



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2017 – 08:48 pm GMT

PDB ID : 1KLK
Title : CRYSTAL STRUCTURE OF PNEUMOCYSTIS CARINII DIHYDROFOLATE REDUCTASE TERNARY COMPLEX WITH PT653 AND NADPH
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Deposited on : 2001-12-12
Resolution : 2.30 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

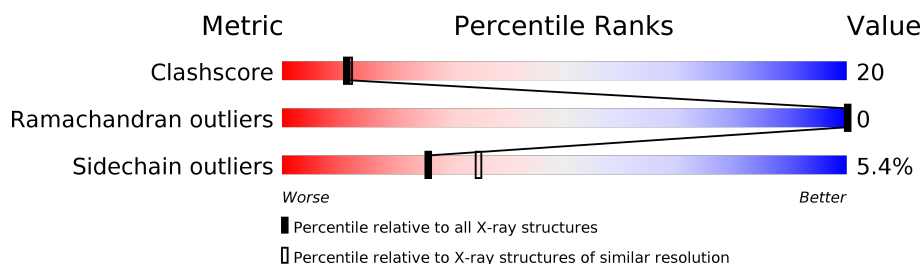
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	206	

2 Entry composition [i](#)

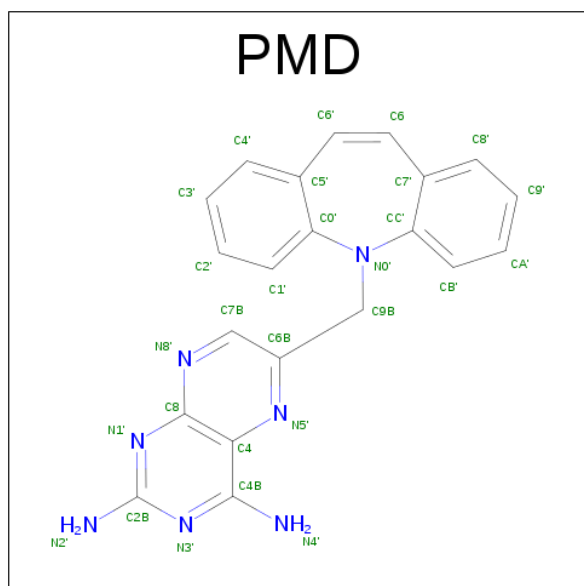
There are 4 unique types of molecules in this entry. The entry contains 1770 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Dihydrofolate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	S	0	0	0
			1661	1072	283	300	6			

- Molecule 2 is [N-(2,4-DIAMINOPTERIDIN-6-YL)-METHYL]-DIBENZ[B,F]AZEPINE (three-letter code: PMD) (formula: C₂₁H₁₇N₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			28	21	7		

- Molecule 3 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: C₂₁H₃₀N₇O₁₇P₃).



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

- Molecule 4 is water.

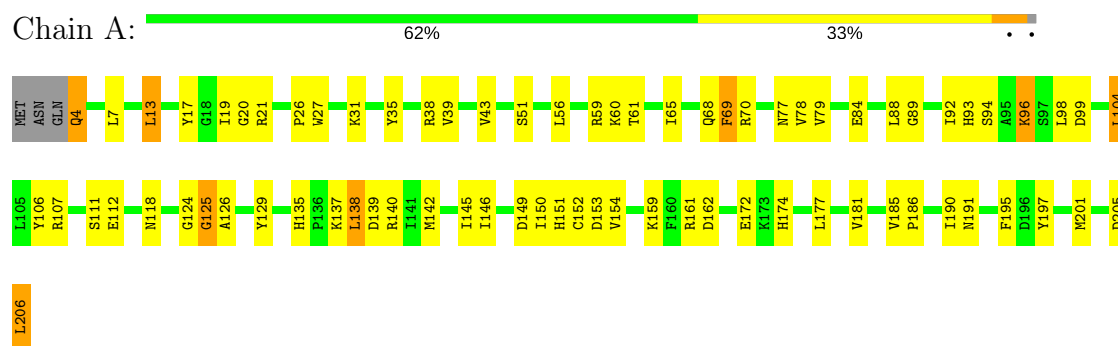
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	33	Total	O	0	0
			33	33		

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.

Note EDS was not executed.

- Molecule 1: Dihydrofolate reductase



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	36.84Å 43.14Å 59.32Å 90.00° 95.03° 90.00°	Depositor
Resolution (Å)	8.00 – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.196 , 0.289	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1770	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PMD, NDP

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	0.45	0/1703	0.77	2/2297 (0.1%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	206	LEU	CA-C-O	16.04	153.79	120.10
1	A	125	GLY	N-CA-C	-5.40	99.61	113.10

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	1661	0	1667	65	0
2	A	28	0	17	3	0
3	A	48	0	25	10	0
4	A	33	0	0	3	0
All	All	1770	0	1709	68	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 20.

