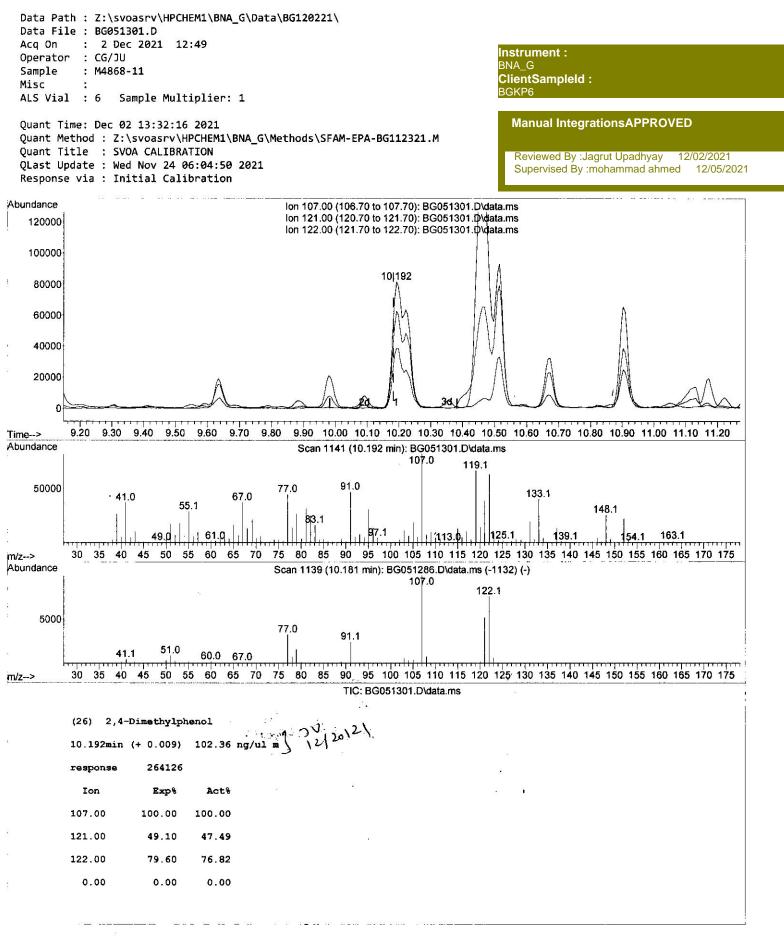


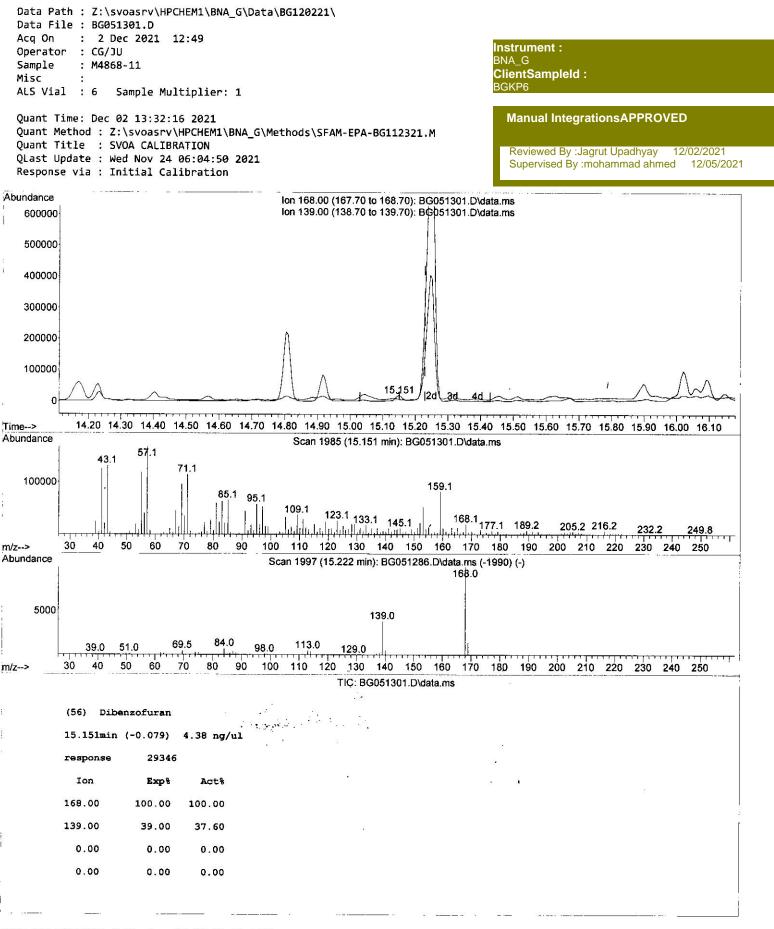
SFAM-EPA-BG112321.M Thu Dec 02 13:34:53 2021

