

wwPDB X-ray Structure Validation Summary Report (i)

May 15, 2020 – 05:19 am BST

PDB ID : 4U7W

Title : The crystal structure of the terminal R domain from the myxalamid PKS-

NRPS biosynthetic pathway

Authors: Tsai, S.C.; Keasling, J.D.; Luo, R.; Barajas, J.F.; Phelan, R.M.; Schaub, A.J.;

Kliewer, J.

Deposited on : 2014-07-31

Resolution : 1.90 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 $\frac{\text{https://www.wwpdb.org/validation/2017/XrayValidationReportHelp}}{\text{with specific help available everywhere you see the (i) symbol.}$

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)

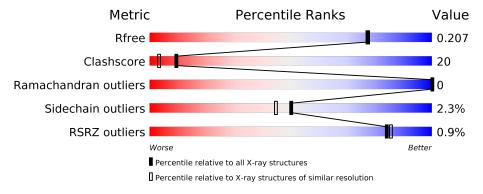
Validation Pipeline (wwPDB-VP) : 2.11

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 1.90 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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				
1	A	420	64%	27%	• 7%		
1	В	420	59%	34%	• 6%		

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
3	ACT	A	1602	-	-	X	-
3	ACT	В	1602	-	-	X	-



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6640 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 MxaA.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	A	390	Total 3001	C 1892				0	0	0
1	В	394	Total 3033	C 1912				0	0	0

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1094	MSE	-	initiating methionine	UNP Q93TX2
A	1095	GLY	_	expression tag	UNP Q93TX2
A	1096	SER	-	expression tag	UNP Q93TX2
A	1097	SER	-	expression tag	UNP Q93TX2
A	1098	HIS	-	expression tag	UNP Q93TX2
A	1099	HIS	-	expression tag	UNP Q93TX2
A	1100	HIS	-	expression tag	UNP Q93TX2
A	1101	HIS	-	expression tag	UNP Q93TX2
A	1102	HIS	-	expression tag	UNP Q93TX2
A	1103	HIS	-	expression tag	UNP Q93TX2
A	1104	SER	-	expression tag	UNP Q93TX2
A	1105	SER	-	expression tag	UNP Q93TX2
A	1106	GLY	-	expression tag	UNP Q93TX2
A	1107	LEU	-	expression tag	UNP Q93TX2
A	1108	VAL	-	expression tag	UNP Q93TX2
A	1109	PRO	-	expression tag	UNP Q93TX2
A	1110	ARG	-	expression tag	UNP Q93TX2
A	1111	GLY	-	expression tag	UNP Q93TX2
A	1112	SER	-	expression tag	UNP Q93TX2
A	1113	HIS	-	expression tag	UNP Q93TX2
A	1114	MSE	_	expression tag	UNP Q93TX2
В	1094	MSE	-	initiating methionine	UNP Q93TX2
В	1095	GLY	-	expression tag	UNP Q93TX2
В	1096	SER	-	expression tag	UNP Q93TX2
В	1097	SER	-	expression tag	UNP Q93TX2

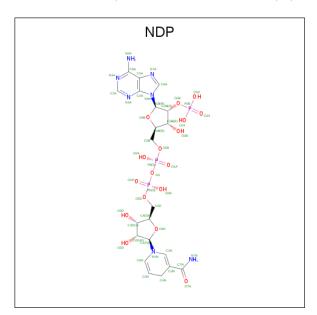
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Chain	Residue	Modelled	Actual	Comment	Reference
В	1098	HIS	-	expression tag	UNP Q93TX2
В	1099	HIS	-	expression tag	UNP Q93TX2
В	1100	HIS	1	expression tag	UNP Q93TX2
В	1101	HIS	1	expression tag	UNP Q93TX2
В	1102	HIS	-	expression tag	UNP Q93TX2
В	1103	HIS	-	expression tag	UNP Q93TX2
В	1104	SER	-	expression tag	UNP Q93TX2
В	1105	SER	-	expression tag	UNP Q93TX2
В	1106	GLY	-	expression tag	UNP Q93TX2
В	1107	LEU	1	expression tag	UNP Q93TX2
В	1108	VAL	-	expression tag	UNP Q93TX2
В	1109	PRO	-	expression tag	UNP Q93TX2
В	1110	ARG	-	expression tag	UNP Q93TX2
В	1111	GLY	-	expression tag	UNP Q93TX2
В	1112	SER	-	expression tag	UNP Q93TX2
В	1113	HIS	=	expression tag	UNP Q93TX2
В	1114	MSE	-	expression tag	UNP Q93TX2

