

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM030921\  
 Data File : BM029057.D  
 Acq On : 09 Mar 2021 21:56  
 Operator : CG/JU  
 Sample : M1603-01  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 31-NORTH

Quant Time: Mar 09 23:40:30 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM022321.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Mar 09 10:08:02 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min)     |
|-----------------------------|--------|------|----------|--------|-------|--------------|
| -----                       |        |      |          |        |       |              |
| Internal Standards          |        |      |          |        |       |              |
| 1) 1,4-Dichlorobenzene-d4   | 7.669  | 152  | 8981     | 20.00  | ng    | 0.00         |
| 21) Naphthalene-d8          | 10.457 | 136  | 33663    | 20.00  | ng    | # 0.00       |
| 39) Acenaphthene-d10        | 14.333 | 164  | 18609    | 20.00  | ng    | 0.00         |
| 64) Phenanthrene-d10        | 17.080 | 188  | 38937    | 20.00  | ng    | 0.00         |
| 76) Chrysene-d12            | 21.256 | 240  | 40347    | 20.00  | ng    | 0.00         |
| 86) Perylene-d12            | 23.409 | 264  | 43092    | 20.00  | ng    | -0.01        |
| System Monitoring Compounds |        |      |          |        |       |              |
| 5) 2-Fluorophenol           | 5.240  | 112  | 60651    | 111.94 | ng    | 0.00         |
| 7) Phenol-d6                | 6.863  | 99   | 70009    | 97.82  | ng    | 0.00         |
| 23) Nitrobenzene-d5         | 8.839  | 82   | 45501    | 72.92  | ng    | 0.00         |
| 42) 2,4,6-Tribromophenol    | 15.833 | 330  | 23157    | 105.00 | ng    | 0.00         |
| 45) 2-Fluorobiphenyl        | 12.945 | 172  | 97778    | 69.95  | ng    | 0.00         |
| 79) Terphenyl-d14           | 19.703 | 244  | 141439   | 64.73  | ng    | 0.00         |
| Target Compounds            |        |      |          |        |       |              |
| 10) Phenol                  | 6.887  | 94   | 1543     | 2.17   | ng    | Qvalue<br>90 |
| -----                       |        |      |          |        |       |              |

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

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