

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG102921\  
 Data File : BG050784.D  
 Acq On : 29 Oct 2021 19:22  
 Operator : CG/JU  
 Sample : M4406-02  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 CHRT-20310

Quant Time: Oct 30 01:38:24 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG100621.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Oct 14 10:56:22 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.247  | 152  | 31482    | 20.000  | ng    | -0.01    |
| 21) Naphthalene-d8          | 11.073 | 136  | 130567   | 20.000  | ng    | #-0.01   |
| 39) Acenaphthene-d10        | 14.869 | 164  | 83859    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.613 | 188  | 192997   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.919 | 240  | 185918   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 25.333 | 264  | 184729   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.779  | 112  | 278222   | 132.168 | ng    | -0.01    |
| 7) Phenol-d6                | 7.389  | 99   | 335766   | 110.168 | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 9.416  | 82   | 297401   | 94.290  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 16.355 | 330  | 120199   | 150.274 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.494 | 172  | 544333   | 97.692  | ng    | -0.01    |
| 79) Terphenyl-d14           | 20.204 | 244  | 910817   | 95.473  | ng    | 0.00     |

|                  |        |
|------------------|--------|
| Target Compounds | Qvalue |
| -----            |        |

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

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