

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM110323\  
 Data File : BM042583.D  
 Acq On : 03 Nov 2023 22:48  
 Operator : MA/JU  
 Sample : 05220-04DL 5X  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 Z-7(0-5)DL

Quant Time: Nov 03 23:27:11 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM103023.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Nov 03 15:52:20 2023  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 11/04/2023  
 Supervised By :mohammad ahmed 11/04/2023

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.916  | 152  | 55378    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.733 | 136  | 203179   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.569 | 164  | 126211   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.321 | 188  | 277249   | 20.000 | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.503 | 240  | 288797   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12            | 23.921 | 264  | 318973   | 20.000 | ng    | 0.00     |
|                             |        |      |          |        |       |          |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.451  | 112  | 80827    | 23.555 | ng    | 0.01     |
| 7) Phenol-d6                | 7.081  | 99   | 102180   | 21.713 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.116  | 82   | 70823    | 16.249 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.063 | 330  | 41996    | 25.486 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.186 | 172  | 146719   | 15.642 | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.933 | 244  | 244732   | 14.369 | ng    | 0.00     |
|                             |        |      |          |        |       |          |
| Target Compounds            |        |      |          |        |       | Qvalue   |
| 13) 1,4-Dichlorobenzene     | 7.951  | 146  | 139999   | 34.594 | ng    | 98       |
| 49) Acenaphthylene          | 14.298 | 152  | 74364    | 6.344  | ng    | 99       |
| 58) Fluorene                | 15.616 | 166  | 21899    | 2.340  | ng    | # 94     |
| 71) Phenanthrene            | 17.362 | 178  | 203726   | 14.061 | ng    | 99       |
| 72) Anthracene              | 17.457 | 178  | 120335   | 8.385  | ng    | 100      |
| 75) Fluoranthene            | 19.380 | 202  | 548908   | 32.252 | ng    | 99       |
| 78) Pyrene                  | 19.745 | 202  | 547582   | 27.613 | ng    | 99       |
| 81) Benzo(a)anthracene      | 21.486 | 228  | 306593   | 16.686 | ng    | 98       |
| 83) Chrysene                | 21.539 | 228  | 264209   | 14.155 | ng    | 96       |
| 87) Indeno(1,2,3-cd)pyrene  | 26.450 | 276  | 148342   | 7.814  | ng    | # 93     |
| 88) Benzo(b)fluoranthene    | 23.174 | 252  | 318119   | 17.901 | ng    | 97       |
| 89) Benzo(k)fluoranthene    | 23.221 | 252  | 105087m  | 5.307  | ng    |          |
| 90) Benzo(a)pyrene          | 23.815 | 252  | 252211   | 14.416 | ng    | 96       |
| 91) Dibenzo(a,h)anthracene  | 26.462 | 278  | 41393    | 6.296  | ng    | # 86     |
| 92) Benzo(g,h,i)perylene    | 27.244 | 276  | 152681   | 9.125  | ng    | 97       |
| -----                       |        |      |          |        |       |          |

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

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