Page: 1



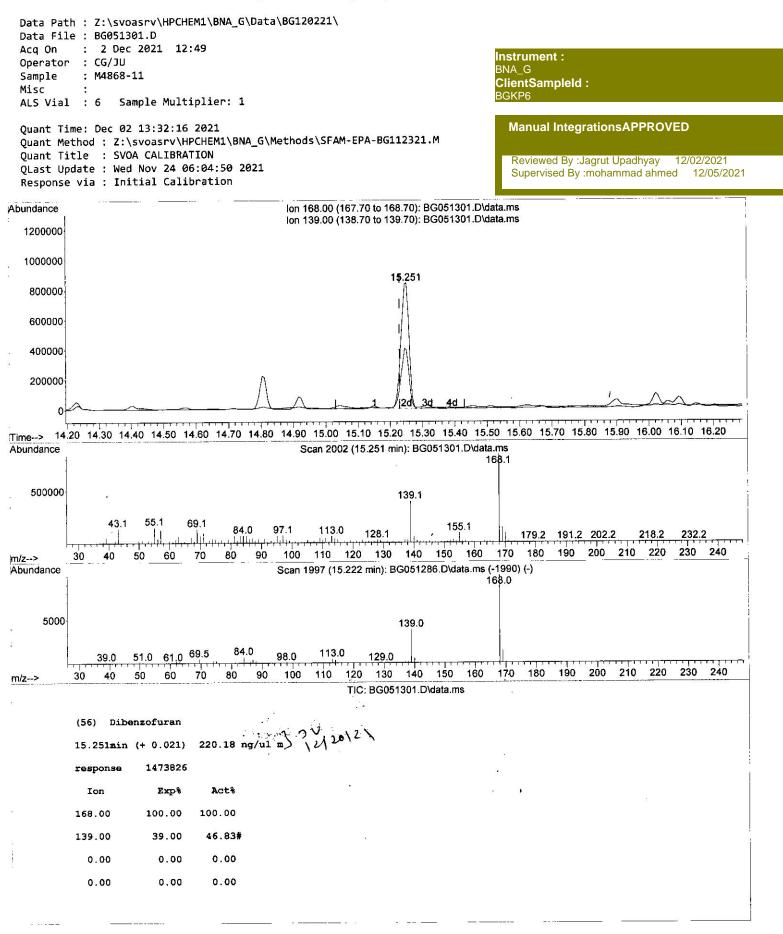
SFAM-EPA-BG112321.M Thu Dec 02 13:35:09 2021

Page: 1



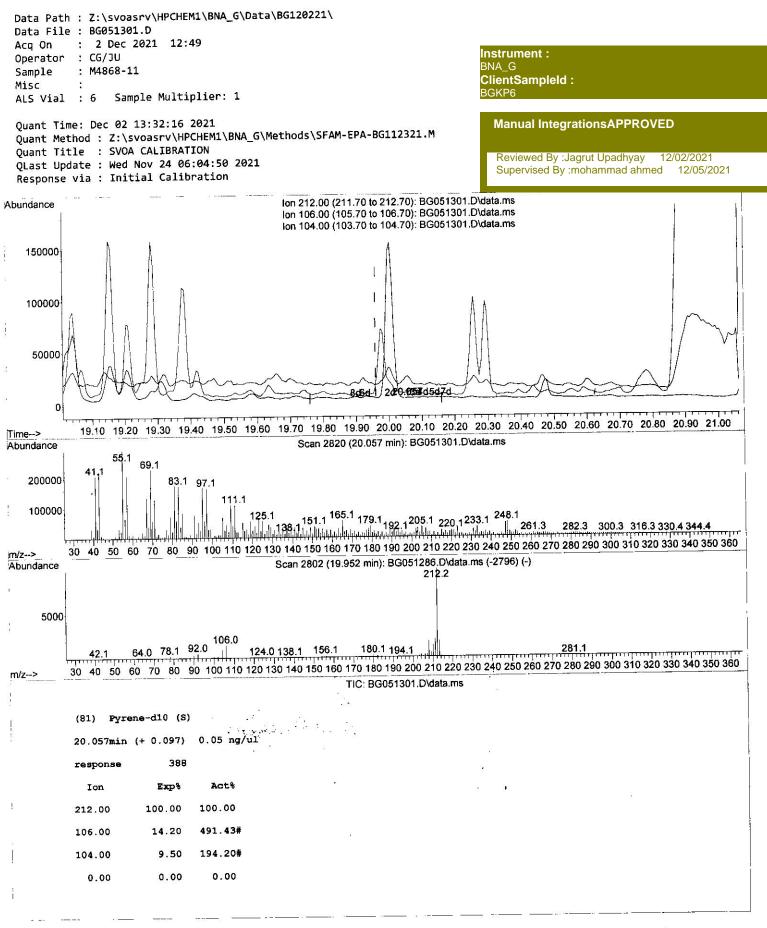
SFAM-EPA-BG112321.M Thu Dec 02 13:38:51 2021

