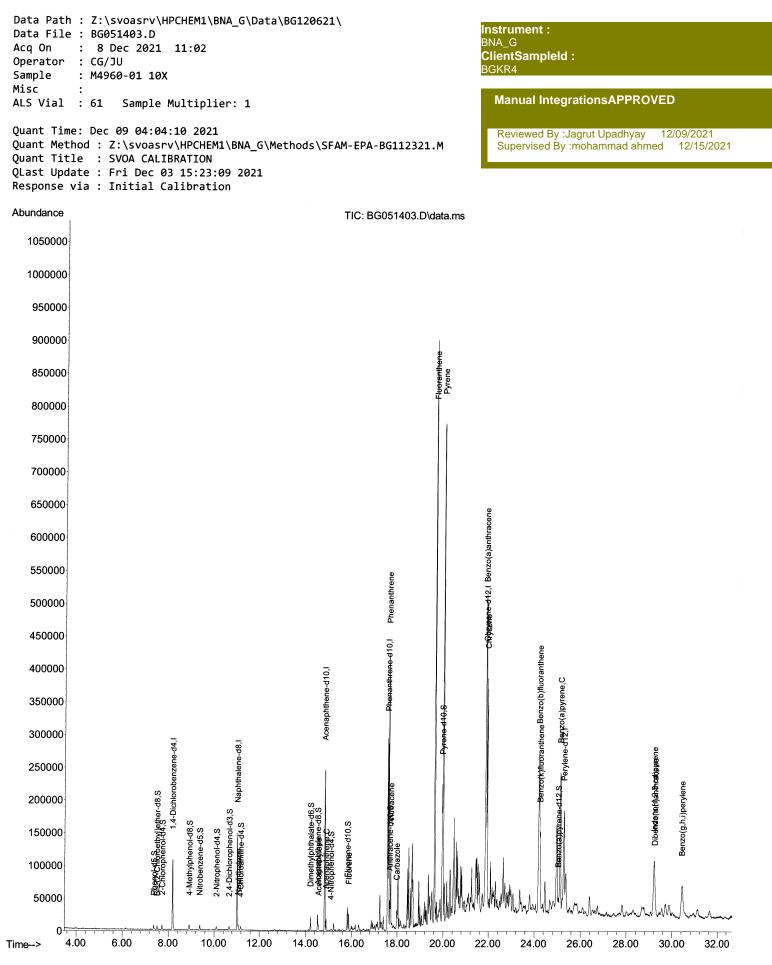
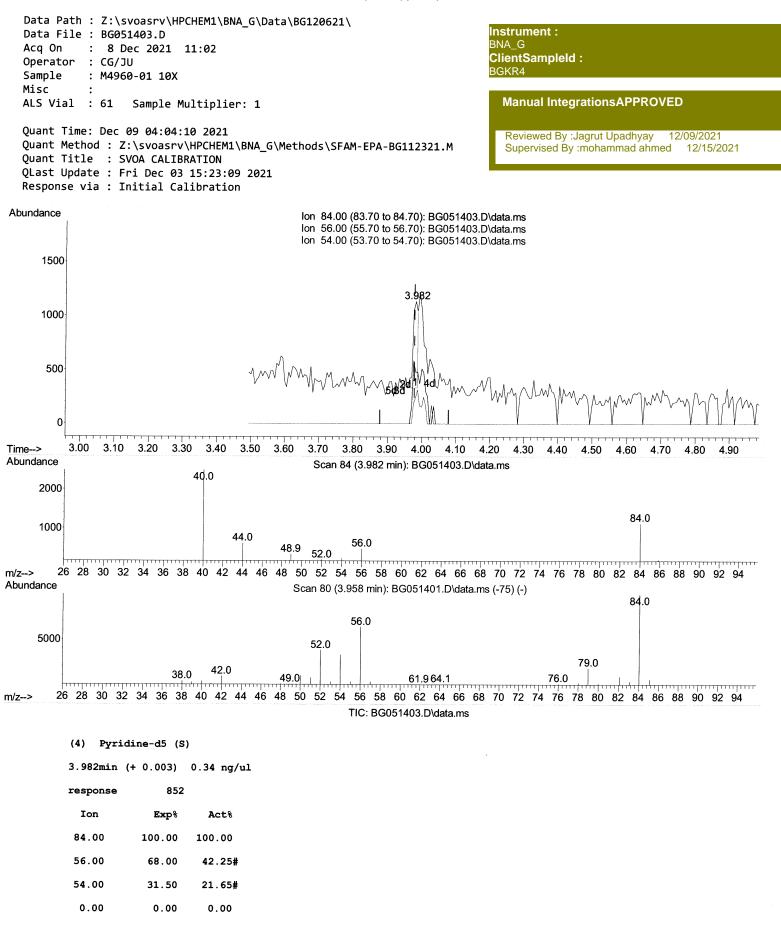
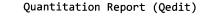
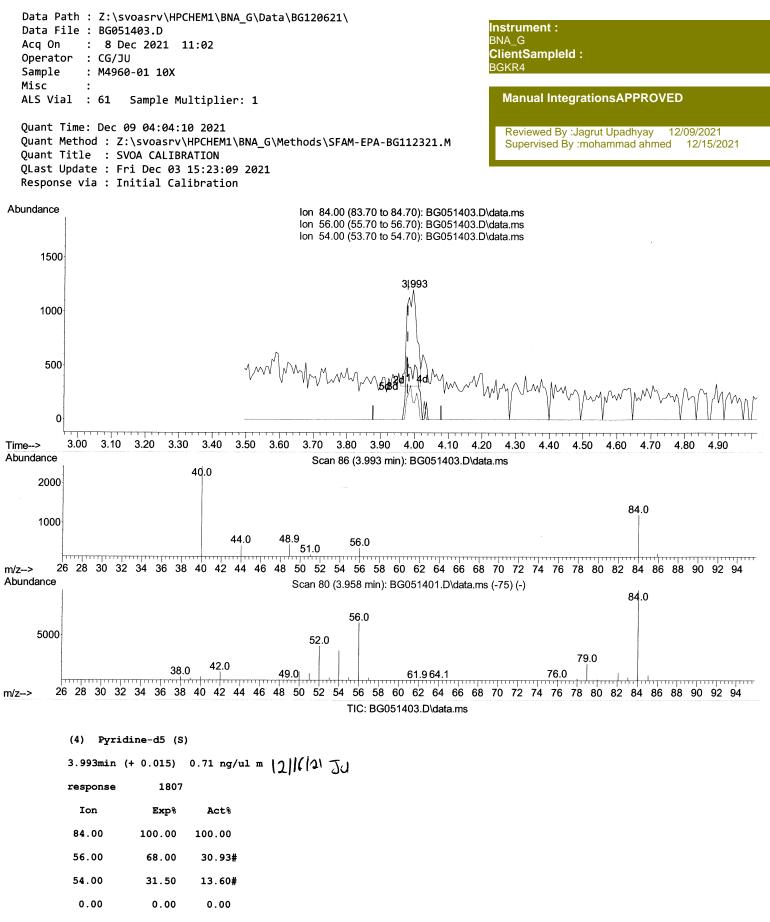
(QT Reviewed)



