

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080822\
 Data File : BF129652.D
 Acq On : 09 Aug 2022 02:44
 Operator : CG\JU
 Sample : N3963-18
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RVE-222

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 08/09/2022
 Supervised By :Jagrut Upadhyay 08/09/2022

Quant Time: Aug 09 04:09:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 16:27:34 2022
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) | |
|-----------------------------|--------|------|----------|---------|-------|----------|-----------|
| Internal Standards | | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 6.839 | 152 | 127539 | 20.000 | ng | -0.01 | |
| 21) Naphthalene-d8 | 8.122 | 136 | 479383 | 20.000 | ng | -0.01 | |
| 39) Acenaphthene-d10 | 9.880 | 164 | 206405 | 20.000 | ng | -0.01 | |
| 64) Phenanthrene-d10 | 11.369 | 188 | 293617 | 20.000 | ng | -0.01 | |
| 76) Chrysene-d12 | 14.015 | 240 | 273327 | 20.000 | ng | -0.01 | |
| 86) Perylene-d12 | 15.492 | 264 | 199556 | 20.000 | ng | -0.01 | |
| System Monitoring Compounds | | | | | | | |
| 5) 2-Fluorophenol | 5.475 | 112 | 895518 | 114.381 | ng | 0.00 | |
| 7) Phenol-d6 | 6.492 | 99 | 1096308 | 111.403 | ng | -0.01 | |
| 23) Nitrobenzene-d5 | 7.404 | 82 | 693800 | 79.005 | ng | -0.02 | |
| 42) 2,4,6-Tribromophenol | 10.674 | 330 | 201751 | 101.667 | ng | -0.01 | |
| 45) 2-Fluorobiphenyl | 9.198 | 172 | 1177683 | 88.264 | ng | -0.01 | |
| 79) Terphenyl-d14 | 12.957 | 244 | 1018457 | 70.297 | ng | -0.01 | |
| Target Compounds | | | | | | | |
| 71) Phenanthrene | 11.392 | 178 | 43014 | 2.684 | ng | | Qvalue 97 |
| 75) Fluoranthene | 12.586 | 202 | 128993 | 7.943 | ng | | 98 |
| 78) Pyrene | 12.816 | 202 | 128271 | 5.612 | ng | | 98 |
| 81) Benzo(a)anthracene | 14.004 | 228 | 74170 | 3.972 | ng | | 98 |
| 83) Chrysene | 14.039 | 228 | 85041 | 4.833 | ng | | 98 |
| 87) Indeno(1,2,3-cd)pyrene | 16.968 | 276 | 39203 | 2.917 | ng | | 97 |
| 88) Benzo(b)fluoranthene | 15.057 | 252 | 103835m | 7.844 | ng | | |
| 89) Benzo(k)fluoranthene | 15.080 | 252 | 37457m | 2.888 | ng | | |
| 90) Benzo(a)pyrene | 15.427 | 252 | 60695 | 5.648 | ng | # | 94 |
| 92) Benzo(g,h,i)perylene | 17.421 | 276 | 39840 | 3.565 | ng | | 97 |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080822\
 Data File : BF129652.D
 Acq On : 09 Aug 2022 02:44
 Operator : CG\JU
 Sample : N3963-18
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
ClientSampleId :
 RVE-222

Quant Time: Aug 09 04:09:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 16:27:34 2022
 Response via : Initial Calibration

Manual Integrations
APPROVED
 Reviewed By :Christian Giraldo 08/09/2022
 Supervised By :Jagrut Upadhyay 08/09/2022