Page: 1

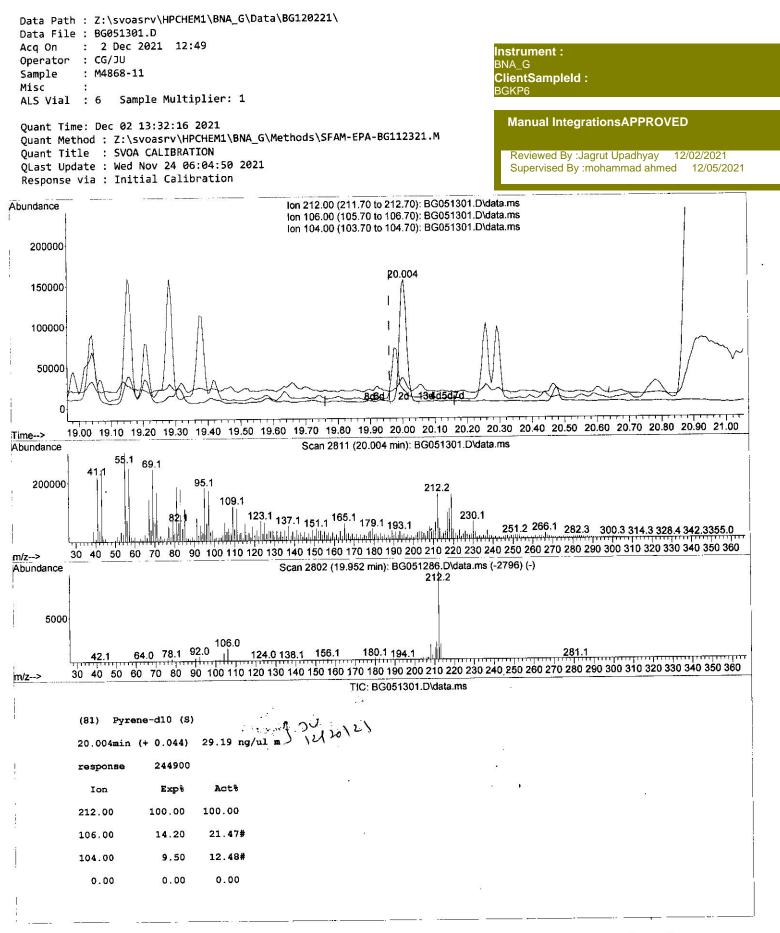


SFAM-EPA-BG112321.M Thu Dec 02 13:39:04 2021

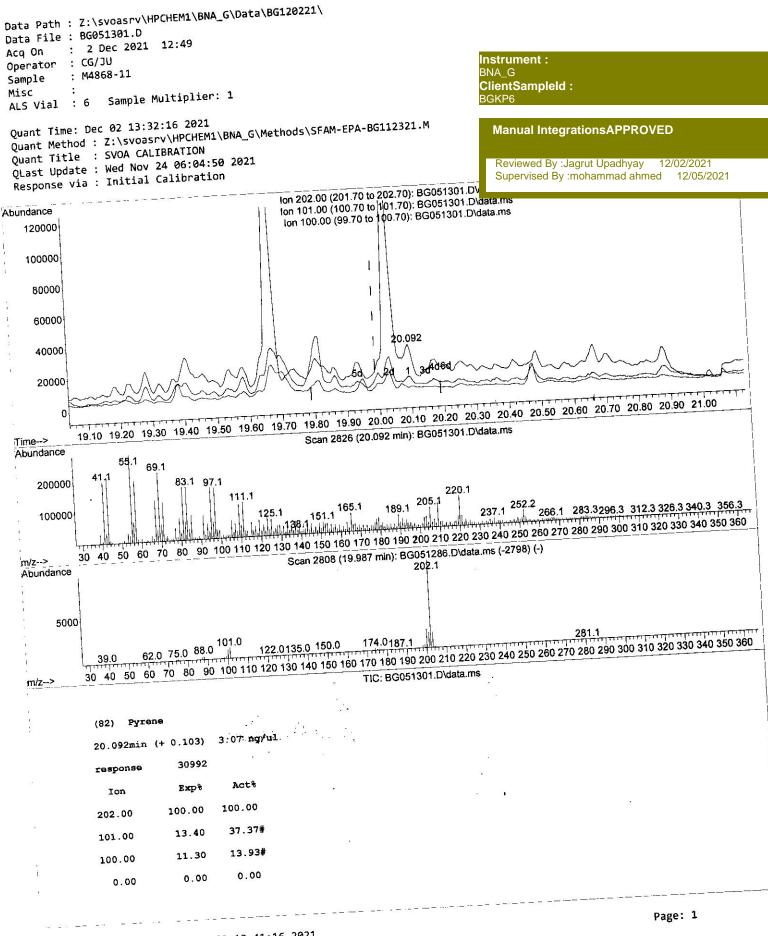
Page: 1



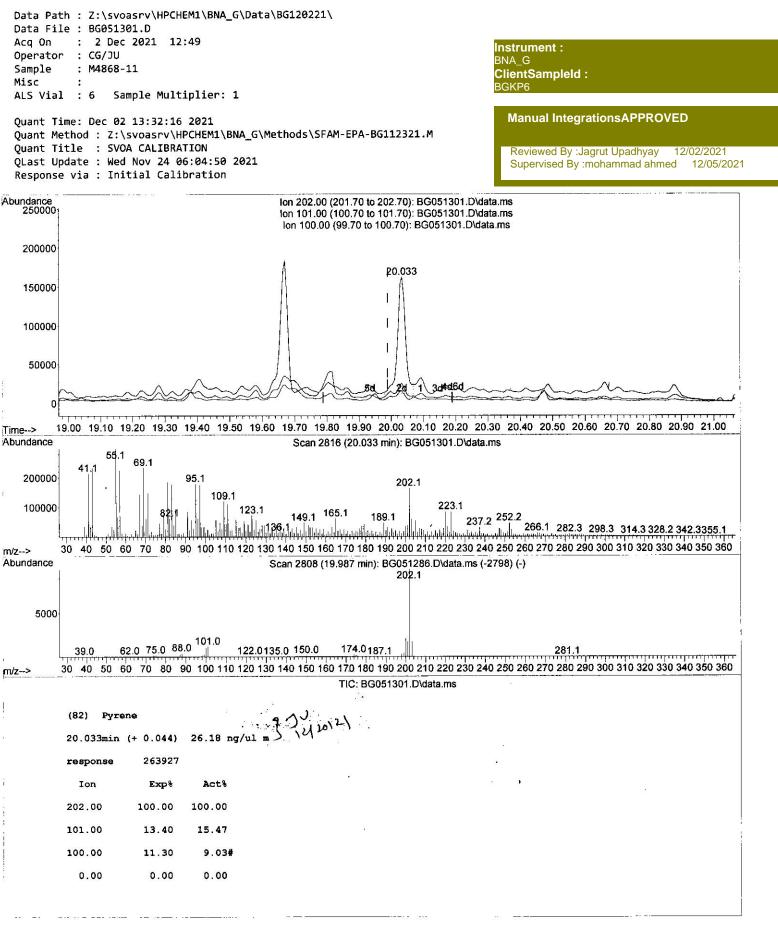
SFAM-EPA-BG112321.M Thu Dec 02 13:40:51 2021