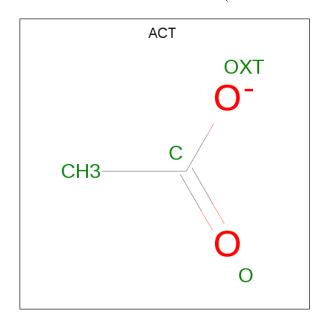
• Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
9	Λ	1	Total	С	N	О	Р	0	0
	2 A	1	48	21	7	17	3	0	
9	D	1	Total	С	N	О	Р	0	0
2	2 B	B 1		48	21	7	17	3	



 \bullet Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $\mathrm{C_2H_3O_2}).$



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

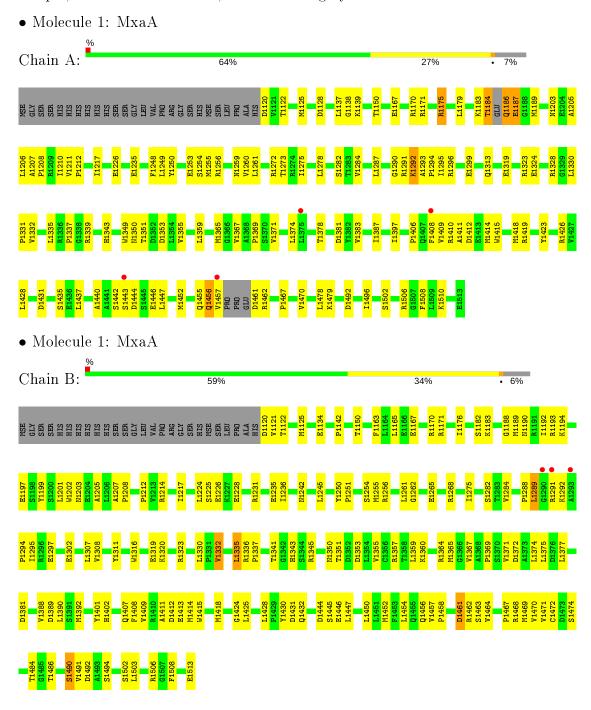
• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	269	Total O 269 269	0	0
4	В	233	Total O 233 233	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.





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	51.11Å 159.05Å 51.18Å	Danagitan
a, b, c, α , β , γ	90.00° 106.14° 90.00°	Depositor
Resolution (Å)	41.82 - 1.90	Depositor
Resolution (A)	41.82 - 1.90	EDS
% Data completeness	99.9 (41.82-1.90)	Depositor
(in resolution range)	98.5 (41.82-1.90)	EDS
R_{merge}	0.21	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) > 1$	$3.36 \; ({\rm at} \; 1.89 {\rm \AA})$	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
P. P.	0.158 , 0.203	Depositor
R, R_{free}	0.163 , 0.207	DCC
R_{free} test set	1995 reflections (3.24%)	wwPDB-VP
Wilson B-factor (Å ²)	23.2	Xtriage
Anisotropy	0.391	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.30 , 32.6	EDS
L-test for twinning ²	$< L >=0.41, < L^2>=0.23$	Xtriage
Estimated twinning fraction	0.287 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6640	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.60% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



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: NDP, ACT

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	Boı	nd lengths	Bond angles		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z >5	
1	A	0.42	1/3045~(0.0%)	0.54	1/4116 (0.0%)	
1	В	0.38	0/3081	0.53	0/4170	
All	All	0.40	$1/6126 \ (0.0\%)$	0.54	1/8286 (0.0%)	

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(\mathbf{\mathring{A}})$	$\mathbf{Ideal}(\mathbf{\mathring{A}})$
1	A	1294	PRO	N-CD	5.28	1.55	1.47

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
1	A	1293	ALA	C-N-CD	5.51	139.97	128.40

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	3001	0	3034	98	1
1	В	3033	0	3062	142	1
2	A	48	0	24	1	0
2	В	48	0	24	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	4	0	3	2	0
3	В	4	0	3	2	0
4	A	269	0	0	9	0
4	В	233	0	0	5	0
All	All	6640	0	6150	241	2

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 20.

