

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM050525\  
 Data File : BM050098.D  
 Acq On : 05 May 2025 13:40  
 Operator : RC/JU  
 Sample : Q1907-01DL 25X  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 CO-8R-WCDL

Quant Time: May 05 14:15:28 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM042825.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Apr 28 18:09:16 2025  
 Response via : Initial Calibration

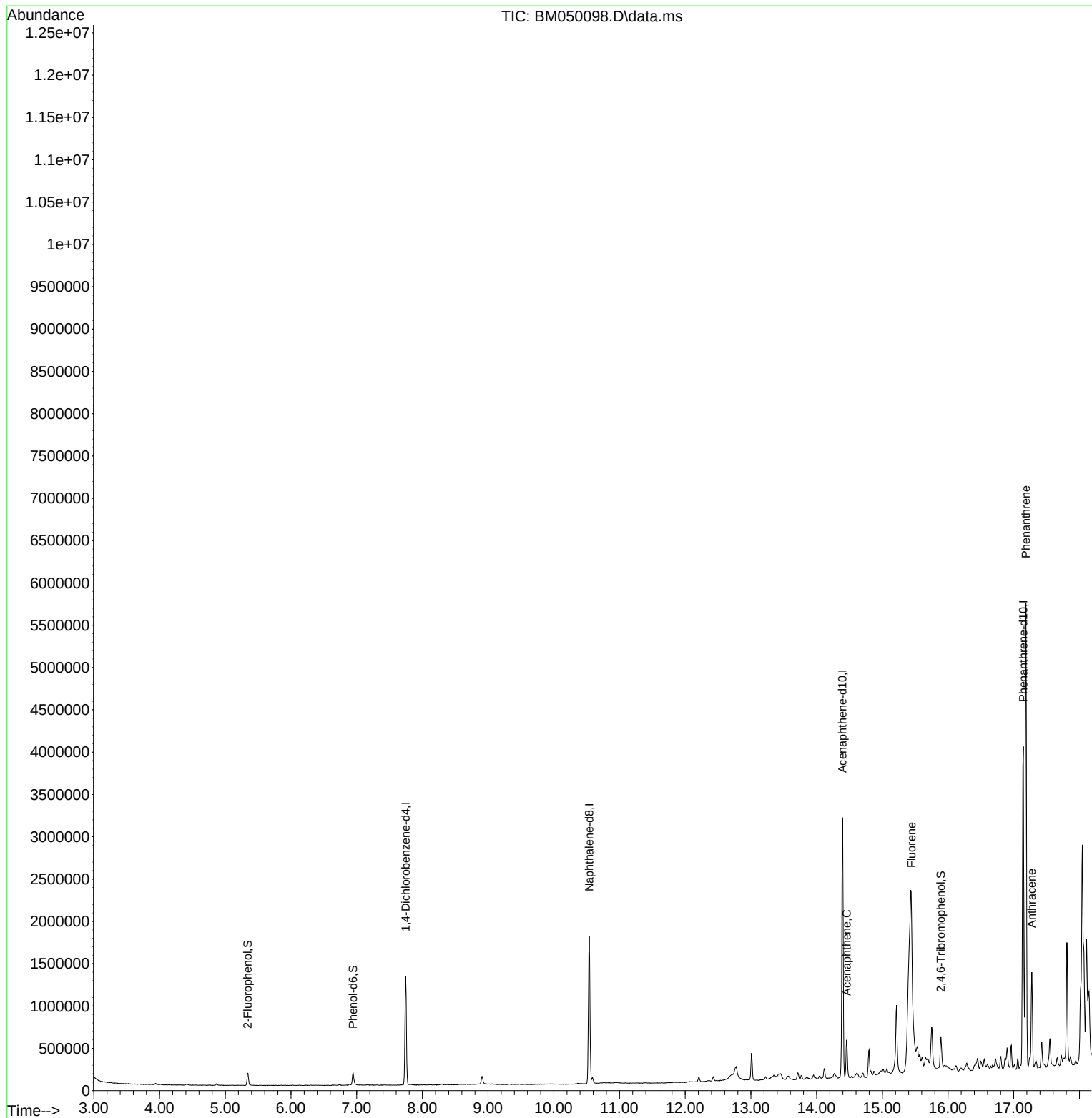
| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.745  | 152  | 373100   | 20.000 | ng    | -0.02    |
| 21) Naphthalene-d8          | 10.539 | 136  | 1493906  | 20.000 | ng    | -0.02    |
| 39) Acenaphthene-d10        | 14.392 | 164  | 1094766  | 20.000 | ng    | -0.02    |
| 64) Phenanthrene-d10        | 17.145 | 188  | 2314689  | 20.000 | ng    | -0.01    |
| 76) Chrysene-d12            | 21.386 | 240  | 2008283  | 20.000 | ng    | -0.01    |
| 86) Perylene-d12            | 24.380 | 264  | 1568873  | 20.000 | ng    | -0.02    |
|                             |        |      |          |        |       |          |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.340  | 112  | 75333    | 3.536  | ng    | -0.02    |
| 7) Phenol-d6                | 6.945  | 99   | 102248   | 3.818  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.904  | 82   | 59788    | 1.972  | ng    | -0.02    |
| 42) 2,4,6-Tribromophenol    | 15.892 | 330  | 54481    | 3.366  | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 13.010 | 172  | 173403   | 1.874  | ng    | -0.02    |
| 79) Terphenyl-d14           | 19.768 | 244  | 257693   | 1.899  | ng    | -0.01    |
|                             |        |      |          |        |       |          |
| Target Compounds            |        |      |          |        |       | Qvalue   |
| 52) Acenaphthene            | 14.457 | 154  | 159950   | 2.719  | ng    | 99       |
| 58) Fluorene                | 15.445 | 166  | 246811   | 3.081  | ng    | 99       |
| 71) Phenanthrene            | 17.186 | 178  | 3350459  | 25.664 | ng    | 100      |
| 72) Anthracene              | 17.274 | 178  | 739955   | 5.601  | ng    | 99       |
| 75) Fluoranthene            | 19.203 | 202  | 3935448  | 25.676 | ng    | 99       |
| 78) Pyrene                  | 19.568 | 202  | 3614643  | 25.631 | ng    | 100      |
| 81) Benzo(a)anthracene      | 21.368 | 228  | 1410304  | 10.198 | ng    | 99       |
| 83) Chrysene                | 21.427 | 228  | 1303463  | 10.184 | ng    | 98       |
| 87) Indeno(1,2,3-cd)pyrene  | 27.756 | 276  | 524922   | 4.490  | ng    | 98       |
| 88) Benzo(b)fluoranthene    | 23.433 | 252  | 1202702  | 11.788 | ng    | 98       |
| 89) Benzo(k)fluoranthene    | 23.491 | 252  | 422321m  | 4.092  | ng    |          |
| 90) Benzo(a)pyrene          | 24.238 | 252  | 902585   | 9.445  | ng    | 97       |
| 92) Benzo(g,h,i)perylene    | 28.809 | 276  | 519885   | 5.699  | ng    | 99       |
| -----                       |        |      |          |        |       |          |

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

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