SFAM-EPA-BG112321.M Thu Dec 02 13:41:07 2021



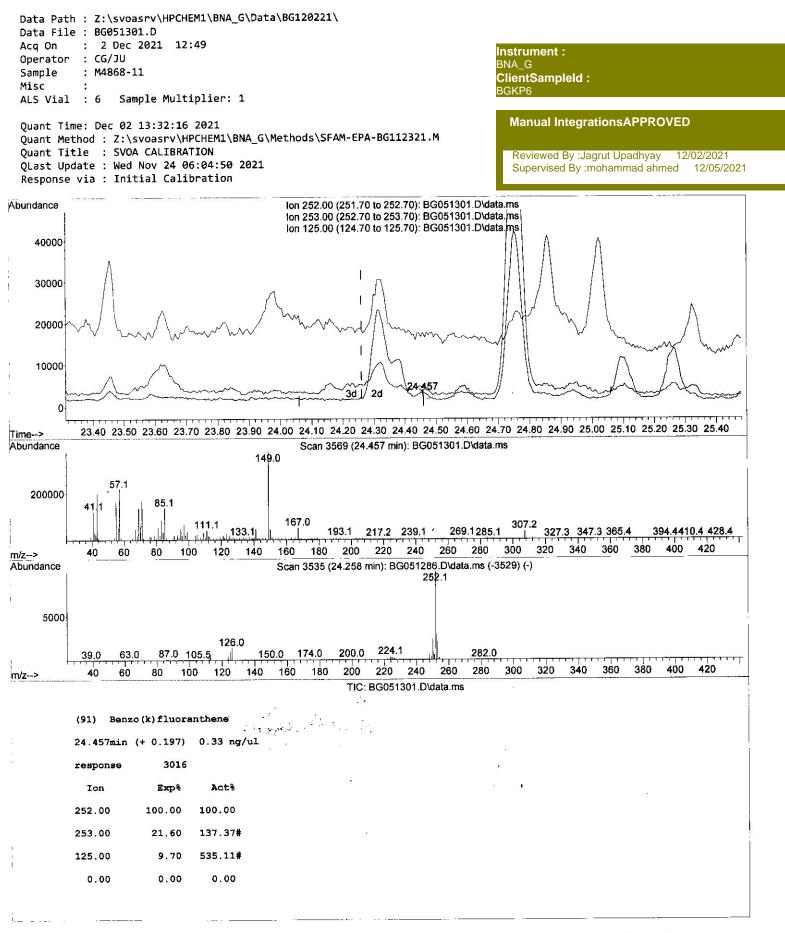
SFAM-EPA-BG112321.M Thu Dec 02 13:41:16 2021



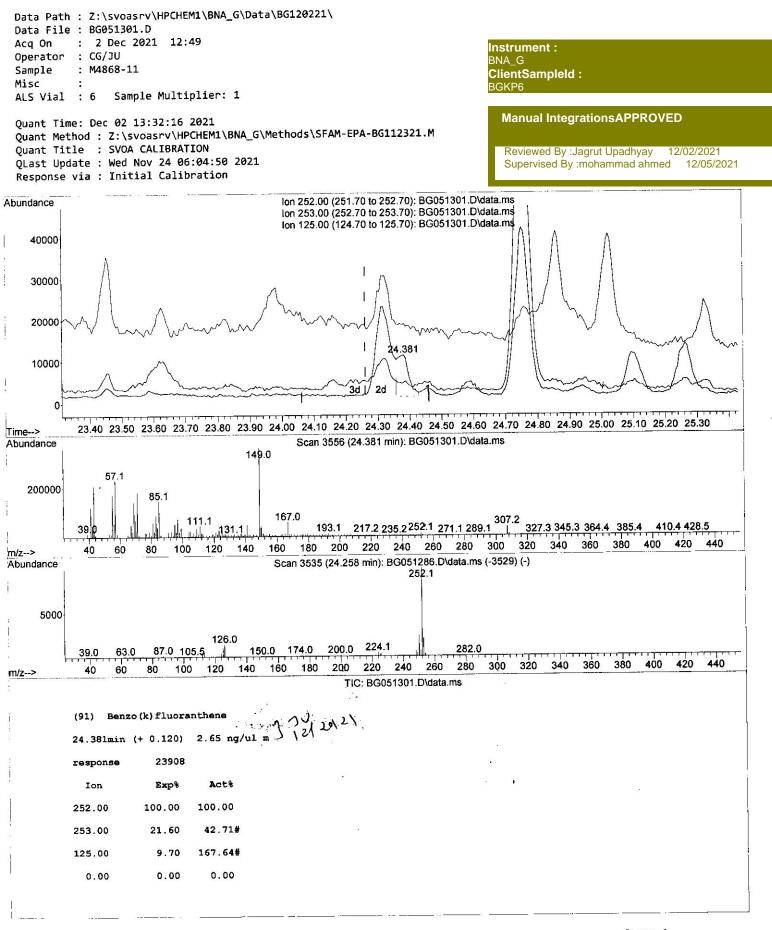
SFAM-EPA-BG112321.M Thu Dec 02 13:41:38 2021

