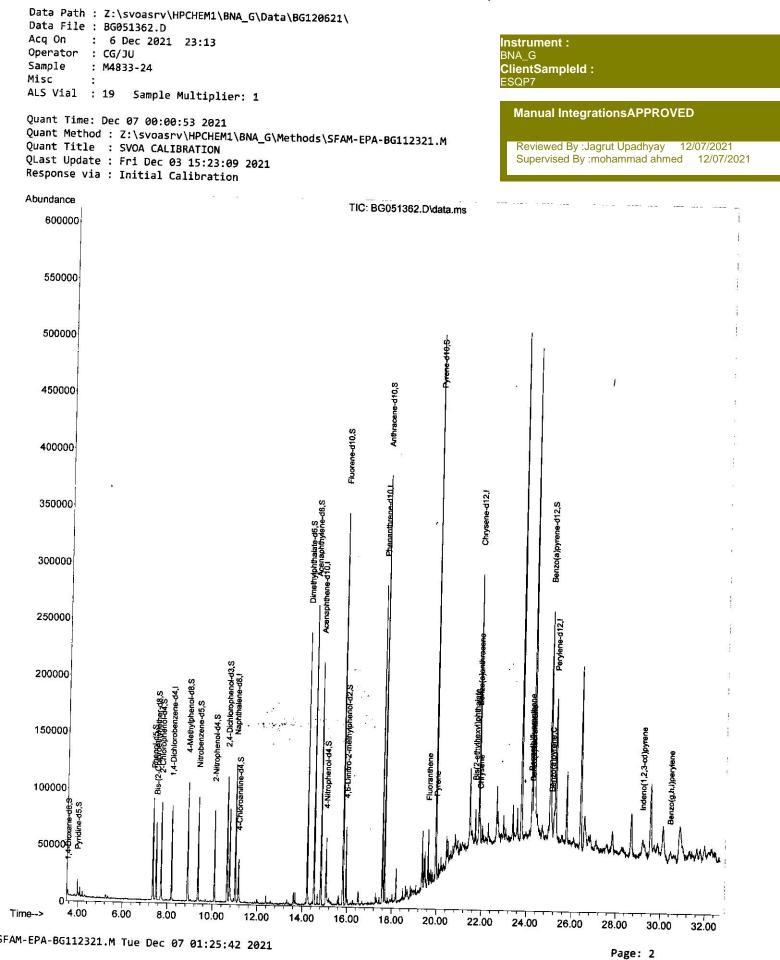
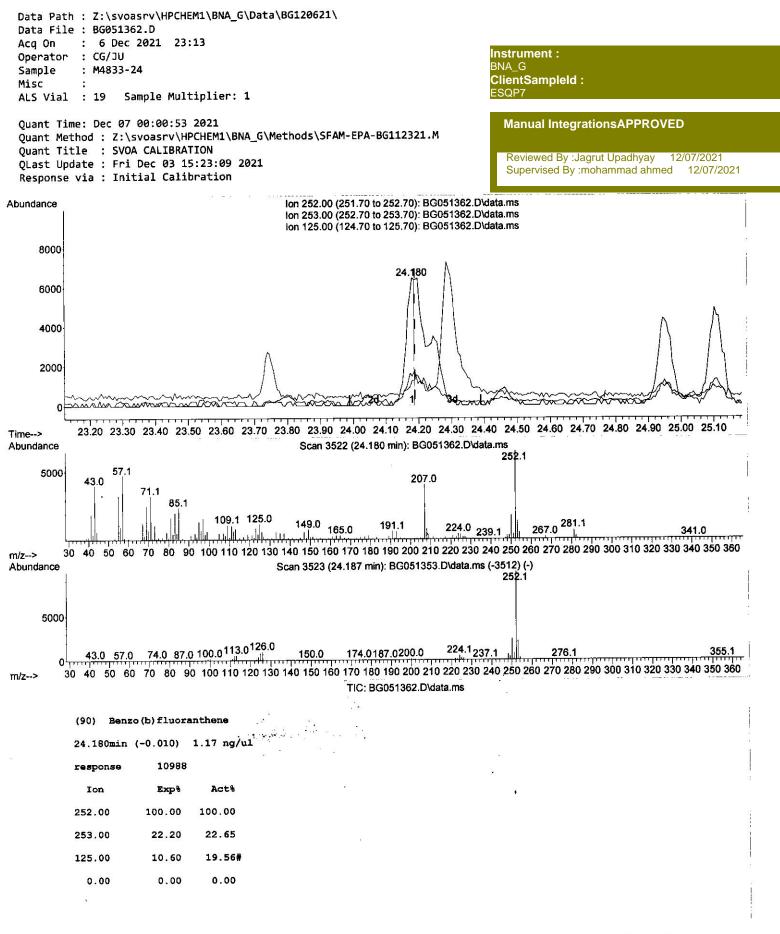
Quantitation Report