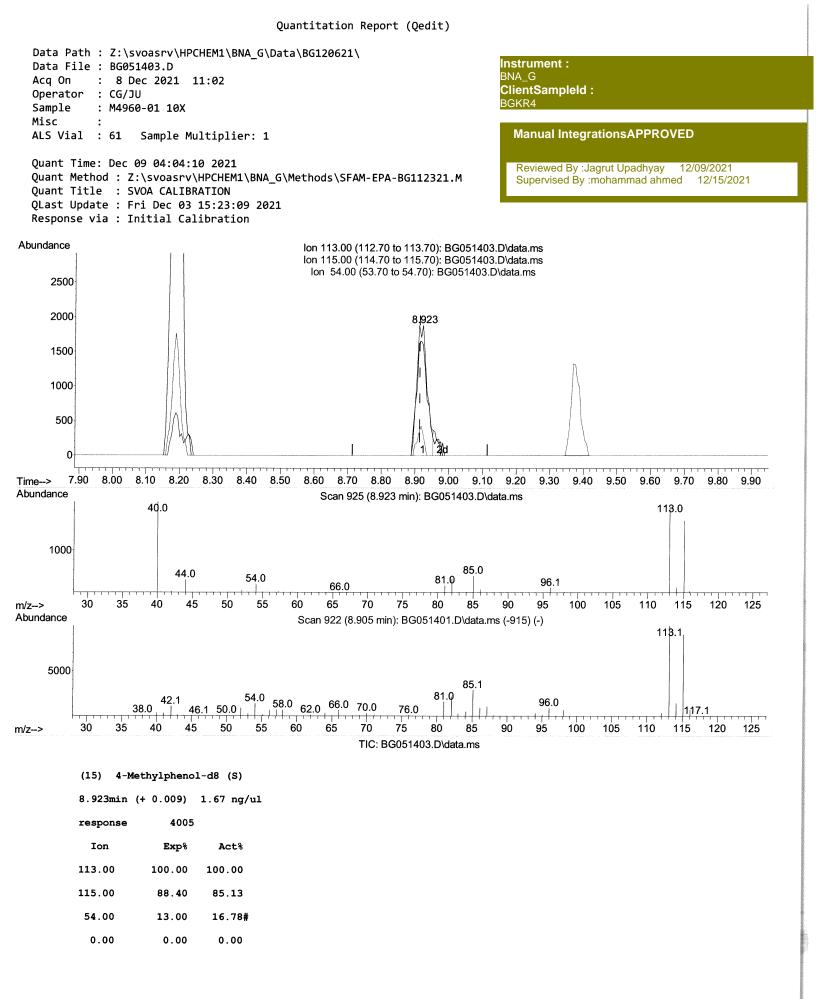


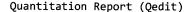


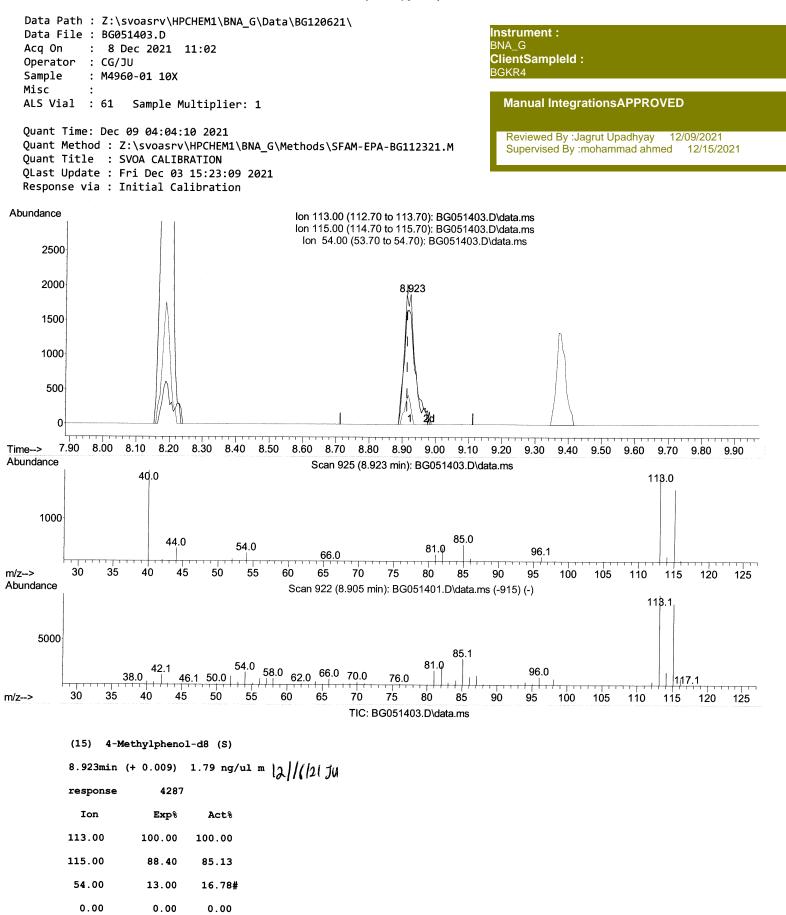