The worst 5 of 241 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{aligned} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{aligned}$	$egin{array}{c} { m Clash} \\ { m overlap} \ ({ m \AA}) \end{array}$
1:B:1245:LEU:HD23	1:B:1255:MSE:HG2	1.42	1.01
1:A:1292:LYS:HE2	1:A:1292:LYS:HA	1.39	1.01
1:B:1407:GLN:NE2	1:B:1491:VAL:H	1.61	0.96
1:A:1291:ARG:NH1	1:A:1295:ILE:HG12	1.84	0.91
1:A:1282:SER:O	1:A:1337:PRO:HD2	1.69	0.91

All (2) 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	$egin{array}{c} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{array}$	$egin{array}{c} ext{Clash} \ ext{overlap } (ext{Å}) \end{array}$
1:B:1225:SER:OG	1:B:1513:GLU:OE2[1_454]	2.04	0.16
1:A:1211:VAL:CG1	1:A:1292:LYS:CG[1_556]	2.18	0.02

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	384/420 (91%)	373 (97%)	11 (3%)	0	100	100	
1	В	392/420 (93%)	380 (97%)	12 (3%)	0	100	100	

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	entiles
All	All	776/840 (92%)	753 (97%)	23 (3%)	0	100	100

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	325/341 (95%)	316 (97%)	9 (3%)	43 36
1	В	329/341 (96%)	323 (98%)	6 (2%)	59 55
All	All	654/682 (96%)	639 (98%)	15 (2%)	50 45

5 of 15 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1292	LYS
1	A	1456	GLN
1	В	1456	GLN
1	A	1187	GLU
1	В	1335	LEU

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

Mol	Chain	Res	Type
1	В	1407	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 carbohydrates in this entry.

5.6 Ligand geometry (i)

4 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	es Link	Bond lengths			Bond angles		
MIOI	Type		nes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NDP	A	1601	-	45,52,52	3.76	18 (40%)	53,80,80	1.36	3 (5%)
2	NDP	В	1601	-	45,52,52	3.74	17 (37%)	53,80,80	1.55	4 (7%)
3	ACT	В	1602	-	1,3,3	2.23	1 (100%)	0,3,3	0.00	-
3	ACT	A	1602	-	1,3,3	2.64	1 (100%)	0,3,3	0.00	-

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	NDP	A	1601	-	-	7/30/77/77	0/5/5/5
2	NDP	В	1601	-	-	6/30/77/77	0/5/5/5

The worst 5 of 37 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(\operatorname{\AA})$
2	A	1601	NDP	O4B-C1B	13.68	1.60	1.41
2	В	1601	NDP	O4B-C1B	13.07	1.59	1.41
2	В	1601	NDP	C6N-C5N	11.36	1.53	1.33
2	A	1601	NDP	C6N-C5N	11.18	1.53	1.33
2	В	1601	NDP	O4D-C1D	7.71	1.60	1.42

The worst 5 of 7 bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	В	1601	NDP	C1B-N9A-C4A	-6.67	114.92	126.64
2	A	1601	NDP	N3A-C2A-N1A	-5.76	119.67	128.68
2	В	1601	NDP	N3A-C2A-N1A	-5.47	120.14	128.68
2	A	1601	NDP	C1B-N9A-C4A	-5.22	117.48	126.64
2	В	1601	NDP	C5A-C6A-N6A	3.47	125.63	120.35

There are no chirality outliers.

5 of 13 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1601	NDP	C5D-O5D-PN-O1N
2	В	1601	NDP	O4D-C4D-C5D-O5D
2	В	1601	NDP	O4D-C1D-N1N-C6N
2	A	1601	NDP	O4D-C1D-N1N-C6N
2	В	1601	NDP	C3D-C4D-C5D-O5D

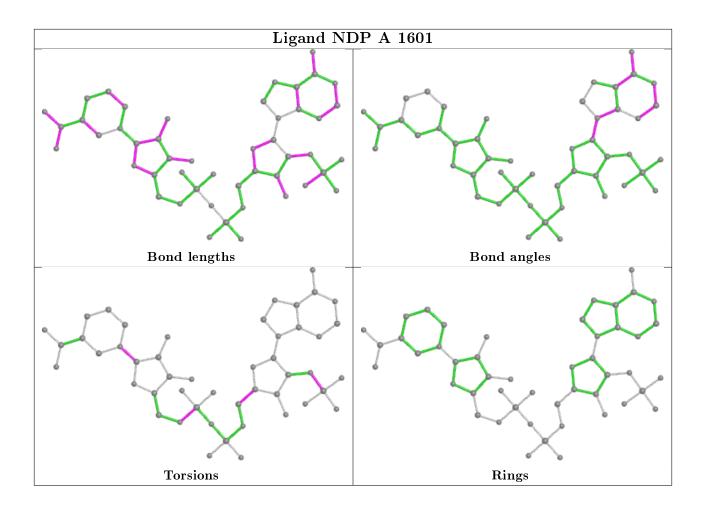
There are no ring outliers.

