



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 02:35 PM GMT

PDB ID : 1E5P
Title : CRYSTAL STRUCTURE OF APHRODISIN, A SEX PHEROMONE FROM FEMALE HAMSTER
Authors : Vincent, F.; Brown, K.; Spinelli, S.; Cambillau, C.; Tegoni, M.
Deposited on : 2000-07-28
Resolution : 1.63 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

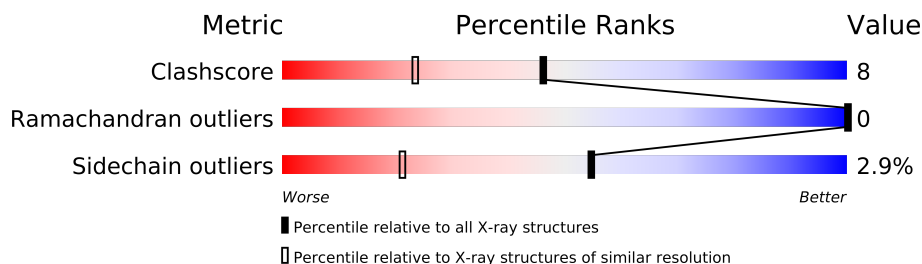
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.63 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	1525 (1.66-1.62)
Ramachandran outliers	78287	1490 (1.66-1.62)
Sidechain outliers	78261	1490 (1.66-1.62)

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	151	
1	B	151	
1	C	151	
1	D	151	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5493 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 APHRODISIN.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	Se	12	40	0
			1206	767	194	237	4	4			
1	B	148	Total	C	N	O	S	Se	28	41	0
			1186	755	192	232	4	3			
1	C	147	Total	C	N	O	S	Se	36	46	0
			1184	753	192	232	4	3			
1	D	149	Total	C	N	O	S	Se	29	39	0
			1214	776	194	236	5	3			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	73	GLU	GLN	CONFLICT	UNP P09465
A	45	MSE	MET	MODIFIED RESIDUE	UNP P09465
A	99	MSE	MET	MODIFIED RESIDUE	UNP P09465
A	108	MSE	MET	MODIFIED RESIDUE	UNP P09465
B	73	GLU	GLN	CONFLICT	UNP P09465
B	45	MSE	MET	MODIFIED RESIDUE	UNP P09465
B	99	MSE	MET	MODIFIED RESIDUE	UNP P09465
B	108	MSE	MET	MODIFIED RESIDUE	UNP P09465
C	73	GLU	GLN	CONFLICT	UNP P09465
C	45	MSE	MET	MODIFIED RESIDUE	UNP P09465
C	99	MSE	MET	MODIFIED RESIDUE	UNP P09465
C	108	MSE	MET	MODIFIED RESIDUE	UNP P09465
D	73	GLU	GLN	CONFLICT	UNP P09465
D	45	MSE	MET	MODIFIED RESIDUE	UNP P09465
D	99	MSE	MET	MODIFIED RESIDUE	UNP P09465
D	108	MSE	MET	MODIFIED RESIDUE	UNP P09465

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	179	Total 179	O 179	0	0
2	B	187	Total 187	O 187	0	0
2	C	148	Total 148	O 148	0	0
2	D	189	Total 189	O 189	0	0

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: APHRODISIN

Chain A: 



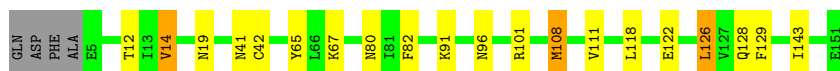
- Molecule 1: APHRODISIN

Chain B: 



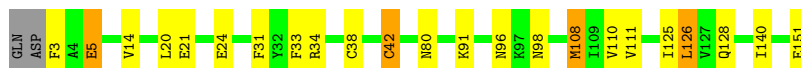
- Molecule 1: APHRODISIN

Chain C: 



- Molecule 1: APHRODISIN

Chain D: 



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 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.05Å 48.92Å 123.55Å 90.00° 102.63° 90.00°	Depositor
Resolution (Å)	15.00 – 1.63	Depositor
% Data completeness (in resolution range)	96.0 (15.00-1.63)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.60	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.194 , 0.234	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5493	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

5 Model quality i

5.1 Standard geometry i

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	0.65	0/1253	1.06	2/1688 (0.1%)
1	B	0.73	2/1211 (0.2%)	1.04	3/1634 (0.2%)
1	C	0.65	0/1214	1.00	2/1638 (0.1%)
1	D	0.82	1/1266 (0.1%)	1.12	4/1708 (0.2%)
All	All	0.72	3/4944 (0.1%)	1.06	11/6668 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	151	GLU	CB-CG	-14.98	1.23	1.52
1	B	46[A]	GLU	CD-OE1	7.71	1.34	1.25
1	B	73[A]	GLU	CD-OE1	-5.99	1.19	1.25

