

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120121\  
 Data File : BP008182.D  
 Acq On : 01 Dec 2021 20:26  
 Operator : CG/JU  
 Sample : M4721-03 2X  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SAMPLE-3

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 12/03/2021  
 Supervised By :mohammad ahmed 12/05/2021

Quant Time: Dec 01 21:16:52 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP112521.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Dec 01 12:06:05 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min)  |
|-----------------------------|--------|------|----------|--------|-------|-----------|
| -----                       |        |      |          |        |       |           |
| Internal Standards          |        |      |          |        |       |           |
| 1) 1,4-Dichlorobenzene-d4   | 7.799  | 152  | 116024   | 20.000 | ng    | 0.00      |
| 21) Naphthalene-d8          | 10.593 | 136  | 433622   | 20.000 | ng    | 0.00      |
| 39) Acenaphthene-d10        | 14.446 | 164  | 270515   | 20.000 | ng    | 0.00      |
| 64) Phenanthrene-d10        | 17.204 | 188  | 578847   | 20.000 | ng    | 0.00      |
| 76) Chrysene-d12            | 21.310 | 240  | 606883   | 20.000 | ng    | # 0.00    |
| 86) Perylene-d12            | 23.704 | 264  | 661967   | 20.000 | ng    | 0.00      |
| System Monitoring Compounds |        |      |          |        |       |           |
| 5) 2-Fluorophenol           | 5.393  | 112  | 307707   | 42.701 | ng    | 0.00      |
| 7) Phenol-d6                | 6.981  | 99   | 398891   | 41.006 | ng    | 0.00      |
| 23) Nitrobenzene-d5         | 8.958  | 82   | 269332   | 28.244 | ng    | 0.00      |
| 42) 2,4,6-Tribromophenol    | 15.945 | 330  | 155989   | 45.114 | ng    | 0.00      |
| 45) 2-Fluorobiphenyl        | 13.063 | 172  | 516127   | 27.714 | ng    | 0.00      |
| 79) Terphenyl-d14           | 19.775 | 244  | 801501   | 27.601 | ng    | 0.00      |
| Target Compounds            |        |      |          |        |       |           |
| 75) Fluoranthene            | 19.222 | 202  | 135747   | 3.297  | ng    | Qvalue 99 |
| 81) Benzo(a)anthracene      | 21.292 | 228  | 86993    | 2.163  | ng    | # 95      |
| 83) Chrysene                | 21.345 | 228  | 116558   | 3.045  | ng    | # 92      |
| 88) Benzo(b)fluoranthene    | 22.980 | 252  | 106821m  | 2.677  | ng    |           |
| 90) Benzo(a)pyrene          | 23.598 | 252  | 73498    | 2.155  | ng    | # 81      |
| -----                       |        |      |          |        |       |           |

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

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