4 monomers are involved in 7 short contacts:

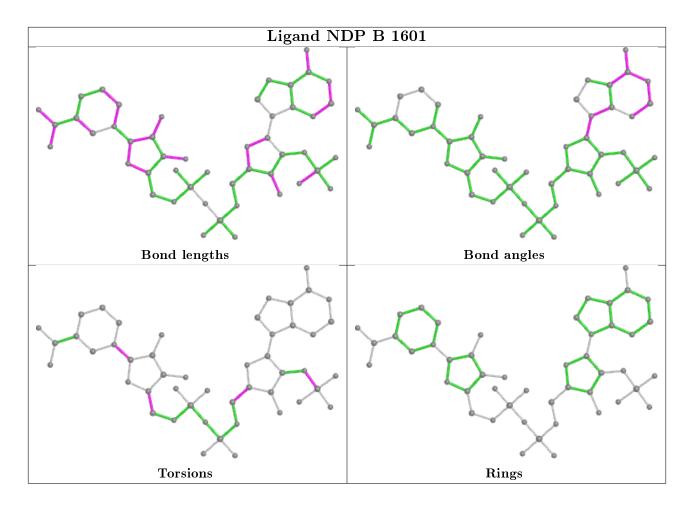
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1601	NDP	1	0
2	В	1601	NDP	2	0
3	В	1602	ACT	2	0
3	A	1602	ACT	2	0

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 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	$\begin{array}{c cccc} \textbf{Analysed} & <& RSRZ> & \#RSRZ>2 \\ \end{array}$		$OWAB(\AA^2)$	Q < 0.9		
1	A	381/420 (90%)	-0.46	4 (1%) 82	84	13, 24, 52, 100	0
1	В	385/420 (91%)	-0.40	3 (0%) 86	87	13, 26, 55, 92	0
All	All	766/840 (91%)	-0.43	7 (0%) 84	85	13, 25, 54, 100	0

The worst 5 of 7 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1408	PHE	3.1
1	A	1443	SER	3.0
1	A	1457	VAL	2.9
1	В	1293	ALA	2.6
1	В	1290	GLY	2.2

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.

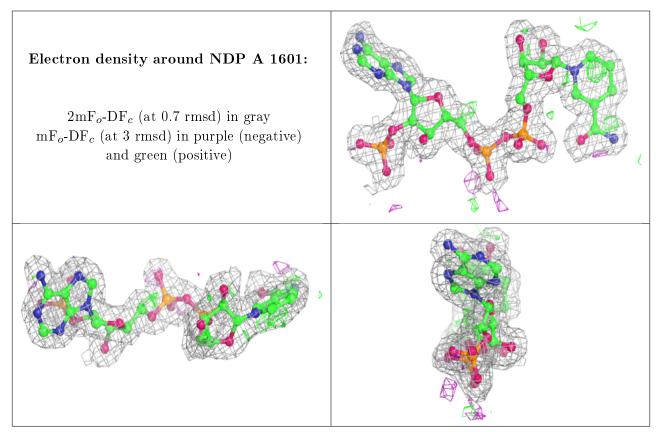
6.4 Ligands (i)

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.

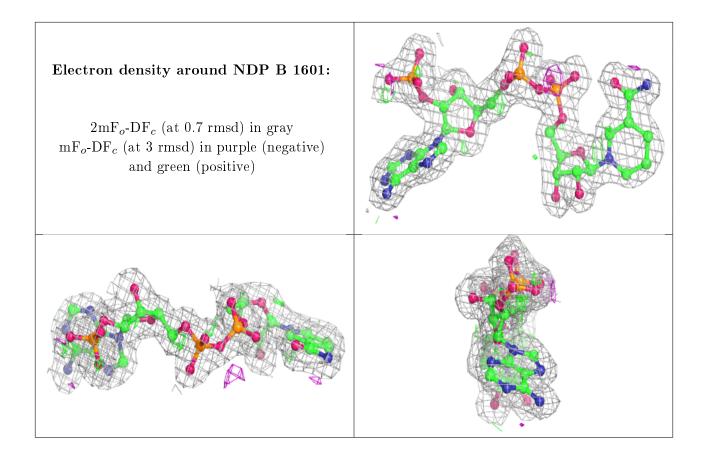


M	ol Type	Chain	Res	Atoms	RSCC	RSR	${f B-factors(\AA^2)}$	Q<0.9
3	ACT	A	1602	4/4	0.94	0.09	22,25,26,26	0
3	ACT	В	1602	4/4	0.95	0.07	20,22,23,24	0
2	NDP	A	1601	48/48	0.98	0.08	12,16,21,23	0
2	NDP	В	1601	48/48	0.98	0.08	14,18,22,23	0

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.

