



wwPDB X-ray Structure Validation Summary Report

Feb 28, 2014 – 10:38 PM GMT

PDB ID : 1E6W
Title : RAT BRAIN 3-HYDROXYACYL-COA DEHYDROGENASE BINARY
COMPLEX WITH NADH AND ESTRADIOL
Authors : Read, J.A.; Powell, A.J.; Brady, R.L.
Deposited on : 2000-08-23
Resolution : 1.70 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>

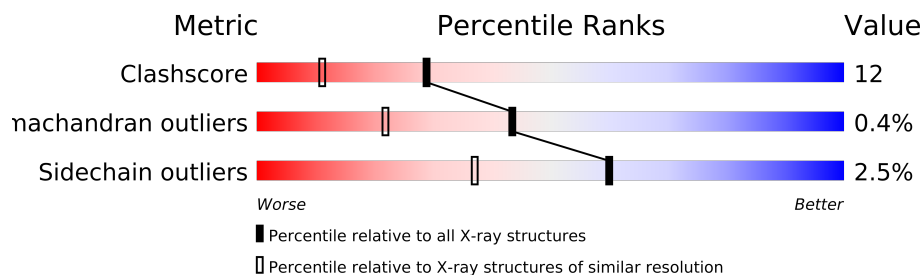
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
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 1.70 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	79885	2929 (1.70-1.70)
Ramachandran outliers	78287	2878 (1.70-1.70)
Sidechain outliers	78261	2878 (1.70-1.70)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	260	
1	B	260	
1	C	260	
1	D	260	

2 Entry composition i

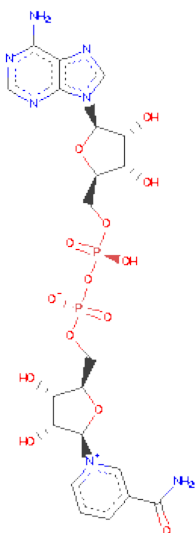
There are 4 unique types of molecules in this entry. The entry contains 8106 atoms, of which 0 are hydrogen and 0 are deuterium.

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 SHORT CHAIN 3-HYDROXYACYL-COA DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	0	8	0
			1841	1160	323	351	7			
1	B	247	Total	C	N	O	S	0	11	0
			1848	1167	328	347	6			
1	C	248	Total	C	N	O	S	0	5	0
			1820	1147	324	342	7			
1	D	255	Total	C	N	O	S	0	2	0
			1856	1172	329	349	6			

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



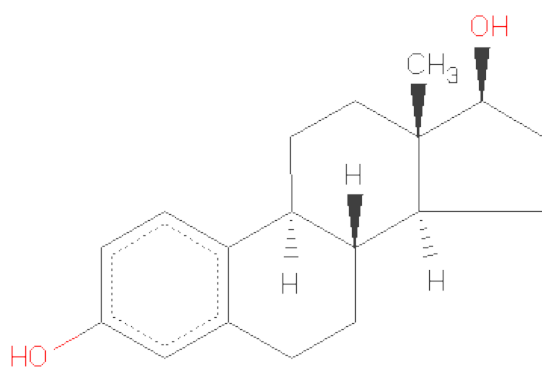
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is ESTRADIOL (three-letter code: EST) (formula: $C_{18}H_{24}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	L	1	Total	C	O	0	0
			20	18	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	146	Total	O	0	0
			146	146		
4	B	128	Total	O	0	0
			128	128		
4	C	143	Total	O	0	0
			143	143		
4	D	128	Total	O	0	0
			128	128		

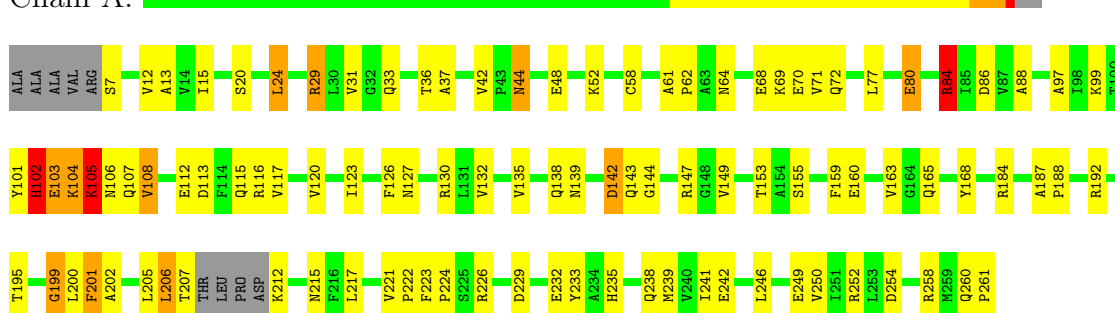
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

