

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM072321\  
 Data File : BM031109.D  
 Acq On : 24 Jul 2021 03:57  
 Operator : CG/JU  
 Sample : M2970-17DL 5X  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BGED0DL

Manual Integrations  
 APPROVED

mohammad  
 7/26/2021 1:26:31 PM

Quant Time: Jul 24 04:33:28 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM072221.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Jul 24 04:20:53 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|------|--------|----------|
| -----                       |        |      |          |      |        |          |
| Internal Standards          |        |      |          |      |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.639  | 152  | 1897     | 0.40 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.401 | 136  | 7332     | 0.40 | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.265 | 164  | 4861     | 0.40 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.009 | 188  | 10291    | 0.40 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.193 | 240  | 9008     | 0.40 | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.365 | 264  | 7804     | 0.40 | ng/ul  | # 0.00   |
| System Monitoring Compounds |        |      |          |      |        |          |
| 3) 1,4-Dioxane-d8           | 3.217  | 96   | 1236     | 0.62 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 11.993 | 152  | 606      | 0.05 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 19.035 | 212  | 1448     | 0.05 | ng/ul  | 0.00     |
| Target Compounds            |        |      |          |      |        |          |
|                             |        |      |          |      | Qvalue |          |
| 15) Phenanthrene            | 17.051 | 178  | 3106     | 0.10 | ng/ul  | 96       |
| 16) Anthracene              | 17.141 | 178  | 1421     | 0.05 | ng/ul  | 95       |
| 19) Fluoranthene            | 19.069 | 202  | 12647    | 0.35 | ng/ul  | 96       |
| 20) Pyrene                  | 19.432 | 202  | 9466     | 0.26 | ng/ul  | 96       |
| 21) Benzo(a)anthracene      | 21.178 | 228  | 7559     | 0.21 | ng/ul  | 94       |
| 22) Chrysene                | 21.227 | 228  | 6456     | 0.17 | ng/ul  | 98       |
| 24) Benzo(b)fluoranthene    | 22.720 | 252  | 7750m    | 0.21 | ng/ul  |          |
| 25) Benzo(k)fluoranthene    | 22.759 | 252  | 3513m    | 0.09 | ng/ul  |          |
| 26) Benzo(a)pyrene          | 23.270 | 252  | 3909     | 0.12 | ng/ul# | 71       |
| 27) Indeno(1,2,3-cd)pyrene  | 25.516 | 276  | 2376     | 0.06 | ng/ul# | 99       |
| 28) Dibenzo(a,h)anthracene  | 25.526 | 278  | 816      | 0.02 | ng/ul# | 42       |
| -----                       |        |      |          |      |        |          |

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

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