



Emidio Camaioni is an Associate Professor in Medicinal Chemistry at the *Dipartimento di Scienze Farmaceutiche* (Department of Pharmaceutical Sciences), University of Perugia (Italy).

Academic Track Record

In 1992 he received his degree with honours (*cum laude*) in *Chimica e Tecnologia Farmaceutiche* (Chemistry and Drug Technologies) from Faculty of Pharmacy at the University of Camerino (Italy).

From 1993 to 1996 he attended a three years Doctoral Course in *Chimica del Farmaco* (Drug Chemistry) under the supervision of Prof. Gloria Cristalli. On July 8th, 1997 at the University of Pisa, he received his PhD discussing an experimental thesis entitled *Agonisti ed Antagonisti del Recettore Adenosinico* (Agonists and Antagonists for Adenosine Receptors).

From 01/10/1996 to 07/12/1997 he spent a postdoctoral experience conducting research activity in medicinal chemistry at the National Institutes of Health, Molecular Recognition Section, LBC (Bethesda, MD, USA) under the supervision of Dr. Kenneth A. Jacobson.

From 1998 to 1999 he got a post-doctoral fellowship doing medicinal chemistry research at *Dipartimento di Scienze Chimiche* - University of Camerino under the supervision of Prof. Gloria Cristalli and then at *Istituto di Chimica Farmaceutica* – University of Perugia, under the supervision of Prof. Roberto Pellicciari.

In November 1999 he became Assistant Professor (D.R. n. 3462-20/10/1999 - SSD CHIM/08 Medicinal Chemistry) at the *Dipartimento di Chimica e Tecnologia del Farmaco*, University of Perugia (Italy).

In January 2005 he became Associate Professor (D.R. n. 2918-30/12/2004) in Medicinal Chemistry (SSD CHIM/08 – 3D1, D.M. 336-29/07/2011) at the *Dipartimento di Chimica e Tecnologia del Farmaco* (from 2014, Department of Pharmaceutical Sciences), University of Perugia (Italy), where he is currently working. From January 2008 he is Associate Professor confirmed (D.R. 891-22/04/2008). He was a member of the board of the PhD course *Dottorato in Chimica e Tecnologia del Farmaco a Profilo Nazionale ed Internazionale* (from 2005 up to 2013); of the PhD course *Dottorato in Scienze Chimiche e Farmaceutiche* (from 2013 up to 2016) and from 2014 is a board member of the PhD course *Dottorato in Scienze Farmaceutiche*.

Teaching Activities

From 2001 up to 2008 he taught a practical course of *Analisi dei Medicinali II* (Drug Analysis II, 12 CFU) for the degree in *Controllo Qualità nel Settore Industriale, Farmaceutico ed Alimentare* (CQSIFA), Faculty of Pharmacy.

From 2008 up to 2012 he taught a practical course of *Analisi dei Farmaci II* (Drug Analysis II, 8 CFU) for the degree in *Chimica e Tecnologia Farmaceutiche* (Chemistry and Drug Technologies), Faculty of Pharmacy.

From 2013, he is teaching a practical course of *Analisi Chimico-Farmaceutica II* (Drug Analysis II, 8 CFU) for the degree in *Chimica e Tecnologia Farmaceutiche* (Chemistry and Drug Technologies), Department of Pharmaceutical Sciences (DSF). From 2018 he is also teaching the theoretical course of *Prodotti Nutraceutici* (Nutraceutical Products, 2 CFU) for the degree in *Scienze della Alimentazione e della Nutrizione Umana* (Human Feeding and Nutrition Sciences, DSF).

He usually supervises students in carrying out the work of experimental thesis in medicinal chemistry and in nutraceutical sciences. He has been a member of numerous committees of the exams.

Scientific Research Fields

His research interests spanned in various medicinal, analytical and computational chemistry fields. Indeed, his research activities are, particularly, focused on design and synthesis of biological active heterocyclic derivatives. More in detail: i) design and synthesis of nucleosides and nucleotides as antiviral, antitumor and ligands for purinergic receptors; ii) design and synthesis of new modulators of glutamate receptors; iii) design, synthesis and characterization of novel steroidal ligands for the nuclear FXR receptor; iv) design and synthesis of novel inhibitors of the PARP family of proteins as neuroprotecting and/or anticancer agents; v) design and synthesis of novel enzyme inhibitors involved in the catabolism of the tryptophan (kynurenine pathway): Recent explorative new fields: i) drug repurposing and ii) circular economy in the development of natural pesticides.

The research activities is highlighted by a high number of peer reviewed international publications, book chapters, patent and abstracts presented at various national and international symposia (listed below). Bibliometric indices from Web of Science (update 09/07/2021): Hirsch index, 30; number of citations over 3000 (average citations per item: 30.4).

Patent (1)

1 J. L. Boyer, T. K. Harden, K. A. Jacobson, **E. Camaioni**. Preparation and use of nucleotide bisphosphate as P2Y antagonists. PCT Int. Appl. 1998, 63pp WO 981843, Chem. Abst. 1998, 129:4815.

Book (1)

- 1 **E. Camaioni.** Compendio Teorico-Pratico di Analisi Farmaceutica. ISBN-10: 8831973223; Bertoni Ed. (2018) pp.280.

