



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 30, 2022 – 06:04 PM JST

PDB ID : 7F7L  
Title : Crystal structure of AKR4C17 bound with NADPH  
Authors : Li, H.; Yang, Y.; Hu, Y.; Chen, C.-C.; Huang, J.-W.; Min, J.; Dai, L.; Guo, R.-T.  
Deposited on : 2021-06-30  
Resolution : 2.25 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	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.29
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)	:	2.29

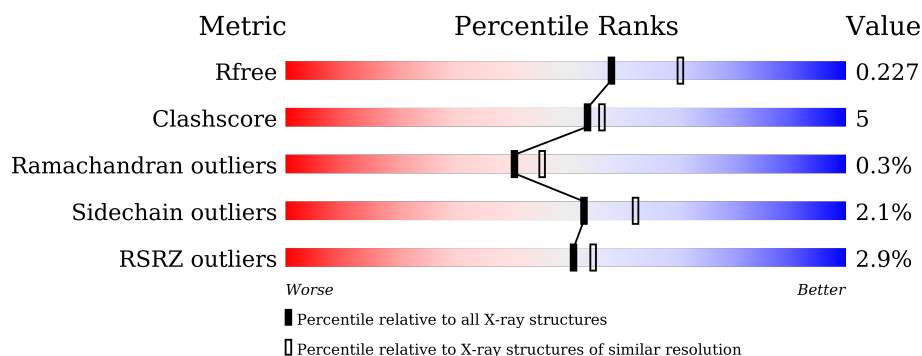
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.25 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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	321	<div> <div>3%</div> <div> <div></div> <div>84%</div> <div>11%</div> <div>..</div> </div> </div>
1	B	321	<div> <div>2%</div> <div> <div></div> <div>86%</div> <div>9%</div> <div>..</div> </div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5335 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 AKR4-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	309	Total	C	N	O	S	0	1	0
			2444	1561	423	453	7			
1	B	310	Total	C	N	O	S	0	0	0
			2442	1560	421	453	8			

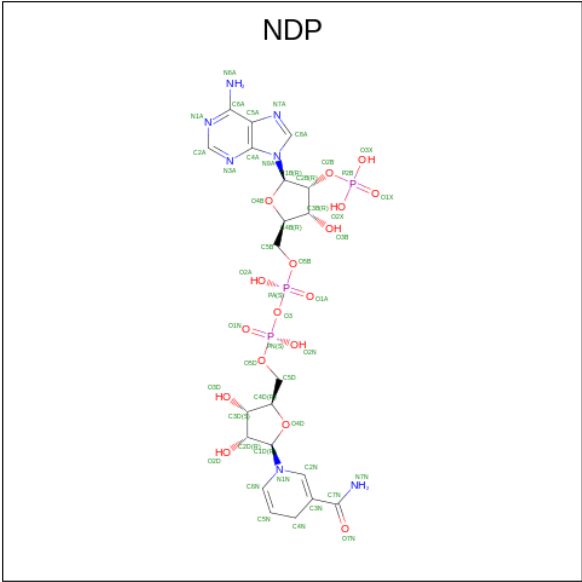
There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	GLY	-	expression tag	UNP A0A5J6VLZ7
A	-9	ALA	-	expression tag	UNP A0A5J6VLZ7
A	-8	GLY	-	expression tag	UNP A0A5J6VLZ7
A	-7	ALA	-	expression tag	UNP A0A5J6VLZ7
A	-6	GLY	-	expression tag	UNP A0A5J6VLZ7
A	-5	ALA	-	expression tag	UNP A0A5J6VLZ7
A	-4	GLY	-	expression tag	UNP A0A5J6VLZ7
A	-3	ALA	-	expression tag	UNP A0A5J6VLZ7
A	-2	GLY	-	expression tag	UNP A0A5J6VLZ7
A	-1	ALA	-	expression tag	UNP A0A5J6VLZ7
A	0	GLY	-	expression tag	UNP A0A5J6VLZ7
B	-10	GLY	-	expression tag	UNP A0A5J6VLZ7
B	-9	ALA	-	expression tag	UNP A0A5J6VLZ7
B	-8	GLY	-	expression tag	UNP A0A5J6VLZ7
B	-7	ALA	-	expression tag	UNP A0A5J6VLZ7
B	-6	GLY	-	expression tag	UNP A0A5J6VLZ7
B	-5	ALA	-	expression tag	UNP A0A5J6VLZ7
B	-4	GLY	-	expression tag	UNP A0A5J6VLZ7
B	-3	ALA	-	expression tag	UNP A0A5J6VLZ7
B	-2	GLY	-	expression tag	UNP A0A5J6VLZ7
B	-1	ALA	-	expression tag	UNP A0A5J6VLZ7
B	0	GLY	-	expression tag	UNP A0A5J6VLZ7

- Molecule 2 is COBALT (II) ION (three-letter code: CO) (formula: Co).

