

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF120920\  
 Data File : BF122626.D  
 Acq On : 9 Dec 2020 23:29  
 Operator : JU/CG  
 Sample : L4970-13  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 4S1-A-D-COMP

Manual Integrations  
 APPROVED

mohammad  
 12/10/2020 2:43:27 PM

Quant Time: Dec 10 02:52:51 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF120520.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Dec 07 12:08:49 2020  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min)  |
|-----------------------------|--------|------|----------|--------|-------|-----------|
| -----                       |        |      |          |        |       |           |
| Internal Standards          |        |      |          |        |       |           |
| 1) 1,4-Dichlorobenzene-d4   | 6.839  | 152  | 184206   | 20.00  | ng    | 0.00      |
| 21) Naphthalene-d8          | 8.122  | 136  | 735147   | 20.00  | ng    | 0.00      |
| 39) Acenaphthene-d10        | 9.874  | 164  | 394213   | 20.00  | ng    | 0.00      |
| 64) Phenanthrene-d10        | 11.363 | 188  | 740089   | 20.00  | ng    | 0.00      |
| 76) Chrysene-d12            | 13.998 | 240  | 664557   | 20.00  | ng    | -0.01     |
| 86) Perylene-d12            | 15.462 | 264  | 550110   | 20.00  | ng    | 0.00      |
| System Monitoring Compounds |        |      |          |        |       |           |
| 5) 2-Fluorophenol           | 5.475  | 112  | 1143208  | 98.25  | ng    | 0.02      |
| 7) Phenol-d6                | 6.469  | 99   | 1263664  | 81.89  | ng    | -0.01     |
| 23) Nitrobenzene-d5         | 7.404  | 82   | 1035820  | 70.59  | ng    | 0.00      |
| 42) 2,4,6-Tribromophenol    | 10.669 | 330  | 565255   | 109.54 | ng    | 0.00      |
| 45) 2-Fluorobiphenyl        | 9.198  | 172  | 1819760  | 62.44  | ng    | 0.00      |
| 79) Terphenyl-d14           | 12.951 | 244  | 2221318  | 58.51  | ng    | 0.00      |
| Target Compounds            |        |      |          |        |       |           |
| 10) Phenol                  | 6.481  | 94   | 36718    | 2.30   | ng    | Qvalue 95 |
| 32) Benzoic acid            | 7.892  | 122  | 17184m   | 2.07   | ng    |           |
| 50) Dimethylphthalate       | 9.592  | 163  | 119371   | 3.79   | ng    | 100       |
| -----                       |        |      |          |        |       |           |

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

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