Book chapters (7)

- 1 G. Cristalli, **E. Camaioni**, R. Volpini, S. Vittori. Selective A2 adenosine receptor agonists with potent inhibitory activity on platelet aggregation. From Adenosine and Adenine Nucleotides: from Molecular Biology to Integrative Physiology, Kluwer Academic Publishers, L. Belardinelli and A. Pelleg Eds.(1995) 140-148.
- 2 G. Cristalli, **E. Camaioni**, E. Di Francesco, R. Volpini, S. Vittori. Chemical and pharmacological profile of selective adenosine receptor agonists. From Perspectives in Receptor Research, Elsevier Science Publishers, Amsterdam, Pharmacochimistry Library, D. Giardinà, S. Piergentili, and M. Pigni Eds,(1996) 165-180.
- 3 K. A. Jacobson, Y.-C. Kim, **E. Camaioni**, A. M. van Rhee. Structure-activity relationships of P2 receptor agonists and antagonists. Chapter 4 in the P2 Nucleotide Receptors, in the series The receptors; Humana Press: Clifton, NJ, J. T. Turner, G. Weisman, J. Fedan Eds, (1997). 81-107.
- 4 K. A. Jacobson, C. Hoffmann, Y. C. Kim, **E. Camaioni**, E. Nandan, S. Y. Jang, D. P. Guo, X. D. Ji, K. I. Von, S. Moro, A. U. Ziganshin, A. Rychkov, B. F. King, S. G. S. S. Brown Wildman, G. Burnstock, J. L. Boyer, A. Mohanram, T. K. Harden. Molecular recognition in P2 receptors: Ligand development aided by molecular modeling and mutagenesis. Progr. in Brain Res. Illes P.; Zimmermann H. Eds, (1999) 120, 119-132.
- 5 R. Pellicciari, A. Laura, **E. Camaioni**, G. Costantino, M. Marinozzi. Heterocycles as companions on rote to drug discovery. From Seminars in Organic Synthesis, XXVII Summer School "A. Corbella", SCI Ed. (2002) pp 111-134.
- 6 R. Pellicciari, **E. Camaioni**, G. Costantino. Life or Death Decisions. The case of Poly(ADP-Ribose)Polymerase (PARP) as therapeutic target for brain ischemia. Chapter 3 in Progress in Medicinal Chemistry (2004), 42, 125-169.
- 7 B. Natalini, R. Sardella, A. Macchiarulo, M. Marinozzi, **E. Camaioni**, R. Pellicciari. Mechanistic Aspects and Applications of Chiral Ligand-Exchange Chromatography. From Advances in Chromatography, Boca Raton, FL, United States (2011), 49, 71-134.

Abstracts on peer reviewed journals (16)

- 1 T. K. Harden, R. K. Palmer, **E. Camaioni**, J. R. Schachter, K. A. Jacobson, R. A. Nicholas, J. L. Boyer. The P2Y₁ receptor: agonist selectivity, evolution of a high affinity and second messenger selectivity. Drug Dev Res. 1998, 43 (1), 4.
- 2 S. Costanzi, **E. Camaioni**, R. Volpini, S. Vittori, G. Lupidi, G. Cristalli. Synthesis and biological evaluation of adenosine deaminase inhibitors. Drug Dev Res. 1998, 43 (1), 16.
- 3 **E. Camaioni**, J. L. Boyer, A. Mohanram, T. K. Harden, K. A. Jacobson. Deoxyadenosine bisphosphate derivatives as potent antagonists at P2Y₁ receptors. Drug Dev Res. 1998, 43 (1), 26.
- 4 R. Volpini, S. Vittori, S. Costanzi, **E. Camaioni**, G. Baraldi, K. A. Jakcobson, G. Cristalli. Novel trisubstituted adenosine-uronamides as potential agonist at A₃ adenosine receptor. Drug Dev Res. 1998, 43 (1), 31.
- 5 S. Vittori, **E. Camaioni**, R. Volpini, K.-N. Klotz, G. Cristalli. Synthesis and adenosine receptor affinity of (8-bromo-9-substituted purines. Drug Dev Res. 1998, 43 (1), 31.
- 6 S. Moro, D. Guo, **E. Camaioni**, J. L. Boyer, T. K. Harden, K. A. Jacobson. Human P2Y₁receptor: molecular modeling and site-directed mutagenesis as tools to identify agonist and antagonist recognition sites. Drug Dev Res. 1998, 43 (1), 32.
- 7 A. U. Ziganshin, A. V. Rychkov, L. E. Ziganshina, Y.-C. Kim, **E. Camaioni**, G. Burnstock, K. A. Jacobson. Antagonistic profiles of new derivatives of pyridoxalphosphate-6-azophenyl-2', 4'-disulphonic acid. Naunyn Sch. Arch. Pharm. 358(1), R132, 1998.
- 8 G. Costantino, **E. Camaioni**, A. Macchiarulo, R. Pellicciari. Molecular modelling approaches towards the design of new poly(ADP-ribose) polymerase (PARP) inhibitors. Journal of Neurochemistry. 2001, 76 (Suppl 1), 5-6.
- 9 G. Costantino, M. Belenkin, A. Macchiarulo, **E. Camaioni**, R. Pellicciari. Molecular determinants for metabotropic glutamate receptor activation and blockade. Molecular dynamics and electrostatic potential analysis studies. Drug Future 2002, 27 (Suppl. A), 80.
- 10 R. Pellicciari, **E. Camaioni**, A. Macchiarulo, G. Costantino. PARP-1 and its inhibitors: molecular modeling studies. Medical Science Monitor 2003, 9 (Suppl 1), 55.
- 11 R. Pellicciari, **E. Camaioni**, F. Moroni, A. Chiarugi, G. Costantino. Novel PARP inhibitors: synthesis and molecular modeling studies. Medical Science Monitor 2005, 11 (Suppl 1), 23-24.
- 12 A. Antonyan, G. Lupidi, **E. Camaioni**, S. Nardanyan, S. Sharoyan. Inhibition of DPPiV by flavonoids as an alternative strategy for the treatment of type-2 diabetes. FEBS J. 2006, 273, 288.
- 13 **E. Camaioni**, P. Sabbatini, G. Costantino, D. Bellocchi, A. Chiarugi, F. Moroni, A. Wood, R. Pellicciari. Design, Molecular Modeling Studies, Synthesis and Preliminary Evaluation of Novel PARP-1 Inhibitors on the Road of Antiischemic Agents. Drug Future 2006, 31 (Suppl. A), 177.

