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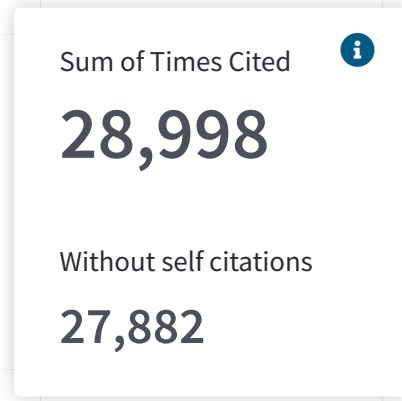
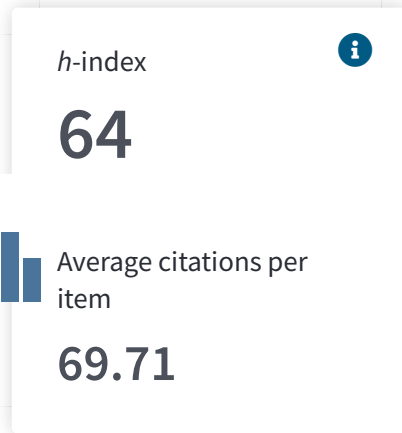
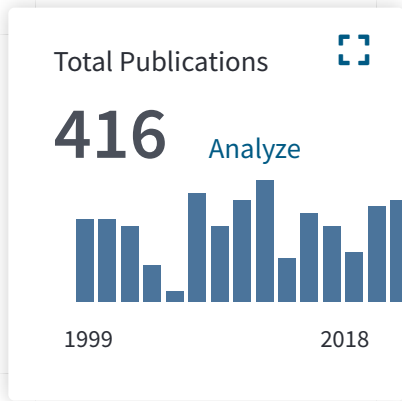


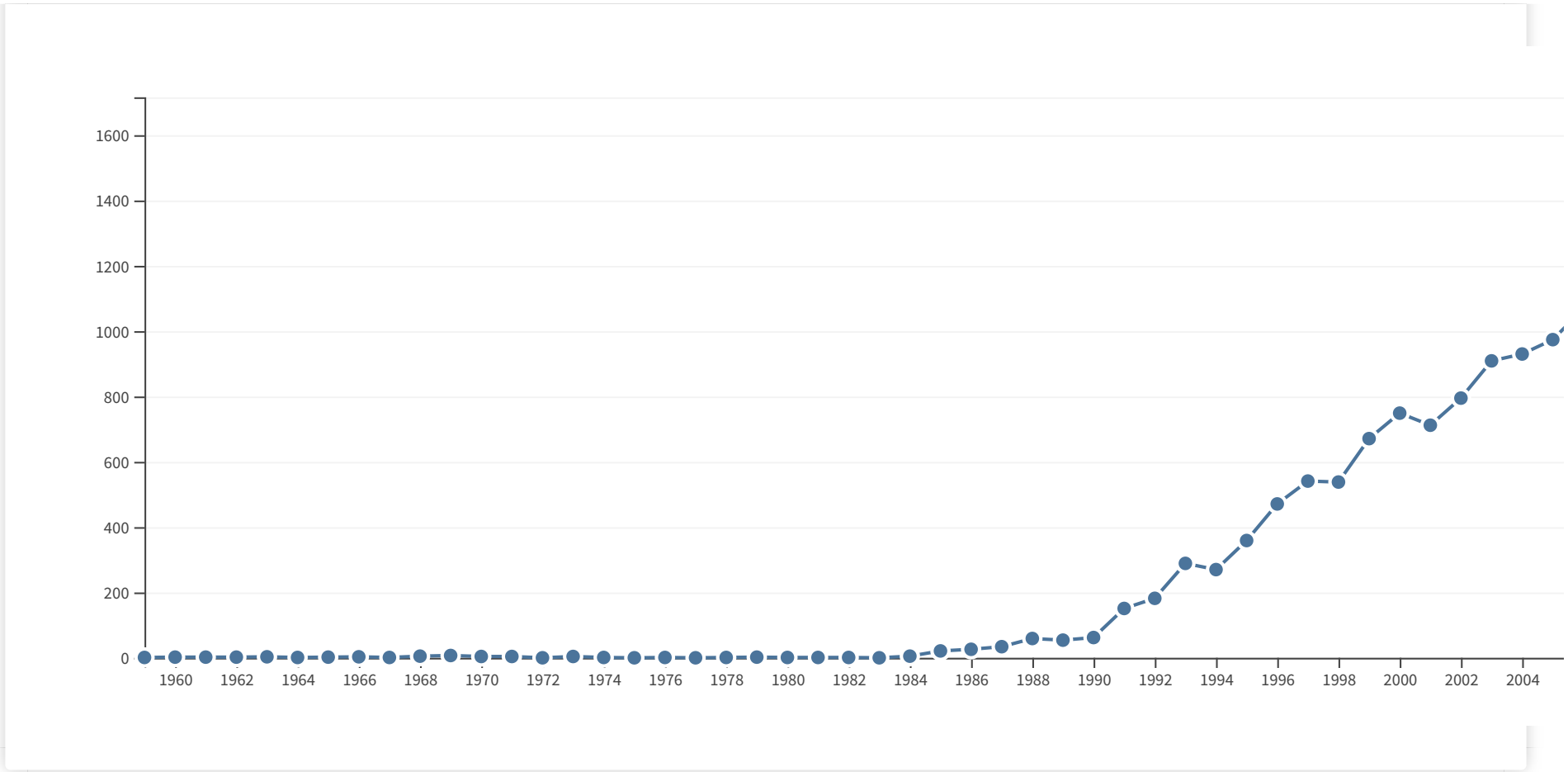
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JOURNAL OF THE AMERICAN CHEMICAL SOCIETY Volume: 117 Issue: 19 Pages: 5179-5197 Published: MAY 17 1995

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By: Case, DA; Cheatham, TE; Darden, T; et al.

JOURNAL OF COMPUTATIONAL CHEMISTRY Volume: 26 Issue: 16 Pages: 1668-1688 Published: DEC 2005

359 398 360 369 185 4317 287.80

3. **ATOMIC CHARGES DERIVED FROM SEMIEMPIRICAL METHODS**

By: BESLER, BH; MERZ, KM; KOLLMAN, PA

JOURNAL OF COMPUTATIONAL CHEMISTRY Volume: 11 Issue: 4 Pages: 431-439 Published: MAY 1990

104 125 117 134 53 2451 81.70

4. **Prediction of drug absorption using multivariate statistics**

By: Egan, WJ; Merz, KM; Baldwin, JJ

JOURNAL OF MEDICINAL CHEMISTRY Volume: 43 Issue: 21 Pages: 3867-3877 Published: OCT 19 2000

31 28 34 42 28 355 17.75

5. **A second generation force field for the simulation of proteins, nucleic acids, and organic molecules (vol 117, pg 5179, 1995)**

	By: Cornell, WD; Cieplak, P; Bayly, CI; et al. JOURNAL OF THE AMERICAN CHEMICAL SOCIETY Volume: 118 Issue: 9 Pages: 2309-2309 Published: MAR 6 1996	22	27	24	27	5	308	12.83
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By: HOOPS, SC; ANDERSON, KW; MERZ, KM

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8262-8270 Published: OCT 23 1991

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