

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071522\  
 Data File : BF129247.D  
 Acq On : 15 Jul 2022 19:53  
 Operator : CG\JU  
 Sample : N3697-01  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PIPELINE-PIGGING-TK2045L

Quant Time: Jul 16 01:42:51 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF071522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jul 15 12:55:49 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-------------------------------|--------|------|----------|---------|-------|----------|
| -----                         |        |      |          |         |       |          |
| Internal Standards            |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 6.898  | 152  | 232505   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8            | 8.181  | 136  | 929447   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10          | 9.939  | 164  | 472753   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10          | 11.427 | 188  | 766744   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12              | 14.068 | 240  | 456583   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12              | 15.562 | 264  | 430281   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |         |       |          |
| 5) 2-Fluorophenol             | 5.522  | 112  | 1619814  | 111.264 | ng    | 0.00     |
| 7) Phenol-d6                  | 6.522  | 99   | 1902463  | 101.038 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 7.457  | 82   | 1485760  | 85.449  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol      | 10.733 | 330  | 546626   | 124.574 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 9.257  | 172  | 2202377  | 78.926  | ng    | 0.00     |
| 79) Terphenyl-d14             | 13.016 | 244  | 1635615  | 67.014  | ng    | 0.00     |
| Target Compounds              |        |      |          |         |       |          |
|                               |        |      |          |         |       | Qvalue   |
| 20) 3+4-Methylphenols         | 7.292  | 107  | 85578    | 5.404   | ng    | # 74     |
| 27) 2,4-Dimethylphenol        | 7.839  | 122  | 50407    | 3.902   | ng    | # 92     |
| 37) 2-Methylnaphthalene       | 8.892  | 142  | 88960    | 3.099   | ng    | 99       |
| 38) 1-Methylnaphthalene       | 8.992  | 142  | 62747    | 2.261   | ng    | 99       |
| 71) Phenanthrene              | 11.451 | 178  | 81412    | 2.091   | ng    | # 92     |
| 84) Bis(2-ethylhexyl)phtha... | 14.039 | 149  | 62640    | 2.858   | ng    | 94       |
| -----                         |        |      |          |         |       |          |

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

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