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	108	MSE	CA-CB-CG	7.58	126.19	113.30
1	D	151	GLU	CA-CB-CG	6.55	127.81	113.40
1	B	34	ARG	NE-CZ-NH2	-5.78	117.41	120.30
1	D	34	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	C	108	MSE	CA-CB-CG	5.76	123.10	113.30
1	C	101[A]	ARG	NE-CZ-NH2	5.58	123.09	120.30
1	B	56[A]	GLN	CA-CB-CG	5.50	125.49	113.40
1	D	5[A]	GLU	CB-CA-C	-5.44	99.52	110.40
1	A	134[A]	LYS	CB-CG-CD	-5.37	97.64	111.60
1	A	32	TYR	CB-CG-CD2	-5.29	117.83	121.00
1	B	108	MSE	CA-CB-CG	5.19	122.12	113.30

There are no chirality outliers.

There are no planarity outliers.

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	1206	0	1004	20	0
1	B	1186	0	963	24	0
1	C	1184	0	953	18	0
1	D	1214	0	1034	16	0
2	A	179	0	0	2	0
2	B	187	0	0	5	0
2	C	148	0	0	2	0
2	D	189	0	0	1	0
All	All	5493	0	3954	74	0

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 8.

All (74) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:79[A]:ASN:HD22	1:A:99[A]:MSE:HE3	1.28	0.98
1:C:91:LYS:HE2	1:C:118:LEU:HD23	1.48	0.91
1:C:118:LEU:HD22	1:C:122:GLU:HB3	1.58	0.86
1:B:14[A]:VAL:HG13	1:B:111:VAL:HB	1.60	0.83
1:D:14[A]:VAL:HG22	1:D:111:VAL:HB	1.69	0.74
1:D:80:ASN:HD21	1:D:98:ASN:HD22	1.37	0.71
1:C:80:ASN:HB3	1:C:96:ASN:HD21	1.55	0.71
1:B:128[A]:GLN:NE2	1:D:128[A]:GLN:NE2	2.39	0.70
1:B:6:LEU:HD13	1:B:45[A]:MSE:HE3	1.76	0.68
1:A:137:VAL:HG13	1:A:140:ILE:HD12	1.75	0.68
1:C:14[A]:VAL:HG13	1:C:111:VAL:HB	1.77	0.67
1:C:65[A]:TYR:CE2	1:C:67[A]:LYS:HE2	2.28	0.67
1:B:14[A]:VAL:HG21	1:B:126:LEU:HD11	1.75	0.67
1:B:36:ILE:HG21	1:B:45[A]:MSE:HE2	1.77	0.66
1:B:80:ASN:HB3	1:B:96:ASN:HD21	1.62	0.64
1:A:80:ASN:HB3	1:A:96:ASN:HD21	1.62	0.63
1:D:80:ASN:ND2	1:D:98:ASN:HD22	1.98	0.62
1:B:128[A]:GLN:HE22	1:D:125:ILE:HA	1.64	0.62
1:D:80:ASN:HB3	1:D:96:ASN:HD21	1.64	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:14[A]:VAL:HG22	1:A:111:VAL:HB	1.82	0.62
1:D:38:CYS:SG	1:D:42[B]:CYS:HB3	2.39	0.61
1:A:79[A]:ASN:HD22	1:A:99[A]:MSE:CE	2.10	0.61
1:B:9[A]:LYS:HE3	2:B:2184:HOH:O	2.01	0.61
1:C:14[B]:VAL:HG22	1:C:111:VAL:HB	1.82	0.60
1:A:90[A]:ASP:OD2	1:A:91:LYS:NZ	2.35	0.60
1:C:65[A]:TYR:HE2	1:C:67[A]:LYS:HE2	1.67	0.59
1:A:14[B]:VAL:HG13	1:A:111:VAL:HB	1.84	0.59
1:C:19[A]:ASN:HB3	2:C:2016:HOH:O	2.03	0.58
1:C:14[A]:VAL:HG21	1:C:126:LEU:HD11	1.86	0.57
1:B:4:ALA:HB1	1:B:6:LEU:HD12	1.86	0.56
1:B:125:ILE:HA	1:D:128[A]:GLN:HE22	1.71	0.56
1:B:4:ALA:HA	2:B:2001:HOH:O	2.06	0.55
1:A:99[A]:MSE:HE2	2:A:2066:HOH:O	2.06	0.55
1:A:128[A]:GLN:NE2	1:C:128:GLN:OE1	2.41	0.53
1:A:9[A]:LYS:HE2	1:A:11[A]:TYR:OH	2.08	0.53
1:C:14[B]:VAL:HG21	1:C:126:LEU:HD11	1.92	0.52
1:A:82:PHE:HB3	1:A:96:ASN:HD22	1.75	0.51
