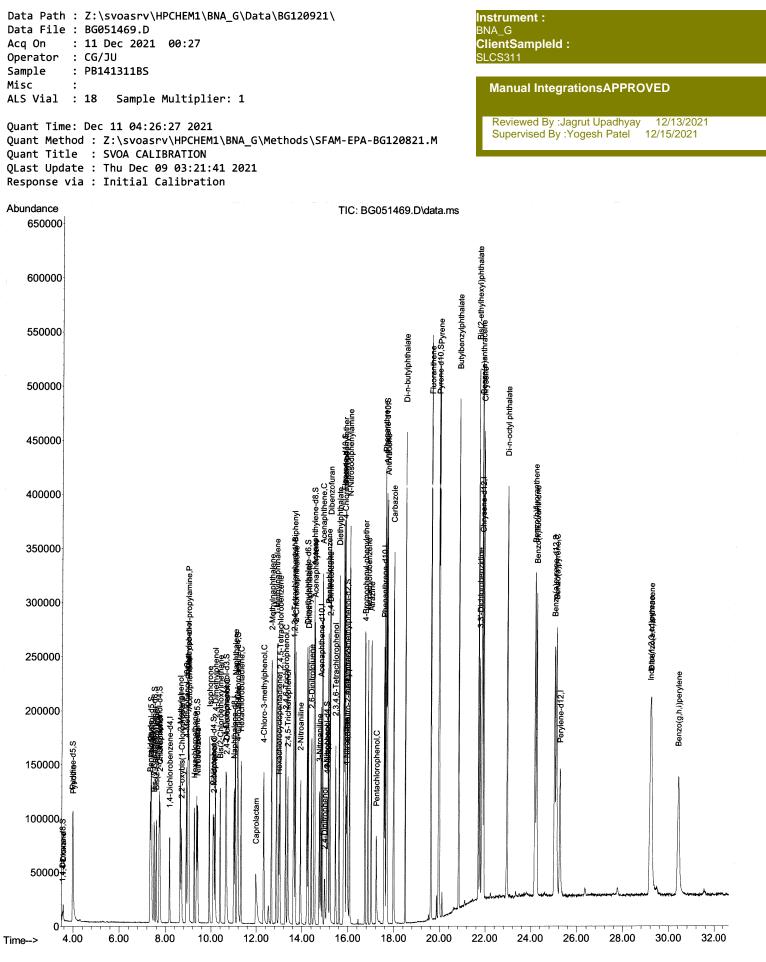
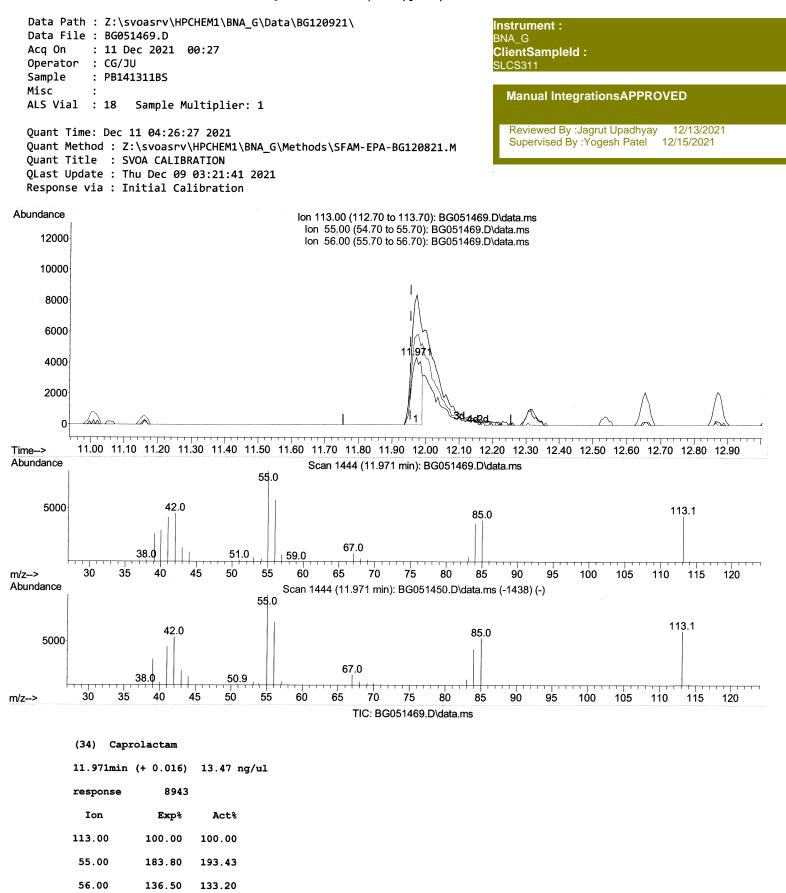
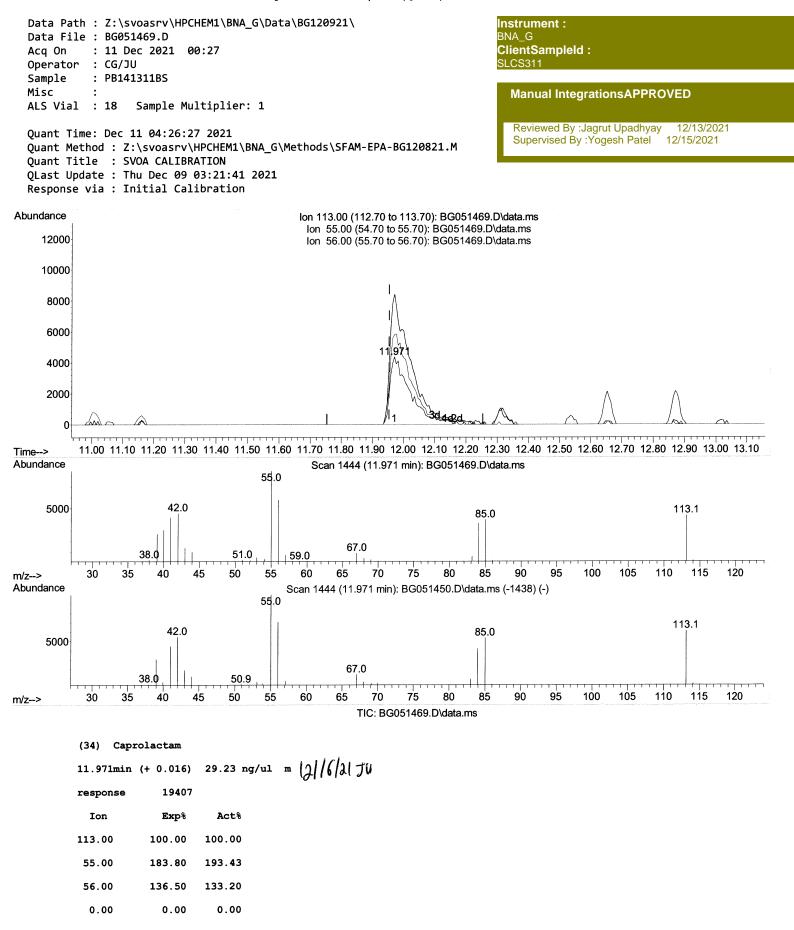
(QT Reviewed)



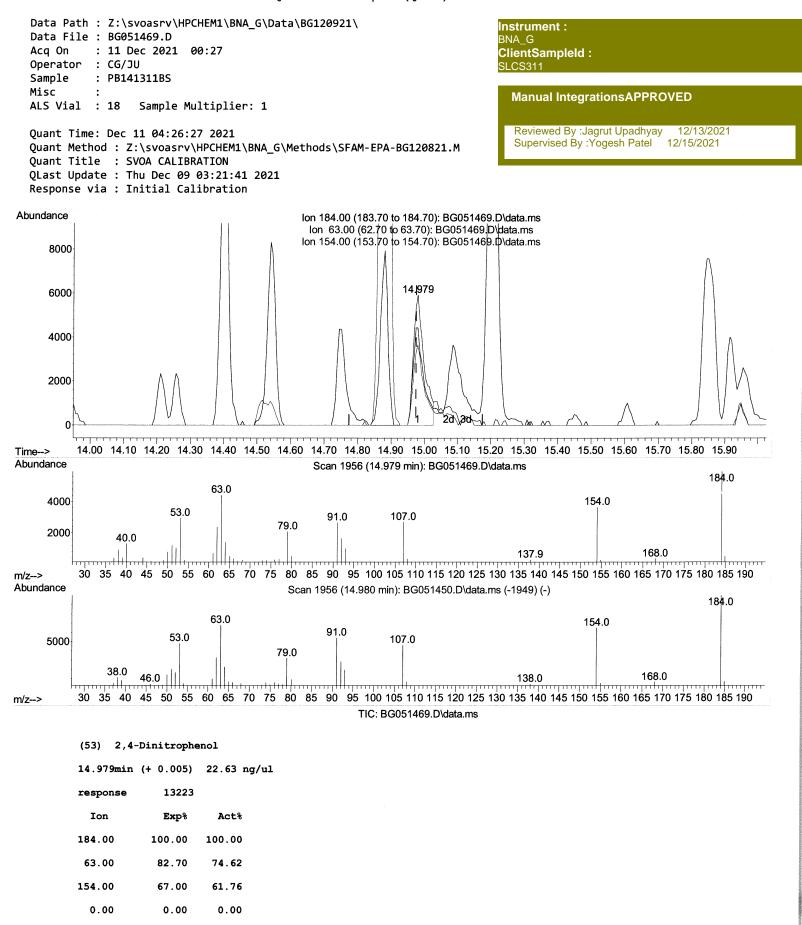


0.00 0.00 0.00

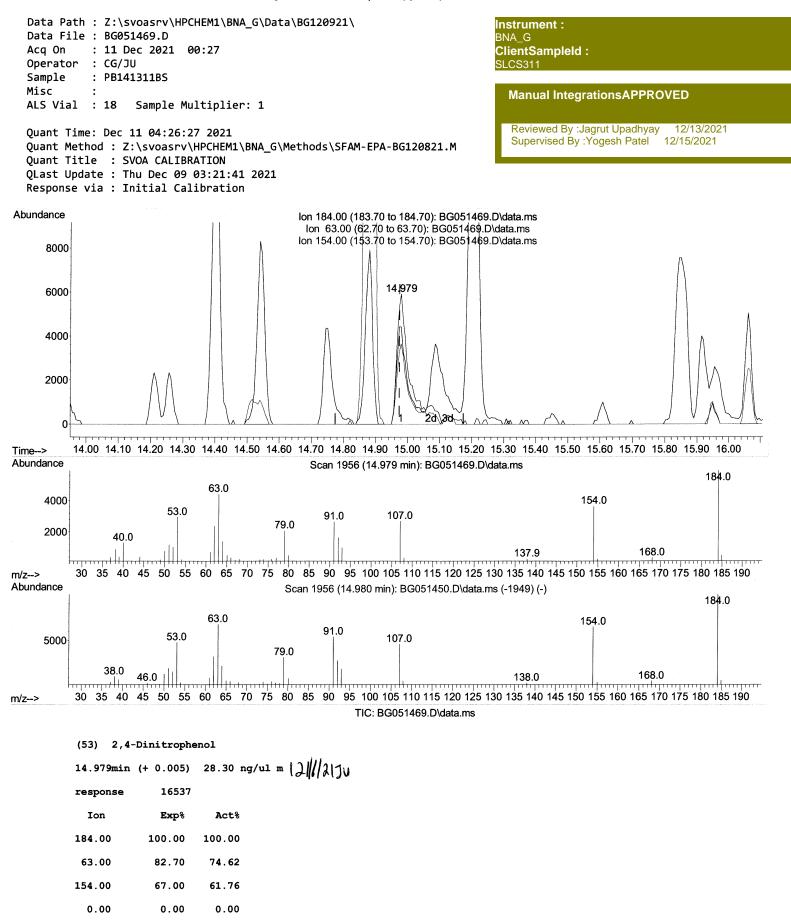


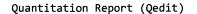


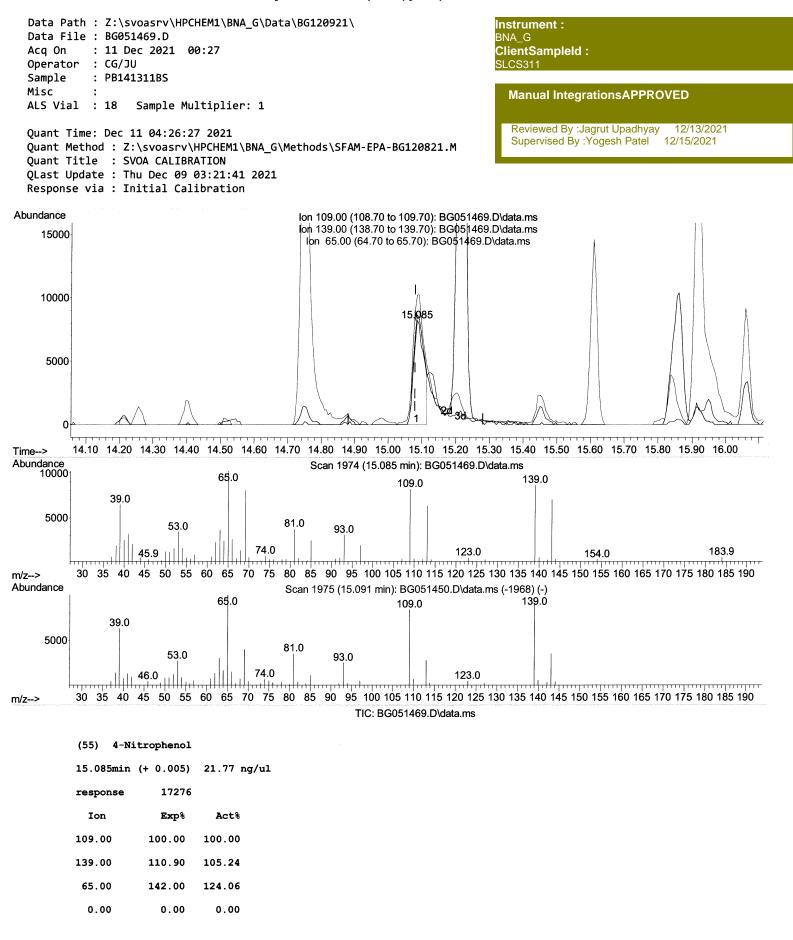




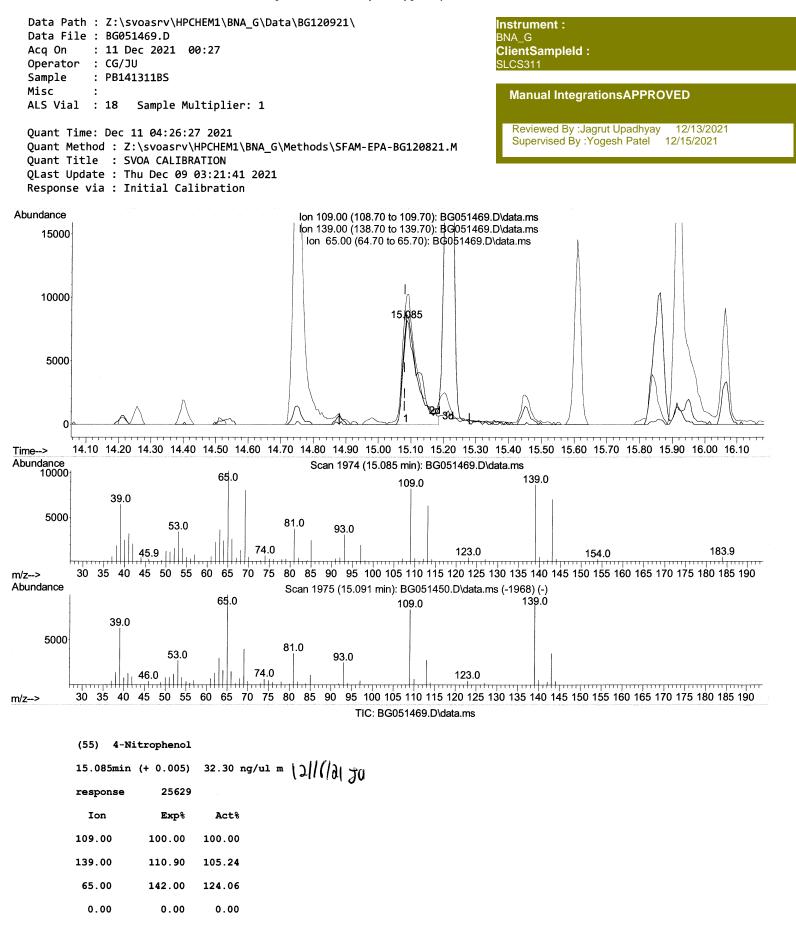




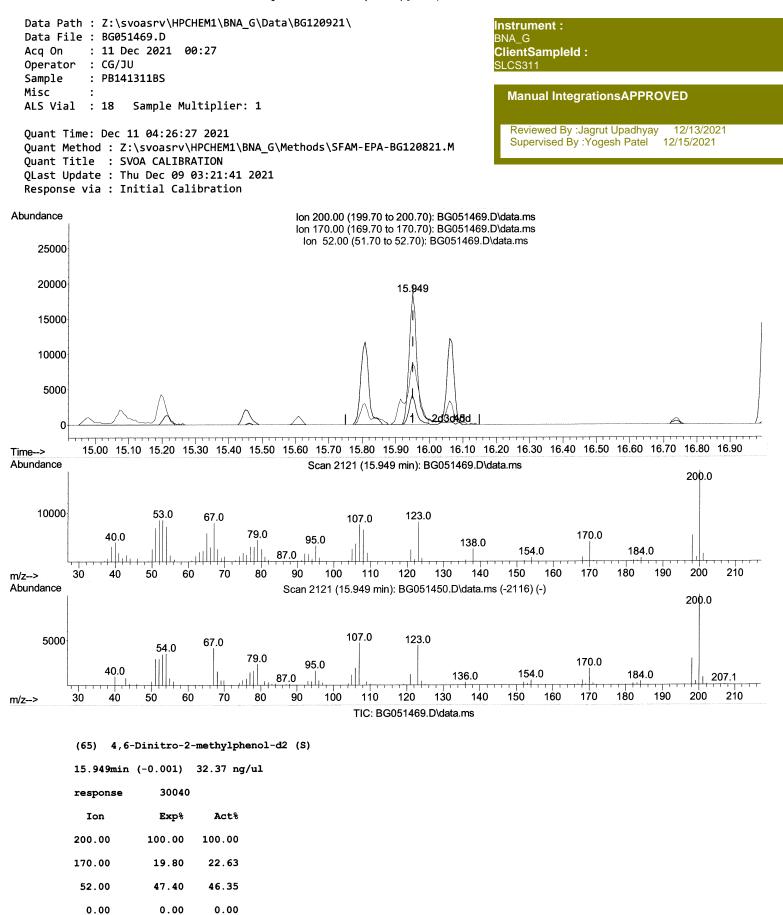


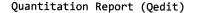


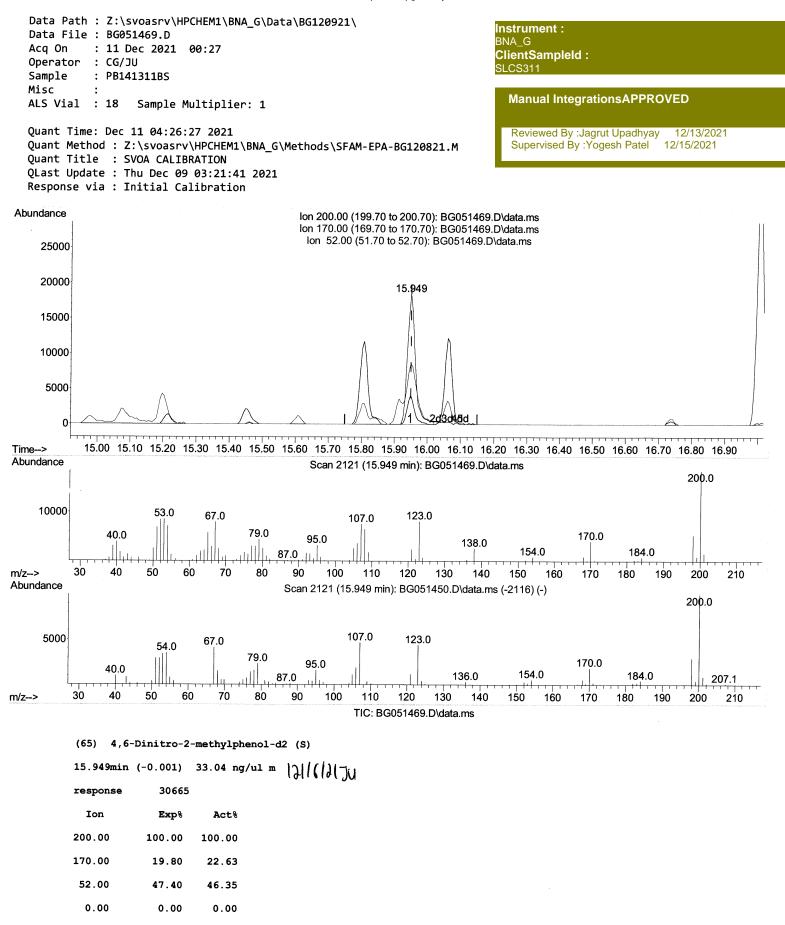




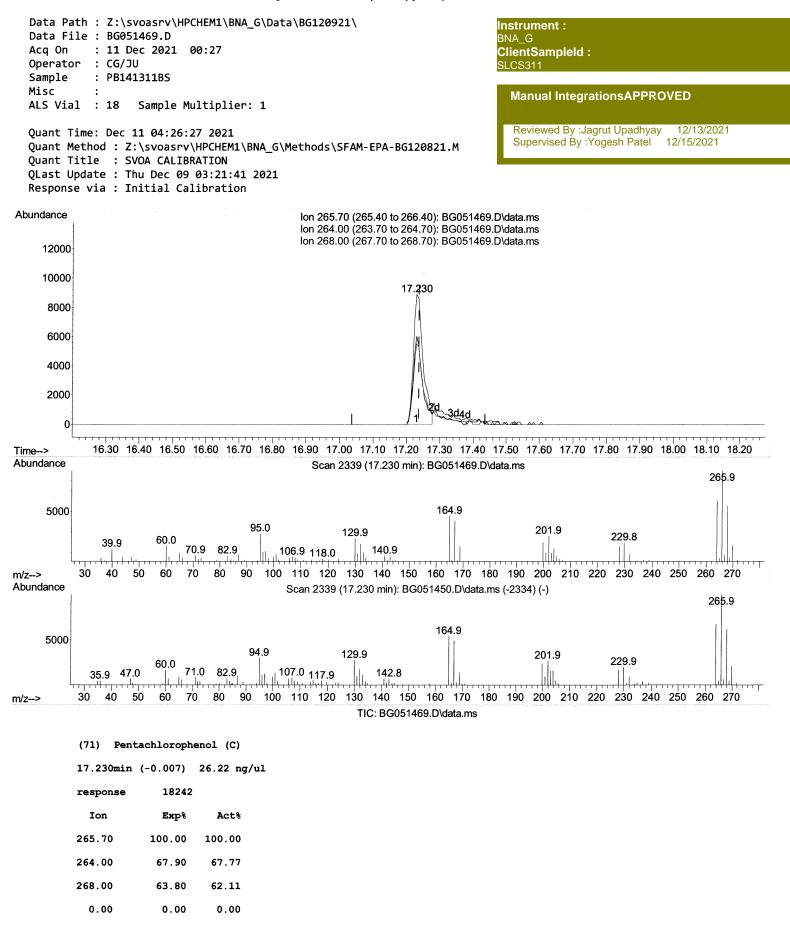




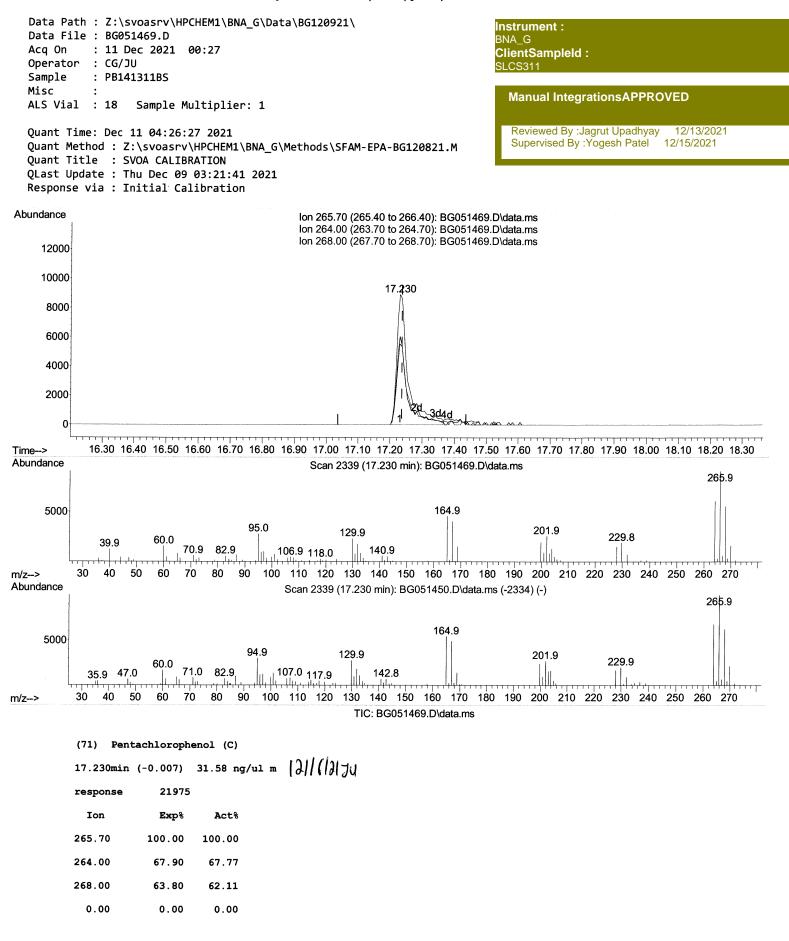












Data Path : Z:\svoasrv\HPCHEM1	BNA_G\Da	ta\BG1	120921\		Instrument :		
Data File : BG051469.D					BNA_G		
Acq On : 11 Dec 2021 00:27					ClientSampleId :		
Operator : CG/JU					SLCS311		
Sample : PB141311BS Misc :							