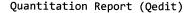


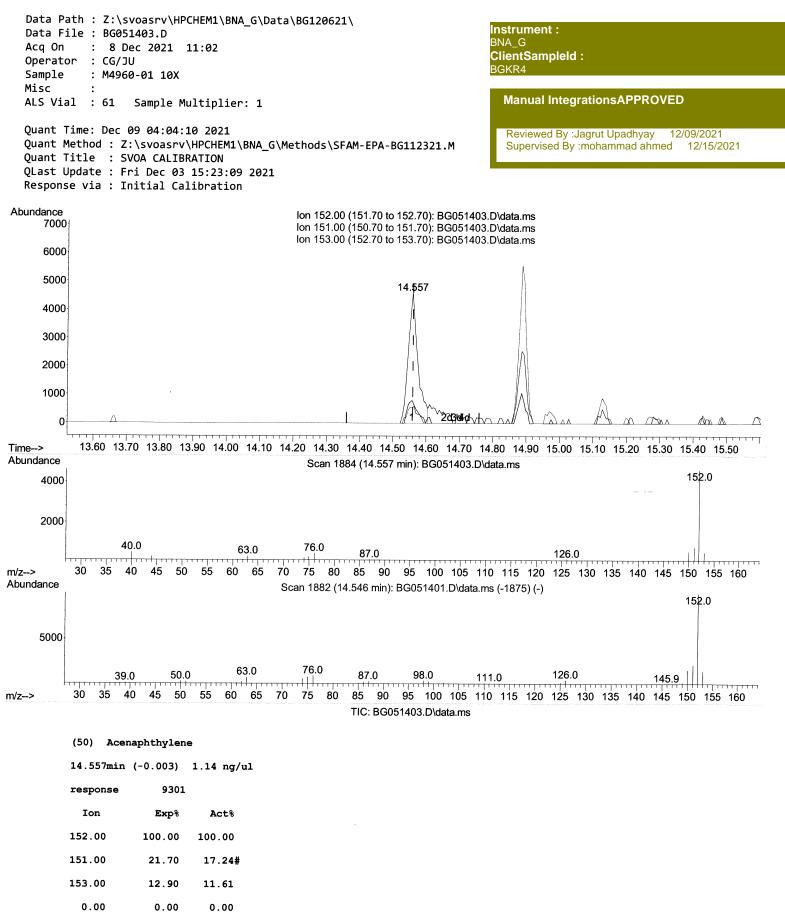
SFAM-EPA-BG112321.M Thu Dec 09 04:21:42 2021

