

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM060222\  
 Data File : BM035377.D  
 Acq On : 03 Jun 2022 00:09  
 Operator : CG/JU  
 Sample : N3099-06  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 20220429-SPLP-AMENDED-COMPOSITE-I

Quant Time: Jun 03 02:16:00 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM060222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jun 02 16:16:23 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min)     |
|-----------------------------|--------|------|----------|---------|-------|--------------|
| -----                       |        |      |          |         |       |              |
| Internal Standards          |        |      |          |         |       |              |
| 1) 1,4-Dichlorobenzene-d4   | 7.851  | 152  | 170626   | 20.000  | ng    | 0.00         |
| 21) Naphthalene-d8          | 10.645 | 136  | 633813   | 20.000  | ng    | #-0.01       |
| 39) Acenaphthene-d10        | 14.486 | 164  | 423254   | 20.000  | ng    | -0.01        |
| 64) Phenanthrene-d10        | 17.233 | 188  | 943467   | 20.000  | ng    | #-0.01       |
| 76) Chrysene-d12            | 21.415 | 240  | 1066140  | 20.000  | ng    | #-0.01       |
| 86) Perylene-d12            | 23.774 | 264  | 1182643  | 20.000  | ng    | -0.01        |
| System Monitoring Compounds |        |      |          |         |       |              |
| 5) 2-Fluorophenol           | 5.434  | 112  | 510524   | 62.873  | ng    | 0.00         |
| 7) Phenol-d6                | 7.034  | 99   | 368444   | 33.218  | ng    | 0.00         |
| 23) Nitrobenzene-d5         | 9.010  | 82   | 928414   | 93.010  | ng    | 0.00         |
| 42) 2,4,6-Tribromophenol    | 15.980 | 330  | 853328   | 125.656 | ng    | -0.01        |
| 45) 2-Fluorobiphenyl        | 13.116 | 172  | 2590632  | 87.911  | ng    | 0.00         |
| 79) Terphenyl-d14           | 19.862 | 244  | 4887531  | 91.332  | ng    | 0.00         |
| Target Compounds            |        |      |          |         |       |              |
| 32) Benzoic acid            | 9.963  | 122  | 14776    | 2.105   | ng    | Qvalue<br>84 |
| -----                       |        |      |          |         |       |              |

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

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