1:D:31:PHE:HZ	1:D:110:VAL:HG21	1.76	0.51
1:A:6:LEU:HD13	1:A:45:MSE:SE	2.60	0.51
1:A:97[A]:LYS:HE3	2:A:2134:HOH:O	2.11	0.50
1:D:21[A]:GLU:OE1	1:D:24[A]:GLU:OE1	2.30	0.50
1:A:128[A]:GLN:OE1	1:C:129:PHE:CD1	2.65	0.49
1:B:14[A]:VAL:HG13	1:B:111:VAL:CB	2.39	0.48
1:D:31:PHE:CE1	1:D:33[B]:PHE:CE1	3.00	0.48
1:D:96:ASN:HB3	1:D:108:MSE:HG3	1.96	0.48
1:B:5:GLU:HG2	1:B:87:ILE:HG21	1.96	0.47
1:B:82:PHE:HB3	1:B:96:ASN:HD22	1.80	0.47
1:B:97[A]:LYS:HE2	2:B:2139:HOH:O	2.14	0.47
1:D:14[A]:VAL:CG2	1:D:111:VAL:HB	2.41	0.47
1:A:69:ASN:C	1:A:69:ASN:HD22	2.18	0.46
1:C:96:ASN:HB3	1:C:108:MSE:HG3	1.97	0.46
1:A:128[A]:GLN:NE2	1:C:128:GLN:HB2	2.31	0.46
1:C:82:PHE:HB3	1:C:96:ASN:HD22	1.81	0.45
1:B:96:ASN:HB3	1:B:108:MSE:HG3	1.97	0.45
1:D:126[B]:LEU:CD1	1:D:140:ILE:HD13	2.46	0.45
1:B:9[A]:LYS:HE2	1:B:11:TYR:OH	2.17	0.45
1:A:36:ILE:HG23	1:A:45:MSE:HG3	1.98	0.45
1:C:12:THR:HG21	1:C:143:ILE:HD11	1.98	0.44
1:A:69:ASN:ND2	1:A:71:THR:H	2.16	0.43
1:B:6:LEU:HD13	1:B:45[A]:MSE:HE1	2.00	0.42
1:C:111:VAL:HG11	1:C:118:LEU:HD21	2.03	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:5:GLU:HG3	2:B:2005:HOH:O	2.20	0.41
1:A:79[A]:ASN:ND2	1:A:99[A]:MSE:HE3	2.12	0.41
1:B:137:VAL:HG11	2:B:2072:HOH:O	2.19	0.41
1:B:53:THR:O	1:B:54[A]:ASN:HB2	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	152/151 (101%)	148 (97%)	4 (3%)	0	100	100
1	B	147/151 (97%)	142 (97%)	5 (3%)	0	100	100
1	C	147/151 (97%)	143 (97%)	4 (3%)	0	100	100
1	D	153/151 (101%)	149 (97%)	4 (3%)	0	100	100
All	All	599/604 (99%)	582 (97%)	17 (3%)	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	136/130 (105%)	133 (98%)	3 (2%)	64	35
1	B	131/130 (101%)	127 (97%)	4 (3%)	52	19
1	C	132/130 (102%)	127 (96%)	5 (4%)	44	14
1	D	137/130 (105%)	130 (95%)	7 (5%)	33	7
All	All	536/520 (103%)	517 (96%)	19 (4%)	55	16

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	CYS
1	A	69	ASN
1	A	89[A]	SER
1	B	14[A]	VAL
1	B	14[B]	VAL
1	B	42[A]	CYS
1	B	126	LEU
1	C	14[A]	VAL
1	C	14[B]	VAL
1	C	41[A]	ASN
1	C	42[A]	CYS
1	C	126	LEU
1	D	3	PHE
1	D	5[A]	GLU
1	D	20	LEU
1	D	42[A]	CYS
1	D	42[B]	CYS
1	D	126[A]	LEU
1	D	126[B]	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41[A]	ASN
1	A	69	ASN
1	A	79[A]	ASN
1	A	96	ASN
1	A	139	ASN
1	A	142[A]	ASN
1	B	96	ASN
1	B	98	ASN
1	B	128[A]	GLN
1	B	131[A]	HIS
1	B	139	ASN
1	B	142[A]	ASN
1	C	19[A]	ASN
1	C	96	ASN
1	C	123	ASN
1	C	139	ASN
1	C	142[A]	ASN
1	D	80	ASN

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Mol	Chain	Res	Type
1	D	96	ASN
1	D	128[A]	GLN
1	D	131	HIS
1	D	139	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 ⓘ

There are no ligands in this entry.

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.