

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM111323\  
 Data File : BM042727.D  
 Acq On : 14 Nov 2023 07:38  
 Operator : MA/JU  
 Sample : 05016-04  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 A49A2

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 11/15/2023  
 Supervised By :mohammad ahmed 11/19/2023

Quant Time: Nov 14 15:05:04 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM111323.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 14 00:02:28 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.834  | 152  | 37290    | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.651 | 136  | 152590   | 20.000 | ng/u1  | # 0.00   |
| 38) Acenaphthene-d10          | 14.492 | 164  | 100276   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.245 | 188  | 223499   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.427 | 240  | 221524   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 23.792 | 264  | 268361   | 20.000 | ng/u1  | -0.01    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.187  | 96   | 3125     | 2.630  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.634  | 84   | 18313    | 6.195  | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.010  | 99   | 22888    | 6.307  | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.181  | 67   | 86042    | 33.739 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.357  | 132  | 63493    | 22.673 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.557  | 113  | 43300    | 15.153 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.040  | 128  | 41233    | 30.821 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.745  | 143  | 45106    | 28.821 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.275 | 165  | 76862    | 26.413 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.828 | 131  | 84827    | 22.959 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.916 | 166  | 319590   | 33.675 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.192 | 160  | 326992   | 32.202 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.786 | 143  | 3230m    | 3.012  | ng/u1  | 0.04     |
| 60) Fluorene-d10              | 15.486 | 176  | 265734   | 33.813 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.633 | 200  | 45517    | 26.886 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.345 | 188  | 462579   | 36.931 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.639 | 212  | 619804   | 42.727 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.639 | 264  | 628189   | 39.134 | ng/u1  | -0.01    |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 16) Acetophenone              | 8.693  | 105  | 8607     | 1.791  | ng/u1  | 96       |
| -----                         |        |      |          |        |        |          |

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

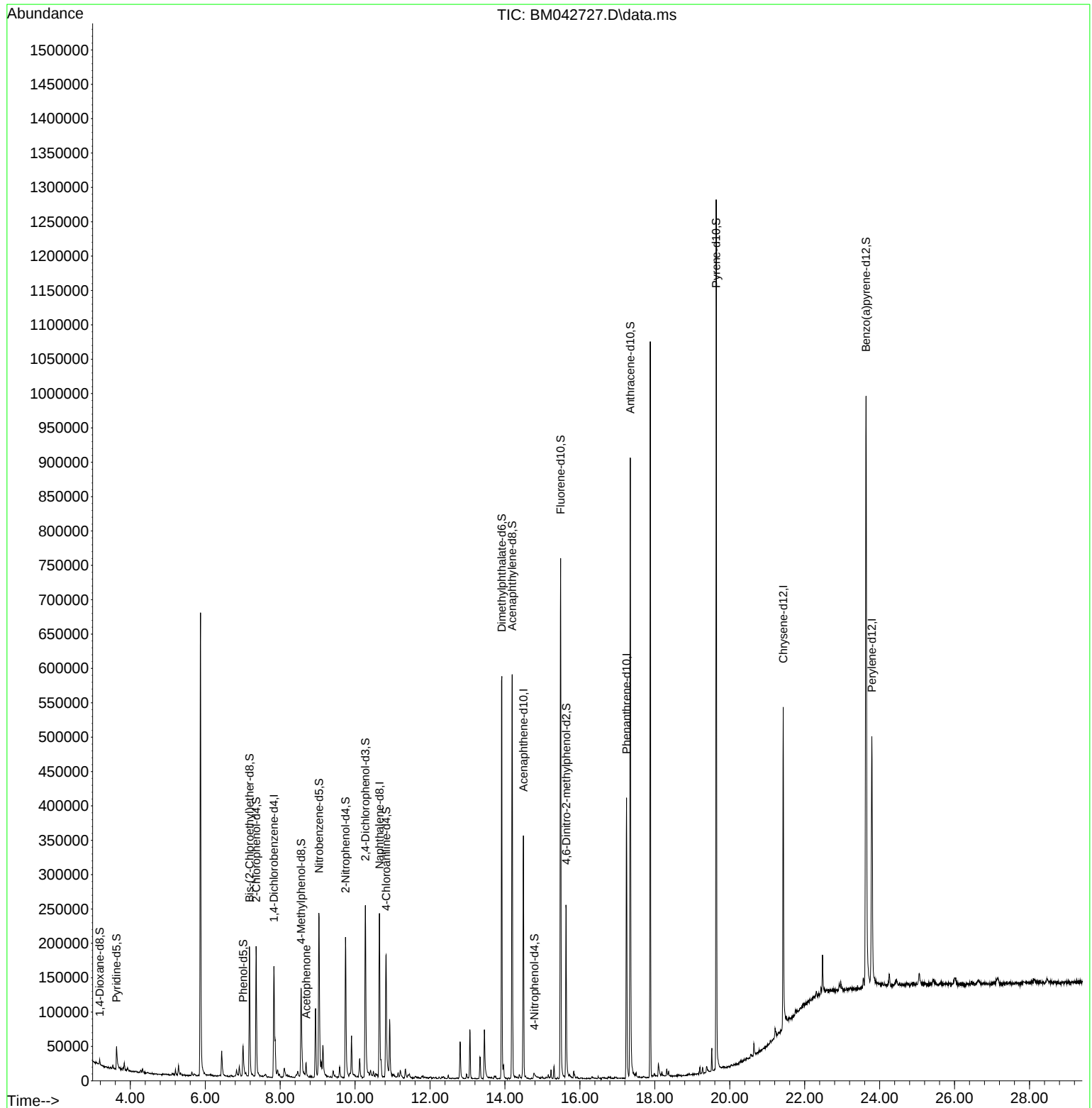
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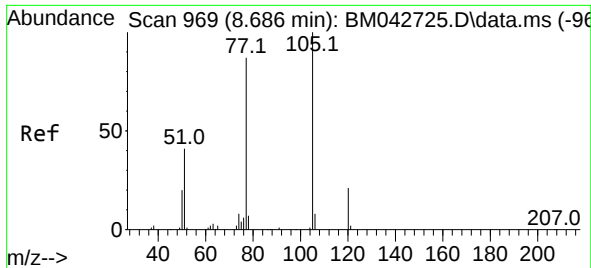
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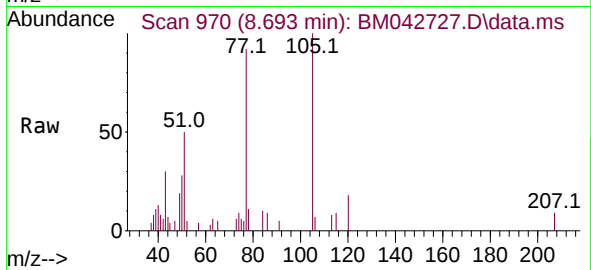
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#16  
Acetophenone  
Concen: 1.791 ng/ul  
RT: 8.693 min Scan# 91  
Delta R.T. 0.000 min  
Lab File: BM042727.D  
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|           |       |       |       |
|-----------|-------|-------|-------|
| Tgt Ion:  | 105   | Resp: | 860   |
| Ion Ratio | Lower | Upper |       |
| 105       | 100   |       |       |
| 77        | 91.5  | 71.6  | 107.4 |
| 51        | 49.6  | 35.4  | 53.2  |

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