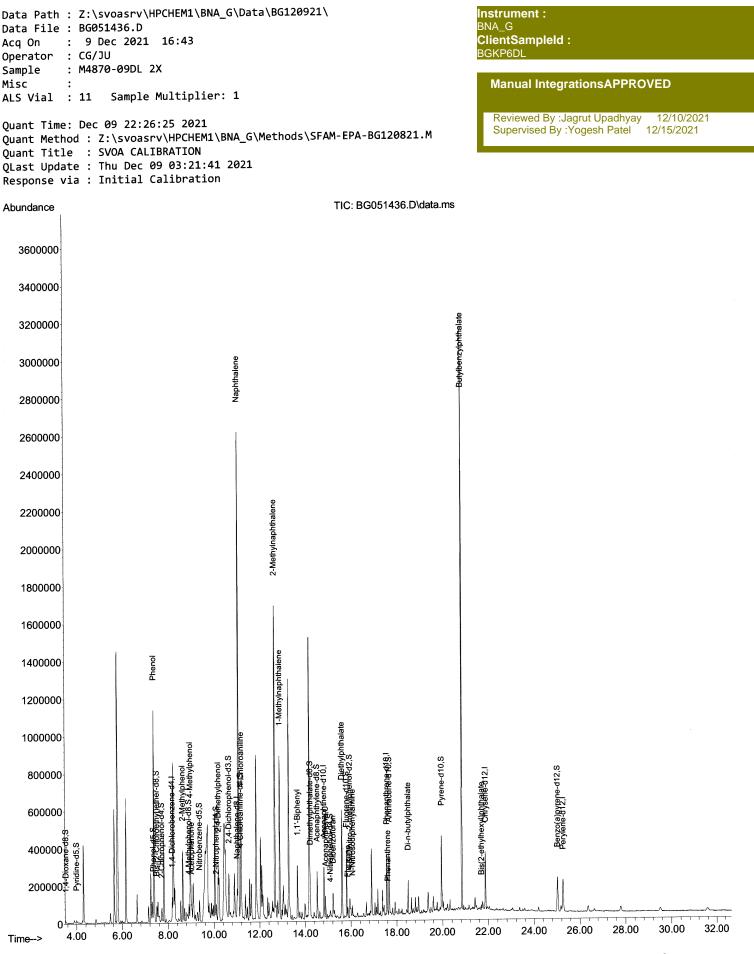
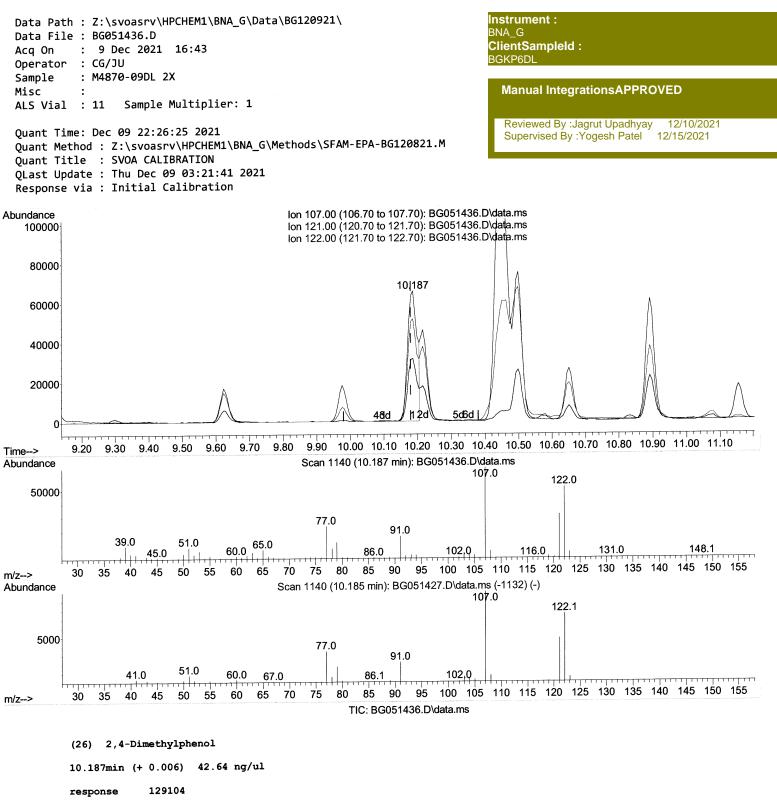
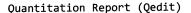
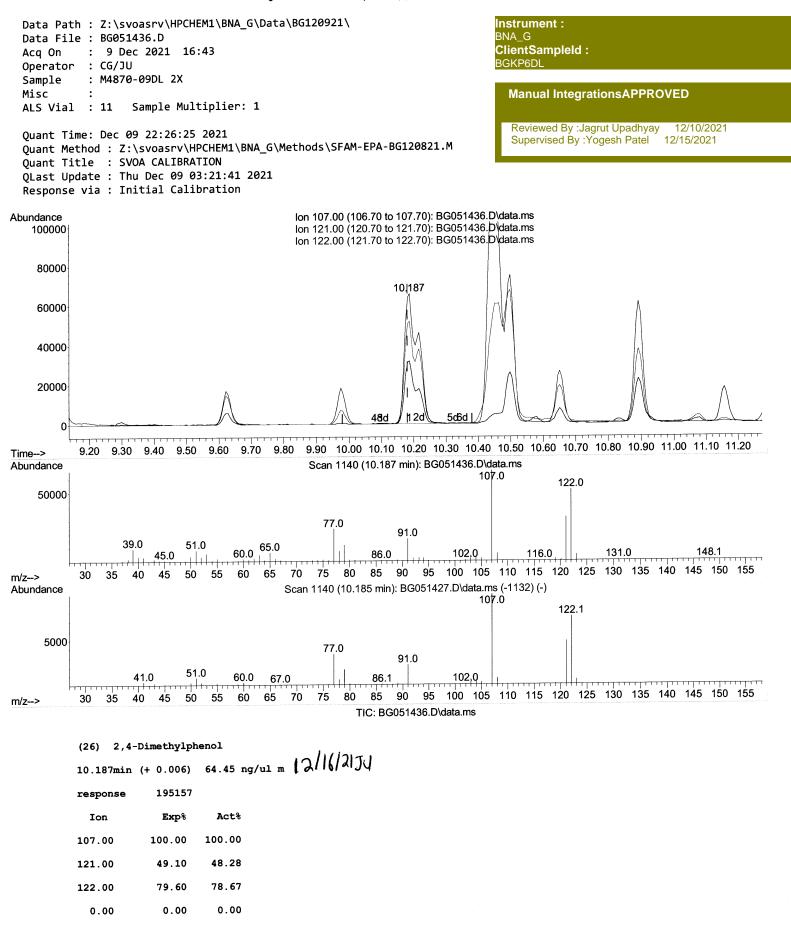
(QT Reviewed)

