

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092920\  
 Data File : BF121734.D  
 Acq On : 29 Sep 2020 16:48  
 Operator : JU/CG  
 Sample : L4172-10  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 B-7(12-14)

Manual Integrations  
 APPROVED

mohammad  
 10/1/2020 1:29:49 PM

Quant Time: Sep 29 17:43:45 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Sep 29 11:12:14 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 6.78  | 152  | 140220   | 20.00 | ng    | 0.00     |
| 21) Naphthalene-d8        | 8.07  | 136  | 528204   | 20.00 | ng    | 0.00     |
| 39) Acenaphthene-d10      | 9.82  | 164  | 256189   | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10      | 11.31 | 188  | 449484   | 20.00 | ng    | 0.00     |
| 76) Chrysene-d12          | 13.95 | 240  | 322058   | 20.00 | ng    | 0.00     |
| 86) Perylene-d12          | 15.39 | 264  | 269302   | 20.00 | ng    | 0.00     |

|                             |       |     |        |       |    |       |
|-----------------------------|-------|-----|--------|-------|----|-------|
| System Monitoring Compounds |       |     |        |       |    |       |
| 5) 2-Fluorophenol           | 5.41  | 112 | 753985 | 79.91 | ng | 0.00  |
| 7) Phenol-d6                | 6.42  | 99  | 963907 | 77.87 | ng | 0.00  |
| 23) Nitrobenzene-d5         | 7.35  | 82  | 625871 | 63.62 | ng | -0.01 |
| 42) 2,4,6-Tribromophenol    | 10.61 | 330 | 230352 | 72.34 | ng | 0.00  |
| 45) 2-Fluorobiphenyl        | 9.15  | 172 | 984032 | 58.17 | ng | 0.00  |
| 79) Terphenyl-d14           | 12.89 | 244 | 828979 | 52.15 | ng | 0.00  |

|                                |       |     |         |        |    |     |
|--------------------------------|-------|-----|---------|--------|----|-----|
| Target Compounds               |       |     |         | Qvalue |    |     |
| 31) Naphthalene                | 8.09  | 128 | 176726  | 6.338  | ng | 99  |
| 37) 2-Methylnaphthalene        | 8.78  | 142 | 48208   | 2.638  | ng | 99  |
| 38) 1-Methylnaphthalene        | 8.88  | 142 | 34688   | 2.040  | ng | 98  |
| 49) Acenaphthylene             | 9.68  | 152 | 64638   | 2.603  | ng | 99  |
| 50) Dimethylphthalate          | 9.54  | 163 | 154004  | 8.294  | ng | 97  |
| 52) Acenaphthene               | 9.86  | 154 | 152111  | 9.698  | ng | 99  |
| 55) Dibenzofuran               | 10.03 | 168 | 130143  | 5.892  | ng | 98  |
| 58) Fluorene                   | 10.37 | 166 | 163307  | 9.668  | ng | 100 |
| 71) Phenanthrene               | 11.34 | 178 | 1440443 | 58.818 | ng | 100 |
| 72) Anthracene                 | 11.38 | 178 | 374714  | 14.827 | ng | 99  |
| 73) Carbazole                  | 11.54 | 167 | 201067  | 8.816  | ng | 97  |
| 75) Fluoranthene               | 12.53 | 202 | 1791747 | 68.537 | ng | 99  |
| 78) Pyrene                     | 12.76 | 202 | 1495480 | 63.555 | ng | 99  |
| 81) Benzo(a)anthracene         | 13.94 | 228 | 765216  | 37.557 | ng | 98  |
| 83) Chrysene                   | 13.97 | 228 | 666815  | 32.319 | ng | 97  |
| 84) Bis(2-ethylhexyl)phthalate | 13.93 | 149 | 28081   | 2.209  | ng | 99  |
| 87) Indeno(1,2,3-cd)pyrene     | 16.82 | 276 | 306031  | 17.450 | ng | 94  |
| 88) Benzo(b)fluoranthene       | 14.98 | 252 | 784317m | 43.566 | ng |     |
| 89) Benzo(k)fluoranthene       | 15.00 | 252 | 186602m | 11.319 | ng |     |
| 90) Benzo(a)pyrene             | 15.33 | 252 | 478546  | 31.271 | ng | 95  |
| 91) Dibenzo(a,h)anthracene     | 16.82 | 278 | 77938m  | 5.339  | ng |     |
| 92) Benzo(g,h,i)perylene       | 17.25 | 276 | 263635  | 18.371 | ng | 98  |

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

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