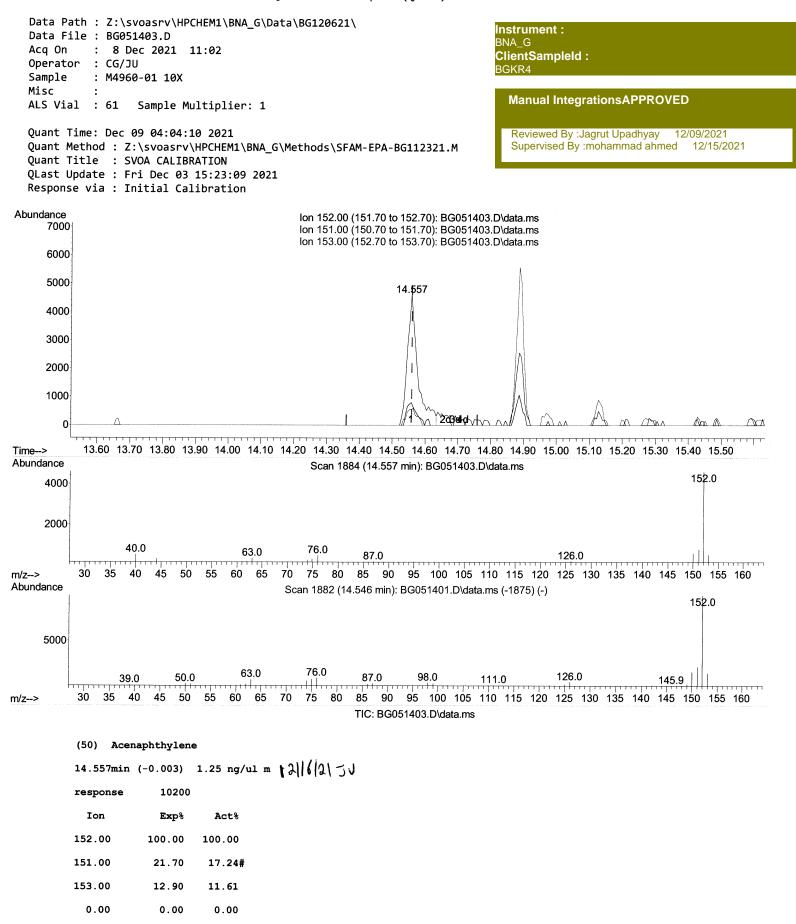




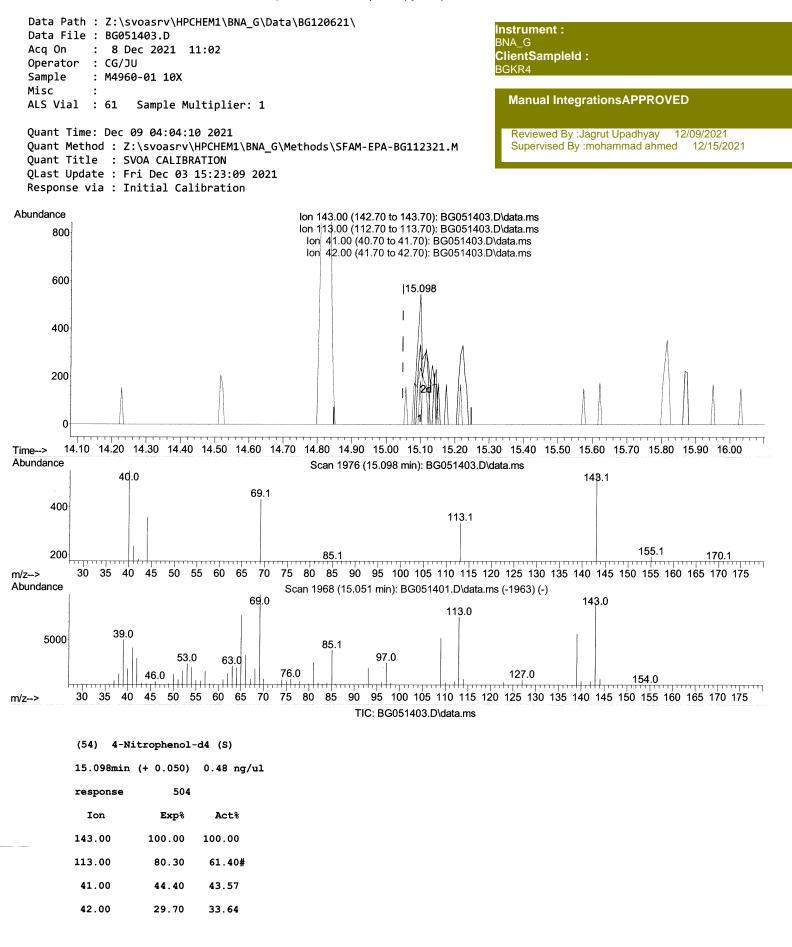




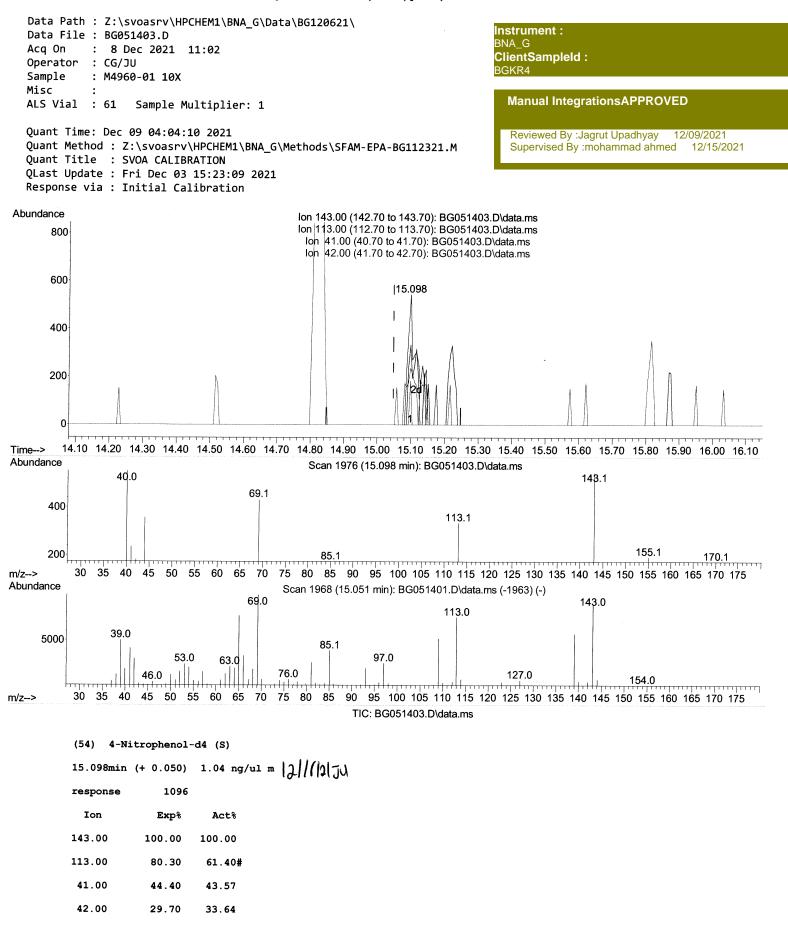