Page: 1

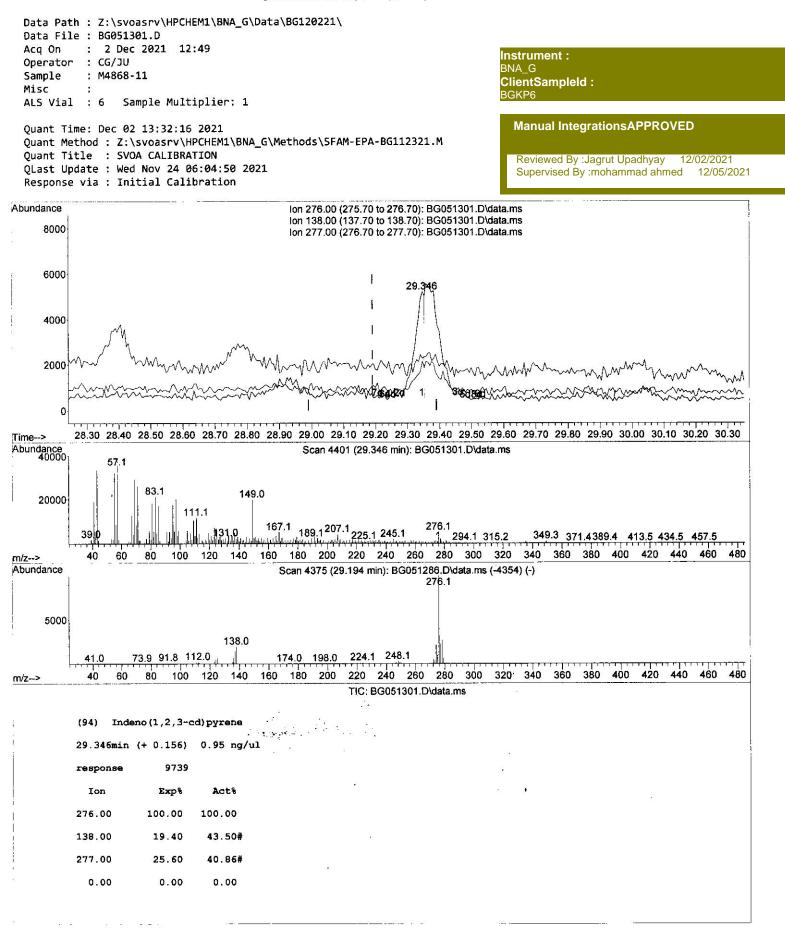
٩.



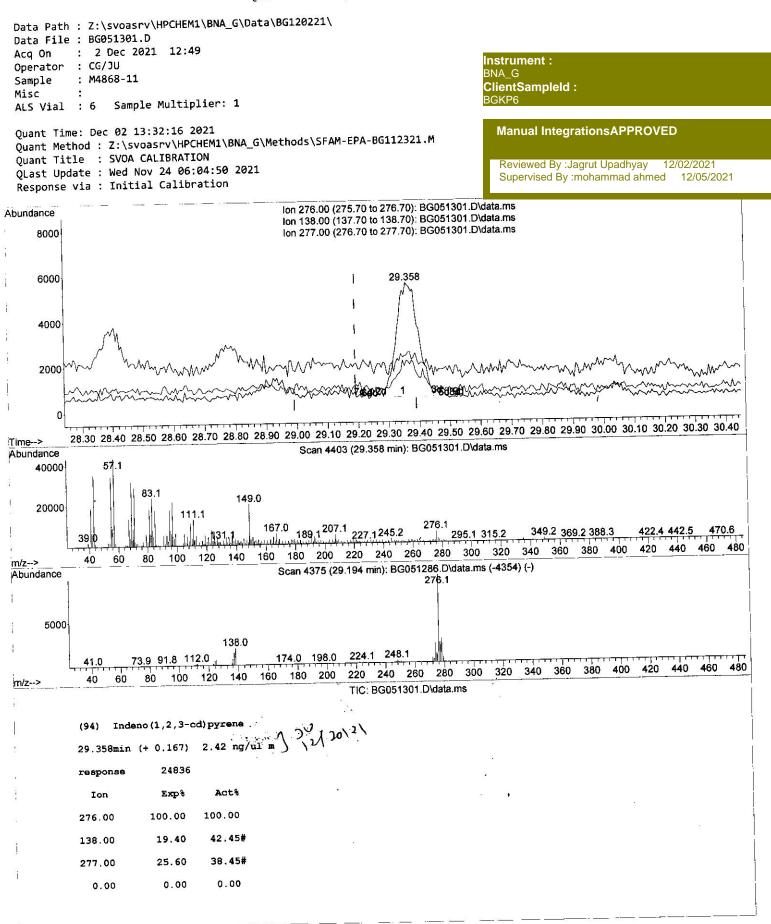
SFAM-EPA-BG112321.M Thu Dec 02 13:45:30 2021



