

Hany W Darwish

List of Publications by Year in descending order

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Version: 2024-02-01

73
papers

1,202
citations

279701

23
h-index

434063

31
g-index

76
all docs

76
docs citations

76
times ranked

883
citing authors

#	ARTICLE	IF	CITATIONS
1	Three different spectrophotometric methods manipulating ratio spectra for determination of binary mixture of Amlodipine and Atorvastatin. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2011, 83, 140-148.	2.0	87
2	Fluorescence spectroscopic and molecular docking studies of the binding interaction between the new anaplastic lymphoma kinase inhibitor crizotinib and bovine serum albumin. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2017, 171, 174-182.	2.0	65
3	Three different methods for determination of binary mixture of Amlodipine and Atorvastatin using dual wavelength spectrophotometry. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2013, 104, 70-76.	2.0	49
4	Synthesis, cytotoxic evaluation, and molecular docking studies of novel quinazoline derivatives with benzenesulfonamide and anilide tails: Dual inhibitors of EGFR/HER2. <i>Bioorganic Chemistry</i> , 2020, 95, 103461.	2.0	41
5	Investigation of metabolic degradation of new ALK inhibitor: Entrectinib by LC-MS/MS. <i>Clinica Chimica Acta</i> , 2018, 485, 298-304.	0.5	38
6	LC-MS/MS reveals the formation of iminium and quinone methide reactive intermediates in entrectinib metabolism: In vivo and in vitro metabolic investigation. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , 2018, 160, 19-30.	1.4	37
7	Comparative study between derivative spectrophotometry and multivariate calibration as analytical tools applied for the simultaneous quantitation of Amlodipine, Valsartan and Hydrochlorothiazide. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2013, 113, 215-223.	2.0	35
8	Identification and characterization of in vitro phase I and reactive metabolites of masitinib using a LC-MS/MS method: bioactivation pathway elucidation. <i>RSC Advances</i> , 2017, 7, 4479-4491.	1.7	35
9	Detection and characterization of ponatinib reactive metabolites by liquid chromatography tandem mass spectrometry and elucidation of bioactivation pathways. <i>RSC Advances</i> , 2016, 6, 72575-72585.	1.7	34
10	Identification and characterization of in vivo, in vitro and reactive metabolites of vandetanib using LC-ESI-MS/MS. <i>Chemistry Central Journal</i> , 2018, 12, 99.	2.6	33
11	LC-MS/MS reveals the formation of aldehydes and iminium reactive intermediates in foretinib metabolism: phase I metabolic profiling. <i>RSC Advances</i> , 2017, 7, 36279-36287.	1.7	31
12	A reliable and stable method for the determination of foretinib in human plasma by LC-MS/MS: Application to metabolic stability investigation and excretion rate. <i>European Journal of Mass Spectrometry</i> , 2018, 24, 344-351.	0.5	31
13	LC-MS/MS method for the quantification of masitinib in RLMs matrix and rat urine: application to metabolic stability and excretion rate. <i>Chemistry Central Journal</i> , 2017, 11, 136.	2.6	30
14	Validated LC-MS/MS Method for the Quantification of Ponatinib in Plasma: Application to Metabolic Stability. <i>PLoS ONE</i> , 2016, 11, e0164967.	1.1	29
15	A highly efficient and sensitive LC-MS/MS method for the determination of afatinib in human plasma: application to a metabolic stability study. <i>Biomedical Chromatography</i> , 2016, 30, 1248-1255.	0.8	28
16	Liquid chromatography tandem mass spectrometry method for the quantification of vandetanib in human plasma and rat liver microsomes matrices: metabolic stability investigation. <i>Chemistry Central Journal</i> , 2017, 11, 45.	2.6	28
17	LC-ESI-MS/MS reveals the formation of reactive intermediates in brigatinib metabolism: elucidation of bioactivation pathways. <i>RSC Advances</i> , 2018, 8, 1182-1190.	1.7	28
18	Investigation of the metabolic stability of olmutinib by validated LC-MS/MS: quantification in human plasma. <i>RSC Advances</i> , 2018, 8, 40387-40394.	1.7	28