- 14 **E. Camaioni**, A. Macchiarulo, R. Nuti, R. Pellicciari. Structural and Conformational Aspects of Molecular Recognition by Indoleamine-2,3-Dioxygenase (IDO), a Novel Target for Cancer Therapy. *Drug Future* 2008, 31 (Suppl. A), 304.
- 15 S. Ascitti, **E. Camaioni**, P. Mong, N. Biran, A. Carotti, P. Liscio, G. Akiri, R. Pellicciari, S. A. Aaronson. Therapeutic Potential of a Novel Poly(ADP-ribose) Polymerase Inhibitor, Hydantiq, in Human Pancreatic and Colon Cancers. *Gastroenterology* 2013, 144 (5), 875.
- 16 S. Ascitti, E. Ferretti, S. Valloscuro, M. M. Pires, L. L. Llacuna, D. Esposito, P. Liscio, A. Macchiarulo, **E. Camaioni**, R. Pellicciari, S. A. Aaronson. Novel Wnt-Targeted Therapy Inhibitors for Hepatocellular Carcinoma. *Gastroenterology* Vol. 146, Issue 5, Supplement 1, Page S-690. DDW2014, Chigago (IL, USA), May 3-6, 2014.

Abstracts of Symposium presentations (> 130)

Omissis

Publications in peer reviewed journals (91)

- 1 Cristalli G, Vittori S, Eleuteri A, Volpini R, Cola D, **Camaioni E**, Gariboldi PV, Lupidi G. Synthesis of 1,7-dideazapurine ribonucleosides and deoxyribonucleosides. *Nucleosides Nucleotides*. 1993;12:39-53.
- 2 Cristalli G, Eleuteri A, Vittori S, Volpini R, **Camaioni E**, Lupidi G. Adenosine deaminase inhibitors: structure-activity relationships in 1-deazaadenosine and erythro-9-(2-hydroxy-3-nonyl)adenine analogs. *Drug Dev Res*. 1993;28:253-8.
- 3 Cristalli G, Eleuteri A, Volpini R, Vittori S, **Camaioni E**, Lupidi G. Adenosine deaminase inhibitors: synthesis and structure-activity relationships of 2-hydroxy-3-nonyl derivatives of azoles. *J Med Chem*. 1994;37(1):201-5.
- 4 Cristalli G, Vittori S, Eleuteri A, Volpini R, **Camaioni E**, Lupidi G. Synthesis of 2'-deoxyribonucleoside derivatives of 1-deazapurine. *Nucleosides Nucleotides*. 1994;13:835-48.
- 5 Cristalli G, Volpini R, Vittori S, **Camaioni E**, Monopoli A, Conti A, Dionisotti S, Zocchi C, Ongini E. 2-Alkynyl derivatives of adenosine-5'-N-ethyluronamide: selective A2 adenosine receptor agonists with potent inhibitory activity on platelet aggregation. *J Med Chem*. 1994;37(11):1720-6.
- 6 Monopoli A, Conti A, Dionisotti S, Casati C, **Camaioni E**, Cristalli G, Ongini E. Pharmacology of the highly selective A1 adenosine receptor agonist 2-chloro-N6-cyclopentyladenosine. *Arzneimittelforschung*. 1994;44(12):1305-12.
- 7 **Camaioni E**, Volpini R, Vittori S, Monopoli A, Dionisotti S, Cristalli G. Synthesis and structure-activity relationships of a series of 2-alkynyl-NECA derivatives as selective A2 adenosine receptor agonists. *Res Commun Molecul Pathol P*. 1995;87:95-6.
- 8 Cristalli G, **Camaioni E**, Vittori S, Volpini R, Borea PA, Conti A, Dionisotti S, Ongini E, Monopoli A. 2-Aralkynyl and 2-heteroalkynyl derivatives of adenosine-5'-N-ethyluronamide as selective A2a adenosine receptor agonists. *J Med Chem*. 1995;38(9):1462-72.
- 9 Cristalli G, **Camaioni E**, Vittori S, Volpini R. Platelet aggregation inhibitory activity of selective A2 adenosine receptor agonists. *Nucleosides Nucleotides*. 1995;14:449-53.
- 10 Vittori S, **Camaioni E**, Volpini R, Palù G, Cristalli G. Inhibitory effects of 1-deazaadenosine analogues on HIV replication and adenosine deaminase. *Nucleosides Nucleotides*. 1995;14:603-6.
- 11 Cristalli G, Vittori S, Eleuteri A, Volpini R, **Camaioni E**, Lupidi G, Mahmood N, Bevilacqua F, Palù G. Synthesis and biological evaluation of N6-cycloalkyl derivatives of 1-deazaadenine nucleosides: a new class of anti-human immunodeficiency virus agents. *J Med Chem*. 1995;38(20):4019-25.
- 12 Volpini R, **Camaioni E**, Di Francesco E, Vittori S, Fein T, Schwabe U, Cristalli G. Synthesis and biological activity of 3'-deoxy derivative of 5'-N-methylcarboxamidoadenosine (MECA). **Collect Czech Chem Commun**. 1996;61:33-7.
- 13 **Camaioni E**, Di Francesco E, Volpini R, Vittori S, Cristalli G. Unexpected synthesis of acyclic adenine nucleosides. **Collect Czech Chem Commun**. 1996;61:154-5.
- 14 Vittori S, **Camaioni E**, Di Francesco E, Volpini R, Monopoli A, Dionisotti S, Ongini E, Cristalli G. 2-alkenyl and 2-alkyl derivatives of adenosine and adenosine-5'-N-ethyluronamide: different affinity and selectivity of E- and Z-diastereomers at A2A adenosine receptors. *J Med Chem*. 1996;39(21):4211-7.
- 15 Cristalli G, Volpini R, Vittori S, **Camaioni E**, Rafaiani G, Potenza S, Vita A. Diazepinone nucleosides as inhibitors of cytidine deaminase. *Nucleosides Nucleotides*. 1996;15:1567-80.
- 16 Volpini R, **Camaioni E**, Lupidi G, Vittori S, Cristalli G. Adenosine deaminase inhibitors: synthesis, diastereoisomeric resolution and biological activity of 1-(2-hydroxy-3-nonyl)-1,2,4-triazole-3-carboxamide. *Farmaco*. 1997;52(6-7):429-33..
- 17 Cristalli G, **Camaioni E**, E. Di Francesco, Eleuteri A, Vittori S, Volpini R. Potent and selective ligands for adenosine binding sites. *Nucleosides Nucleotides*. 1997;16:1379-88.
- 18 **Camaioni E**, Di Francesco E, Vittori S, Volpini R, Cristalli G. Adenosine receptor agonists: synthesis and biological evaluation of the diastereoisomers of 2-(3-hydroxy-3-phenyl-1-propyn-1-yl)NECA. *Bioorg Med Chem*. 1997;5(12):2267-75.