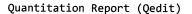


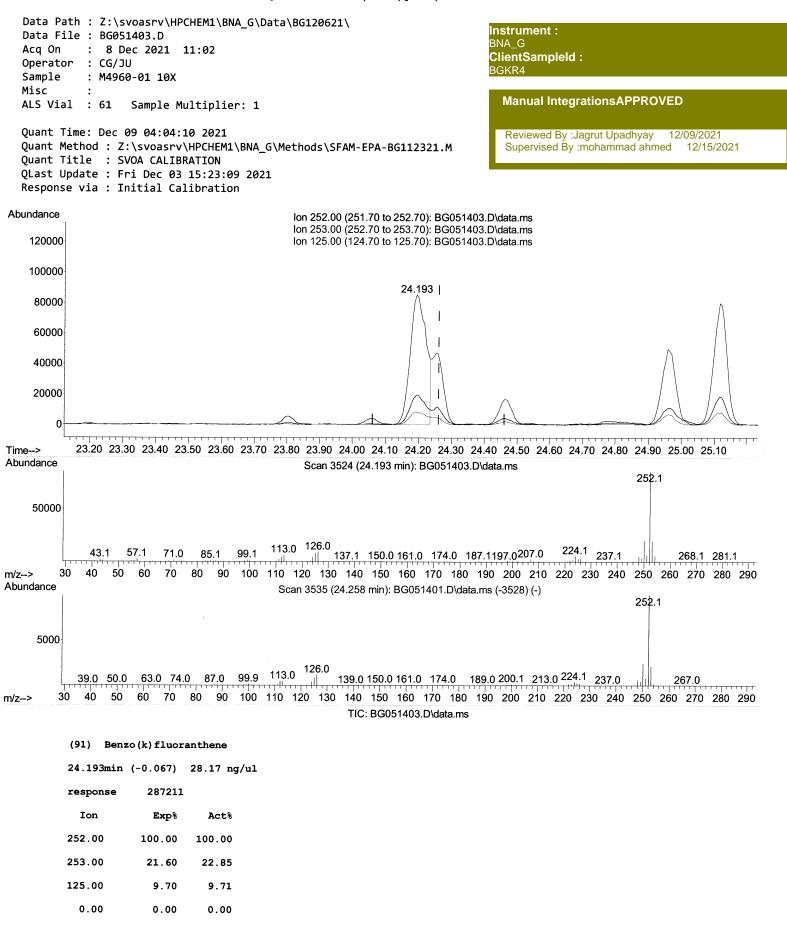




