

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM060925\  
 Data File : BM050253.D  
 Acq On : 09 Jun 2025 22:17  
 Operator : RC/JU  
 Sample : Q2236-03  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 WC-A4-05A-C

Quant Time: Jun 10 01:28:07 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM060525.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 09 11:54:38 2025  
 Response via : Initial Calibration

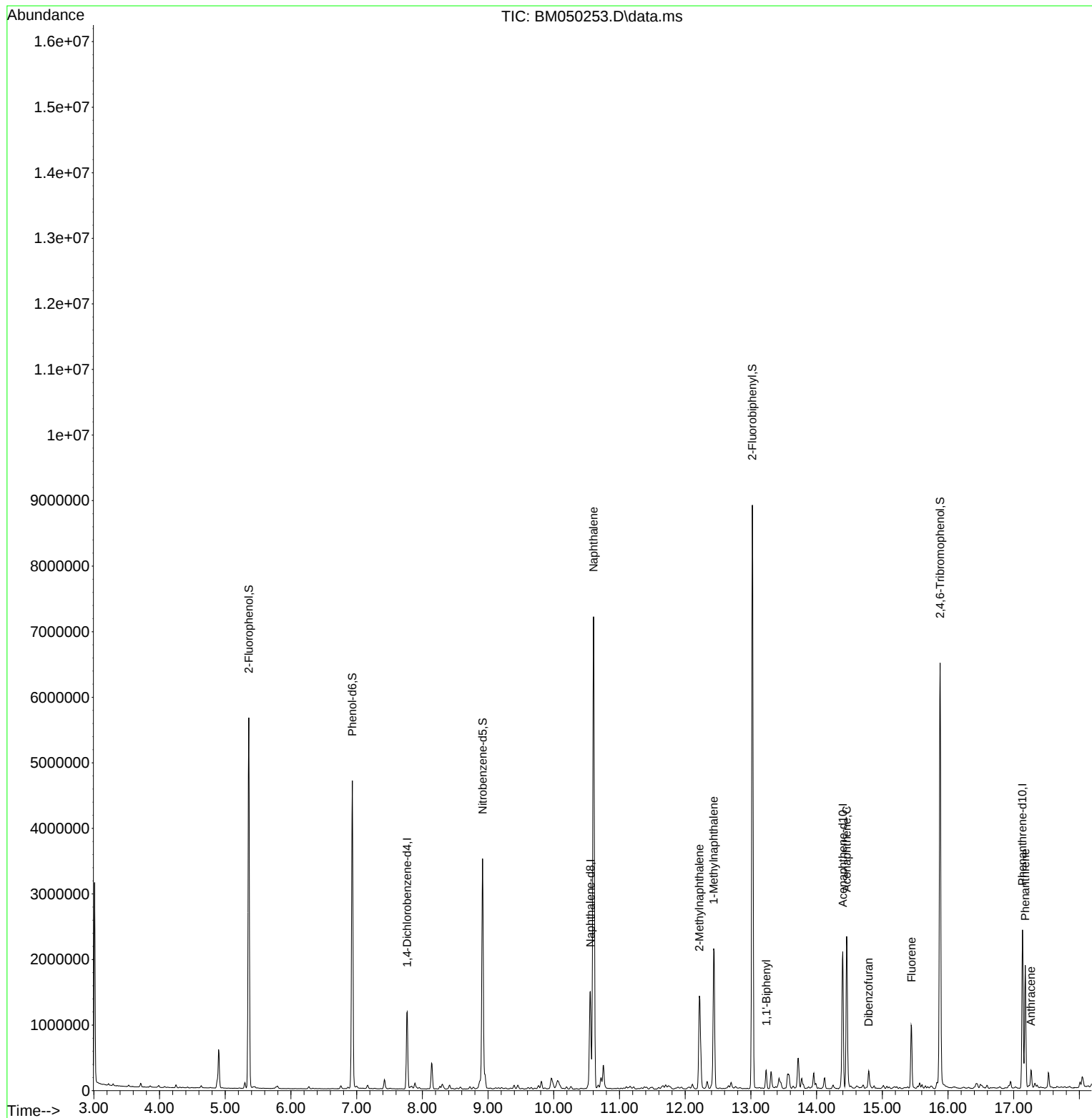
| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.769  | 152  | 339236   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.557 | 136  | 1291741  | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.398 | 164  | 734983   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.133 | 188  | 1365218  | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.362 | 240  | 1409427  | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 24.339 | 264  | 1492062  | 20.000  | ng    | 0.00     |
|                             |        |      |          |         |       |          |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.358  | 112  | 2623964  | 129.029 | ng    | 0.00     |
| 7) Phenol-d6                | 6.934  | 99   | 2974342  | 111.105 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.916  | 82   | 2238948  | 90.145  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.880 | 330  | 1255343  | 150.162 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.022 | 172  | 4616642  | 85.039  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.763 | 244  | 6326070  | 85.056  | ng    | 0.00     |
|                             |        |      |          |         |       |          |
| Target Compounds            |        |      |          |         |       | Qvalue   |
| 31) Naphthalene             | 10.604 | 128  | 6271348  | 93.902  | ng    | 100      |
| 37) 2-Methylnaphthalene     | 12.216 | 142  | 655891   | 16.279  | ng    | 100      |
| 38) 1-Methylnaphthalene     | 12.434 | 142  | 1041234  | 24.349  | ng    | 100      |
| 46) 1,1'-Biphenyl           | 13.234 | 154  | 145035   | 2.635   | ng    | 99       |
| 52) Acenaphthene            | 14.457 | 154  | 812188   | 18.764  | ng    | 98       |
| 55) Dibenzofuran            | 14.792 | 168  | 145848   | 2.303   | ng    | 99       |
| 58) Fluorene                | 15.445 | 166  | 430582   | 8.877   | ng    | 99       |
| 71) Phenanthrene            | 17.174 | 178  | 1073820  | 13.769  | ng    | 99       |
| 72) Anthracene              | 17.263 | 178  | 166472   | 2.134   | ng    | 99       |
| -----                       |        |      |          |         |       |          |

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

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