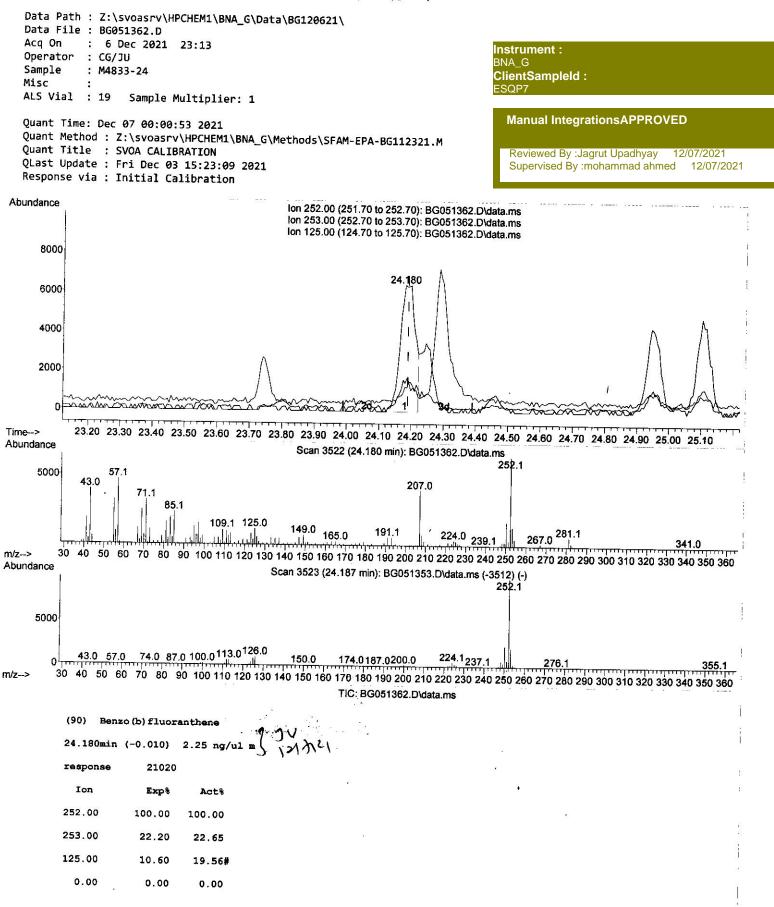
(QT Reviewed)