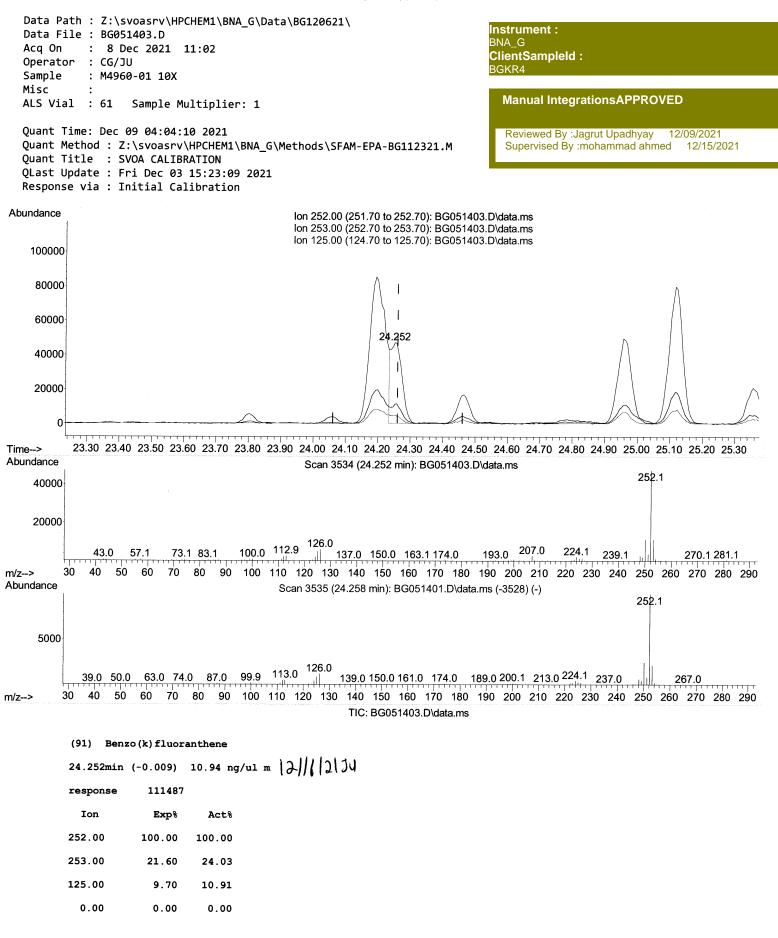


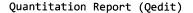
SFAM-EPA-BG112321.M Thu Dec 09 04:23:14 2021