ALS Vial : 18 Sample Multipl	ier: 1				Manual IntegrationsAPPROVED		
Quant Time, Dec 11 04-26-27 202	4				Reviewed By :Jagrut Upadhyay 12/13/2021		
Quant Time: Dec 11 04:26:27 202 Quant Method : Z:\svoasrv\HPCHE		\Methc	ds\SFAM-FP	A-BG120821.M	Supervised By :Yogesh Patel 12/15/2021		
Quant Title : SVOA CALIBRATION		(,					
QLast Update : Thu Dec 09 03:21:41 2021							
Response via : Initial Calibration							
Compound	рт	07	D				
Compound				Conc Units De			
Internal Standards							
 1,4-Dichlorobenzene-d4 	8.187	152	22834	20.000 ng/ul	0.00		
20) Naphthalene-d8	11.007		102355	20.000 ng/ul			
38) Acenaphthene-d10	14.815	164	68461	20.000 ng/ul			
64) Phenanthrene-d10	17.565	188	156164	20.000 ng/ul	0.00		
79) Chrysene-d12	21.865		140607	20.000 ng/ul	0.00		
88) Perylene-d12	25.267	264	137657	20.000 ng/ul	0.00		
System Monitoning Compaunds							
System Monitoring Compounds 3) 1,4-Dioxane-d8	2 524	06	4240	C 111	0.00		
4) Pyridine-d5	3.534 3.969		4249 52614	6.111 ng/uL 26.351 ng/ul	0.00		
7) Phenol-d5	7.359		73452	31.598 ng/ul	0.00 0.00		
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.500		48376	32.452 ng/ul	0.00		
11) 2-Chlorophenol-d4	7.723		53732	32.490 ng/ul	0.00		
15) 4-Methylphenol-d8	8.916		57609	31.547 ng/ul	0.00		
21) Nitrobenzene-d5	9.368		28242	31.808 ng/ul	0.00		
24) 2-Nitrophenol-d4	10.091	143	32003	31.853 ng/ul	0.00		
28) 2,4-Dichlorophenol-d3	10.649	165	53659	32.829 ng/ul	0.00		
31) 4-Chloroaniline-d4	11.160	131	88743	37 .1 20 ng/ul	0.00		
46) Dimethylphthalate-d6	14.216		175639	33.155 ng/ul	0.00		
49) Acenaphthylene-d8	14.515		217102	32.359 ng/ul	0.00		
54) 4-Nitrophenol-d4	15.073		21941	27.510 ng/ul	0.01		
