

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF120324\  
 Data File : BF140716.D  
 Acq On : 03 Dec 2024 17:34  
 Operator : RC/JU  
 Sample : P5045-01DL 2X  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SU-04-11222024DL

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 12/04/2024  
 Supervised By :mohammad ahmed 12/05/2024

Quant Time: Dec 03 17:54:09 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF112124.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Nov 21 15:23:48 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.869  | 152  | 78891    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.151  | 136  | 293060   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.904  | 164  | 153756   | 20.000 | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.398 | 188  | 248875   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 14.057 | 240  | 146599   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12            | 15.580 | 264  | 175946   | 20.000 | ng    | 0.04     |
|                             |        |      |          |        |       |          |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.499  | 112  | 200589   | 43.382 | ng    | 0.00     |
| 7) Phenol-d6                | 6.504  | 99   | 255218   | 41.752 | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 7.428  | 82   | 167951   | 29.314 | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.698 | 330  | 60062    | 36.524 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.222  | 172  | 323041   | 31.304 | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.980 | 244  | 243943   | 25.911 | ng    | -0.01    |
|                             |        |      |          |        |       |          |
| Target Compounds            |        |      |          |        |       | Qvalue   |
| 58) Fluorene                | 10.457 | 166  | 20621    | 2.017  | ng    | 98       |
| 71) Phenanthrene            | 11.422 | 178  | 226837   | 18.961 | ng    | 99       |
| 72) Anthracene              | 11.475 | 178  | 79560    | 6.799  | ng    | 98       |
| 75) Fluoranthene            | 12.616 | 202  | 511490   | 39.379 | ng    | 100      |
| 78) Pyrene                  | 12.845 | 202  | 444891   | 32.821 | ng    | 99       |
| 81) Benzo(a)anthracene      | 14.045 | 228  | 241372   | 24.873 | ng    | 97       |
| 83) Chrysene                | 14.086 | 228  | 177358   | 19.987 | ng    | 98       |
| 87) Indeno(1,2,3-cd)pyrene  | 17.098 | 276  | 112409   | 9.803  | ng    | 96       |
| 88) Benzo(b)fluoranthene    | 15.145 | 252  | 273189m  | 24.720 | ng    |          |
| 89) Benzo(k)fluoranthene    | 15.157 | 252  | 101150m  | 10.459 | ng    |          |
| 90) Benzo(a)pyrene          | 15.516 | 252  | 201814   | 22.459 | ng    | 97       |
| 91) Dibenzo(a,h)anthracene  | 17.104 | 278  | 28510    | 3.036  | ng    | # 85     |
| 92) Benzo(g,h,i)perylene    | 17.557 | 276  | 102663   | 10.714 | ng    | 100      |
| -----                       |        |      |          |        |       |          |

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

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