



wwPDB X-ray Structure Validation Summary Report

Feb 28, 2014 – 10:01 PM GMT

PDB ID : 3CSE
Title : Candida glabrata Dihydrofolate Reductase complexed with NADPH and 2,4-diamino-5-(3-(2,5-dimethoxyphenyl)prop-1-ynyl)-6-ethylpyrimidine(UCP120B)
Authors : Liu, J.; Anderson, A.C.
Deposited on : 2008-04-09
Resolution : 1.60 Å(reported)

This is a wwPDB validation summary 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/ValidationPDFNotes.html>

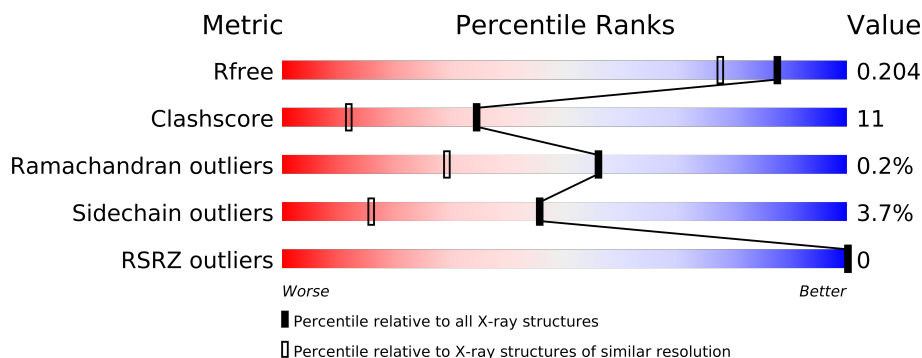
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance



The reported resolution of this entry is 1.60 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}	66092	1872 (1.60-1.60)
Clashscore	79885	2199 (1.60-1.60)
Ramachandran outliers	78287	2126 (1.60-1.60)
Sidechain outliers	78261	2125 (1.60-1.60)
RSRZ outliers	66119	1872 (1.60-1.60)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	227	
1	B	227	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	N22	A	302	X	-
3	N22	B	302	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4222 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

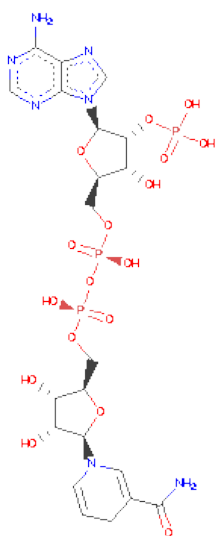
- Molecule 1 is a protein called Dihydrofolate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	225	Total	C	N	O	S	0	0	0
			1846	1180	325	334	7			
1	B	225	Total	C	N	O	S	0	0	0
			1846	1180	325	334	7			

There are 20 discrepancies between the modelled and reference sequences:

