



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:19 PM UTC

PDB ID : 1E5P / pdb_00001e5p
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 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

<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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

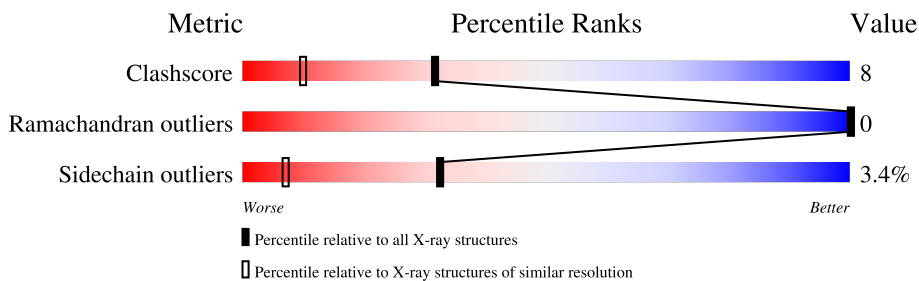
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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	190562	1171 (1.64-1.64)
Ramachandran outliers	187476	1151 (1.64-1.64)
Sidechain outliers	187428	1150 (1.64-1.64)

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 ≥ 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	151	 81% 15% . .
1	B	151	 80% 15% . .
1	C	151	 77% 17% . .
1	D	151	 79% 14% 5% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5493 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 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
B	73	GLU	GLN	conflict	UNP P09465
C	73	GLU	GLN	conflict	UNP P09465
D	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	45	MSE	MET	modified residue	UNP P09465
B	99	MSE	MET	modified residue	UNP P09465
B	108	MSE	MET	modified residue	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	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, DNA and oligosaccharide 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. 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. 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: 



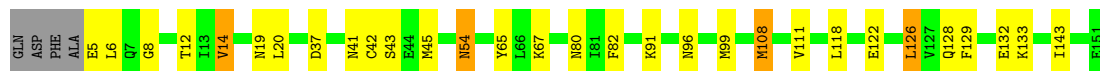
• Molecule 1: APHRODISIN

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 is therefore 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}	0.10	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	0.85	4/1253 (0.3%)	1.33	5/1688 (0.3%)
1	B	0.86	4/1211 (0.3%)	1.35	4/1634 (0.2%)
1	C	0.81	4/1214 (0.3%)	1.30	8/1638 (0.5%)
1	D	0.91	3/1266 (0.2%)	1.42	16/1708 (0.9%)
All	All	0.86	15/4944 (0.3%)	1.35	33/6668 (0.5%)

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	B	0	1
1	C	0	1
1	D	0	1
All	All	0	3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	151	GLU	CB-CG	-9.59	1.23	1.52
1	C	8	GLY	C-N	-7.71	1.22	1.33
1	B	89	SER	C-N	-7.44	1.23	1.33
1	B	45[A]	MSE	SE-CE	-6.26	1.76	1.95
1	D	45[A]	MSE	SE-CE	-5.91	1.77	1.95
1	C	45[A]	MSE	SE-CE	-5.81	1.78	1.95
1	D	5[A]	GLU	CA-CB	5.64	1.62	1.54
1	B	108	MSE	SE-CE	-5.57	1.78	1.95
1	A	45	MSE	SE-CE	-5.52	1.78	1.95
1	C	99[A]	MSE	SE-CE	-5.49	1.78	1.95
1	A	99[A]	MSE	SE-CE	-5.34	1.79	1.95
1	A	99[B]	MSE	SE-CE	-5.34	1.79	1.95
1	C	108	MSE	SE-CE	-5.34	1.79	1.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	99[A]	MSE	SE-CE	-5.24	1.79	1.95
1	A	108	MSE	SE-CE	-5.24	1.79	1.95