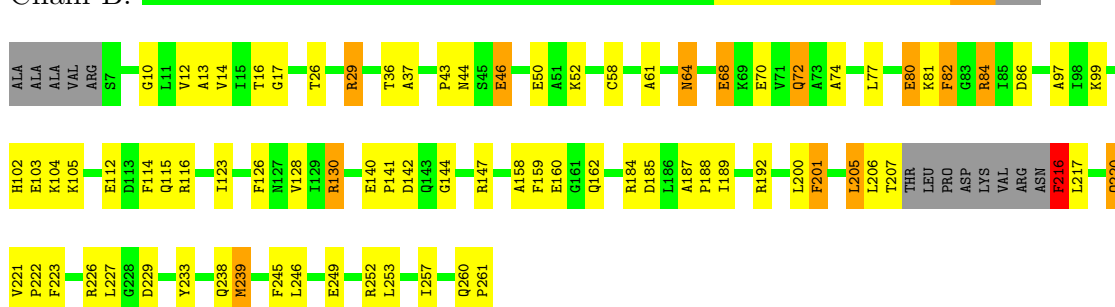
• Molecule 1: SHORT CHAIN 3-HYDROXYACYL-COA DEHYDROGENASE

Chain A:



• Molecule 1: SHORT CHAIN 3-HYDROXYACYL-COA DEHYDROGENASE

Chain B:



• Molecule 1: SHORT CHAIN 3-HYDROXYACYL-COA DEHYDROGENASE

Chain C:



• Molecule 1: SHORT CHAIN 3-HYDROXYACYL-COA DEHYDROGENASE

Chain D: 

ALA	ALA	ALA	VAL	ARG	S7	S20	S25	R29	L30	Q33	T36	D41	V42	P43	M44	S45	E46	G47	E48	T49	E50	A51	L54	N57	K69	E70	V71	K79	E80	R84	I85	D86	V87	A88	C91	A92	A95	K99	T100	Y101	H102	E103	K104	N105	M106
Q107	V108	E112	D113	F114	Q115	R116	M119	I123	F126	R130	L131	V132	V135	M136	G137	Q138	D142	R147	I150	A154	F159	E160	G161	Q162	A167	Y168	S169	K172	I182	A183	R184	D185	P188	I189	G190	I191	R192	T195	P198	G199	L200	F201			
A202	T203	P204	L205	L206	T207	T208	L209	P210	D211	R214	N215	F216	S219	F223	P224	S225	R226	D229	P230	A231	E232	Y233	Q238	M239	N243	P244	F245	L246	V250	I251	R252	L253	D254	G255	A256	I257	R258	F259	Q260	P261					

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	57.60Å 67.43Å 67.50Å 65.22° 73.27° 75.67°	Depositor
Resolution (Å)	30.00 – 1.70	Depositor
% Data completeness (in resolution range)	94.4 (30.00-1.70)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.184 , 0.229	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8106	wwPDB-VP
Average B, all atoms (Å ²)	20.0	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: EST, NAD

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	1.45	9/1906 (0.5%)	2.31	82/2589 (3.2%)
1	B	1.35	3/1928 (0.2%)	2.19	69/2612 (2.6%)
1	C	1.54	8/1869 (0.4%)	2.31	83/2534 (3.3%)
1	D	1.42	9/1895 (0.5%)	2.31	75/2574 (2.9%)
All	All	1.44	29/7598 (0.4%)	2.28	309/10309 (3.0%)

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	9
1	B	0	10
1	C	0	10
1	D	0	10
All	All	0	39

The worst 5 of 29 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	232	GLU	CD-OE2	7.19	1.33	1.25
1	D	25	SER	CA-CB	6.47	1.62	1.52
1	D	255	GLY	N-CA	6.32	1.55	1.46
1	A	101	TYR	CE1-CZ	-6.12	1.30	1.38
1	A	199	GLY	CA-C	6.10	1.61	1.51

The worst 5 of 309 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	84	ARG	CD-NE-CZ	26.25	160.34	123.60
1	D	29	ARG	NE-CZ-NH1	21.82	131.21	120.30
1	A	116	ARG	NE-CZ-NH2	-21.48	109.56	120.30
1	D	252	ARG	NE-CZ-NH1	21.32	130.96	120.30
1	D	226	ARG	NE-CZ-NH2	-20.07	110.27	120.30

There are no chirality outliers.

