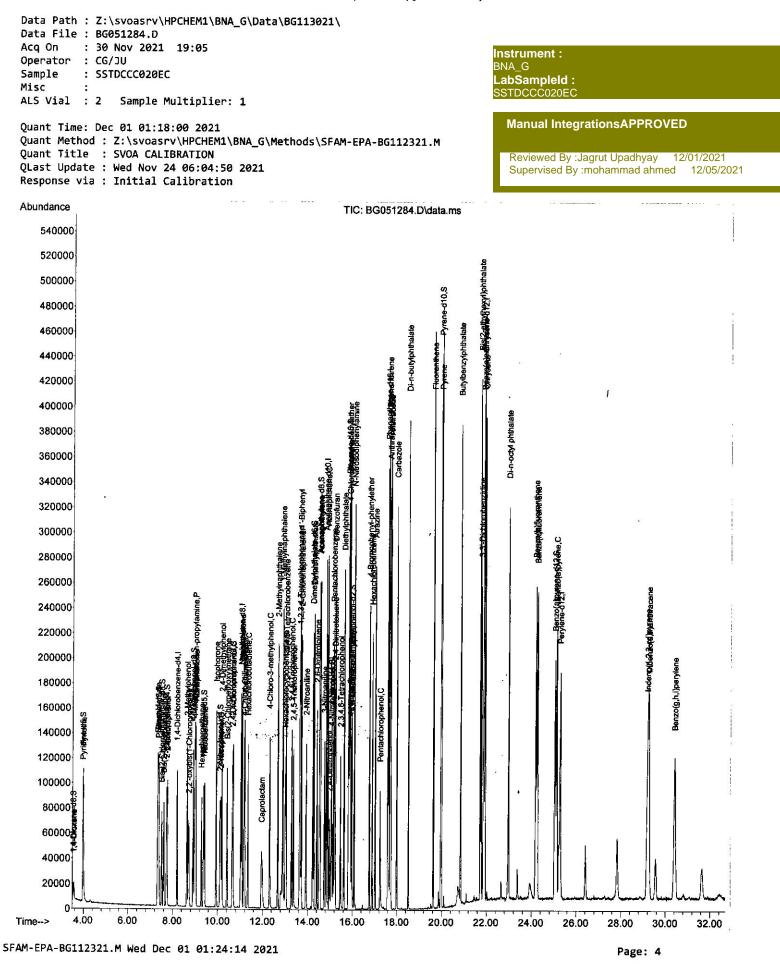
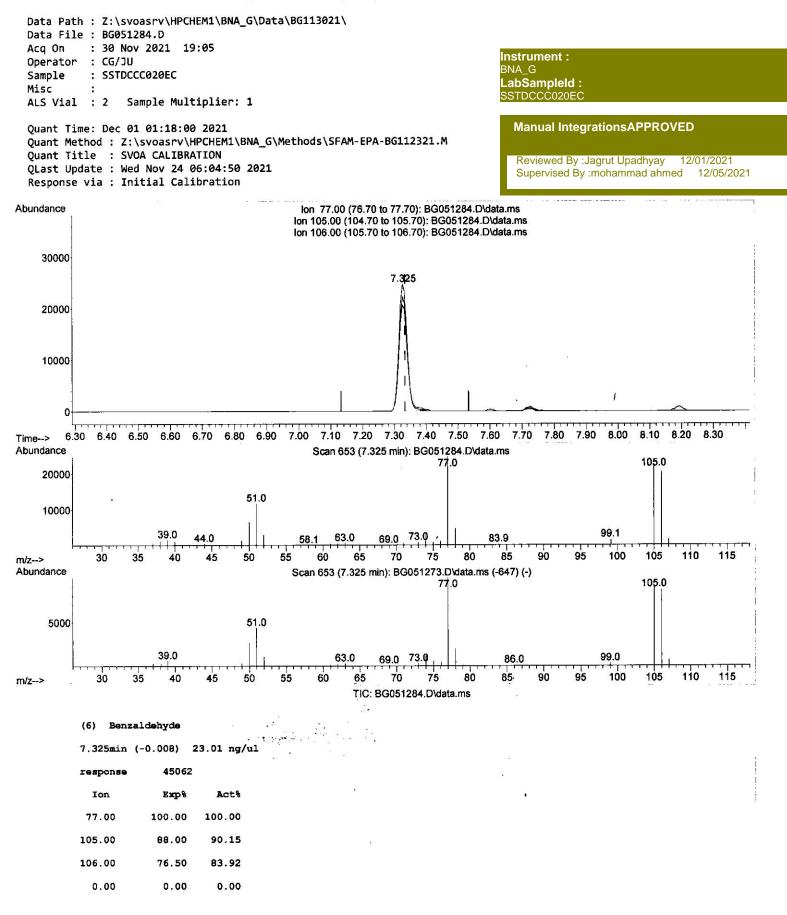
(QT Reviewed)