- 19 Volpini R, **Camaioni E**, Vittori S, Barboni L, Lambertucci C, Cristalli G. Synthesis of new nucleosides by coupling of chloropurines with 2- and 3-deoxy derivatives of N-methyl-D-ribofuranuronamide. *Helv Chim Acta* 1998;81:145-52.
- 20 **Camaioni E**, Boyer JL, Mohanram A, Harden TK, Jacobson KA. Deoxyadenosine bisphosphate derivatives as potent antagonists at P2Y1 receptors. *J Med Chem.* 1998;41(2):183-90.
- 21 Moro S, Guo D, **Camaioni E**, Boyer JL, Harden TK, Jacobson KA. Human P2Y1 receptor: molecular modeling and site-directed mutagenesis as tools to identify agonist and antagonist recognition sites. *J Med Chem.* 1998;41(9):1456-66.
- 22 **Camaioni E**, Costanzi S, Vittori S, Volpini R, Klotz KN, Cristalli G. New substituted 9-alkylpurines as adenosine receptor ligands. *Bioorg Med Chem.* 1998;6(5):523-33.
- 23 Boyer JL, Mohanram A, **Camaioni E**, Jacobson KA, Harden TK. Competitive and selective antagonism of P2Y1 receptors by N6-methyl 2'-deoxyadenosine 3',5'-bisphosphate. *Br J Pharmacol.* 1998;124(1):1-3.
- 24 Barbieri D, Franceschi C, **Camaioni E**, Costanzi S, Vittori S, Volpini R, Cristalli G. Modulation of apoptosis in human lymphocytes by adenosine analogues. *Bioorg Med Chem Lett.* 1998;8(18):2533-8.
- 25 Kim YC, **Camaioni E**, Ziganshin AU, Ji XD, King BF, Wildman SS, Rychkov A, Yoburn J, Kim H, Mohanram A, Harden TK, Boyer JL, Burnstock G, Jacobson KA. Synthesis and Structure-Activity Relationships of Pyridoxal-6-aryloxy-5'-phosphate and Phosphonate Derivatives as P2 Receptor Antagonists. *Drug Dev Res.* 1998;45(2):52-66.
- 26 Volpini R, **Camaioni E**, Costanzi S, Vittori S, Cristalli G. Synthesis of new 3'-deoxyribonucleosides employing the acid-catalyzed fusion method. *Helv Chim Acta* 1998;81:2326-31.
- 27 Cristalli G, **Camaioni E**, Costanzi S, Vittori S, Volpini R, Klotz KN. Characterization of Potent Ligands at Human Recombinant Adenosine Receptors. *Drug Dev Res.* 1998;45:176-81.
- 28 Ziganshin AU, Rychkov AV, Ziganshina LE, Kim YC, **Camaioni E**, Burnstock G, Jacobson KA. Some 6-phenylazopyridoxal phosphate derivatives influence P2-purinoreceptor-mediated effects. *Pharmaceutical Chem. J.* 1998;32:399-401.
- 29 Klotz KN, **Camaioni E**, Volpini R, Kachler S, Vittori S, Cristalli G. 2-Substituted N-ethylcarboxamidoadenosine derivatives as high-affinity agonists at human A3 adenosine receptors. *Naunyn Schmiedebergs Arch Pharmacol.* 1999;360(2):103-8.
- 30 Nandan E, **Camaioni E**, Jang SY, Kim YC, Cristalli G, Herdewijn P, Secrist JA 3rd, Tiwari KN, Mohanram A, Harden TK, Boyer JL, Jacobson KA. Structure-activity relationships of bisphosphate nucleotide derivatives as P2Y1 receptor antagonists and partial agonists. *J Med Chem.* 1999;42(9):1625-38.
