

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092520\  
 Data File : BF121696.D  
 Acq On : 25 Sep 2020 15:22  
 Operator : JU/CG  
 Sample : L4102-06  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PE-1B-(3.0-3.5)

Manual Integrations  
 APPROVED

mohammad  
 9/28/2020 11:10:31 AM

Quant Time: Sep 25 15:56:23 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Sep 25 14:15:19 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 6.80  | 152  | 140184   | 20.00 | ng    | 0.00     |
| 21) Naphthalene-d8        | 8.08  | 136  | 546268   | 20.00 | ng    | 0.00     |
| 39) Acenaphthene-d10      | 9.83  | 164  | 286443   | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10      | 11.32 | 188  | 508446   | 20.00 | ng    | 0.00     |
| 76) Chrysene-d12          | 13.96 | 240  | 426832   | 20.00 | ng    | 0.00     |
| 86) Perylene-d12          | 15.41 | 264  | 344163   | 20.00 | ng    | 0.00     |

| System Monitoring Compounds | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 5) 2-Fluorophenol           | 5.42  | 112  | 847900   | 89.89  | ng    | 0.00     |
| 7) Phenol-d6                | 6.43  | 99   | 1108573  | 89.58  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.36  | 82   | 760235   | 74.72  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.63 | 330  | 359119   | 100.87 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.16  | 172  | 1348351  | 71.29  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.92 | 244  | 1434877  | 68.10  | ng    | 0.00     |

| Target Compounds               | R.T.  | QIon | Response | Conc   | Units | Qvalue |
|--------------------------------|-------|------|----------|--------|-------|--------|
| 50) Dimethylphthalate          | 9.55  | 163  | 269767   | 12.994 | ng    | 99     |
| 71) Phenanthrene               | 11.35 | 178  | 88819    | 3.206  | ng    | 96     |
| 75) Fluoranthene               | 12.54 | 202  | 200995   | 6.797  | ng    | 98     |
| 78) Pyrene                     | 12.77 | 202  | 188925   | 6.058  | ng    | 98     |
| 81) Benzo(a)anthracene         | 13.95 | 228  | 95455    | 3.535  | ng    | 99     |
| 83) Chrysene                   | 13.99 | 228  | 101345   | 3.706  | ng    | 97     |
| 84) Bis(2-ethylhexyl)phthalate | 13.94 | 149  | 54800    | 3.252  | ng    | 99     |
| 87) Indeno(1,2,3-cd)pyrene     | 16.83 | 276  | 47827    | 2.134  | ng    | 97     |
| 88) Benzo(b)fluoranthene       | 14.99 | 252  | 120443m  | 5.235  | ng    |        |
| 90) Benzo(a)pyrene             | 15.35 | 252  | 71897    | 3.676  | ng    | 98     |
| 92) Benzo(g,h,i)perylene       | 17.26 | 276  | 50340    | 2.745  | ng    | 98     |

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

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