# Robustness of SPME Arrow Immersion Sampling: Polycyclic Aromatic Hydrocarbons in Drinking Water

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#### **Abstract**

Analyzing polycyclic aromatic hydrocarbons (PAHs) in drinking water becomes challenging especially for hard water (high mineral content water) and/or when sodium thiosulfate being added for treating chlorine content in drinking water during analysis. In this article, repeatability for determination of PAHs in drinking water using SPME Arrow (PDMS) immersion technique with gas chromatography mass spectrometer (GC/MS) was carried out. Single SPME Arrow had been used in this experiment to determine its durability. The repeatability (n=100) which characterized as %RSD was less than 10.93% for all analytes, ranged between 1.74 % - 10.93 %. The relative peak area of the 100<sup>th</sup> injection against the 1<sup>st</sup> injection peak area for all analytes ranged between 0.600 to 0.947; indicating SPME Arrow fiber is dutiful for at least 100 injections with immersion technique. In conclusion, fully automated SPME Arrow immersion technique with GC/MS analysis is an excellent option to analyze PAH compounds in drinking water; with higher throughput and robustness.



#### **Keywords**

Polycyclic aromatic hydrocarbons (PAHs), solid phase microextraction (SPME), SPME Arrow, immersion, PDMS

#### Introduction

Robustness of SPME and SPME Arrow used for immersion sampling is much lesser compared headspace sampling. During immersion extraction, SPME Arrow phase is fully contacted with sample matrix; e.g. drinking water and tap water which containing minerals and residual chlorine. This will affect SPME and SPME Arrow performance. In this poster, repeatability for determination of PAHs in drinking water using SPME Arrow immersion technique with GC/MS was carried out. Automated SPME Arrow workflow was optimized and one (1) SPME Arrow (PDMS) was used throughout the analysis to determine SPME Arrow durability.

#### **Chemicals**

- Pak-Mix 16 in Cyclohexane (Neochema GmbH)
- Internal Standards Mix 25 in Acetone (Dr. Ehrenstorfer GmbH)
- Benzo[a]pyrene D12 in Cyclohexane (Dr. Ehrenstorfer GmbH)
- Sodium Thiosulfate (Sigma Aldrich)
- Acetone ( > 99.5% Reagent grade) (Sigma Aldrich)
- Water (HPLC grade) (J.T. Baker)

#### **Test Solution**

Switzerland tap water (drinkable, 250mg/L total minerals) spiked with

- 16 PAH compounds (100.8306 pg/mL for each PAH)
- 4 deuterated PAH compounds (136.4374 pg/mL for each deuterated PAH)

#### Instrumentations

- PAL RTC 120 system with 1.1 mm Ø 100 µm PDMS SPME Arrow (ARR11-P-100/20-P1)
- Shimadzu GC-2010 with Restek Rtxi-5MS Column (30 m X 0.25 mmID X 0.25 µm)
  - Inlet Temperature: 280°C
  - GC Carrier: Helium at 1.0 mL/min (Splitless)
  - GC Oven: Initial 35°C (5 minutes); 40°C/min to 150°C; 20°C/min to 250°C; 10°C/min to 305°C (22 minutes)
- GCMS-QP2010
  - Interface Temperature: 310°C
  - Source Temperature: 250°C
  - Mode: EI with SIM

Module. Post-conditioned at 250°C for 5 minutes.



Post-conditioning SPME Arrow at Conditioning Module.

Adding reducing agent for water samples containing residual chlorine.
 Intensive cleaning to eliminate minerals build up at SPME Arrow.

Figure 1: Automated SPME Arrow workflow modified for PAHs in water analysis.

#### **Analysis**

•15 mL Test Solution in 20 mL vial

Replicates of 100 samples analyzed using one (1) SPME Arrow

#### **Results**



Table 2: Reproducibility (%RSD) of relative peak area (corrected with corresponding ISTD) for each analyte (n=100).

SYSTEM

Ingenious sample handling

| <u>Analyte</u><br>Name/ISTD | <u>%RSD of</u><br><u>Relative Peak</u><br><u>Area</u> | <u>Analyte</u><br>Name/ISTD | <u>%RSD of</u><br><u>Relative Peak</u><br><u>Area</u> |
|-----------------------------|---|-----------------------------|---|
| N/Ace-d10                   | 4.68  | BaAnt/Chr-d12               | 2.27  |
| Acy/Ace-d10                 | 3.35  | Chr/Chr-d12                 | 2.15  |
| Ace/Ace-d10                 | 1.81  | BbFI/BaPyr-d12              | 4.17  |
| F/Ace-d10                   | 2.86  | BkFI/BaPyr-d12              | 4.45  |
| P/P-d10                     | 1.74  | BaPyr/BaPyr-d12             | 2.35  |
| Ant/P-d10                   | 3.87  | lpyr/BaPyr-d12              | 9.54  |
| FI/P-d10                    | 2.25  | DBahAnt/BaPyr-d12           | 10.93   |
| Pyr/P-d10                   | 2.75  | BghiPer/BaPyr-d12           | 7.60  |