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

- Molecule 3 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: C<sub>21</sub>H<sub>30</sub>N<sub>7</sub>O<sub>17</sub>P<sub>3</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	158	Total	O	0	0
			158	158		
5	B	184	Total	O	0	0
			184	184		



- Molecule 1: AKR4-2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.43Å 77.43Å 452.75Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	24.82 – 2.25 24.99 – 2.25	Depositor EDS
% Data completeness (in resolution range)	88.6 (24.82-2.25) 84.2 (24.99-2.25)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.68 (at 2.26Å)	Xtriage
Refinement program	PHENIX 1.17.1	Depositor
R, $R_{free}$	0.181 , 0.226 0.182 , 0.227	Depositor DCC
$R_{free}$ test set	1870 reflections (5.49%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.0	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 39.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5335	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.82% 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: NDP, CO, SO4

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	0.43	0/2507	0.59	0/3410
1	B	0.43	0/2504	0.58	0/3405
All	All	0.43	0/5011	0.58	0/6815

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

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	2444	0	2420	27	0
1	B	2442	0	2426	25	0
2	A	1	0	0	0	0
3	A	48	0	26	4	0
3	B	48	0	26	6	0
4	B	10	0	0	0	0
5	A	158	0	0	4	0
5	B	184	0	0	0	0
All	All	5335	0	4898	52	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 5.



All (52) 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:84:HIS:HB2	1:B:115:ARG:HG3	1.45	0.96
1:A:87:GLU:OE2	5:A:501:HOH:O	2.01	0.78
1:B:207:SER:HA	3:B:403:NDP:O1A	1.84	0.76
1:A:37:LYS:HG2	1:A:67:VAL:HG11	1.70	0.73
1:A:229:LYS:NZ	5:A:503:HOH:O	2.24	0.71
1:A:37:LYS:HE2	1:A:67:VAL:HG21	1.73	0.70
1:A:33:TYR:CZ	1:A:37:LYS:HE3	2.29	0.67
1:B:33:TYR:CZ	1:B:37:LYS:HD2	2.31	0.65
1:B:115:ARG:HD2	1:B:128:ILE:HB	1.81	0.62
1:B:1:MET:HG2	1:B:3:ARG:H	1.64	0.61
1:A:220:GLU:HB2	1:A:223:ILE:HD12	1.82	0.61
1:A:24:ASP:HB3	1:A:25:PRO:CD	2.32	0.60
1:A:117:LYS:HG2	1:A:128:ILE:HD11	1.84	0.58
1:A:40:TYR:CE1	1:A:263:LEU:HG	2.41	0.55
1:A:33:TYR:OH	1:A:37:LYS:HE3	2.07	0.55
1:B:33:TYR:CE2	1:B:37:LYS:HD2	2.43	0.54
1:B:212:TRP:CZ3	1:B:213:MET:HG2	2.44	0.52
1:A:14:PRO:HB2	1:A:42:HIS:HB2	1.93	0.50
1:B:84:HIS:CB	1:B:115:ARG:HG3	2.31	0.50
1:A:258:ARG:HD2	3:A:402:NDP:C6A	2.41	0.49
1:B:28:VAL:HG11	1:B:51:ASN:HB2	1.93	0.49
1:B:54:GLU:OE2	1:B:54:GLU:N	2.39	0.49
1:B:202:TYR:HB2	3:B:403:NDP:O4D	2.12	0.49
1:B:25:PRO:HB3	1:B:54:GLU:OE1	2.12	0.49
1:A:253:SER:HB2	3:A:402:NDP:H8A	1.95	0.49
1:B:25:PRO:O	1:B:54:GLU:HG2	2.13	0.48
1:A:202:TYR:CZ	3:A:402:NDP:H41N	2.49	0.47
1:B:227:ALA:HB1	1:B:232:LYS:O	2.15	0.46
1:B:225:SER:O	1:B:228:GLU:HG3	2.16	0.46
1:B:202:TYR:CZ	3:B:403:NDP:H41N	2.52	0.45
1:A:80:TRP:CG	1:A:81:ASN:N	2.84	0.45
1:A:23:SER:HB3	1:A:28:VAL:HB	1.98	0.45
1:A:115:ARG:HG3	5:A:525:HOH:O	2.17	0.45
1:B:202:TYR:CE1	3:B:403:NDP:H41N	2.51	0.45
1:B:23:SER:OG	1:B:252:LYS:HE2	2.17	0.45
1:B:222:VAL:O	1:B:226:ILE:HG13	2.17	0.44
1:A:202:TYR:CE1	3:A:402:NDP:H41N	2.52	0.44
1:A:302:HIS:HB3	1:A:310:LEU:HD21	1.99	0.44
1:B:15:SER:HB2	1:B:243:ILE:HD12	2.00	0.43
1:A:69:LYS:HB2	1:A:69:LYS:HE2	1.64	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:ASP:HB3	1:A:25:PRO:HD2	2.00	0.43
1:B:52:GLU:OE2	1:B:77:SER:OG	2.28	0.42
1:A:22:GLN:N	5:A:509:HOH:O	2.46	0.42
1:A:26:GLY:H	1:A:54:GLU:CD	2.19	0.42
1:A:23:SER:CB	1:A:28:VAL:HB	2.49	0.42
1:A:121:ASN:HB3	1:B:270:ILE:O	2.20	0.42
1:B:107:LEU:HA	1:B:150:ALA:O	2.21	0.41
1:A:3:ARG:NH2	1:A:3:ARG:HG2	2.36	0.41
1:B:203:SER:H	3:B:403:NDP:H51N	1.85	0.41
3:B:403:NDP:H3D	3:B:403:NDP:H6N	2.03	0.40
1:A:204:PRO:HB2	1:A:239:LEU:HD23	2.03	0.40
1:B:111:HIS:O	1:B:154:SER:HB3	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	308/321 (96%)	301 (98%)	5 (2%)	2 (1%)	25	25
1	B	308/321 (96%)	301 (98%)	7 (2%)	0	100	100
All	All	616/642 (96%)	602 (98%)	12 (2%)	2 (0%)	41	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	211	THR
1	A	24	ASP