- 31 Vittori S, **Camaioni E**, Costanzi S, Volpini R, Cristalli G. Coupling of 2,6-dichloropurine and 2,6-dichlorodeazapurines with ribose and ribose modified sugars. *Nucleosides Nucleotides.* 1999;18:587-90.
- 32 Vittori S, **Camaioni E**, Costanzi S, Volpini R, Klotz KN, Cristalli G. Synthesis and receptor affinity of polysubstituted adenosines. *Nucleosides Nucleotides.* 1999;18(4-5):739-40.
- 33 Vittori S, **Camaioni E**, Costanzi S, Volpini R, Lupidi G, Cristalli G. Structure-activity relationships of adenosine deaminase inhibitors. *Nucleosides Nucleotides.* 1999;18:741-2.
- 34 Volpini R, **Camaioni E**, Costanzi S, Vittori S, Klotz KN, Cristalli G. Synthesis of di- and tri-substituted adenosine derivatives and their affinities at human adenosine receptor subtypes. *Nucleosides Nucleotides.* 1999;18(11-12):2511-20.
- 35 Ziganshin AU, Zaitsev AP, Zaitseva IP, Kim YC, **Camaioni E**, Burnstock G, Jacobson KA. Effects of new arylazido compounds of pyridoxal phosphate on ecto-ATP-ase activity of guinea pig tissues. *Pharm Chem J.* 2000;34(5):226-228.
- 36 Cristalli G, Costanzi S, Lambertucci C, Lupidi G, Vittori S, Volpini R, **Camaioni E**. Adenosine deaminase: functional implications and different classes of inhibitors. *Med Res Rev.* 2001;21(2):105-28.
- 37 Costantino G, Maltoni K, Marinozzi M, **Camaioni E**, Prezeau L, Pin JP, Pellicciari R. Synthesis and biological evaluation of 2-(3'-(1H-tetrazol-5-yl) bicyclo[1.1.1]pent-1-yl)glycine (S-TBPG), a novel mGlu1 receptor antagonist. *Bioorg Med Chem.* 2001;9(2):221-7.
- 38 Natalini B, **Camaioni E**, Colagioia S, Pellicciari R. Predicting bile acids intestinal absorption with RP-HPLC on a dynamically coated artificial membrane. *Chromatographia* 2001;53:453-5.
- 39 Pellicciari R, Costantino G, Marinozzi M, Macchiarulo A, **Camaioni E**, Natalini B. Metabotropic glutamate receptors: structure and new subtype-selective ligands. *Farmaco.* 2001;56(1-2):91-4.
- 40 Costantino G, Macchiarulo A, **Camaioni E**, Pellicciari R. Modeling of poly(ADP-ribose)polymerase (PARP) inhibitors. Docking of ligands and quantitative structure-activity relationship analysis. *J Med Chem.* 2001;44(23):3786-94.
- 41 Pellicciari R, Marinozzi M, **Camaioni E**, del Carmen Nunez M, Costantino G, Gasparini F, Giorgi G, Macchiarulo A, Subramanian N. Spiro[2.2]pentane as a dissymmetric scaffold for conformationally constrained analogues of glutamic acid: focus on racemic 1-aminospiro[2.2]pentyl-1,4-dicarboxylic acids. *J Org Chem.* 2002;67(16):5497-507.
- 42 Pellicciari R, Fiorucci S, **Camaioni E**, Clerici C, Costantino G, Maloney PR, Morelli A, Parks DJ, Willson TM. 6 α -ethyl-chenodeoxycholic acid (6-ECDCA), a potent and selective FXR agonist endowed with anticholestatic activity. *J Med Chem.* 2002;45(17):3569-72.