60) Fluorene-d10 65) 4,6-Dinitro-2-methylph	15.808	176 200	156836	33.258 ng/ul	> 0.00 12/16121 JU		
73) Anthracene-d10	17.664		240247	32.880 ng/ul	0.00		
81) Pyrene-d10	19.944		292447	34.603 ng/ul	0.00		
92) Benzo(a)pyrene-d12	25.026		252421	35.551 ng/ul	0.00		
				0.			
Target Compounds				-	value		
2) 1,4-Dioxane	3.569	88	8877	11.441 ng/uL	88		
5) Pyridine	3.992	79	54128	25.971 ng/ul	96		
6) Benzaldehyde 8) Phenol	7.324 7.388	77	50207	33.975 ng/ul	98		
10) Bis(2-Chloroethyl)ether	7.588	94 93	73782 55290	31.010 ng/ul	99 99		
12) 2-Chlorophenol	7.752	128	52182	30.339 ng/ul 30.802 ng/ul	98		
13) 2-Methylphenol	8.646	108	53890	30.422 ng/ul	98		
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.704	45	82888	30.226 ng/ul	94		
16) Acetophenone	9.022	105	85829	30.352 ng/ul	95		
17) N-Nitroso-di-n-propyla	8.992	70	51710	30.503 ng/ul	99		