### 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	264/264 (100%)	258 (98%)	6 (2%)	50	59
1	B	264/264 (100%)	259 (98%)	5 (2%)	57	66
All	All	528/528 (100%)	517 (98%)	11 (2%)	53	62

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	149	ARG
1	A	202	TYR
1	A	216	ASN
1	A	228	GLU
1	A	229	LYS
1	A	297	SER
1	B	202	TYR
1	B	207	SER
1	B	213	MET
1	B	228	GLU
1	B	291	PHE

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

Mol	Chain	Res	Type
1	B	22	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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	SO4	B	401	-	4,4,4	0.20	0	6,6,6	0.36	0
3	NDP	B	403	-	45,52,52	1.09	4 (8%)	53,80,80	1.42	9 (16%)
4	SO4	B	402	-	4,4,4	0.14	0	6,6,6	0.26	0
3	NDP	A	402	-	45,52,52	1.06	3 (6%)	53,80,80	1.18	5 (9%)

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
3	NDP	B	403	-	-	5/30/77/77	0/5/5/5
3	NDP	A	402	-	-	7/30/77/77	0/5/5/5

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	NDP	C6N-C5N	3.16	1.39	1.33
3	B	403	NDP	C5A-C4A	2.95	1.48	1.40
3	B	403	NDP	C6N-C5N	2.88	1.38	1.33
3	A	402	NDP	C5A-C4A	2.75	1.48	1.40
3	B	403	NDP	P2B-O2B	2.28	1.63	1.59
3	B	403	NDP	O4B-C1B	2.09	1.44	1.41
3	A	402	NDP	O4B-C1B	2.03	1.43	1.41