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19	Investigation of metabolic stability of the novel ALK inhibitor brigatinib by liquid chromatography tandem mass spectrometry. <i>Clinica Chimica Acta</i> , 2018, 480, 180-185.	0.5	27
20	Solid-state potentiometric sensor for the rapid assay of the biologically active biogenic amine (tyramine) as a marker of food spoilage. <i>Food Chemistry</i> , 2021, 346, 128911.	4.2	27
21	Support vector regression and artificial neural network models for stability indicating analysis of mebeverine hydrochloride and sulphiride mixtures in pharmaceutical preparation: A comparative study. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2012, 86, 515-526.	2.0	25
22	LC-MS/MS reveals the formation of reactive ortho-quinone and iminium intermediates in saracatinib metabolism: Phase I metabolic profiling. <i>Clinica Chimica Acta</i> , 2018, 482, 84-94.	0.5	25
23	Identification of reactive intermediate formation and bioactivation pathways in Abemaciclib metabolism by LC-MS/MS: <i>in vitro</i> metabolic investigation. <i>Royal Society Open Science</i> , 2019, 6, 181714.	1.1	24
24	Different approaches in Partial Least Squares and Artificial Neural Network models applied for the analysis of a ternary mixture of Amlodipine, Valsartan and Hydrochlorothiazide. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2014, 122, 744-750.	2.0	23
25	Multivariate Versus Classical Univariate Calibration Methods for Spectrofluorimetric Data: Application to Simultaneous Determination of Olmesartan Medoxamil and Amlodipine Besylate in their Combined Dosage Form. <i>Journal of Fluorescence</i> , 2013, 23, 79-91.	1.3	19
26	Spectroscopic and molecular docking studies of the binding of the angiotensin II receptor blockers (ARBs) azilsartan, eprosartan and olmesartan to bovine serum albumin. <i>Journal of Luminescence</i> , 2018, 203, 616-628.	1.5	19
27	Application of smart spectrophotometric methods and artificial neural network for the simultaneous quantitation of olmesartan medoxamil, amlodipine besylate and hydrochlorothiazide in their combined pharmaceutical dosage form. <i>Chemistry Central Journal</i> , 2013, 7, 22.	2.6	16
28	Micellar enhanced synchronous spectrofluorimetric method for determination of dasatinib in tablets, human plasma and urine: application to <i>in vitro</i> drug release and content uniformity test. <i>RSC Advances</i> , 2015, 5, 54471-54480.	1.7	16
29	<i>In silico</i> and <i>in vitro</i> metabolism of ribociclib: a mass spectrometric approach to bioactivation pathway elucidation and metabolite profiling. <i>RSC Advances</i> , 2020, 10, 22668-22683.	1.7	16
30	Rapid validated liquid chromatographic method coupled with Tandem mass spectrometry for quantification of nintedanib in human plasma. <i>Tropical Journal of Pharmaceutical Research</i> , 2016, 15, 2467.	0.2	13
31	Belizatinib: Novel reactive intermediates and bioactivation pathways characterized by LC-MS/MS. <i>Journal of Pharmaceutical and Biomedical Analysis</i> , 2019, 171, 132-147.	1.4	13
32	Advanced stability indicating chemometric methods for quantitation of amlodipine and atorvastatin in their quinary mixture with acidic degradation products. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2016, 154, 58-66.	2.0	12
33	Response surface methodology for optimization of micellar-enhanced spectrofluorimetric method for assay of foretinib in bulk powder and human urine. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2021, 257, 119811.	2.0	12
34	Comparative ANNs with Different Input Layers and GA-PLS Study for Simultaneous Spectrofluorimetric Determination of Melatonin and Pyridoxine HCl in the Presence of Melatonin's Main Impurity. <i>Molecules</i> , 2013, 18, 974-996.	1.7	11
35	A new method to determine the new C-Met inhibitor Cabozantinib in dosage form and human plasma via micelle-enhanced spectrofluorimetry. <i>RSC Advances</i> , 2015, 5, 40484-40490.	1.7	11
36	Orthogonal projection to latent structures and first derivative for manipulation of PLSR and SVR chemometric models' prediction: A case study. <i>PLoS ONE</i> , 2019, 14, e0222197.	1.1	11