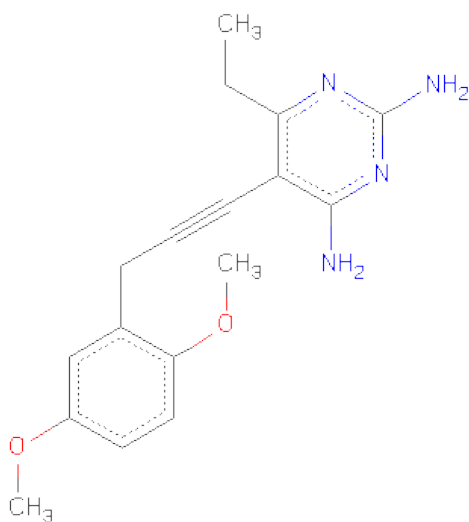
Chain	Residue	Modelled	Actual	Comment	Reference
A	218	LEU	-	EXPRESSION TAG	UNP Q6FPH0
A	219	GLU	-	EXPRESSION TAG	UNP Q6FPH0
A	220	HIS	-	EXPRESSION TAG	UNP Q6FPH0
A	221	HIS	-	EXPRESSION TAG	UNP Q6FPH0
A	222	HIS	-	EXPRESSION TAG	UNP Q6FPH0
A	223	HIS	-	EXPRESSION TAG	UNP Q6FPH0
A	224	HIS	-	EXPRESSION TAG	UNP Q6FPH0
A	225	HIS	-	EXPRESSION TAG	UNP Q6FPH0
A	226	HIS	-	EXPRESSION TAG	UNP Q6FPH0
A	227	HIS	-	EXPRESSION TAG	UNP Q6FPH0
B	218	LEU	-	EXPRESSION TAG	UNP Q6FPH0
B	219	GLU	-	EXPRESSION TAG	UNP Q6FPH0
B	220	HIS	-	EXPRESSION TAG	UNP Q6FPH0
B	221	HIS	-	EXPRESSION TAG	UNP Q6FPH0
B	222	HIS	-	EXPRESSION TAG	UNP Q6FPH0
B	223	HIS	-	EXPRESSION TAG	UNP Q6FPH0
B	224	HIS	-	EXPRESSION TAG	UNP Q6FPH0
B	225	HIS	-	EXPRESSION TAG	UNP Q6FPH0
B	226	HIS	-	EXPRESSION TAG	UNP Q6FPH0
B	227	HIS	-	EXPRESSION TAG	UNP Q6FPH0

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDEPHOSPHATE (three-letter code: NDP) (formula: C₂₁H₃₀N₇O₁₇P₃).



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

- Molecule 3 is 5-[3-(2,5-DIMETHOXYPHENYL)PROP-1-YN-1-YL]-6-ETHYLPYRIMIDIN E-2,4-DIAMINE (three-letter code: N22) (formula: $C_{17}H_{20}N_4O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			23	17	4	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			23	17	4	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	194	Total	O	0	0
			194	194		
4	B	194	Total	O	0	0
			194	194		

4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	42.69Å 42.69Å 230.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.03 – 1.60 40.03 – 1.60	Depositor EDS
% Data completeness (in resolution range)	93.2 (40.03-1.60) 93.2 (40.03-1.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.181 , 0.232 0.172 , 0.204	Depositor DCC
R_{free} test set	2558 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 39.1	EDS
Estimated twinning fraction	0.479 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 50234 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4222	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.73% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NDP, N22

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	1.63	17/1891 (0.9%)	1.57	21/2556 (0.8%)
1	B	1.65	14/1891 (0.7%)	1.51	22/2556 (0.9%)
All	All	1.64	31/3782 (0.8%)	1.54	43/5112 (0.8%)

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	A	0	1
1	B	0	3
All	All	0	4

The worst 5 of 31 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	191	GLU	CD-OE1	8.18	1.34	1.25
1	A	89	GLU	CD-OE2	7.36	1.33	1.25
1	A	127	TYR	CD1-CE1	6.91	1.49	1.39
1	B	89	GLU	CG-CD	6.76	1.62	1.51
1	A	207	TYR	CG-CD1	6.58	1.47	1.39

The worst 5 of 43 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	128	ARG	NE-CZ-NH2	15.88	128.24	120.30
1	A	37	ARG	NE-CZ-NH1	14.56	127.58	120.30
1	A	37	ARG	NE-CZ-NH2	-13.44	113.58	120.30
1	A	128	ARG	NE-CZ-NH1	-13.03	113.78	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	99	ARG	NE-CZ-NH1	10.54	125.57	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	83	GLY	Peptide
1	B	185	GLY	Peptide
1	B	81	PHE	Peptide
1	B	83	GLY	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1846	0	1846	45	0
1	B	1846	0	1846	35	3
2	A	48	0	26	1	0
2	B	48	0	26	1	0
3	A	23	0	20	2	0
3	B	23	0	20	2	0
4	A	194	0	0	21	5
4	B	194	0	0	18	1
All	All	4222	0	3784	82	6

