4 Bromoacetanilide Ir Spectrum Analysis

4-Bromoacetanilide IR Spectrum Analysis: A Comprehensive Guide

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Dr. Eleanor Vance holds a PhD in Analytical Chemistry from Stanford University and has over 15 years of experience in spectroscopic techniques, with a particular focus on infrared (IR) spectroscopy and its application in organic compound characterization. Her research has been extensively published in peer-reviewed journals, including several publications directly related to the analysis of substituted anilides using IR spectroscopy. Her expertise makes her uniquely qualified to provide a detailed analysis of 4-bromoacetanilide IR spectrum analysis.

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Editor: Dr. Robert Miller, PhD, Organic Chemistry

Dr. Robert Miller, PhD in Organic Chemistry from MIT, brings over 20 years of experience in organic synthesis and characterization to the editorial process. His expertise in functional group analysis and spectroscopic interpretation enhances the article's scientific rigor and ensures clarity for a wide range of readers. His in-depth understanding of organic molecules, specifically substituted aromatic amides, provides valuable oversight for the accuracy and completeness of the 4-bromoacetanilide IR spectrum analysis presented.

1. Historical Context of IR Spectroscopy and its Application to 4-Bromoacetanilide

Infrared spectroscopy, a cornerstone of analytical chemistry, has a rich history dating back to the early 19th century. The development of Fourier Transform Infrared (FTIR) spectroscopy in the latter half of the 20th century revolutionized the field, allowing for faster and more sensitive analysis. This advancement significantly impacted the analysis of organic compounds like 4-bromoacetanilide. Early applications focused on qualitative identification based on characteristic absorption bands. The advent of computational chemistry further enhanced the interpretation of IR spectra, enabling more precise predictions and comparisons with experimental data obtained from 4-bromoacetanilide

2. Understanding the 4-Bromoacetanilide Molecule

4-Bromoacetanilide is an aromatic compound consisting of a benzene ring substituted with a bromine atom at the para position (4-position) and an acetamide group (-NHCOCH3). This specific arrangement of functional groups dictates the characteristic absorption bands observed in its IR spectrum. The presence of the bromine atom influences the electron density distribution within the molecule, subtly affecting the vibrational frequencies. A thorough understanding of 4-bromoacetanilide's molecular structure is crucial for accurately interpreting its IR spectrum.

3. Key Absorption Bands in the 4-Bromoacetanilide IR Spectrum

The 4-bromoacetanilide IR spectrum analysis reveals several key absorption bands that provide crucial information about its functional groups. These include:

N-H stretching: A sharp peak typically observed around 3300-3500 cm⁻¹, indicating the presence of the amide N-H bond. The exact position can be slightly affected by hydrogen bonding.

C=O stretching: A strong absorption band usually appearing around 1650-1700 $\rm cm^{-1}$, characteristic of the carbonyl group (C=O) in the acetamide moiety. The position and intensity can vary depending on intermolecular interactions.

C-N stretching: A medium intensity band usually observed around $1250-1350~cm^{-1}$, associated with the C-N stretching vibration in the amide group.

Aromatic C-H stretching: Several bands appear in the 3000-3100 cm⁻¹ region, representative of the aromatic C-H stretching vibrations of the benzene ring.

C-Br stretching: A weak to medium absorption band might be observed in the 500-600 cm⁻¹ region, corresponding to the C-Br stretching vibration. This band's intensity can vary depending on the instrument and sample preparation.

Careful analysis of the positions, intensities, and shapes of these absorption bands provides conclusive evidence for the presence of all functional groups expected in 4-bromoacetanilide and allows for confirmation of its molecular structure via 4-bromoacetanilide IR spectrum analysis.

