

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF110724\  
 Data File : BF140284.D  
 Acq On : 07 Nov 2024 19:18  
 Operator : RC/JU  
 Sample : P4700-01 5X  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 MH-8

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 11/08/2024  
 Supervised By :mohammad ahmed 11/11/2024

Quant Time: Nov 08 01:30:56 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF110524.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Nov 05 16:47:56 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.875  | 152  | 137866   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.157  | 136  | 483639   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.916  | 164  | 221709   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.404 | 188  | 369560   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 14.057 | 240  | 272097   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12            | 15.557 | 264  | 205210   | 20.000 | ng    | 0.00     |
|                             |        |      |          |        |       |          |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.487  | 112  | 109364   | 13.146 | ng    | 0.00     |
| 7) Phenol-d6                | 6.493  | 99   | 138699   | 12.962 | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 7.434  | 82   | 82502    | 8.481  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.704 | 330  | 25880    | 11.809 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.234  | 172  | 152726   | 11.000 | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.992 | 244  | 151212   | 9.104  | ng    | 0.00     |
|                             |        |      |          |        |       |          |
| Target Compounds            |        |      |          |        |       | Qvalue   |
| 71) Phenanthrene            | 11.428 | 178  | 53705    | 3.045  | ng    | 98       |
| 75) Fluoranthene            | 12.622 | 202  | 136964   | 7.427  | ng    | 98       |
| 78) Pyrene                  | 12.851 | 202  | 133585   | 5.946  | ng    | 99       |
| 81) Benzo(a)anthracene      | 14.045 | 228  | 64442    | 3.690  | ng    | 94       |
| 83) Chrysene                | 14.080 | 228  | 47937    | 2.932  | ng    | 96       |
| 88) Benzo(b)fluoranthene    | 15.121 | 252  | 53104m   | 4.278  | ng    |          |
| 90) Benzo(a)pyrene          | 15.492 | 252  | 34362    | 3.358  | ng    | # 85     |
| 92) Benzo(g,h,i)perylene    | 17.515 | 276  | 22485    | 2.224  | ng    | # 91     |
| -----                       |        |      |          |        |       |          |

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

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