

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\  
 Data File : BF117875.D  
 Acq On : 19 Nov 2019 22:25  
 Operator : JU  
 Sample : K5665-01 2X  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 HR-05-111419

Manual Integrations  
 APPROVED

mohammad  
 11/20/2019 1:49:45 PM

Quant Time: Nov 20 01:45:21 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Nov 13 18:36:07 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 6.92  | 152  | 176873   | 20.00 | ng    | -0.01    |
| 21) Naphthalene-d8        | 8.21  | 136  | 614719   | 20.00 | ng    | 0.00     |
| 39) Acenaphthene-d10      | 9.97  | 164  | 321495   | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10      | 11.46 | 188  | 523370   | 20.00 | ng    | -0.01    |
| 76) Chrysene-d12          | 14.11 | 240  | 391370   | 20.00 | ng    | 0.00     |
| 87) Perylene-d12          | 15.62 | 264  | 299342   | 20.00 | ng    | 0.00     |

|                             |       |     |        |       |    |       |
|-----------------------------|-------|-----|--------|-------|----|-------|
| System Monitoring Compounds |       |     |        |       |    |       |
| 5) 2-Fluorophenol           | 5.52  | 112 | 451741 | 45.21 | ng | -0.01 |
| 7) Phenol-d6                | 6.53  | 99  | 569131 | 47.22 | ng | -0.02 |
| 23) Nitrobenzene-d5         | 7.47  | 82  | 392439 | 37.95 | ng | -0.02 |
| 42) 2,4,6-Tribromophenol    | 10.76 | 330 | 116347 | 34.18 | ng | -0.01 |
| 45) 2-Fluorobiphenyl        | 9.29  | 172 | 631233 | 35.22 | ng | 0.00  |
| 79) Terphenyl-d14           | 13.04 | 244 | 619614 | 28.93 | ng | -0.01 |

|                            |       |     |         |        |    |      |
|----------------------------|-------|-----|---------|--------|----|------|
| Target Compounds           |       |     |         | Qvalue |    |      |
| 10) Phenol                 | 6.54  | 94  | 29605   | 2.364  | ng | 84   |
| 31) Naphthalene            | 8.23  | 128 | 88746   | 3.097  | ng | # 72 |
| 37) 2-Methylnaphthalene    | 8.92  | 142 | 123396  | 6.373  | ng | 96   |
| 38) 1-Methylnaphthalene    | 9.02  | 142 | 78716   | 4.300  | ng | # 95 |
| 50) Dimethylphthalate      | 9.67  | 163 | 77200   | 3.424  | ng | # 94 |
| 71) Phenanthrene           | 11.48 | 178 | 267173  | 10.154 | ng | 98   |
| 72) Anthracene             | 11.53 | 178 | 109466  | 4.196  | ng | 98   |
| 75) Fluoranthene           | 12.67 | 202 | 453607  | 13.696 | ng | 99   |
| 78) Pyrene                 | 12.90 | 202 | 541716  | 17.127 | ng | 99   |
| 81) Benzo(a)anthracene     | 14.10 | 228 | 213435  | 8.253  | ng | # 93 |
| 83) Chrysene               | 14.13 | 228 | 236138m | 9.423  | ng |      |
| 86) Indeno(1,2,3-cd)pyrene | 17.15 | 276 | 56821   | 2.664  | ng | 93   |
| 88) Benzo(b)fluoranthene   | 15.17 | 252 | 164770m | 8.472  | ng |      |
| 89) Benzo(k)fluoranthene   | 15.20 | 252 | 40531m  | 2.212  | ng |      |
| 90) Benzo(a)pyrene         | 15.56 | 252 | 114328  | 6.685  | ng | # 89 |
| 92) Benzo(g,h,i)perylene   | 17.62 | 276 | 59793   | 4.078  | ng | 93   |

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