- 43 Costantino G, Macchiarulo A, Entrena-Guadix A, **Camaioni E**, Pellicciari R. Binding mode of 6ECDCA, a potent bile acid agonist of the farnesoid X receptor (FXR). *Bioorg Med Chem Lett*. 2003;13(11):1865-8.
- 44 Chiarugi A, Meli E, Calvani M, Picca R, Baronti R, **Camaioni E**, Costantino G, Marinozzi M, Pellegrini-Giampietro DE, Pellicciari R, Moroni F. Novel isoquinolinone-derived inhibitors of poly(ADP-ribose) polymerase-1: pharmacological characterization and neuroprotective effects in an in vitro model of cerebral ischemia. *J Pharmacol Exp Ther*. 2003;305(3):943-9.
- 45 Pellicciari R, **Camaioni E**, Costantino G, Marinozzi M, Macchiarulo A, Moroni F, Natalini B. Towards new neuroprotective agents: design and synthesis of 4H-thieno[2,3-c] isoquinolin-5-one derivatives as potent PARP-1 inhibitors. *Farmaco*. 2003;58(9):851-8.
- 46 Traversa U, Bombi G, **Camaioni E**, Macchiarulo A, Costantino G, Palmieri C, Caciagli F, Pellicciari R. Rat brain guanosine binding site. Biological studies and pseudo-receptor construction. *Bioorg Med Chem*. 2003;11(24):5417-25.
- 47 Costantino G, Marinozzi M, **Camaioni E**, Natalini B, Sarichelou I, Micheli F, Cavanni P, Faedo S, Noe C, Moroni F, Pellicciari R. Stereoselective synthesis and preliminary evaluation of (+)- and (-)-3-methyl-5-carboxy-thien-2-yl-glycine (3-MATIDA): identification of (+)-3-MATIDA as a novel mGluR1 competitive antagonist. *Farmaco*. 2004;59(2):93-9.
- 48 Pellicciari R, Costantino G, **Camaioni E**, Sadeghpour BM, Entrena A, Willson TM, Fiorucci S, Clerici C, Gioiello A. Bile acid derivatives as ligands of the farnesoid X receptor. Synthesis, evaluation, and structure-activity relationship of a series of body and side chain modified analogues of chenodeoxycholic acid. *J Med Chem*. 2004;47(18):4559-69.
- 49 Natalini B, Sardella R, **Camaioni E**, Natalini S, Pellicciari R. Dominant Factors Affecting The Chromatographic Behavior of Bile Acids. *Chromatographia* 2006;64:343-9.
- 50 Clerici C, Castellani D, Ascitti S, Pellicciari R, Setchell KD, O'Connell NC, Sadeghpour B, **Camaioni E**, Fiorucci S, Renga B, Nardi E, Sabatino G, Clementi M, Giuliano V, Baldoni M, Orlandi S, Mazzocchi A, Morelli A, Morelli O. 3alpha-6alpha-Dihydroxy-7alpha-fluoro-5beta-cholanoate (UPF-680), physicochemical and physiological properties of a new fluorinated bile acid that prevents 17alpha-ethynyl-estradiol-induced cholestasis in rats. *Toxicol Appl Pharmacol*. 2006;214(2):199-208.
- 51 Natalini B, Sardella R, **Camaioni E**, Gioiello A, Pellicciari R. Correlation between CMC and chromatographic index: simple and effective evaluation of the hydrophobic/hydrophilic balance of bile acids. *Anal Bioanal Chem*. 2007;388(8):1681-8.
- 52 Macchiarulo A, Nuti R, Bellocchi D, **Camaioni E**, Pellicciari R. Molecular docking and spatial coarse graining simulations as tools to investigate substrate recognition, enhancer binding and conformational transitions in indoleamine-2,3-dioxygenase (IDO). *Biochim Biophys Acta*. 2007;1774(8):1058-68.
- 53 Pellicciari R, **Camaioni E**, Costantino G, Formentini L, Sabbatini P, Venturoni F, Eren G, Bellocchi D, Chiarugi A, Moroni F. On the way to selective PARP-2 inhibitors. Design, synthesis, and preliminary evaluation of a series of isoquinolinone derivatives. *ChemMedChem*. 2008;3(6):914-23.
- 54 Macchiarulo A, **Camaioni E**, Nuti R, Pellicciari R. Highlights at the gate of tryptophan catabolism: a review on the mechanisms of activation and regulation of indoleamine 2,3-dioxygenase (IDO), a novel target in cancer disease. *Amino Acids*. 2009;37(2):219-29.
- 55 Formentini L, Macchiarulo A, Cipriani G, **Camaioni E**, Rapizzi E, Pellicciari R, Moroni F, Chiarugi A. Poly(ADP-ribose) catabolism triggers AMP-dependent mitochondrial energy failure. *J Biol Chem*. 2009;284(26):17668-76.
- 56 Moroni F, Formentini L, Gerace E, **Camaioni E**, Pellegrini-Giampietro DE, Chiarugi A, Pellicciari R. Selective PARP-2 inhibitors increase apoptosis in hippocampal slices but protect cortical cells in models of post-ischaemic brain damage. *Br J Pharmacol*. 2009;157(5):854-62.
- 57 Natalini B, Sardella R, **Camaioni E**, Macchiarulo A, Gioiello A, Carbone G, Pellicciari R. Derived chromatographic indices as effective tools to study the self-aggregation process of bile acids. *J Pharm Biomed Anal*. 2009;50(4):613-21.
- 58 Natalini B, Sardella R, Giacchè N, Palmiotto S, **Camaioni E**, Marinozzi M, Macchiarulo A, Pellicciari R. Chiral ligand-exchange separation and resolution of extremely rigid glutamate analogs: 1-aminospiro[2.2]pentyl-1,4-dicarboxylic acids. *Anal Bioanal Chem*. 2010;397(5):1997-2011.
- 59 Tentori L, Muzi A, Dorio AS, Scarsella M, Leonetti C, Shah GM, Xu W, **Camaioni E**, Gold B, Pellicciari R, Dantzer F, Zhang J, Graziani G. Pharmacological inhibition of poly(ADP-ribose) polymerase (PARP) activity in PARP-1 silenced tumour cells increases chemosensitivity to temozolomide and to a N3-adenine selective methylating agent. *Curr Cancer Drug Targets*. 2010;10(4):368-83.
- 60 Lupidi G, Scire A, **Camaioni E**, Khalife KH, De Sanctis G, Tanfani F, Damiani E. Thymoquinone, a potential therapeutic agent of *Nigella sativa*, binds to site I of human serum albumin. *Phytomedicine*. 2010;17(10):714-20.
- 61 Pellicciari R, **Camaioni E**, Gilbert AM, Macchiarulo A, Bikker JA, Shah F, Bard J, Costantino G, Gioiello A, Robertson GM, Sabbatini P, Venturoni F, Liscio P, Carotti A, Bellocchi B, Cozzi A, Wood A, Gonzales C, Zaleska MM, Ellingboe JW, Moroni F. Discovery and characterization of novel potent PARP-1 inhibitors endowed with neuroprotective properties: From TIQ-A to HYDAMTIQ. *MedChemComm*. 2011;2:559-65.