SFAM-EPA-BG112321.M Thu Dec 02 13:46:05 2021

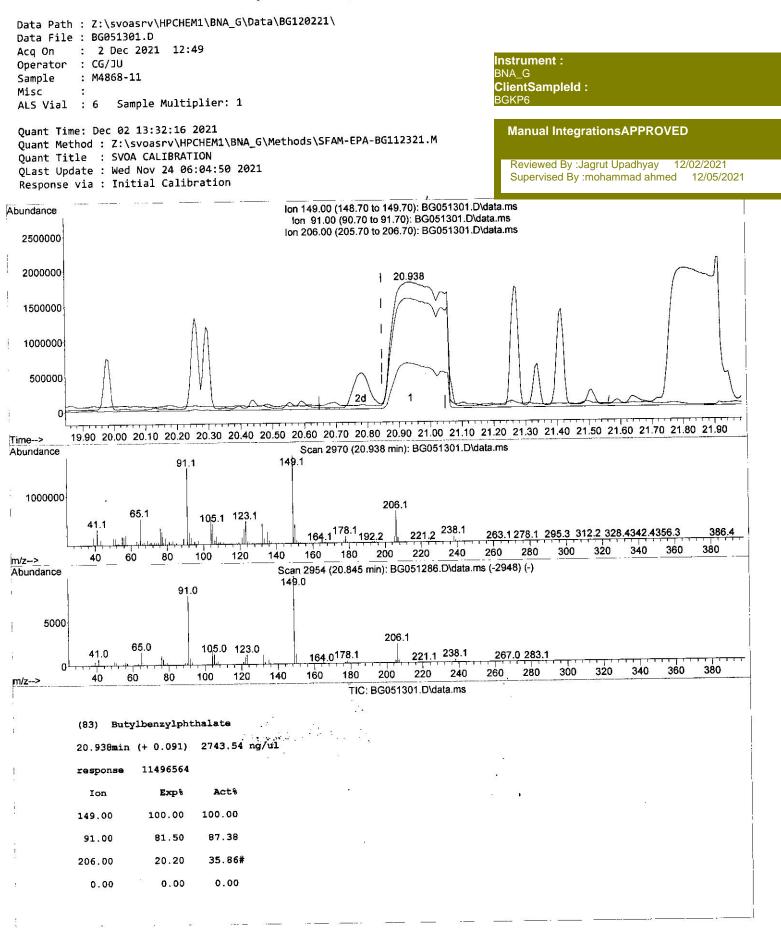


SFAM-EPA-BG112321.M Thu Dec 02 13:46:24 2021

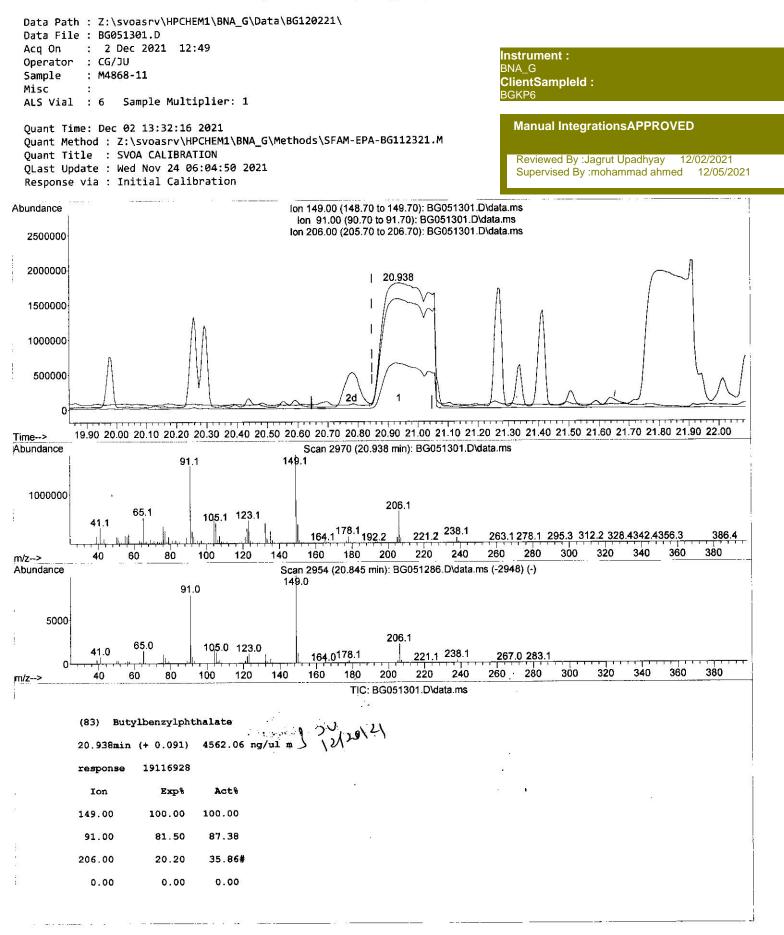


SFAM-EPA-BG112321.M Thu Dec 02 13:46:47 2021

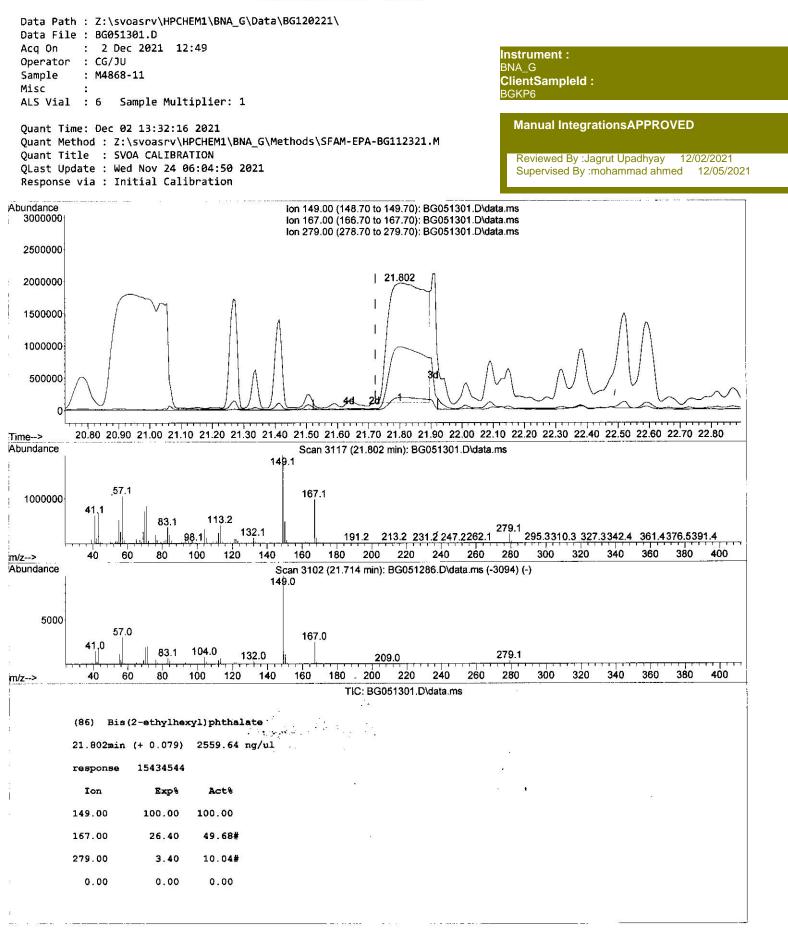




SFAM-EPA-BG112321.M Thu Dec 02 13:56:19 2021

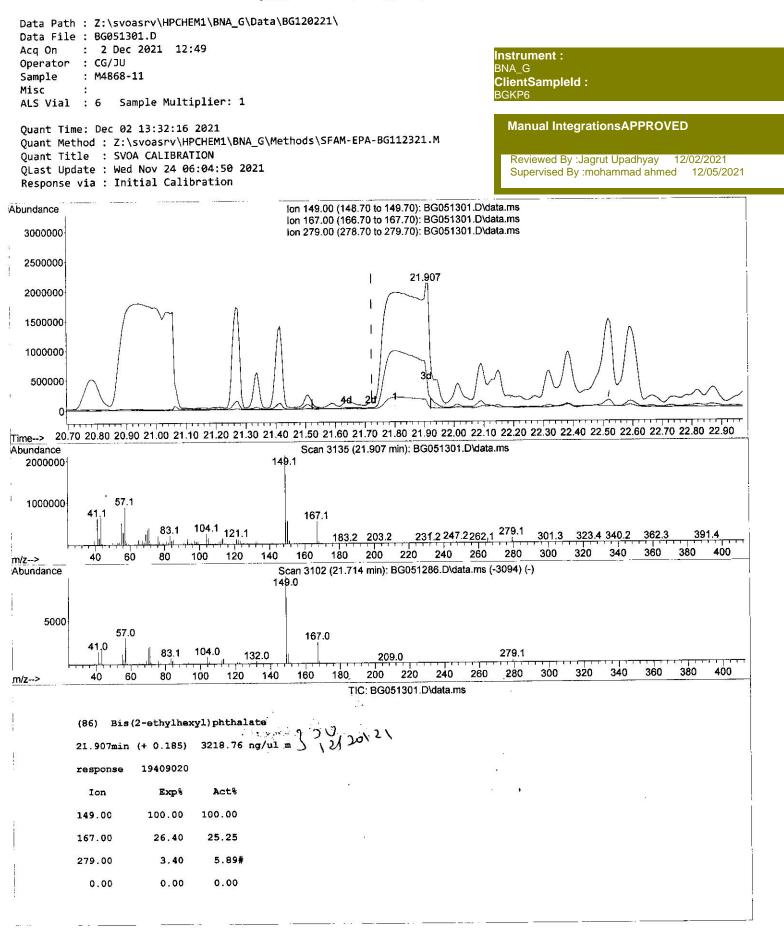


SFAM-EPA-BG112321.M Thu Dec 02 13:56:58 2021



SFAM-EPA-BG112321.M Thu Dec 02 13:57:09 2021

Page: 1



SFAM-EPA-BG112321.M Thu Dec 02 13:57:53 2021

Page: 1

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\ Data File : BG051301.D Acg On : 2 Dec 2021 12:49 Operator : CG/JU BNA_G : M4868-11 Sample Misc BGKP6 ALS Vial : 6 Sample Multiplier: 1 Quant Time: Dec 02 13:32:16 2021 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Compound _____ Internal Standards 20.000 ng/ul 0.00 8,200 152 33082 1) 1,4-Dichlorobenzene-d4 20.000 ng/ul # 0.01 127954 11.044 136 20) Naphthalene-d8 0.01 14.845 164 73407 20.000 ng/ul 38) Acenaphthene-d10 # 0.03 20.000 ng/ul 64) Phenanthrene-d10 17.613 188 145035 20.000 ng/ul # 0.09 21.966 240 138652 79) Chrysene-d12 0.14 25.427 264 142380 20.000 ng/ul 88) Perylene-d12 System Monitoring Compounds 0.00 3832 4.025 ng/uL 3.541 96 3) 1,4-Dioxane-d8 0.00 45152 16.163 ng/ul 3.976 84 4) Pyridine-d5 25.057 ng/ul 0.01 99 81926 7.372 7) Phenol-d5 0.00 7.524 67 59007 28.735 ng/ul 9) Bis-(2-Chloroethyl)eth... 0.00 23.843 ng/ul 11) 2-Chlorophenol-d4 56138 7.736 132 24.752 ng/ul 0.00 8.911 113 65308 15) 4-Methylphenol-d8 0.00 34.550 ng/ul 9.381 128 