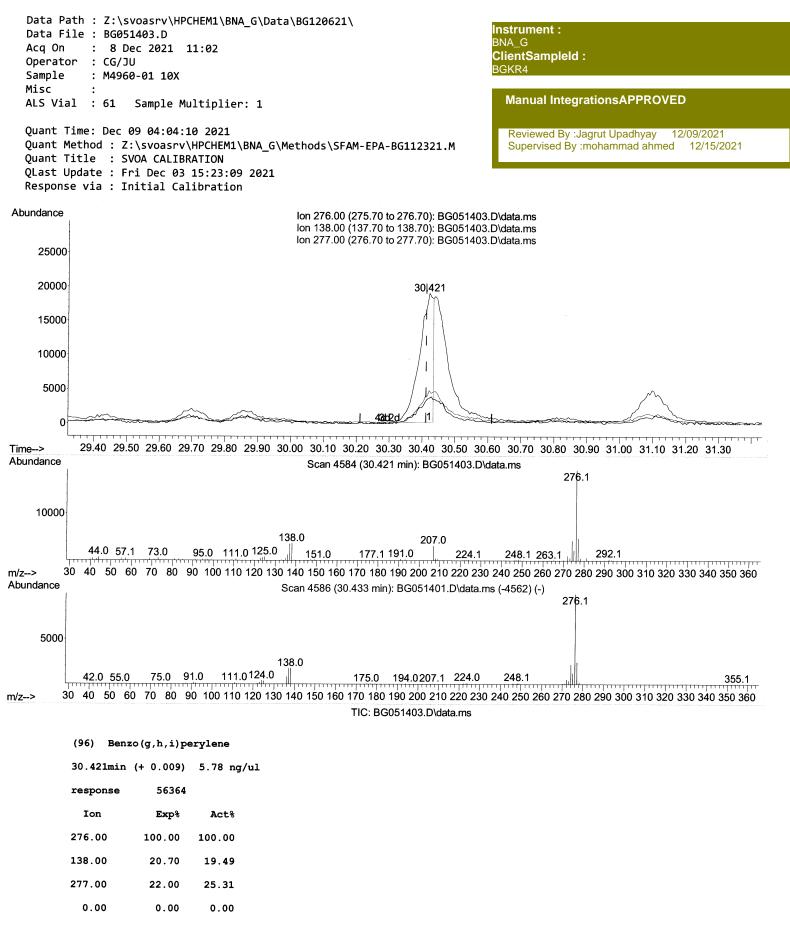




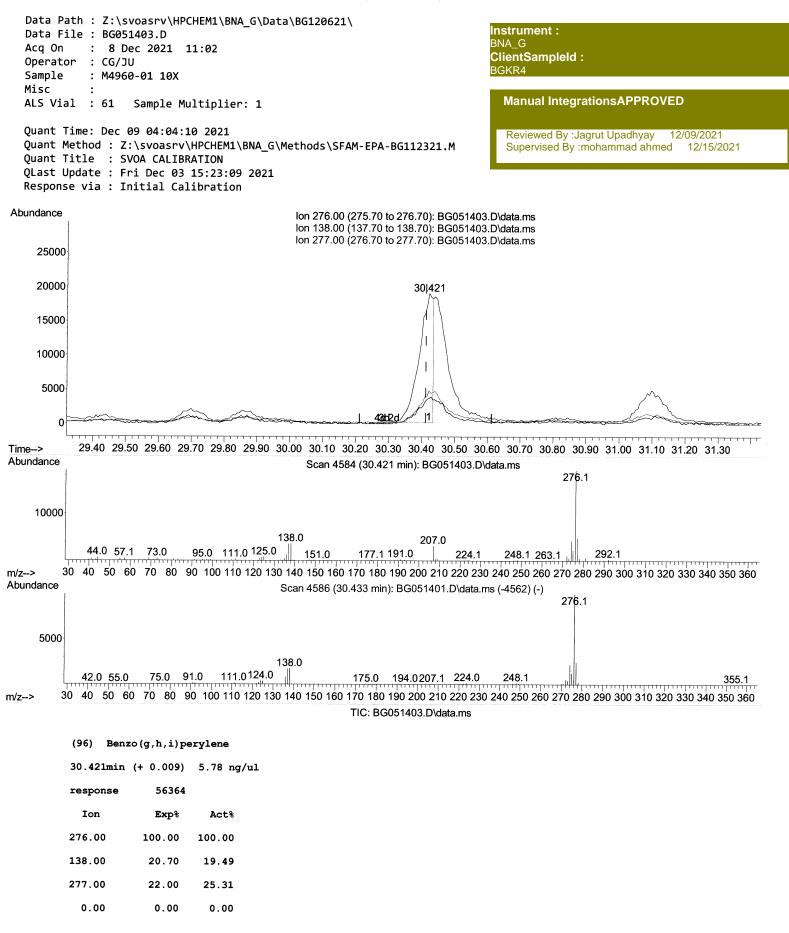


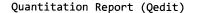


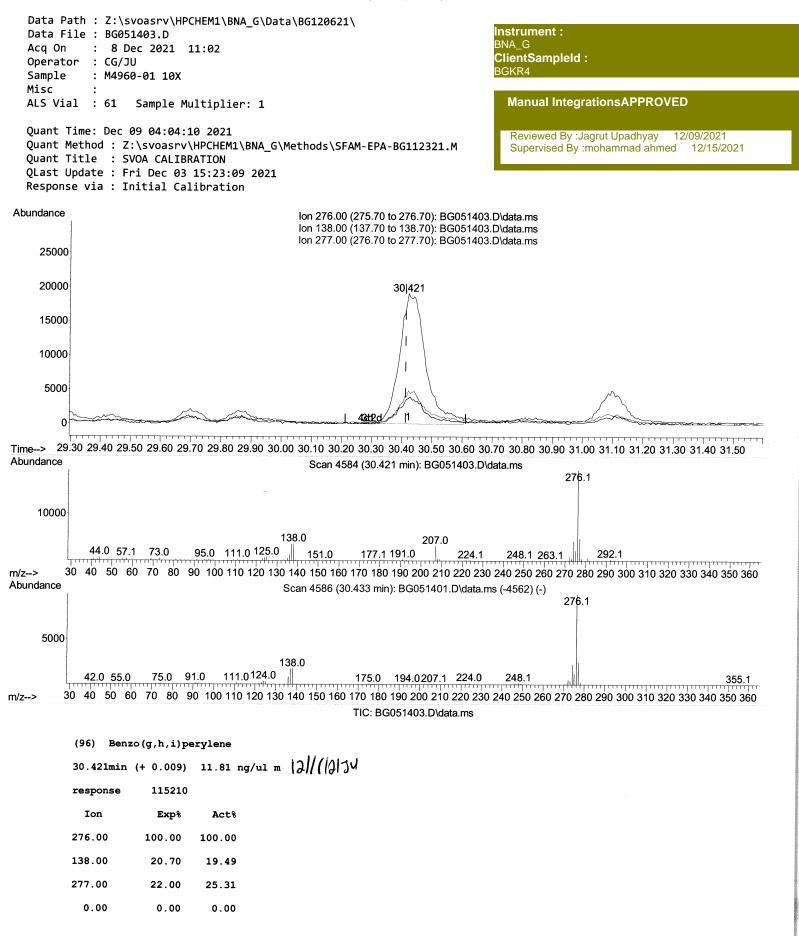












	C		
Data Path : Z:\svoasrv\HPCHEM1	BNA_G\Data\B	G120621\	lu of rum on f
Data File : BG051403.D			Instrument : BNA G
Acq On : 8 Dec 2021 11:02			ClientSampleId :
Operator : CG/JU Sample : M4960-01 10X			BGKR4
Sample : M4960-01 10X Misc :			
ALS Vial : 61 Sample Multipl	ier: 1		Manual IntegrationsAPPROVED
Quant Time: Dec 09 04:04:10 202			Reviewed By :Jagrut Upadhyay 12/09/2021
Quant Method : Z:\svoasrv\HPCHE Quant Title : SVOA CALIBRATION	MI\BNA_G\Met	nods\SFAM-EPA-BG112321.M	Supervised By :mohammad ahmed 12/15/2021
QLast Update : Fri Dec 03 15:23			
Response via : Initial Calibrat			
Compound	R.T. QIO	n Response Conc Units Dev	(Min)
Internal Standards			
1) 1,4-Dichlorobenzene-d4	8.188 15	2 30042 20.000 ng/ul	-0.01
20) Naphthalene-d8	11.020 13	0.	0.00
38) Acenaphthene-d10	14.822 16	•	0.00
64) Phenanthrene-d10	17.577 18	-	0.00
79) Chrysene-d12	21.878 24	0 158319 20.000 ng/ul	0.00
88) Perylene-d12	25.280 26	160964 20.000 ng/ul	0.00
System Monitoring Compounds			
3) 1,4-Dioxane-d8	0.000 9	5 0 0.000 ng/uL	
4) Pyridine-d5	3.993 84	1807m ≥ 0.712 ng/ul>	0.01 ja/16/2174
7) Phenol-d5	7.366 99	9 5563 1.874 ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.507 6	-	0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.730 132	2 3937 1.841 ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.923 113	8 4287m> 1.789 ng/ul>	0.001211/12134
