

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120721\  
 Data File : BP008259.D  
 Acq On : 07 Dec 2021 20:06  
 Operator : CG/JU  
 Sample : M4932-02  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DRUM1-2

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021  
 Supervised By :Sohil Jodhani 12/10/2021

Quant Time: Dec 08 01:10:27 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP112521.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Dec 03 14:39:24 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-------------------------------|--------|------|----------|---------|-------|----------|
| -----                         |        |      |          |         |       |          |
| Internal Standards            |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.781  | 152  | 59393    | 20.000  | ng    | -0.02    |
| 21) Naphthalene-d8            | 10.581 | 136  | 210690   | 20.000  | ng    | -0.02    |
| 39) Acenaphthene-d10          | 14.434 | 164  | 131373   | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10          | 17.198 | 188  | 285426   | 20.000  | ng    | -0.01    |
| 76) Chrysene-d12              | 21.304 | 240  | 352640   | 20.000  | ng    | -0.01    |
| 86) Perylene-d12              | 23.692 | 264  | 426810   | 20.000  | ng    | -0.02    |
|                               |        |      |          |         |       |          |
| System Monitoring Compounds   |        |      |          |         |       |          |
| 5) 2-Fluorophenol             | 5.381  | 112  | 518847   | 140.655 | ng    | 0.00     |
| 7) Phenol-d6                  | 6.975  | 99   | 611520   | 122.805 | ng    | -0.01    |
| 23) Nitrobenzene-d5           | 8.946  | 82   | 482184   | 104.070 | ng    | -0.02    |
| 42) 2,4,6-Tribromophenol      | 15.933 | 330  | 270660   | 161.187 | ng    | -0.02    |
| 45) 2-Fluorobiphenyl          | 13.051 | 172  | 913456   | 100.997 | ng    | -0.02    |
| 79) Terphenyl-d14             | 19.769 | 244  | 1597636  | 94.684  | ng    | -0.01    |
|                               |        |      |          |         |       |          |
| Target Compounds              |        |      |          |         |       | Qvalue   |
| 31) Naphthalene               | 10.628 | 128  | 47819    | 4.431   | ng    | 95       |
| 80) Butylbenzylphthalate      | 20.445 | 149  | 199085   | 19.826  | ng    | 99       |
| 83) Chrysene                  | 21.339 | 228  | 47008    | 2.114   | ng    | 98       |
| 84) Bis(2-ethylhexyl)phtha... | 21.215 | 149  | 327534   | 21.483  | ng    | 100      |
| 85) Di-n-octyl phthalate      | 22.133 | 149  | 104705   | 4.043   | ng    | 98       |
| 88) Benzo(b)fluoranthene      | 22.968 | 252  | 65637m   | 2.551   | ng    |          |
| 89) Benzo(k)fluoranthene      | 23.015 | 252  | 72755    | 2.888   | ng    | 98       |
| 90) Benzo(a)pyrene            | 23.592 | 252  | 54425    | 2.475   | ng    | # 92     |
| -----                         |        |      |          |         |       |          |

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

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