37318 21) Nitrobenzene-d5 27.722 ng/ul 0.00 10.104 143 33777 24) 2-Nitrophenol-d4 27.555 ng/ul 0.00 28) 2,4-Dichlorophenol-d3 10.656 165 56963 17.658 ng/ul 0.00 11.173 131 53413 31) 4-Chloroaniline-d4 27.704 ng/ul 0.00 156478 46) Dimethylphthalate-d6 14.228 166 208284 29.244 ng/ul 0,00 14.540 160 49) Acenaphthylene-d8 30.638 ng/ul 0.01 54) 4-Nitrophenol-d4 15.063 143 28011 0.01 26.045 ng/ul 132473 15.838 176 60) Fluorene-d10 0.03 50 12/20/21 14.769 ng/ul 65) 4,6-Dinitro-2-methylph... 15.985 200 13218 28.441 ng/ul 0.03 197282 17.712 188 73) Anthracene-d10 1 29.191 ng/ul 0.04 212 244900m 20.004 81) Pyrene-d10 210813 0.14 27.724 ng/ul 264 92) Benzo(a)pyrene-d12 25.186 Ovalue Target Compounds 623.448 ng/ul# 82 7.413 94 2111700 8) Phenol 98 214071 84.849 ng/ul 108 13) 2-Methylphenol 8.647 95 364496 ' 135.107 ng/ul 8.981 108 18) 4-Methylphenol 264126m / 102.365 ng/ul 10.192 107 26) 2,4-Dimethylphenol 67 /1032.579 ng/ul# 11.126 128 7189063 30) Naphthalene 58.636 ng/ul 99 178058 32) 4-Chloroaniline 11.197 127 92 1542.546 ng/ul 142 7304919 12.736 36) 2-Methylnaphthalene -784.921 ng/ul# 98 3824199 37) 1-Methylnaphthalene 12.936 142 357.689 ng/ul 93 13.682 154 1961127 43) 1,1'-Biphenyl 8.834 ng/ul# 21 14.569 152 62164 50) Acenaphthylene 14.910 153 206.732 ng/ul 94 959385 12/20121 52) Acenaphthene 220.181 ng/ul 15.251 168 1473826m 56) Dibenzofuran 90 15.638 149 315899 52.640 ng/ul# 59) Diethylphthalate 125.857 ng/ul# 97 166 674806 61) Fluorene 15.897 90 116.360 ng/ul# 178 931809 72) Phenanthrene 17.660 8.838 ng/ul# 52 70286 