Genyuan Li

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	High Dimensional Model Representations. Journal of Physical Chemistry A, 2001, 105, 7765-7777.	1.1	403
2	Practical Approaches To Construct RS-HDMR Component Functions. Journal of Physical Chemistry A, 2002, 106, 8721-8733.	1.1	234
3	Global Sensitivity Analysis for Systems with Independent and/or Correlated Inputs. Journal of Physical Chemistry A, 2010, 114, 6022-6032.	1.1	183
4	Random Sampling-High Dimensional Model Representation (RS-HDMR) and Orthogonality of Its Different Order Component Functions. Journal of Physical Chemistry A, 2006, 110, 2474-2485.	1.1	158
5	Global uncertainty assessments by high dimensional model representations (HDMR). Chemical Engineering Science, 2002, 57, 4445-4460.	1.9	157
6	High Dimensional Model Representations Generated from Low Dimensional Data Samples. I. mp-Cut-HDMR. Journal of Mathematical Chemistry, 2001, 30, 1-30.	0.7	127
7	General formulation of HDMR component functions with independent and correlated variables. Journal of Mathematical Chemistry, 2012, 50, 99-130.	0.7	108
8	Random Samplingâ^'High Dimensional Model Representation (RSâ^'HDMR) with Nonuniformly Distributed Variables:  Application to an Integrated Multimedia/Multipathway Exposure and Dose Model for Trichloroethylene. Journal of Physical Chemistry A, 2003, 107, 4707-4716.	1.1	68
9	Regularized random-sampling high dimensional model representation (RS-HDMR). Journal of Mathematical Chemistry, 2008, 43, 1207-1232.	0.7	59
10	D-MORPH regression: application to modeling with unknown parameters more than observation data. Journal of Mathematical Chemistry, 2010, 48, 1010-1035.	0.7	51
11	Correlation method for variance reduction of Monte Carlo integration in RS-HDMR. Journal of Computational Chemistry, 2003, 24, 277-283.	1.5	38
12	Relationship between sensitivity indices defined by variance- and covariance-based methods. Reliability Engineering and System Safety, 2017, 167, 136-157.	5.1	34
13	Multicut-HDMR with an application to an ionospheric model. Journal of Computational Chemistry, 2004, 25, 1149-1156.	1.5	32
14	Ratio control variate method for efficiently determining high-dimensional model representations. Journal of Computational Chemistry, 2006, 27, 1112-1118.	1.5	32
15	D-MORPH regression for modeling with fewer unknown parameters than observation data. Journal of Mathematical Chemistry, 2012, 50, 1747-1764.	0.7	32
16	High-dimensional model representations generated from low order terms?lp-RS-HDMR. Journal of Computational Chemistry, 2003, 24, 647-656.	1.5	30
17	High efficiency classification of children with autism spectrum disorder. PLoS ONE, 2018, 13, e0192867.	1.1	13
18	High dimensional model representation constructed by support vector regression. I. Independent variables with known probability distributions. Journal of Mathematical Chemistry, 2017, 55, 278-303.	0.7	10

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19	A special singular perturbation method for kinetic model reduction: With application to an H2/O2 oxidation model. Journal of Chemical Physics, 1996, 105, 4065-4075.	1.2	9
20	Analytical HDMR formulas for functions expressed as quadratic polynomials with a multivariate normal distribution. Journal of Mathematical Chemistry, 2014, 52, 2052-2073.	0.7	8
21	Experimental Design of Formulations Utilizing High Dimensional Model Representation. Journal of Physical Chemistry A, 2015, 119, 8237-8249.	1.1	8
22	A lumped model for H2/O2 oxidation in the oscillatory regime. Journal of Chemical Physics, 1995, 102, 7006-7016.	1.2	6
23	A scalable algorithm for molecular property estimation in high dimensional scaffold-based libraries. Journal of Mathematical Chemistry, 2012, 50, 1765-1790.	0.7	5
24	Sparse and nonnegative sparse D-MORPH regression. Journal of Mathematical Chemistry, 2015, 53, 1885-1914.	0.7	4
25	Determination of rate constants for butene isomerization by a temporal inversion method. Journal of Chemical Physics, 1997, 107, 2845-2852.	1.2	2
26	Global Sensitivity Analysis with Mixtures: A Generalized Functional ANOVA Approach. Risk Analysis, 2022, 42, 304-333.	1.5	2
27	Enhancing molecular discovery using descriptorâ€free rearrangement clustering techniques for sparse data sets. AICHE Journal, 2010, 56, 405-418.	1.8	1
28	Molecular discovery by optimal sequential search. Journal of Mathematical Chemistry, 2019, 57, 2110-2141.	0.7	0