

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM061021\  
 Data File : BM030396.D  
 Acq On : 11 Jun 2021 08:00  
 Operator : CG/JU  
 Sample : M2612-37  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 18-B12(70-74)

Quant Time: Jun 11 09:38:07 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM060921.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jun 08 20:51:08 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units | Dev(Min)  |
|-------------------------------|--------|------|----------|---------|-------|-----------|
| -----                         |        |      |          |         |       |           |
| Internal Standards            |        |      |          |         |       |           |
| 1) 1,4-Dichlorobenzene-d4     | 7.839  | 152  | 160552   | 20.000  | ng    | -0.01     |
| 21) Naphthalene-d8            | 10.628 | 136  | 678471   | 20.000  | ng    | #-0.01    |
| 39) Acenaphthene-d10          | 14.463 | 164  | 467458   | 20.000  | ng    | -0.01     |
| 64) Phenanthrene-d10          | 17.198 | 188  | 1080894  | 20.000  | ng    | -0.01     |
| 76) Chrysene-d12              | 21.356 | 240  | 1103749  | 20.000  | ng    | -0.01     |
| 86) Perylene-d12              | 23.633 | 264  | 974968   | 20.000  | ng    | -0.02     |
| System Monitoring Compounds   |        |      |          |         |       |           |
| 5) 2-Fluorophenol             | 5.428  | 112  | 1703644  | 172.501 | ng    | 0.00      |
| 7) Phenol-d6                  | 7.004  | 99   | 2089284  | 146.134 | ng    | 0.00      |
| 23) Nitrobenzene-d5           | 8.998  | 82   | 1336944  | 95.999  | ng    | -0.01     |
| 42) 2,4,6-Tribromophenol      | 15.951 | 330  | 804644   | 136.908 | ng    | 0.00      |
| 45) 2-Fluorobiphenyl          | 13.086 | 172  | 3347728  | 94.179  | ng    | -0.01     |
| 79) Terphenyl-d14             | 19.815 | 244  | 6338740  | 100.871 | ng    | 0.00      |
| Target Compounds              |        |      |          |         |       |           |
| 32) Benzoic acid              | 9.881  | 122  | 50789    | 10.662  | ng    | Qvalue 92 |
| 84) Bis(2-ethylhexyl)phtha... | 21.268 | 149  | 2112     | 4.311   | ng    | # 72      |
| -----                         |        |      |          |         |       |           |

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

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