17.748 178 74) Anthracene 8.192 ng/ul# 20 18.012 167 57187 77) Carbazole 87 278.321 ng/ul# 78) Di-n-butylphthalate 18.547 149 2505250 289404 28.086 ng/ul# 84 202 19.669 80) Fluoranthene 20,033 202 263927m 26.184 ng/ul 82) Pyrene 149 19116928m 4562.060 ng/ul 20.938 83) Butylbenzylphthalate 1 228 87773 9.333 ng/ul# 85) Benzo(a)anthracene 21.943 149 19409020m/3218.764 ng/ul 21.907 86) Bis(2-ethylhexyl)phtha... 91124 75 10.086 ng/ul# 22.013 228 87) Chrysene

SFAM-EPA-BG112321.M Wed Dec 15 03:53:03 2021

Instrument :

ClientSampleId :

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/02/2021 Supervised By :mohammad ahmed 12/05/2021

٩

۰.

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\ Data File : BG051301.D Acq On : 2 Dec 2021 12:49 Operator : CG/JU Sample : M4868-11 Misc : ALS Vial : 6 Sample Multiplier: 1

Quant Time: Dec 02 13:32:16 2021 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration

Compound R.T. QIon Response Conc Units Dev(Min) 89) Di-n-octyl phthalate 23.094 149 10270258 995.667 ng/ul 100 7.237 ng/ul# 90) Benzo(b)fluoranthene 24.311 252 69541 () 1 91) Benzo(k)fluoranthene 24.381 252 23908m/ 2.651 ng/ul 93) Benzo(a)pyrene 25.262 252 36461 3.977 ng/ul# 1 29.358 276 24836m) 94) Indeno(1,2,3-cd)pyrene 2.421 ng/ul 1.751 ng/ul# 96) Benzo(g,h,i)perylene 30.609 276 15115 32 -----

Instrument : BNA_G

ClientSampleId : BGKP6

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/02/2021 Supervised By :mohammad ahmed 12/05/2021

1

2

٩

12/2012

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