

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM081524\  
 Data File : BM047224.D  
 Acq On : 15 Aug 2024 20:53  
 Operator : RC/JU  
 Sample : P3518-13  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 OU4-TS-07-080724

Quant Time: Aug 16 01:20:31 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM081024.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sun Aug 11 23:47:58 2024  
 Response via : Initial Calibration

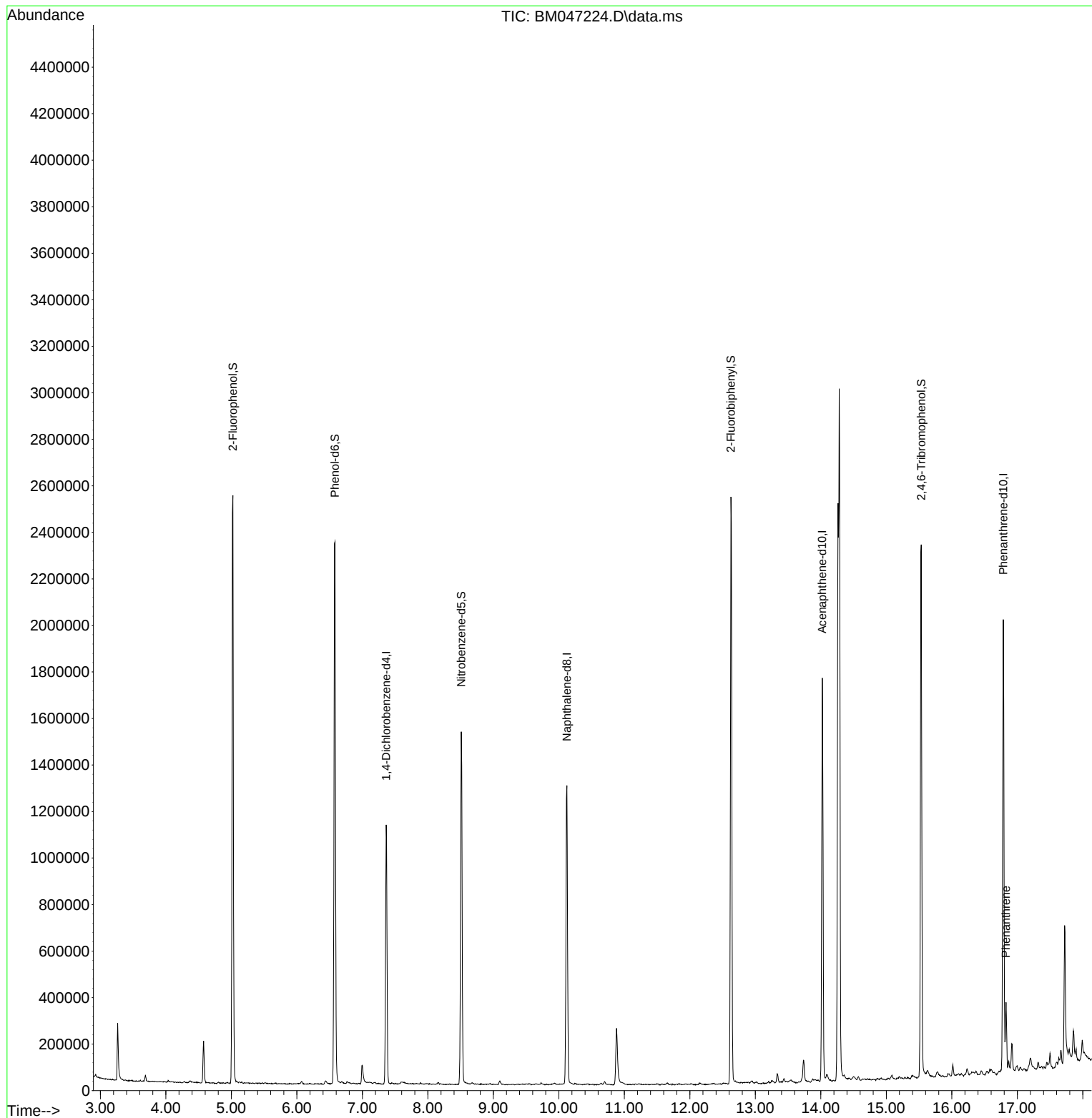
| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.363  | 152  | 301092   | 20.000 | ng    | -0.01    |
| 21) Naphthalene-d8          | 10.122 | 136  | 1058606  | 20.000 | ng    | -0.01    |
| 39) Acenaphthene-d10        | 14.022 | 164  | 606191   | 20.000 | ng    | -0.01    |
| 64) Phenanthrene-d10        | 16.786 | 188  | 1139132  | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 21.021 | 240  | 980146   | 20.000 | ng    | -0.01    |
| 86) Perylene-d12            | 23.715 | 264  | 1128985  | 20.000 | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.022  | 112  | 1062250  | 63.040 | ng    | 0.00     |
| 7) Phenol-d6                | 6.581  | 99   | 1352648  | 58.257 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.510  | 82   | 885015   | 42.096 | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 15.533 | 330  | 418167   | 59.493 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.628 | 172  | 1346893  | 34.273 | ng    | -0.02    |
| 79) Terphenyl-d14           | 19.451 | 244  | 1677203  | 36.666 | ng    | -0.01    |
| Target Compounds            |        |      |          |        |       |          |
|                             |        |      |          |        |       | Qvalue   |
| 71) Phenanthrene            | 16.827 | 178  | 165707   | 2.899  | ng    | 99       |
| 75) Fluoranthene            | 18.863 | 202  | 1091549  | 15.698 | ng    | 99       |
| 78) Pyrene                  | 19.227 | 202  | 835520   | 11.939 | ng    | 99       |
| 81) Benzo(a)anthracene      | 21.004 | 228  | 236001   | 3.627  | ng    | 98       |
| 83) Chrysene                | 21.062 | 228  | 285969   | 4.632  | ng    | 98       |
| 88) Benzo(b)fluoranthene    | 22.874 | 252  | 327999   | 4.892  | ng    | # 97     |
| 90) Benzo(a)pyrene          | 23.592 | 252  | 159252   | 2.742  | ng    | # 91     |
| -----                       |        |      |          |        |       |          |

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

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