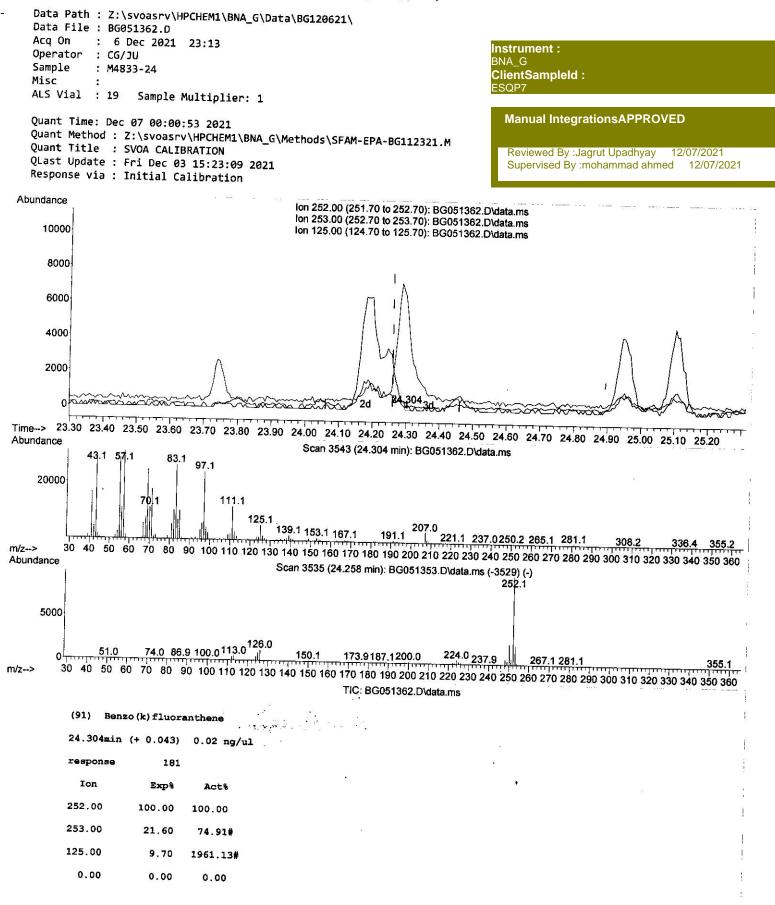


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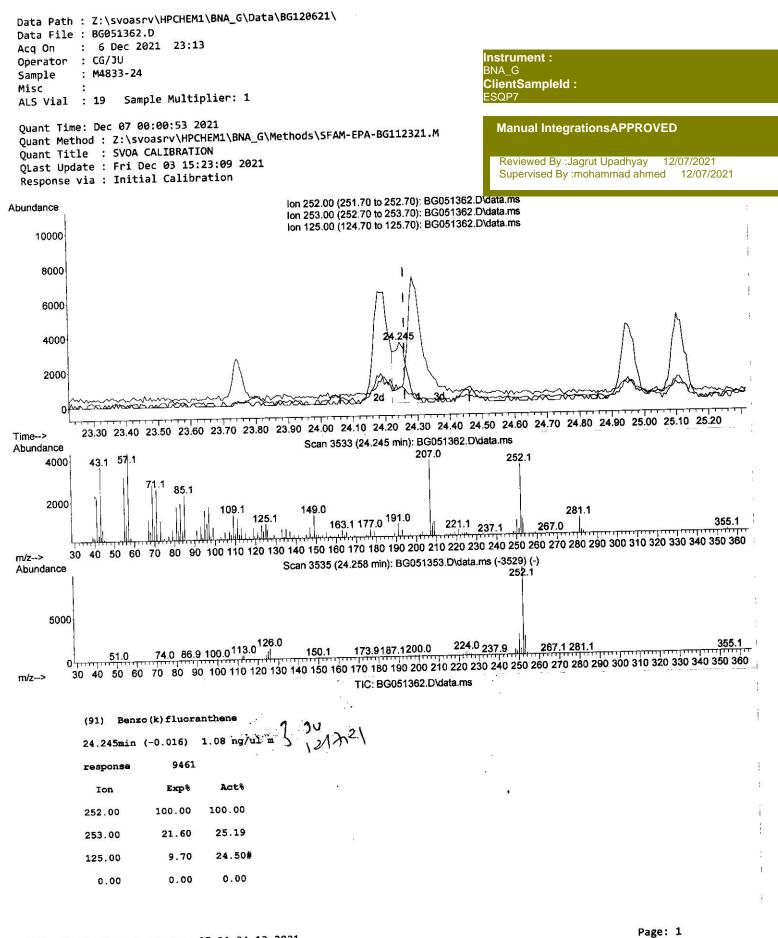