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	403	NDP	O5B-C5B-C4B	3.60	121.38	108.99
3	B	403	NDP	O5B-PA-O1A	-2.99	97.38	109.07
3	A	402	NDP	N3A-C2A-N1A	-2.97	124.03	128.68
3	B	403	NDP	C4A-C5A-N7A	-2.96	106.32	109.40
3	A	402	NDP	C4A-C5A-N7A	-2.70	106.58	109.40
3	B	403	NDP	N3A-C2A-N1A	-2.65	124.53	128.68
3	B	403	NDP	PN-O3-PA	-2.34	124.79	132.83
3	B	403	NDP	O3D-C3D-C4D	2.34	117.81	111.05
3	B	403	NDP	C2A-N1A-C6A	2.32	122.73	118.75
3	A	402	NDP	O5D-C5D-C4D	2.26	116.77	108.99
3	B	403	NDP	O7N-C7N-C3N	-2.12	116.90	120.90
3	A	402	NDP	O3X-P2B-O1X	2.10	118.90	110.68
3	B	403	NDP	C1D-N1N-C2N	-2.10	117.62	121.11
3	A	402	NDP	C2A-N1A-C6A	2.05	122.27	118.75

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	403	NDP	C3B-C4B-C5B-O5B
3	B	403	NDP	O4B-C4B-C5B-O5B
3	B	403	NDP	O4D-C1D-N1N-C2N
3	A	402	NDP	PA-O3-PN-O5D
3	A	402	NDP	O4D-C1D-N1N-C2N
3	A	402	NDP	C4D-C5D-O5D-PN
3	A	402	NDP	C3B-C4B-C5B-O5B
3	A	402	NDP	O4B-C4B-C5B-O5B
3	A	402	NDP	C5D-O5D-PN-O3
3	A	402	NDP	PN-O3-PA-O2A
3	B	403	NDP	PN-O3-PA-O2A
3	B	403	NDP	PA-O3-PN-O2N

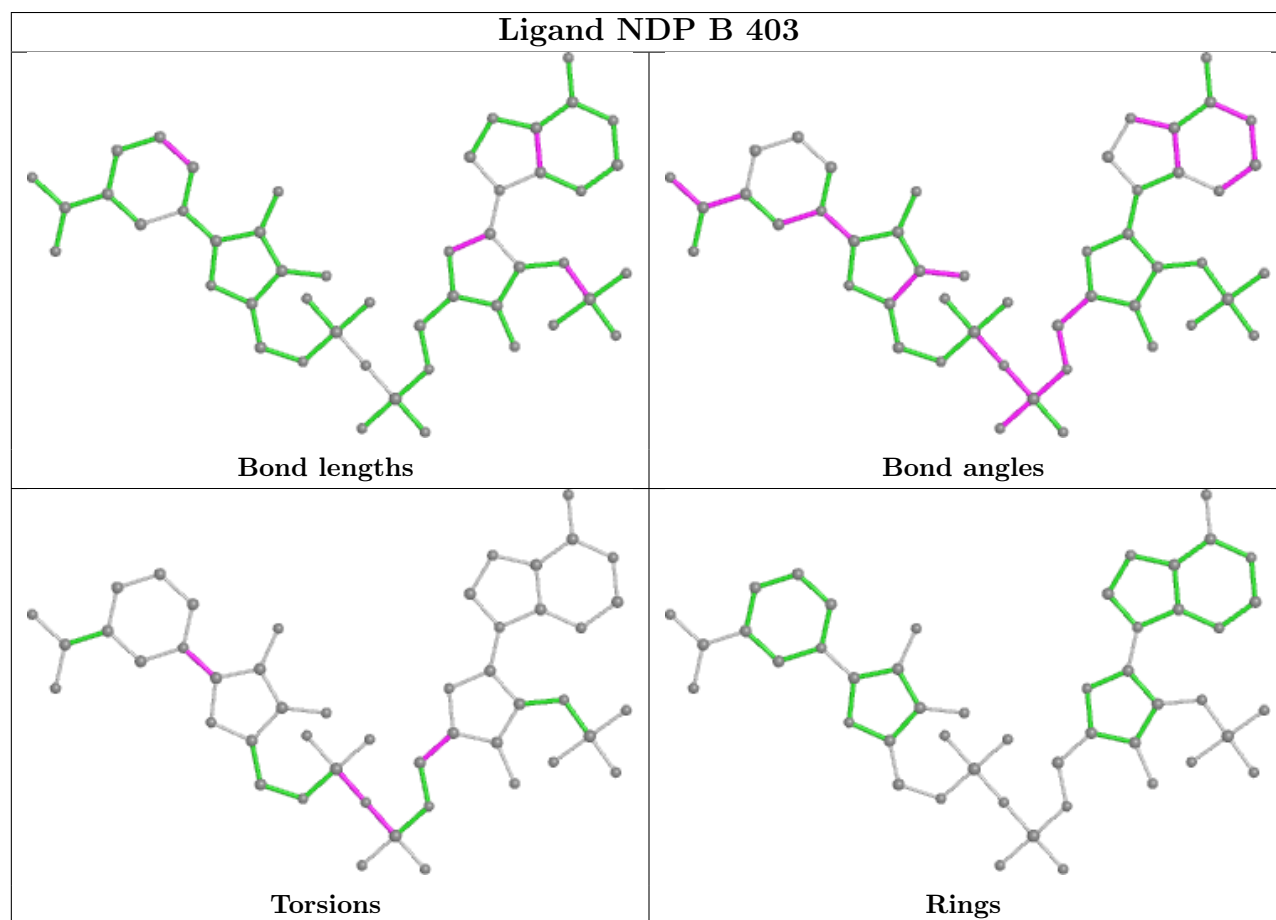
There are no ring outliers.