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37	Micellar Enhanced Spectrofluorimetric Method for the Determination of Ponatinib in Human Plasma and Urine via Cremophor RH 40 as Sensing Agent. <i>International Journal of Analytical Chemistry</i> , 2015, 1-9.	0.4	10
38	Spectrophotometric and computational investigations of charge transfer complexes of chloranilic acid with tyrosine kinase inhibitors and application to development of novel universal 96-microwell assay for their determination in pharmaceutical formulations. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2021, 252, 119482.	2.0	10
39	Synthesis, spectroscopic and computational studies on hydrogen bonded charge transfer complex of duvelisib with chloranilic acid: Application to development of novel 96-microwell spectrophotometric assay. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2022, 264, 120287.	2.0	10
40	Improved partial least squares models for stability-indicating analysis of mebeverine and sulpiride mixtures in pharmaceutical preparation: A comparative study. <i>Drug Testing and Analysis</i> , 2013, 5, 325-333.	1.6	9
41	A Stability-Indicating HPLC-DAD Method for Determination of Stiripentol: Development, Validation, Kinetics, Structure Elucidation and Application to Commercial Dosage Form. <i>Journal of Analytical Methods in Chemistry</i> , 2014, 2014, 1-10.	0.7	9
42	Full spectrum and selected spectrum based multivariate calibration methods for simultaneous determination of betamethasone dipropionate, clotrimazole and benzyl alcohol: Development, validation and application on commercial dosage form. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2016, 169, 50-57.	2.0	9
43	Simultaneous quantitative analysis of olmesartan, amlodipine and hydrochlorothiazide in their combined dosage form utilizing classical and alternating least squares based chemometric methods. <i>Acta Pharmaceutica</i> , 2016, 66, 83-95.	0.9	9
44	Stability-indicating chemometric methods for the determination of tazarotene. <i>Drug Testing and Analysis</i> , 2010, 2, 357-361.	1.6	8
45	Stability indicating spectrophotometric methods for quantitative determination of bromazepam and its degradation product. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2020, 238, 118433.	2.0	8
46	A new spectrofluorimetric assay method for vandetanib in tablets, plasma and urine. <i>Tropical Journal of Pharmaceutical Research</i> , 2016, 15, 2219.	0.2	7
47	Enhanced spectrofluorimetric determination of the multitargeted tyrosine kinase inhibitor, crizotinib, in human plasma via micelle-mediated approach. <i>Tropical Journal of Pharmaceutical Research</i> , 2016, 15, 2209.	0.2	7
48	LC-ESI-MS/MS identification and characterization of ponatinib in vivo phase I and phase II metabolites. <i>Clinica Chimica Acta</i> , 2018, 485, 144-151.	0.5	7
49	Validated liquid chromatography tandem mass spectrometry for simultaneous quantification of foretinib and lapatinib, and application to metabolic stability investigation. <i>RSC Advances</i> , 2019, 9, 19325-19332.	1.7	7
50	Development of innovative artificial neural networks for simultaneous determination of lapatinib and foretinib in human urine by micellar enhanced synchronous spectrofluorimetry. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2020, 238, 118438.	2.0	7
51	Experimental and Computational Evaluation of Chloranilic Acid as an Universal Chromogenic Reagent for the Development of a Novel 96-Microwell Spectrophotometric Assay for Tyrosine Kinase Inhibitors. <i>Molecules</i> , 2021, 26, 744.	1.7	7
52	Development and validation of a stability indicating RP-HPLC-DAD method for the determination of bromazepam. <i>PLoS ONE</i> , 2021, 16, e0244951.	1.1	7
53	A Validated LC-MS/MS Assay for the Simultaneous Quantification of the FDA-Approved Anticancer Mixture (Encorafenib and Binimetinib): Metabolic Stability Estimation. <i>Molecules</i> , 2021, 26, 2717.	1.7	7
54	Irbesartan (a comprehensive profile). <i>Profiles of Drug Substances, Excipients and Related Methodology</i> , 2021, 46, 185-272.	3.5	6

