

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030920\  
 Data File : BF119306.D  
 Acq On : 10 Mar 2020 2:22  
 Operator : CG/JU  
 Sample : L1778-05  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 DN1

Manual Integrations  
 APPROVED

mohammad  
 3/10/2020 2:44:58 PM

Quant Time: Mar 10 10:39:33 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Mar 06 23:38:46 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 6.72  | 152  | 106605   | 20.00 | ng    | 0.00     |
| 21) Naphthalene-d8        | 8.00  | 136  | 399634   | 20.00 | ng    | 0.00     |
| 39) Acenaphthene-d10      | 9.76  | 164  | 217649   | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10      | 11.24 | 188  | 369764   | 20.00 | ng    | 0.00     |
| 76) Chrysene-d12          | 13.89 | 240  | 260361   | 20.00 | ng    | 0.00     |
| 87) Perylene-d12          | 15.32 | 264  | 271217   | 20.00 | ng    | 0.00     |

|                             |       |     |        |       |    |      |
|-----------------------------|-------|-----|--------|-------|----|------|
| System Monitoring Compounds |       |     |        |       |    |      |
| 5) 2-Fluorophenol           | 5.35  | 112 | 614903 | 96.39 | ng | 0.00 |
| 7) Phenol-d6                | 6.38  | 99  | 747077 | 96.30 | ng | 0.00 |
| 23) Nitrobenzene-d5         | 7.29  | 82  | 478306 | 63.19 | ng | 0.00 |
| 42) 2,4,6-Tribromophenol    | 10.55 | 330 | 237584 | 83.44 | ng | 0.00 |
| 45) 2-Fluorobiphenyl        | 9.08  | 172 | 949159 | 64.24 | ng | 0.00 |
| 79) Terphenyl-d14           | 12.82 | 244 | 863206 | 57.08 | ng | 0.00 |

|                            |       |     |         |        |    |      |
|----------------------------|-------|-----|---------|--------|----|------|
| Target Compounds           |       |     |         | Qvalue |    |      |
| 50) Dimethylphthalate      | 9.47  | 163 | 87074   | 5.232  | ng | 98   |
| 52) Acenaphthene           | 9.79  | 154 | 53611   | 4.343  | ng | 99   |
| 55) Dibenzofuran           | 9.96  | 168 | 61707   | 3.349  | ng | 94   |
| 58) Fluorene               | 10.30 | 166 | 63050   | 4.403  | ng | 99   |
| 71) Phenanthrene           | 11.27 | 178 | 1118999 | 55.492 | ng | 99   |
| 72) Anthracene             | 11.32 | 178 | 247433  | 12.281 | ng | 97   |
| 73) Carbazole              | 11.48 | 167 | 128888  | 7.181  | ng | 97   |
| 75) Fluoranthene           | 12.46 | 202 | 1329229 | 62.100 | ng | 98   |
| 78) Pyrene                 | 12.69 | 202 | 1083553 | 51.828 | ng | 99   |
| 80) Butylbenzylphthalate   | 13.29 | 149 | 45766   | 5.201  | ng | 97   |
| 81) Benzo(a)anthracene     | 13.87 | 228 | 591808  | 33.360 | ng | 99   |
| 83) Chrysene               | 13.91 | 228 | 494570  | 27.979 | ng | 97   |
| 86) Indeno(1,2,3-cd)pyrene | 16.73 | 276 | 221140  | 12.920 | ng | 95   |
| 88) Benzo(b)fluoranthene   | 14.91 | 252 | 691971m | 38.707 | ng |      |
| 89) Benzo(k)fluoranthene   | 14.93 | 252 | 191797m | 11.577 | ng |      |
| 90) Benzo(a)pyrene         | 15.26 | 252 | 415564  | 25.756 | ng | 99   |
| 91) Dibenzo(a,h)anthracene | 16.72 | 278 | 60025   | 4.228  | ng | # 88 |
| 92) Benzo(g,h,i)perylene   | 17.15 | 276 | 200406  | 14.287 | ng | 98   |

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