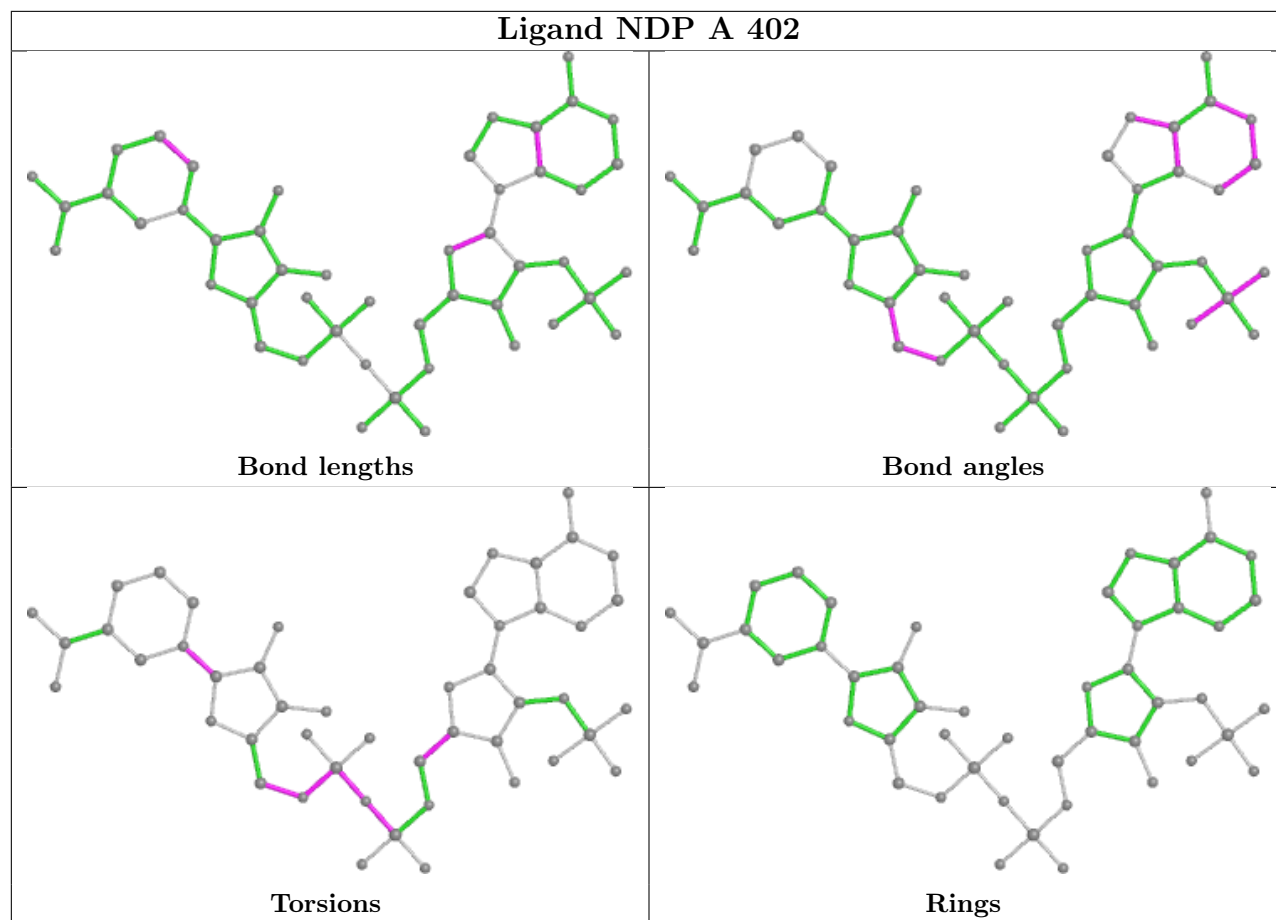
2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	403	NDP	6	0
3	A	402	NDP	4	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](#)

