

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM090424\  
 Data File : BM047443.D  
 Acq On : 04 Sep 2024 18:02  
 Operator : RC/JU  
 Sample : P3750-03  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 346

Quant Time: Sep 05 01:53:08 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM082624.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Aug 26 18:11:40 2024  
 Response via : Initial Calibration

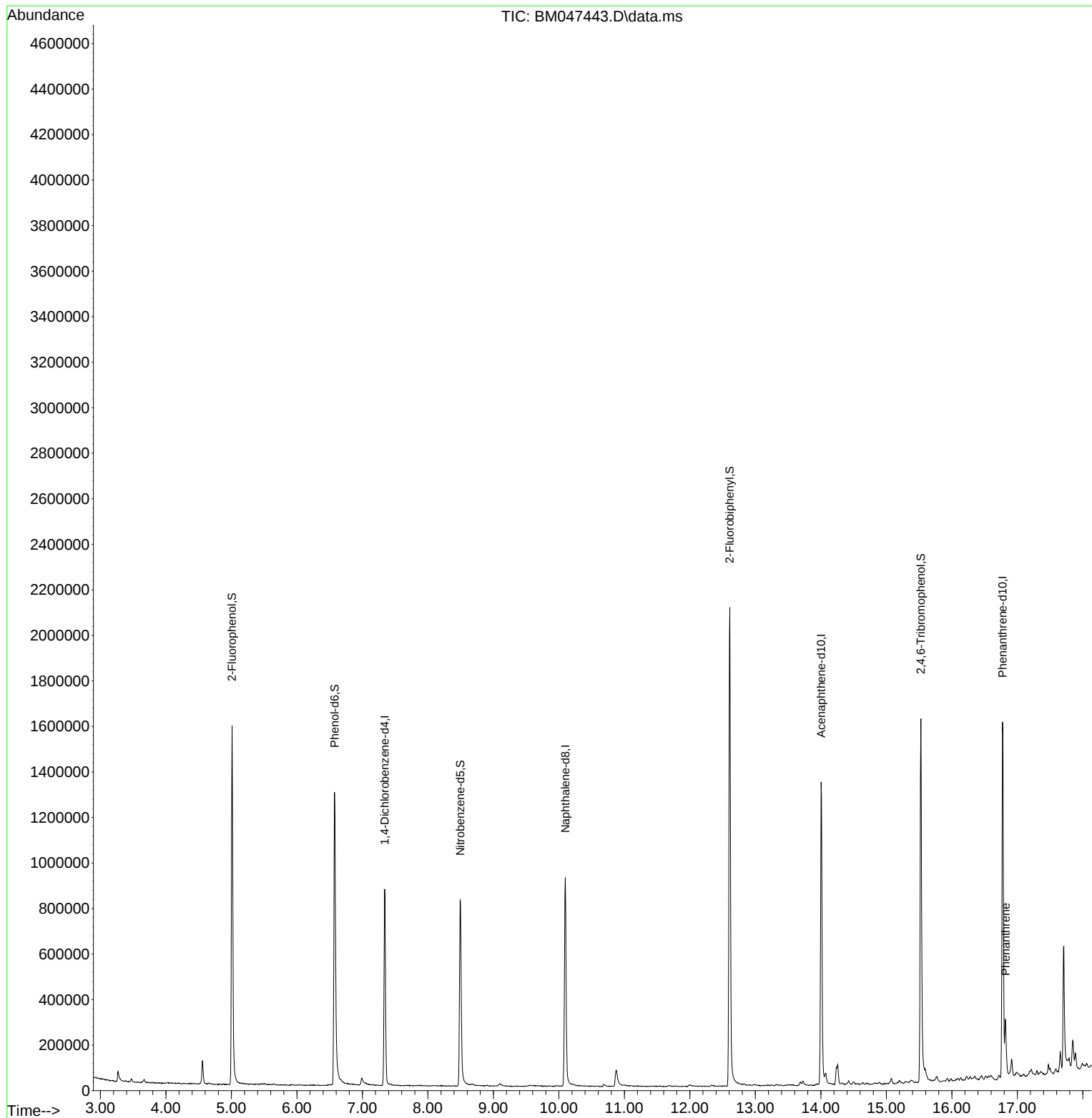
| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.339  | 152  | 256154   | 20.000 | ng    | -0.01    |
| 21) Naphthalene-d8          | 10.098 | 136  | 853340   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.004 | 164  | 534747   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 16.774 | 188  | 1059192  | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 21.021 | 240  | 1017256  | 20.000 | ng    | 0.00     |
| 86) Perylene-d12            | 23.727 | 264  | 1197681  | 20.000 | ng    | 0.01     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.010  | 112  | 747441   | 55.982 | ng    | 0.00     |
| 7) Phenol-d6                | 6.575  | 99   | 929816   | 51.122 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.492  | 82   | 566185   | 36.814 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.527 | 330  | 361000   | 55.955 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.610 | 172  | 1313979  | 37.755 | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.439 | 244  | 1861593  | 32.573 | ng    | 0.00     |
| Target Compounds            |        |      |          |        |       |          |
|                             |        |      |          |        |       | Qvalue   |
| 71) Phenanthrene            | 16.821 | 178  | 160791   | 3.126  | ng    | 100      |
| 75) Fluoranthene            | 18.856 | 202  | 388294   | 6.429  | ng    | 99       |
| 78) Pyrene                  | 19.221 | 202  | 417055   | 5.345  | ng    | 99       |
| 81) Benzo(a)anthracene      | 21.003 | 228  | 219418   | 3.308  | ng    | 96       |
| 83) Chrysene                | 21.062 | 228  | 275577   | 4.416  | ng    | 98       |
| 88) Benzo(b)fluoranthene    | 22.885 | 252  | 317944   | 4.550  | ng    | 97       |
| 90) Benzo(a)pyrene          | 23.603 | 252  | 192683   | 3.189  | ng    | # 96     |
| -----                       |        |      |          |        |       |          |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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