SFAM-EPA-BG112321.M Tue Dec 07 01:23:33 2021



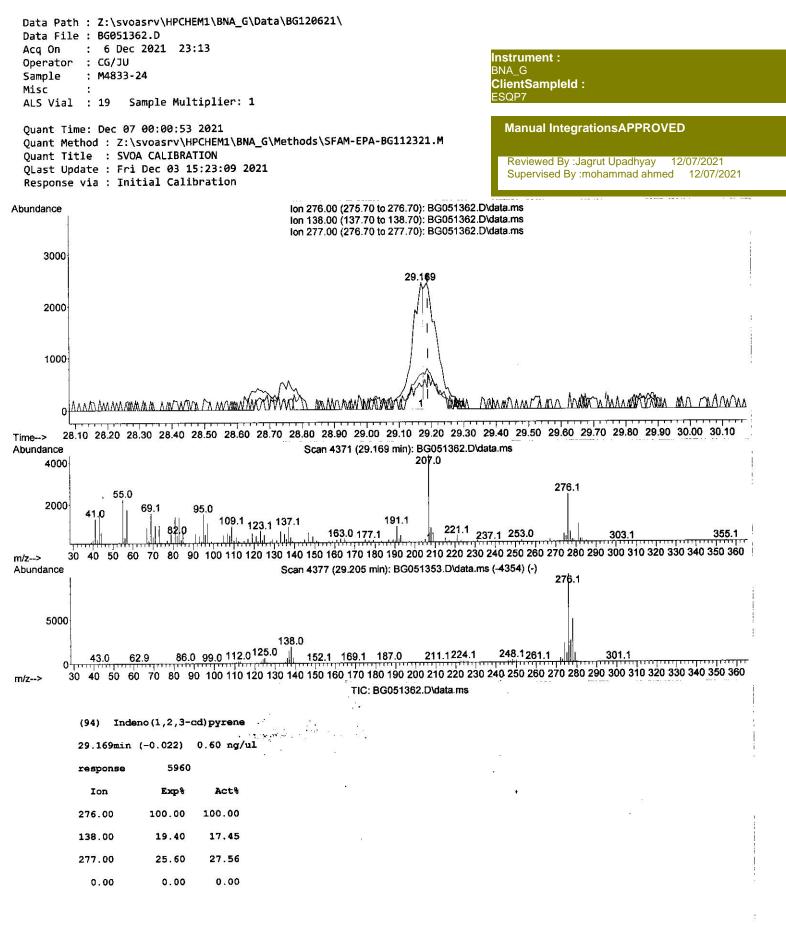


SFAM-EPA-BG112321.M Tue Dec 07 01:24:01 2021



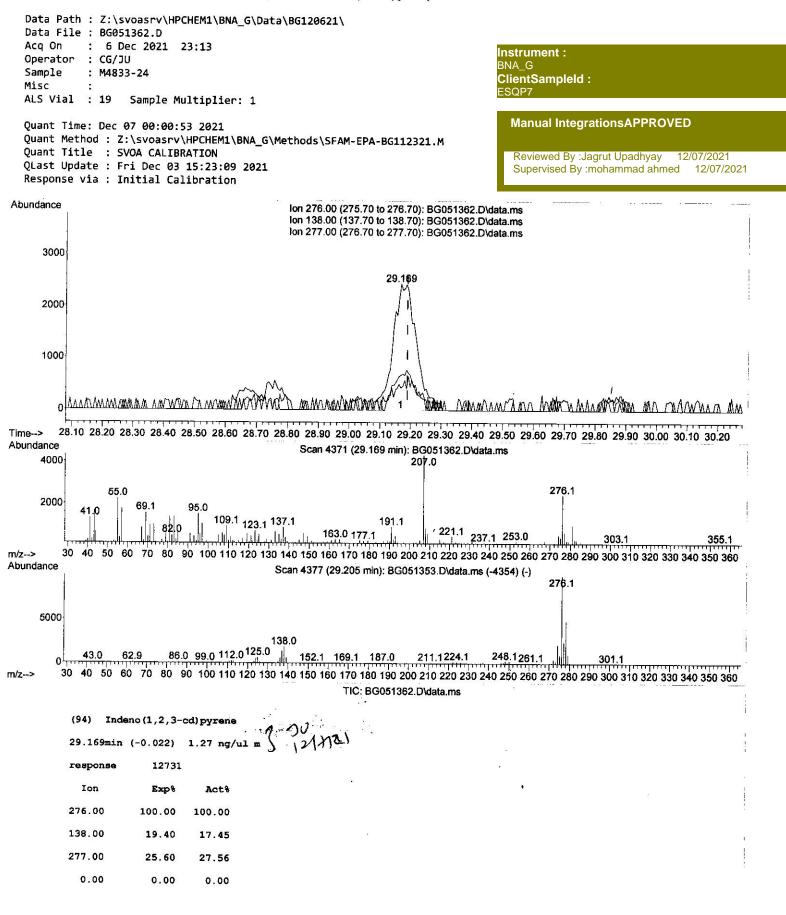
SFAM-EPA-BG112321.M Tue Dec 07 01:24:13 2021