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	5[A]	GLU	CA-C-N	8.30	132.50	120.38
1	D	5[A]	GLU	C-N-CA	8.30	132.50	120.38
1	D	5[A]	GLU	N-CA-C	-7.95	103.39	113.72
1	C	37[A]	ASP	CA-CB-CG	7.81	120.41	112.60
1	D	41[A]	ASN	CA-C-N	7.45	133.03	122.08
1	D	41[A]	ASN	C-N-CA	7.45	133.03	122.08
1	D	151	GLU	CA-CB-CG	6.85	127.81	114.10
1	B	33	PHE	CA-CB-CG	6.65	120.45	113.80
1	D	108	MSE	CA-CB-CG	6.04	126.19	114.10
1	A	41[A]	ASN	CA-C-N	6.01	130.91	122.08
1	A	41[A]	ASN	C-N-CA	6.01	130.91	122.08
1	C	6	LEU	N-CA-C	-6.01	105.60	113.17
1	D	5[A]	GLU	CB-CA-C	-5.97	99.52	109.55
1	A	134[A]	LYS	CB-CG-CD	-5.94	97.64	111.30
1	B	18[B]	ASP	CA-CB-CG	5.92	118.52	112.60
1	C	43	SER	N-CA-C	-5.89	106.05	113.18
1	D	94	PHE	CA-CB-CG	5.71	119.51	113.80
1	D	31	PHE	CA-CB-CG	-5.71	108.09	113.80
1	B	56[A]	GLN	CA-CB-CG	5.70	125.49	114.10
1	C	8	GLY	CA-C-N	5.60	129.51	121.50
1	C	8	GLY	C-N-CA	5.60	129.51	121.50
1	C	54[A]	ASN	O-C-N	-5.59	115.89	121.88
1	D	55	ASN	CA-CB-CG	-5.50	107.10	112.60
1	D	3	PHE	CA-CB-CG	5.47	119.27	113.80
1	B	82	PHE	CA-CB-CG	5.46	119.26	113.80
1	D	96	ASN	OD1-CG-ND2	-5.41	117.19	122.60
1	D	100	ASP	CA-CB-CG	5.39	117.99	112.60
1	A	33	PHE	CA-CB-CG	5.36	119.16	113.80
1	C	20	LEU	N-CA-C	5.18	117.59	111.33
1	A	31[A]	PHE	CA-CB-CG	-5.14	108.66	113.80
1	C	132[A]	GLU	CA-CB-CG	5.08	124.25	114.10
1	D	110	VAL	CA-C-N	-5.07	116.22	123.11
1	D	110	VAL	C-N-CA	-5.07	116.22	123.11

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	90[A]	ASP	Mainchain
1	C	54[A]	ASN	Mainchain
1	D	6[A]	LEU	Mainchain

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	1206	0	1004	20	0
1	B	1186	0	963	23	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	73	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 8.

All (73) 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: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:A:137:VAL:HG13	1:A:140:ILE:HD12	1.75	0.68
1:B:6:LEU:HD13	1:B:45[A]:MSE:HE3	1.76	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:128[A]:GLN:HE22	1:D:125:ILE:HA	1.64	0.62
1:D:80:ASN:ND2	1:D:98:ASN:HD22	1.98	0.62
1:D:80:ASN:HB3	1:D:96:ASN:HD21	1.64	0.62
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:9[A]:LYS:HE2	1:A:11[A]:TYR:OH	2.08	0.53
1:A:128[A]:GLN:NE2	1:C:128:GLN:OE1	2.41	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:69:ASN:C	1:A:69:ASN:HD22	2.18	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: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:B:96:ASN:HB3	1:B:108:MSE:HG3	1.97	0.45
1:C:82:PHE:HB3	1:C:96:ASN:HD22	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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%)	132 (97%)	4 (3%)	37	11
1	B	131/130 (101%)	128 (98%)	3 (2%)	44	17
1	C	132/130 (102%)	125 (95%)	7 (5%)	20	3
1	D	137/130 (105%)	129 (94%)	8 (6%)	18	2
All	All	536/520 (103%)	514 (96%)	22 (4%)	32	4

All (22) 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	A	91	LYS
1	B	14[A]	VAL
1	B	14[B]	VAL
1	B	126	LEU
1	C	5[A]	GLU
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	C	133	LYS
1	D	3	PHE
1	D	5[A]	GLU
1	D	9[A]	LYS
1	D	20	LEU
1	D	42[A]	CYS
1	D	42[B]	CYS
1	D	126[A]	LEU
1	D	126[B]	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	69	ASN
1	A	79[A]	ASN

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Mol	Chain	Res	Type
1	A	96	ASN
1	A	98	ASN
1	A	139	ASN
1	B	35	HIS
1	B	75	GLN
1	B	96	ASN
1	B	116[A]	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	98	ASN
1	C	123	ASN
1	C	142[A]	ASN
1	D	19	ASN
1	D	35	HIS
1	D	80	ASN
1	D	96	ASN
1	D	98	ASN
1	D	128[A]	GLN
1	D	131	HIS
1	D	139	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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.