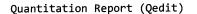


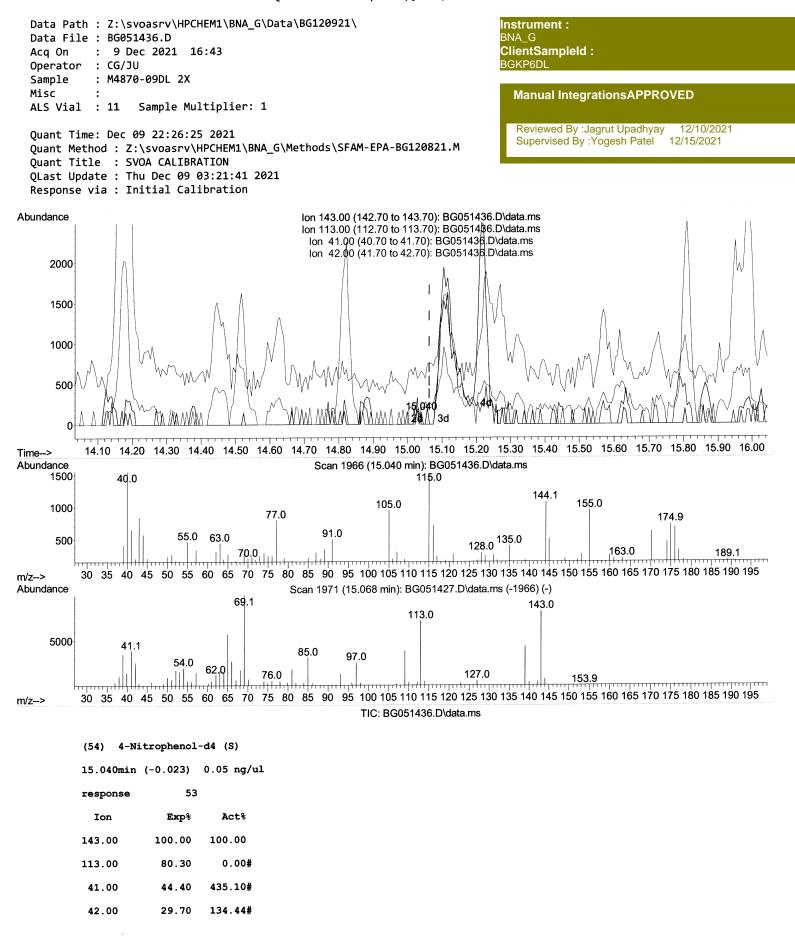


| response | 129104 | | |
|----------|--------|--------|--|
| Ion | Exp% | Act% | |
| 107.00 | 100.00 | 100.00 | |
| 121.00 | 49.10 | 48.28 | |
| 122.00 | 79.60 | 78.67 | |
| 0.00 | 0.00 | 0.00 | |