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SFAM-EPA-BG112321.M Tue Dec 07 01:24:36 2021

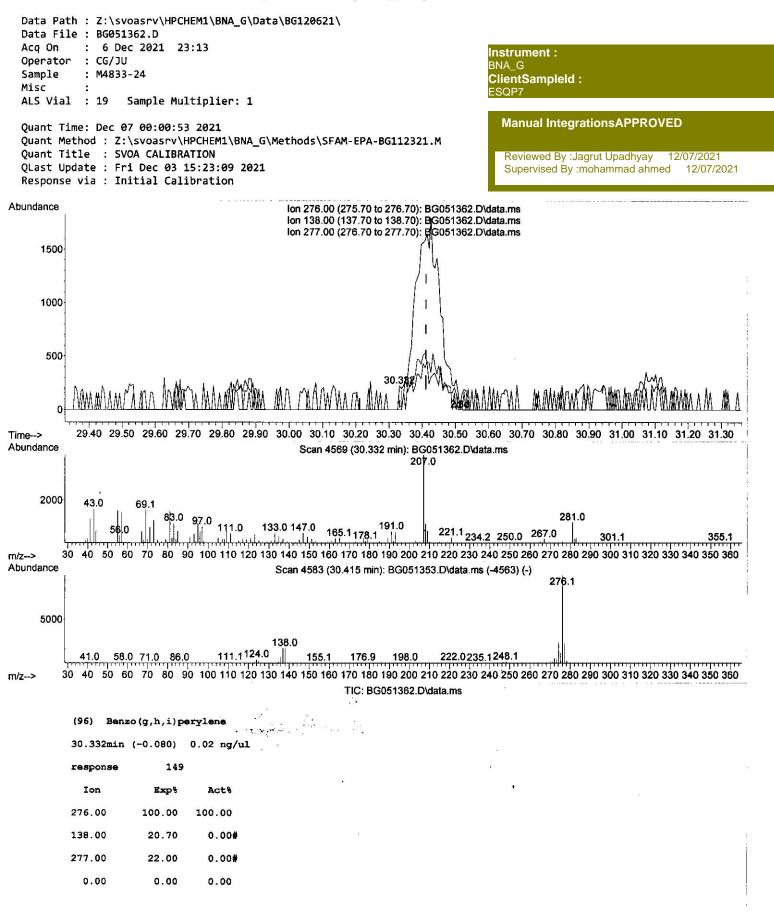
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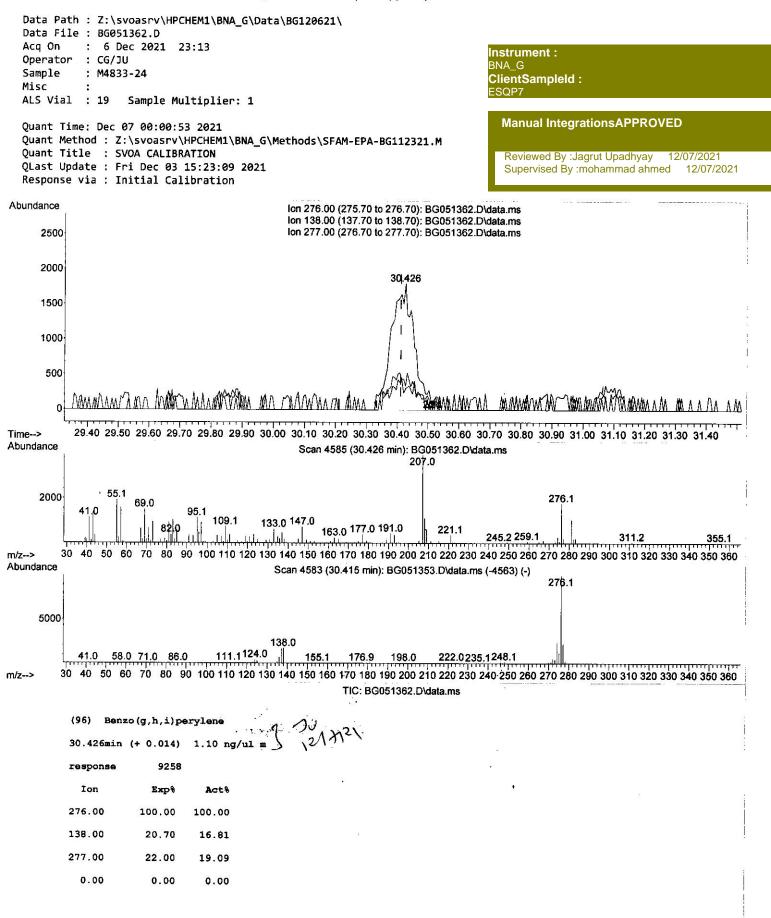
SFAM-EPA-BG112321.M Tue Dec 07 01:24:48 2021

