

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092820\  
 Data File : BF121705.D  
 Acq On : 28 Sep 2020 11:31  
 Operator : JU/CG  
 Sample : L4152-02DL 10X  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 WR-PILE-3-4DL

Manual Integrations  
 APPROVED

mohammad  
 9/29/2020 3:24:08 PM

Quant Time: Sep 28 12:31:51 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Sep 28 11:40:42 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 6.79  | 152  | 163589   | 20.00 | ng    | 0.00     |
| 21) Naphthalene-d8        | 8.07  | 136  | 642801   | 20.00 | ng    | 0.00     |
| 39) Acenaphthene-d10      | 9.83  | 164  | 335819   | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10      | 11.32 | 188  | 663635   | 20.00 | ng    | 0.00     |
| 76) Chrysene-d12          | 13.96 | 240  | 515934   | 20.00 | ng    | 0.00     |
| 86) Perylene-d12          | 15.40 | 264  | 432411   | 20.00 | ng    | 0.00     |

|                             |       |     |        |      |    |       |
|-----------------------------|-------|-----|--------|------|----|-------|
| System Monitoring Compounds |       |     |        |      |    |       |
| 5) 2-Fluorophenol           | 5.41  | 112 | 32957  | 2.99 | ng | 0.00  |
| 7) Phenol-d6                | 6.43  | 99  | 92081  | 6.38 | ng | -0.01 |
| 23) Nitrobenzene-d5         | 7.35  | 82  | 82145  | 6.86 | ng | -0.02 |
| 42) 2,4,6-Tribromophenol    | 10.62 | 330 | 1190   | 0.29 | ng | 0.00  |
| 45) 2-Fluorobiphenyl        | 9.15  | 172 | 171482 | 7.73 | ng | 0.00  |
| 79) Terphenyl-d14           | 12.89 | 244 | 202238 | 7.94 | ng | -0.01 |

|                            |       |     |         |           |        |
|----------------------------|-------|-----|---------|-----------|--------|
| Target Compounds           |       |     |         |           | Qvalue |
| 52) Acenaphthene           | 9.86  | 154 | 93946   | 4.569 ng  | 99     |
| 55) Dibenzofuran           | 10.03 | 168 | 115803  | 4.000 ng  | 93     |
| 58) Fluorene               | 10.37 | 166 | 139589  | 6.304 ng  | 100    |
| 71) Phenanthrene           | 11.35 | 178 | 1696655 | 46.924 ng | 100    |
| 72) Anthracene             | 11.39 | 178 | 505129  | 13.538 ng | 99     |
| 73) Carbazole              | 11.55 | 167 | 180853  | 5.371 ng  | 98     |
| 75) Fluoranthene           | 12.53 | 202 | 2235999 | 57.931 ng | 99     |
| 78) Pyrene                 | 12.76 | 202 | 1893441 | 50.230 ng | 99     |
| 81) Benzo(a)anthracene     | 13.94 | 228 | 1088703 | 33.354 ng | 99     |
| 83) Chrysene               | 13.98 | 228 | 853947  | 25.836 ng | 97     |
| 87) Indeno(1,2,3-cd)pyrene | 16.82 | 276 | 418187  | 14.850 ng | 95     |
| 88) Benzo(b)fluoranthene   | 14.99 | 252 | 938431m | 32.464 ng |        |
| 89) Benzo(k)fluoranthene   | 15.00 | 252 | 329205m | 12.436 ng |        |
| 90) Benzo(a)pyrene         | 15.34 | 252 | 638599  | 25.989 ng | 98     |
| 91) Dibenzo(a,h)anthracene | 16.82 | 278 | 99215m  | 4.232 ng  |        |
| 92) Benzo(g,h,i)perylene   | 17.24 | 276 | 378363  | 16.420 ng | 99     |

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

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