4. Interpretation and Data Analysis of the 4-Bromoacetanilide IR Spectrum

The analysis of a 4-bromoacetanilide IR spectrum involves comparing the observed absorption bands

with known spectral data for similar compounds and utilizing spectral databases. Software packages associated with FTIR instruments often include libraries of reference spectra for comparison. This comparison allows for unambiguous identification of the compound and confirmation of its purity. Furthermore, deviations from expected band positions can offer clues to the presence of impurities or unusual intermolecular interactions. Quantitative analysis, though more challenging with IR, can be performed under controlled conditions to determine the concentration of 4-bromoacetanilide in a mixture.

5. Current Relevance of 4-Bromoacetanilide IR Spectrum Analysis

Despite the emergence of more sophisticated analytical techniques, 4-bromoacetanilide IR spectrum analysis remains highly relevant for several reasons:

Rapid and relatively inexpensive: IR spectroscopy is a relatively fast and cost-effective technique compared to more advanced methods like NMR or mass spectrometry.

Simple sample preparation: IR sample preparation is generally straightforward, requiring minimal sample handling.

Qualitative and quantitative applications: IR spectroscopy can be used for both qualitative identification and quantitative analysis of 4-bromoacetanilide.

Versatile applicability: IR spectroscopy is applicable to a wide range of sample types, including solids, liquids, and gases.

Complementary technique: IR spectroscopy often serves as a complementary technique to other analytical methods, providing valuable supporting data.

6. Applications of 4-Bromoacetanilide and its IR Spectral Analysis

4-Bromoacetanilide, although not widely used as a final product, serves as an important intermediate in the synthesis of various organic compounds. Its IR spectrum analysis is crucial during synthesis and purification to ensure the desired product's identity and purity. This is particularly important in pharmaceutical and agrochemical industries where accurate characterization is essential.

7. Limitations of 4-Bromoacetanilide IR Spectrum Analysis

While powerful, 4-bromoacetanilide IR spectrum analysis has limitations:

Overlapping bands: In complex mixtures, overlapping absorption bands can complicate interpretation.

Sensitivity limitations: IR spectroscopy might not be sensitive enough to detect trace impurities.

Qualitative nature (primarily): While quantitative analysis is possible, it is generally less precise than other methods.

8. Future Trends in IR Spectroscopy and 4-Bromoacetanilide Analysis

Ongoing research focuses on improving the sensitivity and resolution of IR spectroscopy. Coupling IR with other techniques, such as microscopy or chromatography, leads to more comprehensive analysis. Advances in computational chemistry are further enhancing the ability to predict and interpret IR spectra, leading to more accurate 4-bromoacetanilide IR spectrum analysis.

9. Conclusion

4-Bromoacetanilide IR spectrum analysis is a valuable technique with a long history and continued relevance in analytical chemistry. Its straightforward procedure, coupled with its ability to provide crucial information about the compound's functional groups and purity, makes it a valuable tool in both research and industrial settings. While limitations exist, ongoing advancements are enhancing its capabilities and expanding its applications.

FAQs

- 1. What is the primary use of 4-bromoacetanilide? Primarily an intermediate in organic synthesis, not typically used as a final product.
- 2. What are the most important peaks to look for in a 4-bromoacetanilide IR spectrum? N-H stretch, C=O stretch, and aromatic C-H stretches are crucial.
- 3. Can IR spectroscopy quantify 4-bromoacetanilide? Yes, but with less precision than other methods like HPLC or GC.
- 4. What are the potential interferences in the 4-bromoacetanilide IR spectrum? Overlapping peaks from impurities or solvents.
- 5. How does the presence of bromine affect the IR spectrum? It subtly affects the electron distribution, influencing vibrational frequencies.

- 6. What is the difference between FTIR and traditional IR spectroscopy? FTIR is faster and more sensitive due to the use of Fourier Transform mathematics.
- 7. What software is commonly used for IR spectrum analysis? Software packages specific to FTIR instruments, often with spectral libraries.
- 8. What is the typical sample preparation for 4-bromoacetanilide IR analysis? KBr pellet or solution in a suitable solvent.
- 9. Can IR spectroscopy distinguish between isomers of 4-bromoacetanilide? Potentially, depending on the isomer's differing vibrational modes.