## Table 3: Ratio for Peak Area of the 100<sup>th</sup> injection against the 1<sup>st</sup> injection.

| <u>Compound</u> | <u>Ratio</u> | <u>Compound</u> | <u>Ratio</u> | <u>Compound</u> | <u>Ratio</u> |
|-----------------|--------------|-----------------|--------------|-----------------|--------------|
| Ν               | 0.832        | Ant             | 0.854        | BkFl            | 0.826        |
| Acy             | 0.945        | FI              | 0.921        | BaPyr-d12       | 0.700        |
| Ace-d10         | 0.909        | Pyr             | 0.921        | BaPyr           | 0.724        |
| Ace             | 0.839        | BaAnt           | 0.846        | lpyr            | 0.678        |
| F               | 0.858        | Chr-d12         | 0.922        | DBahAnt         | 0.600        |
| P-d10           | 0.947        | Chr             | 0.904        | BghiPer         | 0.710        |
| Р               | 0.893        | BbFl            | 0.772        |                 |              |

## Table 1: Analytes details includes compound name, retention time and quantify ion.

| Compound Name          | <u>Abbrev.</u> | <u>Target/</u><br>ISTD | <u>Group</u> | <u>Retention</u><br><u>Time,</u><br><u>minutes</u> | <u>Quant. Ion,</u><br><u>m/z</u> |
|------------------------|----------------|------------------------|--------------|--|----------------------------------|
| Naphthalene            | Ν              | Target                 | 1            | 9.241  | 128                              |
| Acenaphthylene         | Асу            | Target                 | 1            | 10.688   | 152                              |
| Acenaphthene-d10       | Ace-d10        | ISTD                   | 1            | 10.832   | 164                              |
| Acenaphthene           | Ace            | Target                 | 1            | 10.863   | 153                              |
| Fluorene               | F              | Target                 | 1            | 11.418   | 166                              |
| Phenanthrene-d10       | P-d10          | ISTD                   | 2            | 12.482   | 188                              |
| Phenanthrene           | Р              | Target                 | 2            | 12.508   | 178                              |
| Anthracene             | Ant            | Target                 | 2            | 12.569   | 178                              |
| Fluoranthrene          | FI             | Target                 | 3            | 14.004   | 202                              |
| Pyrene                 | Pyr            | Target                 | 3            | 14.326   | 202                              |
| Benz(a)anthracene      | BaAnt          | Target                 | 3            | 16.258   | 228                              |
| Chrysene-d12           | Chr-d12        | ISTD                   | 3            | 16.275   | 240                              |
| Chrysene               | Chr            | Target                 | 3            | 16.320   | 228                              |
| Benz(b)fluoranthene    | BbFl           | Target                 | 4            | 18.204   | 252                              |
| Benz(k)fluoranthene    | BkFl           | Target                 | 4            | 18.252   | 252                              |
| Benz(a)pyrene-d12      | BaPyr-d12      | ISTD                   | 4            | 18.762   | 264                              |
| Benz(a)pyrene          | BaPyr          | Target                 | 4            | 18.803   | 252                              |
| Indeno(1,2,3-cd)pyrene | lpyr           | Target                 | 4            | 21.349   | 276                              |
| Dibenz(a,h)anthracene  | DBahAnt        | Target                 | 4            | 21.436   | 278                              |
| Benzo(g,h,i)perylene   | BghiPer        | Target                 | 4            | 22.043   | 276                              |

#### Figure 2: TIC Chromatogram.

(A) Full TIC chromatogram.

(B) Zoom in TIC chromatogram at 17.5 minutes to 25.0 minutes.

1-N 2-Acy 3-Ace-d10 4-Ace 5-F 6-P-d10 7-P 8-Ant 9-Fl 10-Pyr 11-BaAnt
12-Chr-d12 13-Chr 14-BbFl 15-BkFl 16-BaPyr-d12 17-BaPyr 18-Ipyr
19-DBahAnt 20-BghiPer

### Conclusions

SPME Arrow is robust to be used for immersion sampling technique with fully automated workflow. In PAHs in drinking water analysis, adequate cleaning procedures for SPME Arrow ensure good reproducibility results with satisfaction GC-MS response obtained for 100 sample injections.



#### **Presented at**

RAFA 2019, November 05-08, Prague, Czech Republic.

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