	0.16	5 (1%) 77 81	26, 38, 59, 85	0
All	All	1122/1146~(97%)	0.19	11 (0%) 82 86	25, 38, 59, 86	0

The worst 5 of 11 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	246	TYR	3.2
1	А	245	LEU	2.9
1	С	250	LYS	2.7
1	А	374	GLY	2.7
1	В	246	TYR	2.6

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 monosaccharides in this entry.

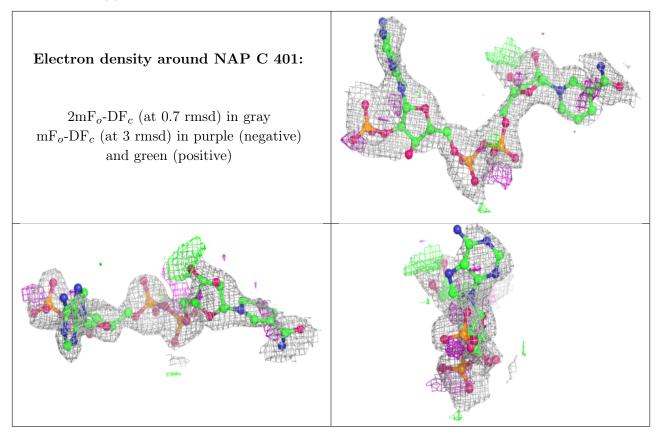
6.4 Ligands (i)

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^{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.



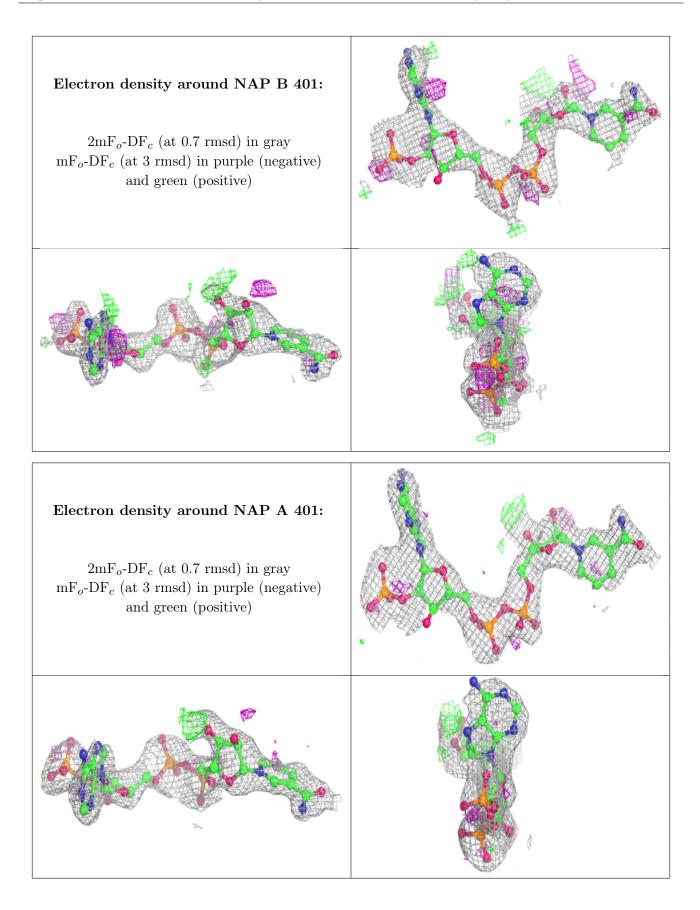
Mol	Type	Chain	Res	Atoms	RSCC	RSR	${f B} ext{-factors}({ m \AA}^2)$	Q < 0.9
2	NAP	С	401	48/48	0.82	0.27	36,73,87,107	0
2	NAP	В	401	48/48	0.83	0.27	40,72,85,103	0
2	NAP	А	401	48/48	0.85	0.26	42,73,88,113	0

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.

