

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062620\  
 Data File : BP002559.D  
 Acq On : 26 Jun 2020 19:37  
 Operator : CG/JU  
 Sample : L3113-01  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DRUM-COMP

Manual Integrations  
 APPROVED

mohammad  
 6/29/2020 3:18:53 PM

Quant Time: Jun 27 00:20:19 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jun 24 13:54:06 2020  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.58  | 152  | 123391   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.35 | 136  | 514381   | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.23 | 164  | 340434   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 16.99 | 188  | 747387   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.10 | 240  | 692738   | 20.00  | ng    | 0.00     |
| 86) Perylene-d12            | 23.30 | 264  | 603089   | 20.00  | ng    | 0.01     |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.20  | 112  | 719148   | 107.85 | ng    | 0.00     |
| 7) Phenol-d6                | 6.78  | 99   | 1040721  | 117.22 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.73  | 82   | 739785   | 85.89  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.73 | 330  | 495107   | 151.90 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.85 | 172  | 1698626  | 83.90  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.58 | 244  | 2466470  | 94.68  | ng    | 0.00     |
| Target Compounds            |       |      |          |        |       |          |
| 50) Dimethylphthalate       | 13.69 | 163  | 172471   | 7.394  | ng    | 100      |
| 71) Phenanthrene            | 17.03 | 178  | 126429   | 3.616  | ng    | 98       |
| 75) Fluoranthene            | 19.02 | 202  | 417164   | 9.996  | ng    | 98       |
| 78) Pyrene                  | 19.37 | 202  | 472210   | 11.129 | ng    | 100      |
| 81) Benzo(a)anthracene      | 21.08 | 228  | 290137   | 7.338  | ng    | 96       |
| 83) Chrysene                | 21.13 | 228  | 407542   | 10.879 | ng    | 97       |
| 87) Indeno(1,2,3-cd)pyrene  | 25.51 | 276  | 168147   | 4.458  | ng    | 97       |
| 88) Benzo(b)fluoranthene    | 22.64 | 252  | 497174m  | 15.287 | ng    |          |
| 89) Benzo(k)fluoranthene    | 22.68 | 252  | 145430m  | 4.706  | ng    |          |
| 90) Benzo(a)pyrene          | 23.20 | 252  | 292823   | 9.608  | ng    | 99       |
| 92) Benzo(g,h,i)perylene    | 26.18 | 276  | 181978   | 5.807  | ng    | 99       |

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

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