

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF110222\  
 Data File : BF130943.D  
 Acq On : 02 Nov 2022 22:13  
 Operator : CG\JU  
 Sample : N5395-11  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 S-8-103122

Manual Integrations  
 APPROVED

Reviewed By : Christian Giraldo 11/03/2022  
 Supervised By : Jagrut Upadhyay 11/03/2022

Quant Time: Nov 03 03:48:23 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Nov 03 03:37:09 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.922  | 152  | 63288    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.210  | 136  | 222969   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 9.969  | 164  | 100682   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.457 | 188  | 144496   | 20.000  | ng    | -0.01    |
| 76) Chrysene-d12            | 14.110 | 240  | 137121   | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 15.621 | 264  | 146622   | 20.000  | ng    | 0.00     |
|                             |        |      |          |         |       |          |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.546  | 112  | 518559   | 142.046 | ng    | 0.00     |
| 7) Phenol-d6                | 6.545  | 99   | 632321   | 133.288 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.487  | 82   | 422332   | 114.214 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.757 | 330  | 114907   | 135.577 | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 9.286  | 172  | 582334   | 87.446  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.051 | 244  | 454166   | 60.694  | ng    | 0.00     |
|                             |        |      |          |         |       |          |
| Target Compounds            |        |      |          |         |       | Qvalue   |
| 20) 3+4-Methylphenols       | 7.316  | 107  | 13634    | 3.017   | ng    | 85       |
| 71) Phenanthrene            | 11.480 | 178  | 15128    | 2.028   | ng    | 98       |
| 75) Fluoranthene            | 12.674 | 202  | 37013    | 4.671   | ng    | 96       |
| 78) Pyrene                  | 12.904 | 202  | 32902    | 2.844   | ng    | 97       |
| 81) Benzo(a)anthracene      | 14.098 | 228  | 22305    | 2.453   | ng    | # 93     |
| 83) Chrysene                | 14.133 | 228  | 23126m   | 2.637   | ng    |          |
| 88) Benzo(b)fluoranthene    | 15.168 | 252  | 29744m   | 3.138   | ng    |          |
| 90) Benzo(a)pyrene          | 15.551 | 252  | 21500    | 2.822   | ng    | # 91     |
| -----                       |        |      |          |         |       |          |

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

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