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	309/321 (96%)	-0.26	10 (3%) 47 50	16, 30, 53, 79	0
1	B	310/321 (96%)	-0.34	8 (2%) 56 59	14, 27, 51, 77	0
All	All	619/642 (96%)	-0.30	18 (2%) 51 55	14, 29, 53, 79	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	212	TRP	4.8
1	A	27	VAL	3.7
1	B	212	TRP	3.7
1	A	211	THR	3.6
1	B	211	THR	3.4
1	A	24	ASP	3.2
1	A	273	ASP	3.2
1	B	209	GLY	3.1
1	A	213	MET	2.9
1	A	310	LEU	2.8
1	B	308	GLY	2.8
1	A	25	PRO	2.6
1	B	214	ASN	2.5
1	A	23	SER	2.4
1	B	213	MET	2.4
1	B	257	GLU	2.3
1	B	50	GLY	2.3
1	A	22	GLN	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](#)

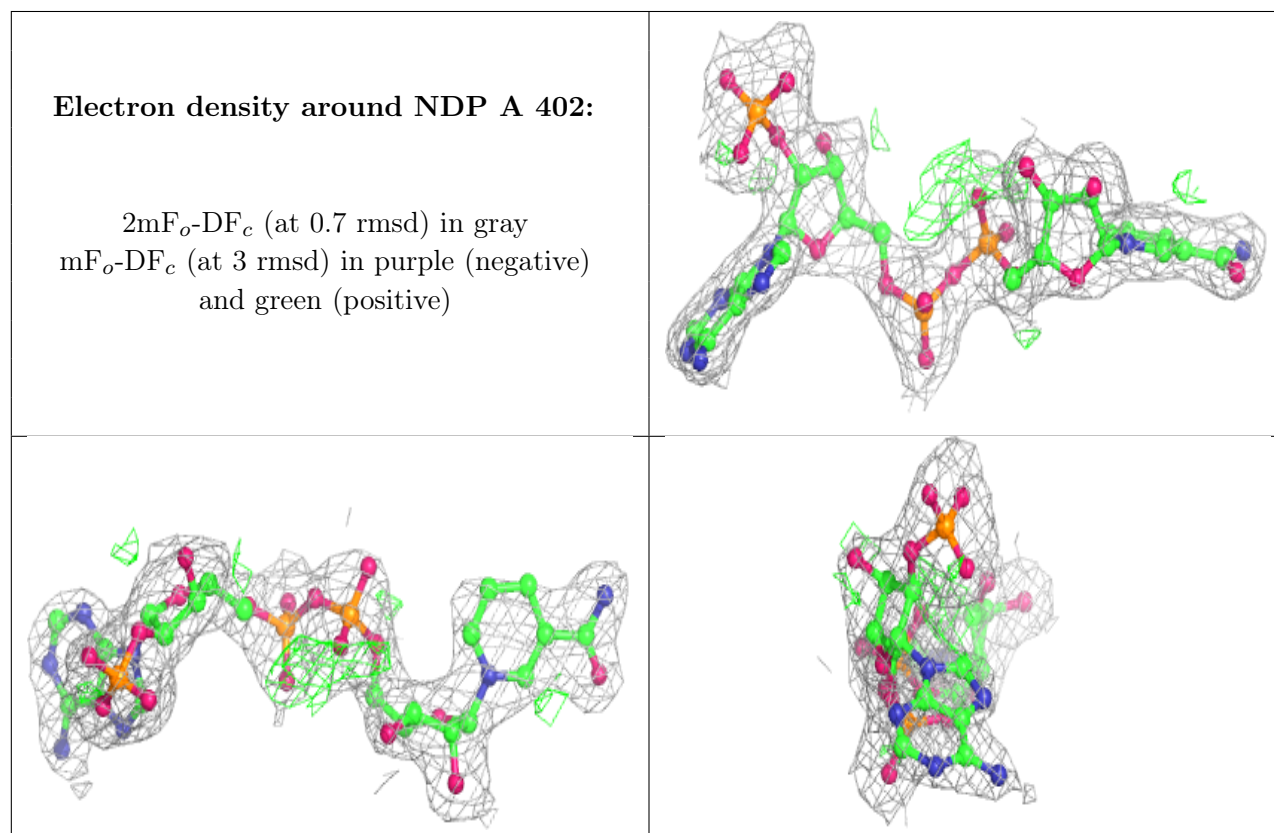
There are no monosaccharides in this entry.