18) 4-Methylphenol	8.980	108	57074	30.670 ng/ul	95		
19) Hexachloroethane	9.262	117	22116	30.192 ng/ul	98		
22) Nitrobenzene	9.409	77	75632	31.310 ng/ul	100		
23) Isophorone	9.926	82	139728	30.129 ng/ul	99		
25) 2-Nitrophenol 26) 2,4-Dimethylphenol	10.126 10.179	139 107	30925	30.738 ng/ul	99		
27) Bis(2-Chloroethoxy)met	10.179	107 93	62593 77587	29.379 ng/ul 30.891 ng/ul	100 98		
29) 2,4-Dichlorophenol	10.402	162	50905	31.766 ng/ul	98 97		
30) Naphthalene	11.060	128	172973	30.773 ng/ul	98		
32) 4-Chloroaniline	11.184	127	67875	28.224 ng/ul	100		
33) Hexachlorobutadiene	11.319	225	34058	31.158 ng/ul	99		
34) Caprolactam	11.971	113		29.226 ng/ul	> [all(laiju		
35) 4-Chloro-3-methylphenol	12.312	107	60535	30.411 ng/ul	97		

Data Path : Z:\svoasrv\HPCHEM1 Data File : BG051469.D Acq On : 11 Dec 2021 00:27 Operator : CG/JU Sample : PB141311BS Misc : ALS Vial : 18 Sample Multip Quant Time: Dec 11 04:26:27 20 Quant Method : Z:\svoasrv\HPCH Quant Title : SVOA CALIBRATIO QLast Update : Thu Dec 09 03:2 Response via : Initial Calibra	lier: 1 21 EM1\BNA_G\Me N 1:41 2021	Instrument : BNA_G ClientSampleId : SLCS311 Manual IntegrationsAPPROVED Reviewed By :Jagrut Upadhyay 12/13/2021 Supervised By :Yogesh Patel 12/15/2021	
Compound		on Response Conc Units Dev(M	lin)
36) 2-Methylnaphthalene	12.653 1		98
37) 1-Methylnaphthalene		42 119628 31.006 ng/ul	97
39) 1,2,4,5-Tetrachloroben		16 66446 31.170 ng/ul	98
40) Hexachlorocyclopentadien		37 30953 27.421 ng/ul	99
41) 2,4,6-Trichlorophenol		96 43217 31.383 ng/ul	96
42) 2,4,5-Trichlorophenol		96 45540 30.890 ng/ul	99
43) 1,1'-Biphenyl 44) 2-Chloronaphthalene		54 158636 30.994 ng/ul	99