All (68) 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:150:ILE:HD11	1:A:195:PHE:HE2	1.39	0.86
1:A:125:GLY:HA3	3:A:207:NDP:O1A	1.83	0.79
3:A:207:NDP:H52A	3:A:207:NDP:H8A	1.68	0.75
1:A:65:ILE:HB	1:A:70:ARG:HD3	1.68	0.75
1:A:21:ARG:HH11	1:A:21:ARG:HG2	1.51	0.73
1:A:142:MET:HE2	1:A:201:MET:HG3	1.71	0.71
1:A:124:GLY:HA2	1:A:129:TYR:CZ	2.24	0.70
1:A:186:PRO:HB2	1:A:190:ILE:HD11	1.74	0.69
1:A:51:SER:OG	1:A:118:ASN:HB2	1.94	0.68
1:A:60:LYS:HD2	3:A:207:NDP:O2N	1.94	0.68
1:A:191:ASN:HA	1:A:195:PHE:O	1.95	0.67
1:A:139:ASP:HB3	1:A:206:LEU:HD11	1.77	0.65
1:A:99:ASP:OD2	1:A:135:HIS:HE1	1.80	0.64
1:A:142:MET:HE1	1:A:201:MET:SD	2.39	0.62
1:A:78:VAL:HG22	1:A:93:HIS:HB2	1.82	0.62
1:A:107:ARG:HH11	1:A:107:ARG:HG2	1.65	0.61
1:A:4:GLN:HA	1:A:106:TYR:OH	2.03	0.59
1:A:153:ASP:OD2	1:A:154:VAL:HG23	2.02	0.59
1:A:26:PRO:HG2	1:A:27:TRP:CE3	2.37	0.59
1:A:56:LEU:HD23	1:A:98:LEU:CD2	2.35	0.56
1:A:21:ARG:NH1	1:A:21:ARG:HG2	2.20	0.56
1:A:174:HIS:HD2	1:A:177:LEU:HD23	1.71	0.56
1:A:19:ILE:O	3:A:207:NDP:H2N	2.06	0.55
1:A:181:VAL:O	1:A:181:VAL:HG12	2.05	0.55
1:A:174:HIS:NE2	1:A:185:VAL:HG23	2.22	0.55
1:A:125:GLY:HA3	3:A:207:NDP:PA	2.46	0.54
2:A:653:PMD:H2'1	3:A:207:NDP:O2D	2.07	0.54
1:A:20:GLY:O	1:A:152:CYS:HB3	2.10	0.52
1:A:124:GLY:HA2	1:A:129:TYR:CE1	2.44	0.52
1:A:145:ILE:O	1:A:197:TYR:HA	2.09	0.52
1:A:77:ASN:N	1:A:77:ASN:HD22	2.09	0.51
1:A:31:LYS:HB3	1:A:197:TYR:OH	2.10	0.51
1:A:146:ILE:N	1:A:146:ILE:HD12	2.26	0.51
1:A:142:MET:HE2	1:A:201:MET:CG	2.39	0.50
1:A:59:ARG:NH2	3:A:207:NDP:O3X	2.44	0.49
1:A:140:ARG:HG2	1:A:142:MET:HE3	1.95	0.49
1:A:159:LYS:HD3	1:A:162:ASP:OD2	2.13	0.49
1:A:38:ARG:HB3	4:A:671:HOH:O	2.12	0.48
1:A:174:HIS:CD2	1:A:177:LEU:HD23	2.49	0.47
1:A:35:TYR:O	1:A:39:VAL:HG23	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:LYS:O	3:A:207:NDP:H2A	2.16	0.46
1:A:84:GLU:OE2	1:A:96:LYS:HD3	2.16	0.46
2:A:653:PMD:N5'	3:A:207:NDP:H42N	2.30	0.46
1:A:89:GLY:HA3	1:A:92:ILE:HG12	1.98	0.46
1:A:17:TYR:OH	1:A:161:ARG:HB2	2.16	0.46
1:A:107:ARG:NH1	1:A:107:ARG:HG2	2.31	0.45
1:A:7:LEU:O	1:A:138:LEU:HD23	2.17	0.45
1:A:13:LEU:HB3	1:A:19:ILE:HG22	1.99	0.45
1:A:43:VAL:O	1:A:43:VAL:HG23	2.17	0.45
1:A:174:HIS:CE1	1:A:185:VAL:O	2.69	0.44
1:A:69:PHE:CD2	2:A:653:PMD:H8'1	2.52	0.44
1:A:68:GLN:HB2	1:A:68:GLN:HE21	1.67	0.43
1:A:126:ALA:HB3	3:A:207:NDP:O1N	2.19	0.43
1:A:174:HIS:HE1	1:A:185:VAL:O	2.02	0.43
1:A:186:PRO:HG3	4:A:670:HOH:O	2.19	0.43
1:A:104:LEU:HD22	1:A:107:ARG:HH22	1.84	0.42
1:A:150:ILE:HG22	1:A:151:HIS:N	2.34	0.42
1:A:98:LEU:HD23	1:A:98:LEU:HA	1.87	0.42
1:A:135:HIS:CD2	1:A:137:LYS:HG2	2.54	0.42
1:A:79:VAL:O	1:A:94:SER:HA	2.21	0.41
1:A:135:HIS:HD2	1:A:137:LYS:HB2	1.85	0.41
1:A:177:LEU:HD13	1:A:201:MET:HB2	2.01	0.41
1:A:142:MET:CE	1:A:201:MET:SD	3.08	0.40
1:A:61:THR:O	1:A:65:ILE:HG13	2.21	0.40
1:A:172:GLU:HG3	1:A:201:MET:HE3	2.02	0.40
1:A:21:ARG:NH1	1:A:21:ARG:CG	2.83	0.40
1:A:4:GLN:HB2	1:A:137:LYS:HE2	2.03	0.40
1:A:111:SER:HA	4:A:677:HOH:O	2.20	0.40