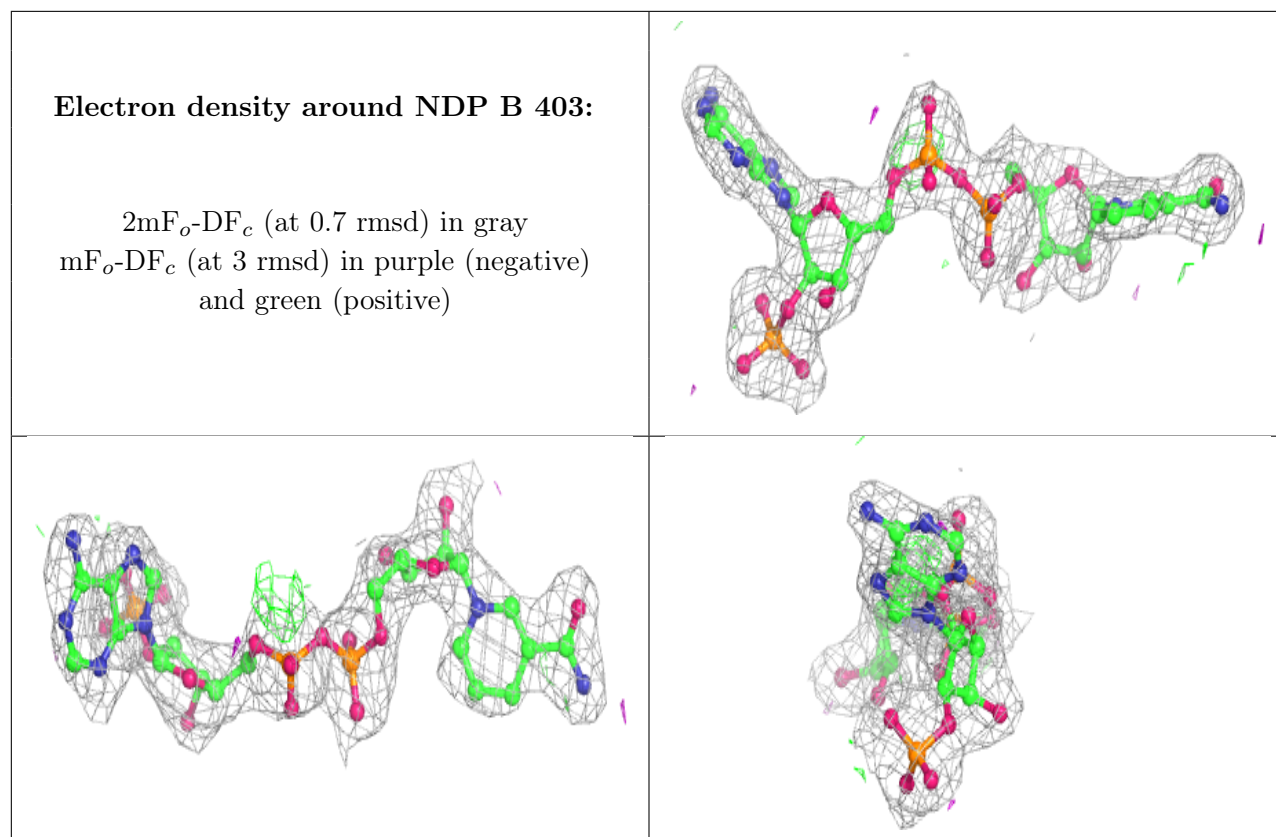
## 6.4 Ligands [i](#)

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<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
3	NDP	A	402	48/48	0.94	0.13	30,45,52,53	0
3	NDP	B	403	48/48	0.94	0.11	19,39,52,61	0
4	SO4	B	402	5/5	0.96	0.12	47,55,70,74	0
4	SO4	B	401	5/5	0.99	0.07	37,41,51,57	0
2	CO	A	401	1/1	0.99	0.10	15,15,15,15	1

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.