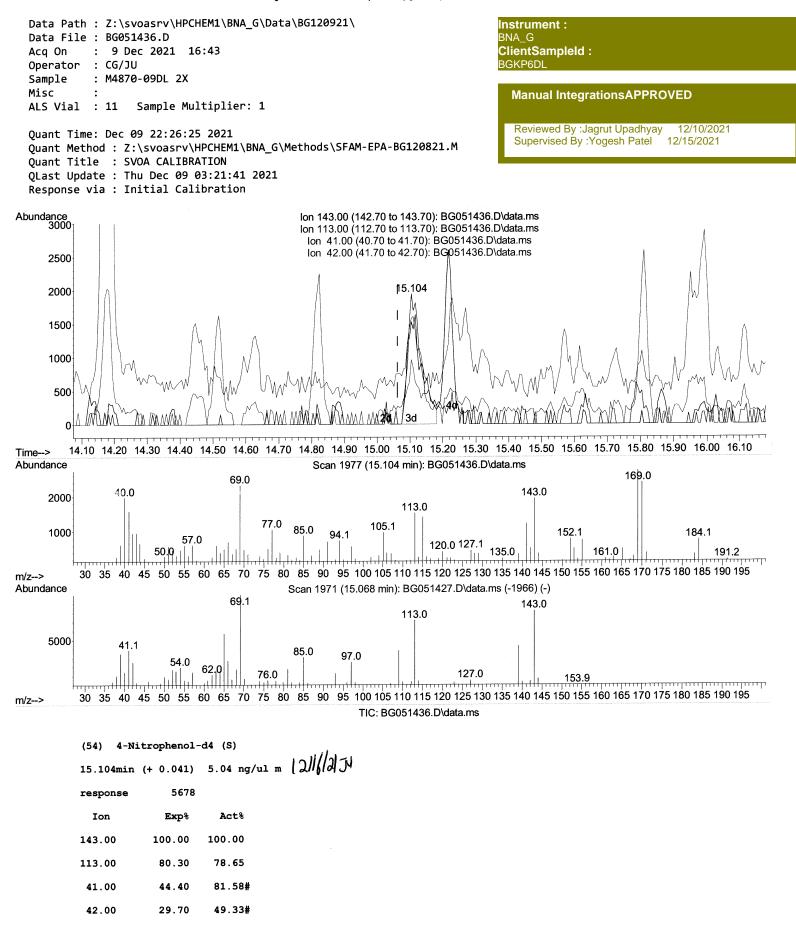












| | | 1001 | 0021 | | | In etrum and a |
|----------------------------------------------|------------------|------------|------------------|--------------------------|----------------------------------------|-----------------------------------------------------------------------------------|
| Data Path : Z:\svoasrv\HPCHEM1\B | NA_G\Data | a/RC17 | 20921/ | | | Instrument : BNA_G |
| Data File : BG051436.D | | | | | | ClientSampleId : |
| Acq On : 9 Dec 2021 16:43 | | | | | | BGKP6DL |
| Operator : CG/JU Sample : M4870-09DL 2X | | | | | | |
| Sample : M4870-09DL 2X Misc : | | | | | | Manual IntegrationsAPPROVED |
| ALS Vial : 11 Sample Multiplic | er: 1 | | | | | Manual mogratione/11 Noveb |
| | | | | | | Deviewed Due loggett Lipselly even 40/40/2024 |
| Quant Time: Dec 09 22:26:25 2021 | | | | | | Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :Yogesh Patel 12/15/2021 |
| Quant Method : Z:\svoasrv\HPCHEM | 1\BNA_G\M | letho | ds\SFAM-EP | A-BG120821.M | М | |
| Quant Title : SVOA CALIBRATION | | | | | | |
| QLast Update : Thu Dec 09 03:21: | 41 2021 | | | | | |
| Response via : Initial Calibrati | on | | | | | |
| Compound | R.T. (| OTon | Response | Conc Units | Dev(Mi | n) |
| Compound | | | | | | |
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 8.183 | 152 | 33529 | 20.000 ng/ | • | 0.00 |
| 20) Naphthalene-d8 | 11.015 | | 145467 | 20.000 ng/ | • | 0.00 |
| 38) Acenaphthene-d10 | 14.822 | | 96663 | 20.000 ng/ | | 0.00 |
| 64) Phenanthrene-d10 | 17.566 | | 208466 | 20.000 ng/ | | 0.00 |
| 79) Chrysene-d12 | 21.873 | | 189228 | 20.000 ng/ | | 0.00 |
| 88) Perylene-d12 | 25.269 | 264 | 187928 | 20.000 ng/ | /ui | 0.00 |
| - · · · · · · · · · · · · · · · · · · · | | | | | | |
| System Monitoring Compounds | 2 520 | 96 | 2232 | 2.186 ng/ | /uL 0 | .00 |
| 3) 1,4-Dioxane-d8 | 3.530 3.994 | 84 | 10585 | 3.610 ng/ | | .03 |
| 4) Pyridine-d5 | 7.360 | 99 | 15240 | 4.465 ng/ | | .00 |
| 7) Phenol-d5 9) Bis-(2-Chloroethyl)eth | 7.507 | 67 | 38324 | 17.508 ng/ | | .00 |
| 11) 2-Chlorophenol-d4 | 7.725 | | 35561 | 14.644 ng/ | | .00 |
| 15) 4-Methylphenol-d8 | 8.912 | | 26995 | 10.067 ng/ | | .00 |
| 21) Nitrobenzene-d5 | 9.370 | | 23773 | 18.839 ng/ | /ul 0 | .00 |