- 62 Natalini B, Sardella R, Gioiello A, Rosatelli E, Ianni F, **Camaioni E**, Pellicciari R. Fast chromatographic determination of the bile salt critical micellar concentration. *Anal Bioanal Chem.* 2011;401(1):267-74.
- 63 Moroni F, Cozzi A, Chiarugi A, Formentini L, **Camaioni E**, Pellegrini-Giampietro DE, Chen Y, Liang S, Zaleska MM, Gonzales C, Wood A, Pellicciari R. Long-lasting neuroprotection and neurological improvement in stroke models with new, potent and brain permeable inhibitors of poly(ADP-ribose) polymerase. *Br J Pharmacol.* 2012;165(5):1487-500.
- 64 Lupidi G, **Camaioni E**, Khalifé H, Avenali L, Damiani E, Tanfani F, Scirè A. Characterization of thymoquinone binding to human α_1 -acid glycoprotein. *J Pharm Sci.* 2012;101(7):2564-73.
- 65 Ekblad T, **Camaioni E**, Schüler H, Macchiarulo A. PARP inhibitors: polypharmacology versus selective inhibition. *FEBS J.* 2013;280(15):3563-75.
- 66 Liscio P, **Camaioni E**, Carotti A, Pellicciari R, Macchiarulo A. From polypharmacology to target specificity: the case of PARP inhibitors. *Curr Top Med Chem.* 2013;13(23):2939-54.
- 67 Antolin AA, Carotti A, Nuti R, Hakkaya A, **Camaioni E**, Mestres J, Pellicciari R, Macchiarulo A. Exploring the effect of PARP-1 flexibility in docking studies. *J Mol Graph Model.* 2013;45:192-201.
- 68 Antonyan A, De A, Vitali LA, Pettinari R, Marchetti F, Gigliobianco MR, Pettinari C, **Camaioni E**, Lupidi G. Evaluation of (arene)Ru(II) complexes of curcumin as inhibitors of dipeptidyl peptidase IV. *Biochimie.* 2014;99:146-52.
- 69 Liscio P, Carotti A, Ascutti S, Karlberg T, Bellocchi D, Llacuna L, Macchiarulo A, Aaronson SA, Schüler H, Pellicciari R, **Camaioni E**. Design, synthesis, crystallographic studies, and preliminary biological appraisal of new substituted triazolo[4,3-b]pyridazin-8-amine derivatives as tankyrase inhibitors. *J Med Chem.* 2014;57(6):2807-12.
- 70 Liscio P, Carotti A, Ascutti S, Ferri M, Pires MM, Valloscuro S, Ziff J, Clark NR, Macchiarulo A, Aaronson SA, Pellicciari R, **Camaioni E**. Scaffold hopping approach on the route to selective tankyrase inhibitors. *Eur J Med Chem.* 2014;87:611-23.
- 71 Marchand JR, Carotti A, Passeri D, Filippini P, Liscio P, **Camaioni E**, Pellicciari R, Gioiello A, Macchiarulo A. Investigating the allosteric reverse signalling of PARP inhibitors with microsecond molecular dynamic simulations and fluorescence anisotropy. *Biochim Biophys Acta.* 2014;1844(10):1765-72.
- 72 Greco FA, Coletti A, **Camaioni E**, Carotti A, Marinozzi M, Gioiello A, Macchiarulo A. The Janus-faced nature of IDO1 in infectious diseases: challenges and therapeutic opportunities. *Future Med Chem.* 2016;8(1):39-54.
- 73 Passeri D, **Camaioni E**, Liscio P, Sabbatini P, Ferri M, Carotti A, Giacchè N, Pellicciari R, Gioiello A, Macchiarulo A. Concepts and Molecular Aspects in the Polypharmacology of PARP-1 Inhibitors. *ChemMedChem.* 2016;11(12):1219-26.
- 74 Carotti A, Ianni F, **Camaioni E**, Pucciarini L, Marinozzi M, Sardella R, Natalini B. N-Decyl-S-trityl-(R)-cysteine, a new chiral selector for "green" ligand-exchange chromatography applications. *J Pharm Biomed Anal.* 2017;144:31-40.
- 75 Coletti A, Greco FA, Dolciami D, **Camaioni E**, Sardella R, Pallotta MT, Volpi C, Orabona C, Grohmann U, Macchiarulo A. Advances in indoleamine 2,3-dioxygenase 1 medicinal chemistry. *Medchemcomm.* 2017;8(7):1378-1392.
- 76 Ferri M, Liscio P, Carotti A, Ascutti S, Sardella R, Macchiarulo A, **Camaioni E**. Targeting Wnt-driven cancers: Discovery of novel tankyrase inhibitors. *Eur J Med Chem.* 2017;142:506-522.
- 77 Ianni F, Scorzoni S, Gentili PL, Michele AD, Frigoli M, **Camaioni E**, Ortica F, Sardella R. Chiral separation of helical chromenes with chloromethyl phenylcarbamate polysaccharide-based stationary phases. *J Sep Sci.* 2017;144:31-40.
- 78 Muzzi M, Gerace E, Buonvicino D, Coppi E, Resta F, Formentini L, Zecchi R, Tigli L, Guasti D, Ferri M, **Camaioni E**, Masi A, Pellegrini-Giampietro DE, Mannaioni G, Bani D, Pugliese AM, Chiarugi A. Dextrampipexole improves bioenergetics and outcome in experimental stroke. *Br J Pharmacol.* 2018;175(2):272-283.
- 79 Buonvicino D, Mazzola F, Zamporlini F, Resta F, Ranieri G, **Camaioni E**, Muzzi M, Zecchi R, Pieraccini G, Dölle C, Calamante M, Bartolucci G, Ziegler M, Stecca B, Raffaelli N, Chiarugi A. Identification of the Nicotinamide Salvage Pathway as a New Toxication Route for Antimetabolites. *Cell Chem Biol.* 2018;25(4):471-82.
- 80 Ianni F, Lisanti A, Marinozzi M, **Camaioni E**, Pucciarini L, Massoli A, Sardella R, Concezzi L, Natalini B. Hydrophobic Amino Acid Content in Onions as Potential Fingerprints of Geographical Origin: The Case of Rossa da Inverno sel. Rojo Duro. *Molecules.* 2018;23(6):E1259.
- 81 Venturoni F, Cerra B, Marinozzi M, **Camaioni E**, Gioiello A, Pellicciari R. BF3 center dot Et2O-Promoted Decomposition of Cyclic alpha-Diazo-beta-Hydroxy Ketones: Novel Insights into Mechanistic Aspects. *Catalysts.* 2018;8(12):e600.
- 82 Ianni F, Sechi P, La Mantia A, Pucciarini L, **Camaioni E**, Cenci Coga BT, Roccaldo S, Natalini B. The Relationships between Somatic Cells and Isoleucine, Leucine and Tyrosine Content in Cow Milk. *Appl Sci.* 2019;9(2), 349.