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 aliphatic and aromatic carboxylic acids, substituted phenols and mineral acids. The bases of
 principal interest are likewise compounds capable of forming ions in water, for example, aliphatic
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 basicity scales in aprotic and water-like solvents are discussed. The role of hydrogen bonding in
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4 Bromoacetanilide Ir Spectrum Analysis: Microscale and Miniscale Organic Chemistry Laboratory Experiments Allen M. Schoffstall, Barbara A. Gaddis, Melvin L. Druelinger, 2000 This work offers ...

IR Spectrum analysis of p-bromoacetanilide - Seth ...

Two moderately intense bands are observed. The appearance of two bands is attributed to Fermi resonance between the fundamental aldehydic C–H stretching and the first overton.

SPECTROSCOPIC ANALYSIS OF ORGANIC COMPOUNDS

IR Spectral Analysis of 4'-Bromoacetanilide • 3300: N-H stretching: N-H bond contains a small atom H, so frequency (ν) and hence wavenumber increases by Hooke's law.

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Interpretation and Data Analysis of the 4-Bromoacetanilide IR Spectrum The analysis of a 4-bromoacetanilide IR spectrum involves comparing the observed absorption bands with known ...

SEM V (H) SPECTROSCOPIC ANALYSIS OF ORGANIC ...

Spectroscopic Analysis of Organic Compounds b-Disubstituted benzene with non-identical substituents For these compounds HA,HB, HC and 1--16 are chemically as well as ...

Practical Organic Chemistry

4-Bromoacetanilide: 1H NMR spectra in DMSO-d 6 The amine proton is in conjugation with the carbonyl group, thus making the N atom partially more electronegative. Hence, b proton ...

Introducing the Concept of Green Synthesis in the ...

A practical, two-step synthesis of 4-bromoacetanilide from aniline through a greener approach is described. Design of the synthetic route is based on the fundamental concepts of green ...

51LC EXP #7 SELECTIVITY IN THE BROMINATION OF ...

Oct 6, $2011 \cdot$ As part of you in-lab work, analyze both the IR and NMR spectra for the bromination products of acetanilide and 4-methylacetanilide. The most important one is the 1 H NMR spectra.

1007 Synthesis of 2,4,6-tribromoaniline from 4-bromoacetanilide

2.14 g (10.0 mmol) 4-bromoacetanilid are dissolved in 7 mL ethanol under stirring and slightly heating in a 250 mL three-neck flask equipped with magnetic stir bar, reflux condenser and ...

A Multistep Synthesis Incorporating a Green

May 4, $2012 \cdot ABSTRACT$: Electrophilic aromatic substitution is a fundamental topic taught in the undergraduate organic chemistry curriculum. A multistep synthesis that includes a safer and ...

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HPLC determination of chlorine in air and water samples ...

The chromatographic peak for 4-bromoacetanilide was identified by comparison of its retention time with that of the authentic substance under the same chromato- graphic conditions, and by ...

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NMR SPECTROSCOPY SEM-4, CC-8 PART-6, PPT-18 - St.

The 1H NMR spectrum of this molecule (Figure 8) is superficially first-order, with a doublet near δ 7.51 for H B and H c (because they couple with H A) and a triplet near δ 7.88 for H A (coupling ...

Table of Characteristic IR Absorptions - vscht.cz

The IR spectrum of octane, plotted as transmission (left) and absorbance (right). As illustrated in the spectrum of octane, even simple organic molecules give rise to complex IR spectra.

51LC EXP #7 SPRING 2012 SELECTIVITY IN THE ...

As part of you in-lab work, analyze both the IR and NMR spectra for the bromination products of acetanilide and 4-methylacetanilide. These are available on the course website.

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4 Bromoacetanilide Ir Spectrum Analysis John C. Gilbert, Stephen F. Martin Techniques and Experiments for Organic Chemistry Addison Ault, 1973 Microscale and Miniscale Organic ...

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