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

The worst 5 of 82 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:190:GLN:HG2	4:A:529:HOH:O	1.17	1.25
1:A:183:LEU:HD12	1:A:200:LEU:CD1	1.84	1.07
1:B:224:HIS:HE1	4:B:447:HOH:O	1.36	1.07
1:A:224:HIS:HE1	4:A:485:HOH:O	1.39	1.03
1:A:219:GLU:OE1	4:A:542:HOH:O	1.80	0.99

The worst 5 of 6 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	Distance(Å)	Clash(Å)
1:B:71:LYS:NZ	4:A:490:HOH:O[1_565]	1.87	0.33
1:B:219:GLU:OE1	4:A:527:HOH:O[1_655]	1.95	0.25
4:B:476:HOH:O	4:B:527:HOH:O[1_545]	2.03	0.17
4:A:460:HOH:O	4:A:478:HOH:O[1_655]	2.10	0.10
1:B:87:LYS:NZ	4:A:542:HOH:O[1_565]	2.16	0.04

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	223/227 (98%)	218 (98%)	5 (2%)	0	100	100
1	B	223/227 (98%)	217 (97%)	5 (2%)	1 (0%)	43	18
All	All	446/454 (98%)	435 (98%)	10 (2%)	1 (0%)	56	29

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	185	GLY

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	205/207 (99%)	197 (96%)	8 (4%)	43	14
1	B	205/207 (99%)	198 (97%)	7 (3%)	49	19
All	All	410/414 (99%)	395 (96%)	15 (4%)	45	16

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

Mol	Chain	Res	Type
1	A	196	LEU
1	A	216	LYS
1	B	145	LEU
1	A	182	GLN
1	B	101	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 15 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	208	GLN
1	A	220	HIS
1	B	136	HIS
1	A	203	GLN
1	B	113	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

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	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NDP	A	301	-	52,52,52	1.59	9 (17%)	80,80,80	2.37	13 (16%)
3	N22	A	302	-	24,24,24	2.64	9 (37%)	32,32,32	4.27	18 (56%)
2	NDP	B	301	-	52,52,52	1.55	10 (19%)	80,80,80	2.35	9 (11%)
3	N22	B	302	-	24,24,24	2.84	15 (62%)	32,32,32	4.18	19 (59%)

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	301	-	-	0/35/77/77	0/3/5/5
3	N22	A	302	-	-	1/12/12/12	0/2/2/2
2	NDP	B	301	-	-	0/35/77/77	0/3/5/5
3	N22	B	302	-	-	0/12/12/12	0/2/2/2

The worst 5 of 43 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	N22	C18-C17	6.07	1.50	1.38
3	B	302	N22	C11-C10	-5.64	1.42	1.47
2	B	301	NDP	O4B-C1B	5.14	1.49	1.41
3	B	302	N22	C6-C1	5.07	1.47	1.42
2	B	301	NDP	PN-O3	-4.60	1.51	1.59

The worst 5 of 59 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	NDP	N3A-C2A-N1A	-16.29	115.09	128.71
2	B	301	NDP	N3A-C2A-N1A	-16.05	115.29	128.71
3	A	302	N22	C3-N4-C5	11.16	124.07	115.99
3	B	302	N22	C18-C13-C14	9.34	129.53	118.43
3	B	302	N22	C3-N4-C5	8.91	122.44	115.99

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	302	N22	C1-C6-C9-C10

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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, 95th 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	A	225/227 (99%)	-0.47	0 100 100	15, 23, 37, 46	0
1	B	225/227 (99%)	-0.45	0 100 100	15, 23, 37, 46	1 (0%)
All	All	450/454 (99%)	-0.46	0 100 100	15, 23, 37, 46	1 (0%)

There are no RSRZ outliers to report.

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	N22	A	302	23/23	0.09	1.46	14,21,33,40	0
3	N22	B	302	23/23	0.08	0.84	15,22,32,38	0
2	NDP	B	301	48/48	0.06	-0.66	15,18,23,26	0
2	NDP	A	301	48/48	0.06	-0.73	14,18,23,25	0

6.5 Other polymers ⓘ

There are no such residues in this entry.