21) Nitrobenzene-d5	9.375 128	8 1795 1.620 ng/ul	0.00
24) 2-Nitrophenol-d4	10.104 143	0,	0.00
28) 2,4-Dichlorophenol-d3	10.668 165	0,	0.01
31) 4-Chloroaniline-d4 46) Dimethylphthalate-d6	11.173 131 14.217 166	0,	0.00
49) Acenaphthylene-d8	14.522 166	0,	0.00 0.00
54) 4-Nitrophenol-d4	15.098 143	1096m > 1.035 mg/ul	- 0.05 (2111)1)
60) Fluorene-d10	15.815 176		0.00
65) 4,6-Dinitro-2-methylph			
73) Anthracene-d10	17.671 188	-	0.00
81) Pyrene-d10	19.951 212	0,	0.00
92) Benzo(a)pyrene-d12	25.045 264	22368 2.602 ng/ul	0.00
Target Compounds		0.97	alue
30) Naphthalene	11.067 128		94
50) Acenaphthylene	14.557 152		
52) Acenaphthene	14.886 153	-	97
61) Fluorene	15.874 166	-	95
72) Phenanthrene	17.619 178	261358 26.042 ng/ul	100
74) Anthracene	17.707 178	0,	98
77) Carbazole	17.983 167	0.	96
80) Fluoranthene	19.622 202	0,	97
82) Pyrene 85) Benzo(a)anthracene	19.986 202 21.855 228	0,	97
87) Chrysene	21.925 228	277392 25.833 ng/ul 249390 24.176 ng/ul	99 98
90) Benzo(b)fluoranthene	24.193 252	287211 26.440 ng/ul	98
91) Benzo(k)fluoranthene	24.252 252	111487m ≻ 10.937 ng/ul >	
93) Benzo(a)pyrene	25.116 252	225764 21.785 ng/ul	96
94) Indeno(1,2,3-cd)pyrene	29.199 276	142388 12.278 ng/ul	97
95) Dibenzo(a,h)anthracene	29.228 278	35767 3.635 ng/ul	93
96) Benzo(g,h,i)perylene	30.421 276	115210m > 11.808 ng/ul =	> [0][(1012)

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

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