| 24) 2-Nitrophenol-d4 | 10.092 | | 24839 | 17.395 ng/ | | .00 |
| 28) 2,4-Dichlorophenol-d3 | 10.651 | | 40739 | 17.538 ng/ | | .00 |
| 31) 4-Chloroaniline-d4 | 11.162 | | 37000 | 10.890 ng/ | | 0.00 |
| 46) Dimethylphthalate-d6 | 14.217 | 166 | 135128 | 18.066 ng | | . 00 |
| 49) Acenaphthylene-d8 | 14.517 | 160 | 181077 | 19.115 ng | | 2.0412116121JU |
| 54) 4-Nitrophenol-d4 | 15.104 | | | > 5.042 ng | | |
| 60) Fluorene-d10 | 15.809 | | 126068 | 18.934 ng | ,, , , , , , , , , , , , , , , , , , , |).00).00 |
| 65) 4,6-Dinitro-2-methylph | 15.956 | | 23682 | 19.117 ng | |). 00 |
| 73) Anthracene-d10 | 17.672 | | 198338 227854 | 20.334 ng, 20.033 ng, | <i>.</i> |).00 |
| 81) Pyrene-d10 | 19.952 25.028 | | 196094 | 20.230 ng | | 0.00 |
| 92) Benzo(a)pyrene-d12 | 25.020 | 204 | 190094 | 20.200 116 | , | |
| Target Compounds | | | | | Qvalu | le |
| 8) Phenol | 7.390 | 94 | 607261 | 173.816 ng | g/ul | 95 |
| 13) 2-Methylphenol | 8.641 | 108 | 130792 | 50.283 ng | | 95 |
| 16) Acetophenone | 9.023 | 105 | 39009 | 9.395 ng | - | 99 |
| 18) 4-Methylphenol | 8.976 | 108 | 196024 | 71.737 ng | | $\frac{97}{1}$ |
| 26) 2,4-Dimethylphenol | 10.187 | 107 | | > 64.452 ng | | [A116121 JU |
| 30) Naphthalene | 11.074 | 128 | 2111723 | 264.348 ng | | 95 |
| 32) 4-Chloroaniline | 11.185 | 127 | 220453 | 64.502 ng | - | 99 100 |
| 36) 2-Methylnaphthalene | 12.660 | 142 | 804474 | 151.015 ng | | 97 |
| 37) 1-Methylnaphthalene | 12.877 | 142 | 369495 | 67.386 ng | - | 98 |
| 43) 1,1'-Biphenyl | 13.653 | 154 | 123925 | 17.148 ng 8.126 ng | _ | 96 |
| 52) Acenaphthene | 14.881 | 153 168 | 49842 70696 | 8.120 ng 8.130 ng | - | 98 |
| 56) Dibenzofuran | 15.216 15.610 | 168 | 317056 | 38.970 ng | | 99 |
| 59) Diethylphthalate | 15.862 | 149 | 25370 | 3.603 ng | - | 92 |
| 61) Fluorene 67) N-Nitrosodiphenylamine | 15.862 | 169 | 23858 | 4.105 ng | | 99 |
| 72) Phenanthrene | 17.613 | 178 | 13874 | 1.235 ng | | 98 |
| 72) Phenantin ene 78) Di-n-butylphthalate | 18.500 | 149 | 132935 | 9.907 ng | | 99 |
| 83) Butylbenzylphthalate | 20.845 | 149 | 920479 | 153.573 ng | | 98 |
| 86) Bis(2-ethylhexyl)phtha | 21.702 | | 12316 | 1.479 ng | g/ul | 96 |
| | | | | | | |
| | | | | | | |

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

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