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	Percentiles	
1	A	201/206 (98%)	183 (91%)	18 (9%)	0	100	100

There are no Ramachandran outliers to report.

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	184/187 (98%)	174 (95%)	10 (5%)	26	35

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLN
1	A	13	LEU
1	A	69	PHE
1	A	88	LEU
1	A	96	LYS
1	A	104	LEU
1	A	112	GLU
1	A	138	LEU
1	A	149	ASP
1	A	205	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	68	GLN
1	A	116	GLN
1	A	127	GLN
1	A	135	HIS
1	A	174	HIS

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

2 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	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NDP	A	207	-	43,52,52	1.57	6 (13%)	49,80,80	1.85	7 (14%)
2	PMD	A	653	-	31,32,32	2.65	15 (48%)	40,46,46	2.27	14 (35%)

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	A	207	-	-	1/30/77/77	0/5/5/5
2	PMD	A	653	-	-	0/4/17/17	0/4/5/5

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	207	NDP	C4N-C5N	-2.50	1.43	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	207	NDP	C5A-N7A	-2.01	1.32	1.39
2	A	653	PMD	C9'-C8'	2.02	1.42	1.38
2	A	653	PMD	C4B-N3'	2.13	1.38	1.33
2	A	653	PMD	C0'-N0'	2.34	1.46	1.43
2	A	653	PMD	C6B-N5'	2.71	1.37	1.32
3	A	207	NDP	C2N-C3N	2.90	1.43	1.34
3	A	207	NDP	C2A-N3A	2.96	1.37	1.32
2	A	653	PMD	C7B-C6B	3.09	1.44	1.39
2	A	653	PMD	C7B-N8'	3.23	1.37	1.31
2	A	653	PMD	C5'-C0'	3.24	1.47	1.41
2	A	653	PMD	CB'-CC'	3.32	1.45	1.39
2	A	653	PMD	C1'-C0'	3.44	1.45	1.39
2	A	653	PMD	C7'-CC'	3.80	1.47	1.41
3	A	207	NDP	O4B-C1B	4.01	1.46	1.41
2	A	653	PMD	C7'-C6	4.04	1.53	1.46
2	A	653	PMD	C9B-C6B	4.29	1.58	1.51
2	A	653	PMD	C5'-C6'	4.51	1.54	1.46
3	A	207	NDP	C6N-C5N	4.66	1.41	1.33
2	A	653	PMD	C9B-N0'	4.79	1.55	1.47
2	A	653	PMD	CC'-N0'	5.20	1.51	1.43

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	207	NDP	N3A-C2A-N1A	-8.16	121.75	128.86
2	A	653	PMD	C9B-N0'-C0'	-6.23	108.00	119.07
2	A	653	PMD	N1'-C2B-N3'	-5.23	119.82	127.46
3	A	207	NDP	O4B-C4B-C5B	-3.21	98.56	109.40
3	A	207	NDP	O4B-C1B-C2B	-2.84	101.63	106.59
2	A	653	PMD	C7B-C6B-N5'	-2.60	119.09	120.80
2	A	653	PMD	C6B-C7B-N8'	-2.50	120.56	123.09
2	A	653	PMD	C9B-N0'-CC'	-2.04	115.44	119.07
2	A	653	PMD	CC'-N0'-C0'	2.04	121.10	116.79
2	A	653	PMD	N2'-C2B-N3'	2.18	120.72	117.24
3	A	207	NDP	O4D-C1D-N1N	2.36	112.81	108.07
2	A	653	PMD	C0'-C5'-C6'	3.08	126.08	122.94
2	A	653	PMD	CC'-C7'-C6	3.11	126.11	122.94
2	A	653	PMD	C7'-CC'-N0'	3.18	120.50	118.90
2	A	653	PMD	C9B-C6B-N5'	3.19	122.13	117.05
2	A	653	PMD	N8'-C8-N1'	3.30	120.31	116.04
2	A	653	PMD	C5'-C0'-N0'	3.44	120.64	118.90
3	A	207	NDP	O2B-C2B-C1B	3.51	123.14	110.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	207	NDP	C4A-C5A-N7A	3.58	112.87	109.41
3	A	207	NDP	O2B-C2B-C3B	3.70	125.33	111.63
2	A	653	PMD	C2B-N1'-C8	5.11	121.13	115.16

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	207	NDP	P2B-O2B-C2B-C3B

There are no ring outliers.

2 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	207	NDP	10	0
2	A	653	PMD	3	0

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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.