5 of 39 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	102	HIS	Mainchain
1	A	105	LYS	Mainchain
1	A	120	VAL	Mainchain
1	A	24	LEU	Mainchain
1	A	7	SER	Mainchain

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	1841	0	1864	54	2
1	B	1848	0	1916	48	4
1	C	1820	0	1872	58	2
1	D	1856	0	1889	42	6
2	A	44	0	26	0	0
2	B	44	0	26	0	0
2	C	44	0	26	2	0
2	D	44	0	26	1	0
3	L	20	0	21	11	0
4	A	146	0	0	15	1
4	B	128	0	0	12	3
4	C	143	0	0	14	3
4	D	128	0	0	11	1
All	All	8106	0	7666	190	11

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 12.

The worst 5 of 190 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:239:MET:HG2	4:D:2109:HOH:O	1.63	0.98
1:A:29:ARG:HE	1:A:238[B]:GLN:NE2	1.62	0.95
1:D:29:ARG:HE	1:D:238[B]:GLN:NE2	1.64	0.94
1:D:29:ARG:HH21	1:D:238[B]:GLN:HE22	0.96	0.94
1:D:29:ARG:HE	1:D:238[B]:GLN:HE21	0.94	0.93

The worst 5 of 11 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:44:ASN:ND2	1:D:49:THR:OG1[1_565]	1.39	0.81
1:B:44:ASN:CG	1:D:49:THR:OG1[1_565]	1.90	0.30
4:B:2037:HOH:O	4:C:2026:HOH:O[1_564]	1.93	0.27
1:D:44:ASN:CB	4:B:2056:HOH:O[1_545]	1.94	0.26
1:B:44:ASN:ND2	1:D:49:THR:CB[1_565]	2.01	0.19

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	255/260 (98%)	245 (96%)	10 (4%)	0	100	100
1	B	254/260 (98%)	241 (95%)	13 (5%)	0	100	100
1	C	249/260 (96%)	237 (95%)	12 (5%)	0	100	100
1	D	255/260 (98%)	239 (94%)	12 (5%)	4 (2%)	14	2
All	All	1013/1040 (97%)	962 (95%)	47 (5%)	4 (0%)	43	22

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	211	ASP
1	D	207	THR

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Mol	Chain	Res	Type
1	D	210	PRO
1	D	206	LEU

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	193/199 (97%)	186 (96%)	7 (4%)	47	22
1	B	199/199 (100%)	197 (99%)	2 (1%)	85	76
1	C	191/199 (96%)	186 (97%)	5 (3%)	59	35
1	D	192/199 (96%)	186 (97%)	6 (3%)	52	27
All	All	775/796 (97%)	755 (97%)	20 (3%)	60	35

5 of 20 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	7	SER
1	C	44	ASN
1	D	104	LYS
1	B	216	PHE
1	B	220	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	44	ASN
1	D	215	ASN
1	C	106	ASN
1	A	143	GLN
1	C	102	HIS

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 ⓘ

5 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	NAD	A	301	-	48,48,48	1.81	12 (25%)	73,73,73	2.71	29 (39%)
2	NAD	B	301	-	48,48,48	1.53	9 (18%)	73,73,73	2.26	23 (31%)
2	NAD	C	301	-	48,48,48	1.88	17 (35%)	73,73,73	2.37	25 (34%)
2	NAD	D	301	-	48,48,48	1.56	8 (16%)	73,73,73	2.25	22 (30%)
3	EST	L	350	-	23,23,23	2.47	3 (13%)	36,36,36	2.79	17 (47%)

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	NAD	A	301	-	-	0/30/62/62	0/3/5/5
2	NAD	B	301	-	-	0/30/62/62	0/3/5/5
2	NAD	C	301	-	-	0/30/62/62	0/3/5/5
2	NAD	D	301	-	-	0/30/62/62	0/3/5/5
3	EST	L	350	-	2/2/5/5	0/0/40/40	0/0/4/4

The worst 5 of 49 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	350	EST	O17-C17	-10.18	1.25	1.43
2	C	301	NAD	O4D-C1D	4.76	1.48	1.41
2	B	301	NAD	C2B-C1B	4.72	1.60	1.53
2	A	301	NAD	C4A-N9A	4.47	1.44	1.37
2	C	301	NAD	C4A-N9A	4.44	1.44	1.37

The worst 5 of 116 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	NAD	C5N-C4N-C3N	-8.84	108.84	120.32
2	A	301	NAD	C6N-C5N-C4N	7.71	131.69	119.44
2	C	301	NAD	C2N-C3N-C4N	7.70	127.03	118.31
3	L	350	EST	C15-C14-C8	7.00	130.15	119.03
2	C	301	NAD	N3A-C2A-N1A	-6.68	123.12	128.71

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	L	350	EST	C14
3	L	350	EST	C17

There are no torsion outliers.

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.