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55	New spectrofluorimetric methods for determination of melatonin in the presence of N-{2-[1-({3-[2-(acetylamino)ethyl]-5-methoxy-1H-indol-2-yl)methyl]-5-methoxy-1H-indol-3-yl]-ethyl}acetamide: a contaminant in commercial melatonin preparations. <i>Chemistry Central Journal</i> , 2012, 6, 36.	2.6	5
56	Development of novel response surface methodology-assisted micellar enhanced synchronous spectrofluorimetric method for determination of vandetanib in tablets, human plasma and urine. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2019, 213, 272-280.	2.0	5
57	Comparative Study of Novel Ratio Spectra and Isoabsorptive Point Based Spectrophotometric Methods: Application on a Binary Mixture of Ascorbic Acid and Rutin. <i>Journal of Analytical Methods in Chemistry</i> , 2016, 2016, 1-12.	0.7	4
58	Characterization of in vivo metabolites in rat urine following an oral dose of masitinib by liquid chromatography tandem mass spectrometry. <i>Chemistry Central Journal</i> , 2018, 12, 61.	2.6	4
59	Five modified classical least squares based models for stability indicating analysis of cyclobenzaprine HCl with its major degradation products: A comparative study. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2018, 204, 598-602.	2.0	4
60	Metabolic Stability Assessment of Larotrectinib Using Liquid Chromatography Tandem Mass Spectrometry. <i>Drug Design, Development and Therapy</i> , 2020, Volume 14, 111-119.	2.0	4
61	Design, Synthesis, Characterization of Novel Ruthenium(II) Catalysts: Highly Efficient and Selective Hydrogenation of Cinnamaldehyde to (E)-3-Phenylprop-2-en-1-ol. <i>Molecules</i> , 2014, 19, 5965-5980.	1.7	3
62	Experimental and computational evaluation of kolliphor RH 40 as a new fluorescence enhancer in development of a micellar-based spectrofluorimetric method for determination of lapatinib in tablets and urine. <i>PLoS ONE</i> , 2020, 15, e0239918.	1.1	3
63	Synthesis, spectroscopic and computational characterization of charge transfer complex of remdesivir with chloranilic acid: Application to development of novel 96-microwell spectrophotometric assay. <i>Journal of Molecular Structure</i> , 2022, 1263, 133104.	1.8	3
64	DNA Binding Test, X-Ray Crystal Structure, Spectral Studies, TG-DTA, and Electrochemistry of [CoX ₂ (dmdpphen)] (Dmdpphen Is 2,9-Dimethyl-4,7-diphenyl-1,10-phenanthroline, X = Cl, and NCS) Complexes. <i>Bioinorganic Chemistry and Applications</i> , 2014, 2014, 1-7.	1.8	2
65	Highly Sensitive Micellar Enhanced Spectrofluorimetric Method for Determination of Mirtazapine in Tablets and Human Urine: Application to In Vitro Drug Release and Content Uniformity Test. <i>Journal of Chemistry</i> , 2016, 2016, 1-8.	0.9	2
66	Determination of Cefoperazone Sodium in Presence of Related Impurities by Improved Classical Least Squares Chemometric Methods: A Comparative Study. <i>Journal of Chemistry</i> , 2016, 2016, 1-8.	0.9	2
67	Reactive intermediates in naquotinib metabolism identified by liquid chromatography-tandem mass spectrometry: phase I metabolic profiling. <i>RSC Advances</i> , 2019, 9, 10211-10225.	1.7	2
68	Novel spectrofluorimetric determination of brigatinib in bulk powder and human urine samples via ion-pair complex formation using eosin Y. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2021, 249, 119210.	2.0	2
69	Crystal structure of (2E)-2-[3-(1H-imidazol-1-yl)-1-phenylpropylidene]-N-(4-methylphenyl)hydrazinecarboxamide, C ₂₀ H ₂₁ N ₅ O. <i>Zeitschrift Fur Kristallographie - New Crystal Structures</i> , 2014, 229, 311-312.	0.1	1
70	Quantification of total mercury in compact fluorescent lamps in Saudi Arabia using inductively coupled plasma mass spectrometry (ICP-MS). <i>Main Group Chemistry</i> , 2016, 15, 117-124.	0.4	1
71	A novel method to determine new potent angiotensin inhibitor, azilsartan, in human plasma via micelle-enhanced spectrofluorimetry using cremophor RH 40. <i>Tropical Journal of Pharmaceutical Research</i> , 2016, 15, 1003.	0.2	1
72	Full spectrum and genetic algorithm-selected spectrum-based chemometric methods for simultaneous determination of azilsartan medoxomil, chlorthalidone, and azilsartan: Development, validation, and application on commercial dosage form. <i>Open Chemistry</i> , 2021, 19, 205-213.	1.0	1

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73	Innovative use of Ir and Ir electron acceptors in the development of three high throughput 96-microwell spectrophotometric assays for crizotinib. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2021, 259, 119884.	2.0	1