

ErgÃ¼n Kasap

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Impact of pressure and composition on the mechanical behavior of In Ga As P and Al In Sb P quaternary alloys. <i>Results in Physics</i> , 2019, 14, 102400.	4.1	10
2	Design studies for the beam position monitor (BPM) front-end electronics of the Turkish accelerator and radiation laboratory in Ankara (TARLA). <i>Turkish Journal of Physics</i> , 2017, 41, 269-276.	1.1	3
3	Torsional potential and nonlinear optical properties of phenyldiazines and phenyltetrazines. <i>Computational and Theoretical Chemistry</i> , 2011, 977, 22-28.	2.5	11
4	Comparing of the host-guest interaction in the Hofmann-1,10-diaminodecane and Hofmann-1,12-diaminododecane-type clathrates. <i>Journal of Inclusion Phenomena and Macrocyclic Chemistry</i> , 2010, 66, 243-249.	1.6	1
5	The structural, electronic and optical properties of $\text{In}_x\text{Ga}_{1-x}\text{P}$ alloys. <i>Physica B: Condensed Matter</i> , 2010, 405, 2357-2361.	2.7	23
6	Ab-initio investigation of structural, electronic and optical properties of $\text{In}_x\text{Ga}_{1-x}\text{As}$, $\text{GaAs}_{1-y}\text{Py}$ ternary and $\text{In}_x\text{Ga}_{1-x}\text{As}_{1-y}\text{Py}$ quaternary semiconductor alloys. <i>Journal of Alloys and Compounds</i> , 2010, 496, 226-233.	5.5	35
7	Conformational Analysis, Dipole Moment and Polarizability of 3-(2-chlorophenyl)thiophene. <i>AIP Conference Proceedings</i> , 2007, , .	0.4	0
8	A theoretical study of molecular structure and potential energy surface for various substituents substituted 3-phenylthiophene. <i>AIP Conference Proceedings</i> , 2007, , .	0.4	0
9	A theoretical study of the linear, nonlinear optical properties and conformational analysis of 3-phenylthiophene and its fluoro derivatives with torsional dependence. <i>Journal of Molecular Structure</i> , 2007, 834-836, 508-515.	3.6	5
10	Investigation of torsional barriers and nonlinear optical (NLO) properties of phenyltriazines. <i>Journal of Molecular Structure</i> , 2007, 834-836, 516-520.	3.6	59
11	Infrared Spectroscopic Study on the Hofmann-dadn-type and the Td-dahxn-type clathrates. <i>Journal of Inclusion Phenomena and Macrocyclic Chemistry</i> , 2007, 58, 49-54.	1.6	1
12	Torsional barriers and nonlinear optical properties of 2-, 3-, 4-phenylpyridine molecules. <i>European Physical Journal D</i> , 2006, 56, 349-358.	0.4	11
13	Infrared Spectroscopic Study of Td-type Piperazinemetal(II) Tetracyanometallate(II) Benzene(1/1) Clathrates: $\text{Cd}(\text{C}_4\text{H}_{10}\text{N}_2)\text{Cd}(\text{CN})_4 \cdot \text{C}_6\text{H}_6$ and $\text{Cd}(\text{C}_4\text{H}_{10}\text{N}_2)\text{Hg}(\text{CN})_4 \cdot 1,25 \text{C}_6\text{H}_6$. <i>Spectroscopy Letters</i> , 2005, 38, 583-594.	1.0	9
14	VIBRATIONAL SPECTROSCOPIC STUDIES ON THE 1,4-DIAMINOBUTANE-Td-TYPE CLATHRATES: $\text{Cd}(\text{dabn})\text{M}(\text{CN})_4 \cdot 1,5\text{C}_6\text{H}_6$ ($\text{M}=\text{Cd}$ or Hg). <i>Spectroscopy Letters</i> , 2002, 35, 811-819.	1.0	5
15	Title is missing!. <i>Journal of Inclusion Phenomena and Macrocyclic Chemistry</i> , 2001, 39, 175-180.	1.6	12
16	Title is missing!. <i>Journal of Inclusion Phenomena and Macrocyclic Chemistry</i> , 2001, 39, 103-108.	1.6	1
17	Infrared spectroscopic study of the Hofmann-dadn-type clathrates: $\text{M}(1,10\text{-diaminodecane})\text{Ni}(\text{CN})_4 \cdot 1,5\text{G}$ ($\text{M}=\text{Co}$, Ni or Cd ; $\text{G}=\text{chlorobenzene}$, 1,2-, 1,3- or 1,4-dichlorobenzene). <i>Vibrational Spectroscopy</i> , 2000, 24, 249-255.	2.2	4
18	An infrared and Raman spectroscopic study on the Hofmann-Td-type 1,4-dioxane clathrates: $\text{M}(\text{NH}_3)_2\text{M}^{\text{2+}}(\text{CN})_4 \cdot 2\text{C}_4\text{H}_8\text{O}_2$ ($\text{M}=\text{Mn}$ or Cd , $\text{M}^{\text{2+}}=\text{Hg}$; $\text{M}=\text{Cd}$, $\text{M}^{\text{2+}}=\text{Cd}$). <i>Journal of Molecular Structure</i> , 1999, 3, 482-483, 81-85.	3.6	8

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19	Investigation of host-guest interactions in the Hofmann-dabn-type clathrates: M(1,4-diaminobutane)Ni(CN) ₄ ·1.5G (M=Co or Ni, G=benzene derivatives). Journal of Molecular Structure, 1999, 482-483, 69-74.	3.6	3
20	Infrared spectroscopic studies of the Hofmann-daon-type clathrates: M(1,8-diaminoctane)Ni(CN) ₄ ·G (M=Co, Ni or Cd; G=1,2-dichlorobenzene or 1,4-dichlorobenzene). Journal of Molecular Structure, 1999, 482-483, 75-79.	3.6	9
21	Title is missing!. Journal of Inclusion Phenomena and Macrocyclic Chemistry, 1999, 33, 285-294.	1.6	3
22	Infrared Spectroscopic Study of the Hofmann-Diam-Type Clathrates: M(1,6-Diaminohexane)Ni(CN) ₄ ·C ₆ H ₆ (M=Ni, Co or Cd). Spectroscopy Letters, 1997, 30, 491-496.	1.0	15
23	Title is missing!. Journal of Inclusion Phenomena and Macrocyclic Chemistry, 1997, 28, 117-124.	1.6	8
24	Title is missing!. Journal of Inclusion Phenomena and Macrocyclic Chemistry, 1997, 28, 259-267.	1.6	17
25	Title is missing!. Journal of Inclusion Phenomena and Macrocyclic Chemistry, 1997, 28, 335-347.	1.6	13
26	Infrared and Raman spectroscopic study of the Hofmann-type clathrates M(1,7-diaminoheptane)Ni(CN) ₄ ·G (M = Ni or Co; G = chlorobenzene, m-xylene or naphthalene). Journal of Molecular Structure, 1997, 408-409, 425-430.	3.6	9
27	Vibrational spectroscopic studies on the en-Td-type benzene clathrates: M(ethylenediamine)M?(CN) ₄ ? _{1/2} C ₆ H ₆ (M=Mn or Cd, M?=Cd or Hg). Journal of Inclusion Phenomena and Macrocyclic Chemistry, 1995, 23, 1-9.	1.6	20
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