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SFAM-EPA-BG112321.M Tue Dec 07 01:25:05 2021



SFAM-EPA-BG112321.M Tue Dec 07 01:25:20 2021

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| Data Path : Z:\svoasrv\HPCHEM1\BN Data File : BG051362.D Acq On : 6 Dec 2021 23:13 | A_G\Data\ | BG120 | 621\ | | _ | |
|--|------------------|--------|-----------------|------------------------------|--------------|--|
| Operator : CG/JU | | | | | | strument : |
| Sample : M4833-24 | | | | | | A_G |
| Misc : | 12 | | | | | entSampleId : |
| ALS Vial : 19 Sample Multiplie | r: 1 | | | | ES | QP7 |
| | | | | | | |
| Quant Time: Dec 07 00:00:53 2021 Quant Method : Z:\svoasrv\HPCHEMI | | othods | SFAM-EPA | -BG112321.M | | Manual IntegrationsAPPROVED |
| Quant Method : Z:\svoasrv\HPCHEMJ | | | | | | |
| Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:0 | 9 2021 | | | | | Reviewed By :Jagrut Upadhyay 12/07/2021 |
| Response via : Initial Calibratio | on | | | | | Supervised By :mohammad ahmed 12/07/2021 |
| Response via : iniciae delle | | | | | Min | |
| Compound | R.T. Ç | Ion | Response | Conc Units Dev(| (MIN) | |
| | | | | | | |
| Internal Standards | | | 225.01 | 20.000 ng/ul | -0.02 | |
| 1) 1,4-Dichlorobenzene-d4 | 8.187 | 152 | 23501 103351 | 20.000 ng/ul | -0.02 | |
| 20) Naphthalene-d8 | 11.013 | | 73607 | 20.000 ng/ul | 0.00 | |
| 38) Acenaphthene-d10 | 14.821 | | 167397 | 20.000 ng/ul | 0.00 | |
| 64) Phenanthrene-d10 | 17.570 | 240 | 141885 | 20.000 ng/ul | 0.00 | |
| 79) Chrysene-d12 | 21.871 25.273 | 264 | 138660 | 20.000 ng/ul | 0.00 | |
| 88) Perylene-d12 | 25.275 | 204 | 190000 | | | |
| - whitening Compounds | | | | | | |
| System Monitoring Compounds | 3,528 | 96 | 2461 | 3.639 ng/uL | -0.02 | |
| 3) 1,4-Dioxane-d8 | 3.969 | 84 | 9627 | 4.851 ng/ul | 0.00 | 17 |
| 4) Pyridine-d5 | 7.353 | 99 | 57674 | 24.831 ng/ul | 0.00 | |
| 7) Phenol-d5 9) Bis-(2-Chloroethyl)eth | 7.500 | 67 | 35382 | 24.255 ng/ul | -0.02 | 1 |
