

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100520\  
 Data File : BF121827.D  
 Acq On : 5 Oct 2020 20:20  
 Operator : JU/CG  
 Sample : L4247-06  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 B-19(4-5)

Manual Integrations  
 APPROVED

mohammad  
 10/7/2020 10:36:16 AM

Quant Time: Oct 06 00:57:36 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Oct 01 14:29:14 2020  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc  | Units | Dev(Min)    |
|-----------------------------|-------|------|----------|-------|-------|-------------|
| 1) 1,4-Dichlorobenzene-d4   | 6.76  | 152  | 113764   | 20.00 | ng    | 0.00        |
| 21) Naphthalene-d8          | 8.05  | 136  | 432195   | 20.00 | ng    | -0.01       |
| 39) Acenaphthene-d10        | 9.80  | 164  | 201958   | 20.00 | ng    | -0.01       |
| 64) Phenanthrene-d10        | 11.29 | 188  | 334082   | 20.00 | ng    | 0.00        |
| 76) Chrysene-d12            | 13.92 | 240  | 248708   | 20.00 | ng    | -0.01       |
| 86) Perylene-d12            | 15.36 | 264  | 242594   | 20.00 | ng    | 0.00        |
| System Monitoring Compounds |       |      |          |       |       |             |
| 5) 2-Fluorophenol           | 5.38  | 112  | 635860   | 83.06 | ng    | 0.00        |
| 7) Phenol-d6                | 6.40  | 99   | 804108   | 80.07 | ng    | -0.01       |
| 23) Nitrobenzene-d5         | 7.33  | 82   | 543758   | 67.55 | ng    | -0.01       |
| 42) 2,4,6-Tribromophenol    | 10.59 | 330  | 201337   | 80.21 | ng    | -0.01       |
| 45) 2-Fluorobiphenyl        | 9.12  | 172  | 838873   | 62.91 | ng    | -0.01       |
| 79) Terphenyl-d14           | 12.87 | 244  | 667598   | 54.38 | ng    | -0.01       |
| Target Compounds            |       |      |          |       |       |             |
| 50) Dimethylphthalate       | 9.52  | 163  | 96329    | 6.581 | ng    | Qvalue # 97 |
| 71) Phenanthrene            | 11.31 | 178  | 84022    | 4.616 | ng    | 97          |
| 75) Fluoranthene            | 12.50 | 202  | 134497   | 6.922 | ng    | 97          |
| 78) Pyrene                  | 12.72 | 202  | 138606   | 7.628 | ng    | 98          |
| 81) Benzo(a)anthracene      | 13.91 | 228  | 71751    | 4.560 | ng    | 97          |
| 83) Chrysene                | 13.94 | 228  | 61333    | 3.849 | ng    | 95          |
| 87) Indeno(1,2,3-cd)pyrene  | 16.77 | 276  | 33773    | 2.138 | ng    | 95          |
| 88) Benzo(b)fluoranthene    | 14.94 | 252  | 87219m   | 5.378 | ng    |             |
| 90) Benzo(a)pyrene          | 15.30 | 252  | 62251    | 4.516 | ng    | # 89        |
| 92) Benzo(g,h,i)perylene    | 17.20 | 276  | 33884    | 2.621 | ng    | 91          |

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

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