44) 2-Chibronaphchaiene 45) 2-Nitroaniline		52 125473 31.245 ng/ul	97
47) Dimethylphthalate	13.922 14.262 1	55 47823 31.506 ng/ul 53 165166 30.936 ng/ul	95 98
48) 2,6-Dinitrotoluene		55 35896 31.773 ng/ul	97
50) Acenaphthylene	14.545 1	0.	98
51) 3-Nitroaniline	14.750 1		95
52) Acenaphthene	14.879 1	0	96
53) 2,4-Dinitrophenol	14.979 1		1211612120
55) 4-Nitrophenol	15.085 10		(01101012)
56) Dibenzofuran	15.214 16		94
57) 2,4-Dinitrotoluene	15.197 10	0.	94
58) 2,3,4,6-Tetrachlorophenol	15.449 23	0.	100
59) Diethylphthalate	15.608 14	S.	99
61) Fluorene	15.861 16	8.	98
62) 4-Chlorophenyl-phenyle		0,	94
63) 4-Nitroaniline	15.914 13	0.	97
66) 4,6-Dinitro-2-methylph 67) N-Nitrosodiphenylamine	15.966 19	0.	99 07
68) 4-Bromophenyl-phenylether		6.	97 94
69) Hexachlorobenzene			97
70) Atrazine	17.006 20		98
71) Pentachlorophenol	17.230 26		(2)/(1)/(1)
72) Phenanthrene	17.612 17		99
74) Anthracene	17.706 17	8 265895 31.569 ng/ul	97
75) 1,2,3,4-Tetrachloroben		6 69512 31.832 ng/uL	96
76) Pentachlorobenzene	15.132 25	0.	99
77) Carbazole	17.982 16	0,	99
78) Di-n-butylphthalate	18.493 14	0.	99
80) Fluoranthene	19.615 20	0	97
82) Pyrene 83) Butylbenzylphthalate	19.979 20 20.831 14	0.	97
84) 3,3'-Dichlorobenzidine		0	98
85) Benzo(a)anthracene	21.754 25 21.848 22	0.	98 99
86) Bis(2-ethylhexyl)phtha			100
87) Chrysene	21.912 22	Q .	99
89) Di-n-octyl phthalate	22.952 14		100
90) Benzo(b)fluoranthene	24.174 25	0.	99
91) Benzo(k)fluoranthene	24.245 25		99
93) Benzo(a)pyrene	25.103 25		98
94) Indeno(1,2,3-cd)pyrene	29.186 27	C .	99
95) Dibenzo(a,h)anthracene	29.221 27	8 251210 31.178 ng/ul	98
96) Benzo(g,h,i)perylene	30.408 27		99

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