- 83 Grispoli L, Karama M, Ianni F, La Mantia A, Pucciarini L, **Camaioni E**, Sardella R, Sechi P, Natalini B, Cenci-Goga BT. The Relationship between *S. aureus* and Branched-Chain Amino Acids Content in Composite Cow Milk. *Animals (Basel)*. 2019;9(11):981.
- 84 Sardella R, **Camaioni E**, Macchiarulo A, Gioiello A, Marinozzi M, Carotti A. Computational studies in enantioselective liquid chromatography: forty years in docking and molecular dynamics based simulations. *Trends in Analytical Chemistry*, 2020;122:115703.
- 85 Mammoli A, Coletti A, Ballarotto M, Riccio A, Carotti A, Grohmann U, **Camaioni E**, Macchiarulo A. New Insights from Crystallographic Data: Diversity of Structural Motifs and Molecular Recognition Properties between Groups of IDO1 Structures. *ChemMedChem*. 2020;15(10):891-9.
- 86 Torres Moreno H, Ianni F, Robles Zepeda RE, López-Romero JC, Vidal-Gutiérrez M, Jacobi Durán MD, Galarini R, **Camaioni E**, Sardella R, Marcotullio MC. Quantitative analysis of cucurbitane-type triterpenes in *Ibervillea sonorae* extracts: Relationship study with their antiproliferative activity. *Steroids*. 2020;161:108676.
- 87 Cerra B, Macchiarulo A, Carotti A, **Camaioni E**, Varfaj I, Sardella R, Gioiello A. Enantioselective HPLC Analysis to Assist the Chemical Exploration of Chiral Imidazolines. *Molecules*. 2020;25(3):640.
- 88 Ferri M, Alunno M, Greco FA, Mammoli A, Saluti G, Carotti A, Sardella R, Macchiarulo A, **Camaioni E**, Liscio P. Fragment based drug design and diversity-oriented synthesis of carboxylic acid isosteres. *Bioorg Med Chem*. 2020;28(22):115731.
- 89 Urru M, Muzzi M, Coppi E, Ranieri G, Buonvicino D, **Camaioni E**, Coppini R, Pugliese AM, Tanaka B, Estacion M, Waxman SG, Dib-Hajj SD, Chiarugi A. Dexpramipexole blocks Nav1.8 sodium channels and provides analgesia in multiple nociceptive and neuropathic pain models. *Pain*. 2020;161(4):831-41.
- 90 Ronchetti R, Moroni G, Carotti A, Gioiello A, **Camaioni E**. Recent advances in urea- and thiourea-containing compounds: focus on innovative approaches in medicinal chemistry and organic synthesis. *RSC Med Chem*. 2021, accepted.
- 91 Mammoli A, Riccio A, Bianconi E, Coletti A, **Camaioni E**, Macchiarulo A. One Key and Multiple Locks: Substrate Binding in Structures of Tryptophan Dioxygenases and Hydroxylases. *ChemMedChem*. 2021, accepted.

Other publications (2)


- 1 Porfiri F, Macerata R, **Camaioni E**. Il rischio chimico nell'ambiente di lavoro calzaturiero. *Wealth Planet* 2015, 36, 14-21.
- 2 Ianni F, Pucciarini L, **Camaioni E**, Alunni G, Sardella R, Natalini B. Branched-chain Amino Acids as Potential Diagnostic and Prognostic Disease Biomarkers. *Int J Clin Res Trials* 2017, 2, 112.

He was also involved in the organizing committee of the following scientific meetings:

- 1) Recent Development in Pharmaceutical Analysis (RDPA2015), Perugia (PG, Italy) June 28 – July 1, 2015.
- 2) National Meeting in Medicinal Chemistry and 10 Young Medicinal Chemists' Symposium (XIV NMMC & 10th NPCF) - Perugia (Italy), September 11-14, 2016.

He is a peer reviewer for PRIN Projects (Italian Ministry of Education, University and Research) and for several international journals.

Perugia, 12/07/2021


Prof. E. Camaioni



DIPARTIMENTO 2018
di ECCELLENZA 2022
progetto DELPHI