

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071020\  
 Data File : BM026741.D  
 Acq On : 10 Jul 2020 19:28  
 Operator : JU/CG  
 Sample : L3217-17  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 E3-41-NHL

Manual Integrations  
 APPROVED

mohammad  
 7/13/2020 1:03:09 PM

Quant Time: Jul 13 09:17:35 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070820.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jul 10 10:38:33 2020  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|-------|------|----------|---------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.33  | 152  | 97178    | 20.00   | ng    | 0.01     |
| 21) Naphthalene-d8          | 10.08 | 136  | 388271   | 20.00   | ng    | 0.00     |
| 39) Acenaphthene-d10        | 13.98 | 164  | 233414   | 20.00   | ng    | 0.00     |
| 64) Phenanthrene-d10        | 16.74 | 188  | 463104   | 20.00   | ng    | 0.00     |
| 76) Chrysene-d12            | 20.96 | 240  | 495696   | 20.00   | ng    | 0.00     |
| 86) Perylene-d12            | 23.03 | 264  | 495741   | 20.00   | ng    | 0.00     |
| System Monitoring Compounds |       |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.08  | 112  | 18954m   | 3.27    | ng    | 0.12     |
| 7) Phenol-d6                | 6.76  | 99   | 33472    | 3.62    | ng    | 0.23     |
| 23) Nitrobenzene-d5         | 8.47  | 82   | 28011    | 3.08    | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.49 | 330  | 12268    | 3.61    | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.59 | 172  | 57530    | 3.28    | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.41 | 244  | 91684    | 3.64    | ng    | 0.00     |
| Target Compounds            |       |      |          |         |       |          |
| 10) Phenol                  | 6.79  | 94   | 471591   | 51.654  | ng    | 98       |
| 17) 2-Methylphenol          | 7.80  | 107  | 122585   | 17.668  | ng    | 99       |
| 20) 3+4-Methylphenols       | 8.13  | 107  | 599436   | 62.842  | ng    | 94       |
| 27) 2,4-Dimethylphenol      | 9.30  | 122  | 110888m  | 19.298  | ng    |          |
| 31) Naphthalene             | 10.13 | 128  | 2648190  | 122.629 | ng    | 100      |
| 32) Benzoic acid            | 9.46  | 122  | 40087    | 13.947  | ng    | 96       |
| 37) 2-Methylnaphthalene     | 11.76 | 142  | 3352580  | 209.575 | ng    | 100      |
| 38) 1-Methylnaphthalene     | 11.98 | 142  | 1650471  | 107.690 | ng    | 99       |
| 46) 1,1'-Biphenyl           | 12.80 | 154  | 881177   | 47.793  | ng    | 98       |
| 52) Acenaphthene            | 14.05 | 154  | 800832   | 55.342  | ng    | 97       |
| 55) Dibenzofuran            | 14.39 | 168  | 739925   | 31.186  | ng    | 97       |
| 58) Fluorene                | 15.05 | 166  | 139423   | 7.097   | ng    | 97       |
| 71) Phenanthrene            | 16.79 | 178  | 140726   | 5.147   | ng    | 97       |

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

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