| 11) 2-Chlorophenol-d4 | 7.717 | 132 | 42143 | 25.197 ng/ul | -0.02 | , , , |
| 15) 4-Methylphenol-d8 | 8.904 | 113 | 42555 | 22.704 ng/ul | 0.00 | |
| 21) Nitrobenzene-d5 | 9.368 | 128 | 22639 | 25.949 ng/ul | 0.00 | |
| 24) 2-Nitrophenol-d4 | 10.091 | 143 | 26296 | 26.720 ng/ul | 0.00 | |
| 28) 2,4-Dichlorophenol-d3 | 10.643 | 165 | 45485 | 27.240 ng/ul | 0.00 0.00 | |
| 31) 4-Chloroaniline-d4 | 11.160 | 131 | 25076 | 10.264 ng/ul | 0.00 | |
| 46) Dimethylphthalate-d6 | 14.216 | 166 | 156149 | 27.570 ng/ul 27.792 ng/ul | 0.00 | |
| 49) Acenaphthylene-d8 | 14.521 | 160 | 198487 | 21.831 ng/ul | 0.00 | |
| 54) 4-Nitrophenol-d4 | 15.056 | 143 | 20014 | 27.967 ng/ul | | |
| 60) Fluorene-d10 | 15.814 | | 142634 | 16.360 ng/ul | | |
| 65) 4,6-Dinitro-2-methylph | 15.949 | 200 | 16899 223331 | 27.895 ng/ul | | |
| 73) Anthracene-d10 | 17.670 | 188 | 251751 | 29.324 ng/ul | | |
| 81) Pyrene-d10 | 19.950 | | 210332 | 28.402 ng/ul | | l de la constante de |
| 92) Benzo(a)pyrene-d12 | 25.038 | 204 | 210332 | | | |
| | | | | | value | |
| Target Compounds | 19.615 | 202 | 29479 | 2.796 ng/ul | | |
| 80) Fluoranthene | 19.979 | | 25327 | 2.455 ng/ul | | |
| 82) Pyrene | 21.854 | | 12659 | 1.315 ng/ul | . 99 | |
| <pre>85) Benzo(a)anthracene 86) Bis(2-ethylhexyl)phtha</pre> | | | 12591 | 2.040 ng/u] | | |
| | 21.918 | - | 15117 • | √ 1.635 ng/u] | 9 | 5 0.1 |
| 87) Chrysene 90) Benzo(b)fluoranthene | 24.180 | | 21020m | 2,246 ng/u | | 01 |
| 91) Benzo(k)fluoranthene | 24.245 | | | 1.077 ng/u | L ~ | 12/2121 |
| 93) Benzo(a)pyrene | 25.103 | 252 | | 1.577 ng/u | L# 9 1 | |
| 94) Indeno(1,2,3-cd)pyrene | 29.169 | 276 | | | L | |
| 96) Benzo(g,h,i)perylene | 30 420 | 5 276 | 9258m | 1.101 ng/u | L | |
| | | | | | | |
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(#) = qualifier out of range (m) = manual integration (+) = signals summed +

SFAM-EPA-BG112321.M Tue Dec 07 01:25:42 2021

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