

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF032922\  
 Data File : BF127576.D  
 Acq On : 29 Mar 2022 20:17  
 Operator : CG\JU  
 Sample : N2123-03DL 5X  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SPLP-LEACHATE-AMENDED-COMP-CS-1DL

Quant Time: Mar 30 08:32:53 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF032122.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Mar 29 07:13:31 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.834  | 152  | 87006    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.116  | 136  | 305932   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.869  | 164  | 160806   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.363 | 188  | 231649   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 14.010 | 240  | 213021   | 20.000 | ng    | # 0.00   |
| 86) Perylene-d12            | 15.492 | 264  | 219262   | 20.000 | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.469  | 112  | 61858    | 11.349 | ng    | 0.01     |
| 7) Phenol-d6                | 6.487  | 99   | 51353    | 6.892  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.398  | 82   | 130547   | 18.240 | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.663 | 330  | 33005    | 19.267 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.186  | 172  | 227262   | 20.637 | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.945 | 244  | 192472   | 14.791 | ng    | 0.00     |
| Target Compounds            |        |      |          |        |       |          |
|                             |        |      |          |        |       | Qvalue   |
| 31) Naphthalene             | 8.139  | 128  | 834729   | 51.392 | ng    | 99       |
| 37) 2-Methylnaphthalene     | 8.828  | 142  | 284290   | 27.220 | ng    | 100      |
| 38) 1-Methylnaphthalene     | 8.928  | 142  | 403750   | 39.096 | ng    | 97       |
| 46) 1,1'-Biphenyl           | 9.292  | 154  | 27065    | 2.228  | ng    | 99       |
| 52) Acenaphthene            | 9.904  | 154  | 127003   | 15.055 | ng    | 99       |
| 58) Fluorene                | 10.422 | 166  | 49081    | 2.740  | ng    | 98       |
| 71) Phenanthrene            | 11.386 | 178  | 109059   | 9.030  | ng    | 98       |
| 73) Carbazole               | 11.604 | 167  | 58518    | 5.106  | ng    | 100      |
| -----                       |        |      |          |        |       |          |

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

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