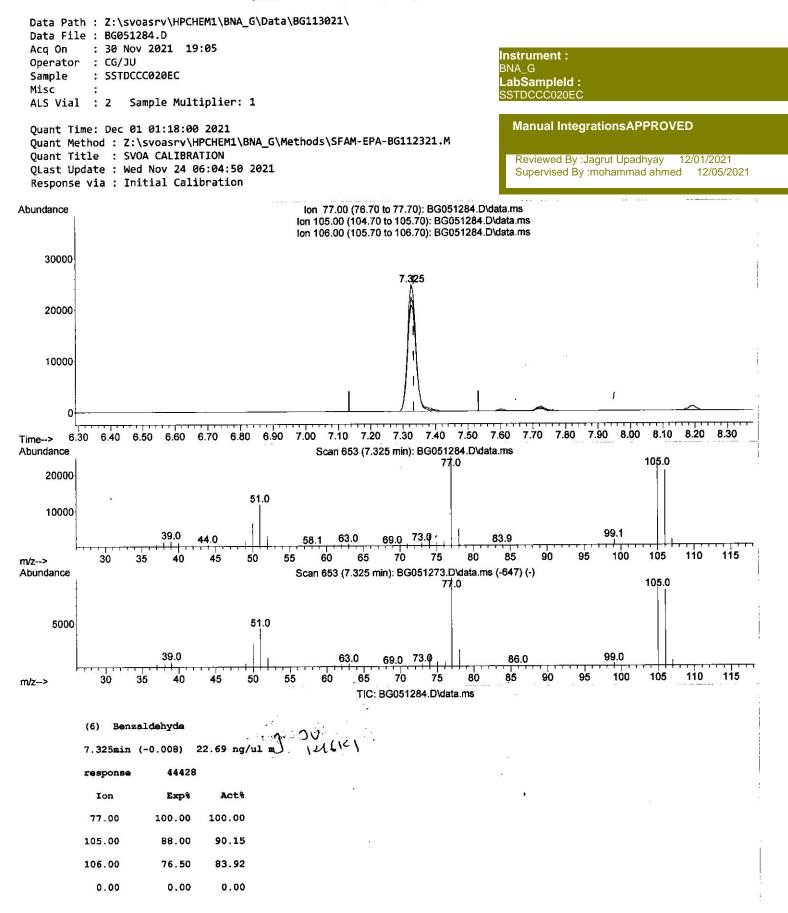




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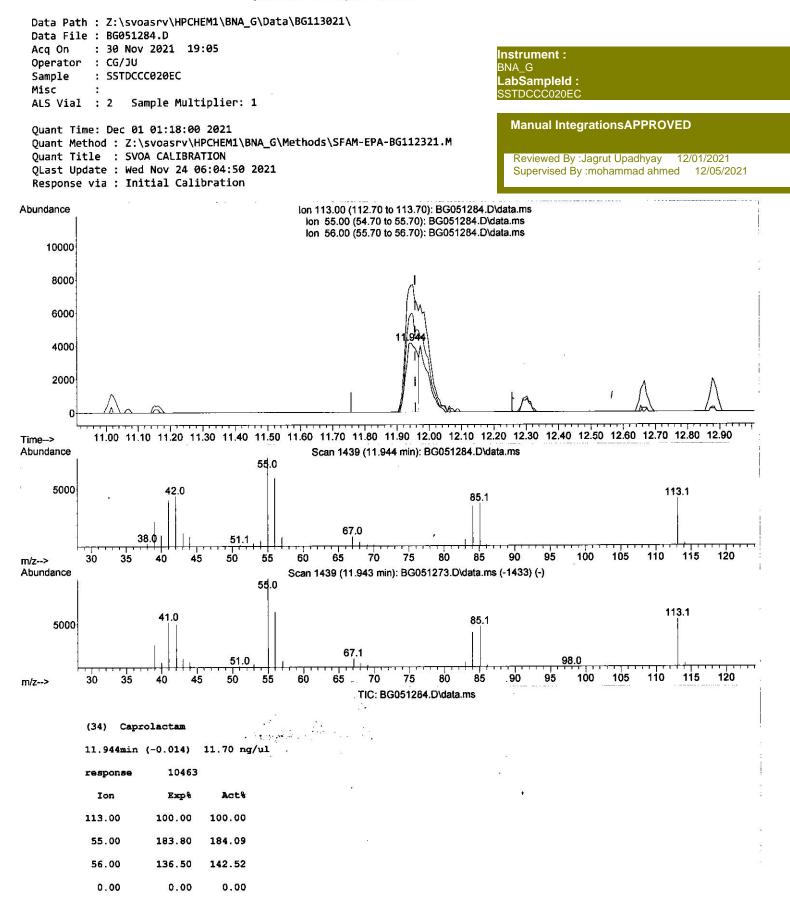
Quantitation Report (Qedit)



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## Quantitation Report (Qedit)

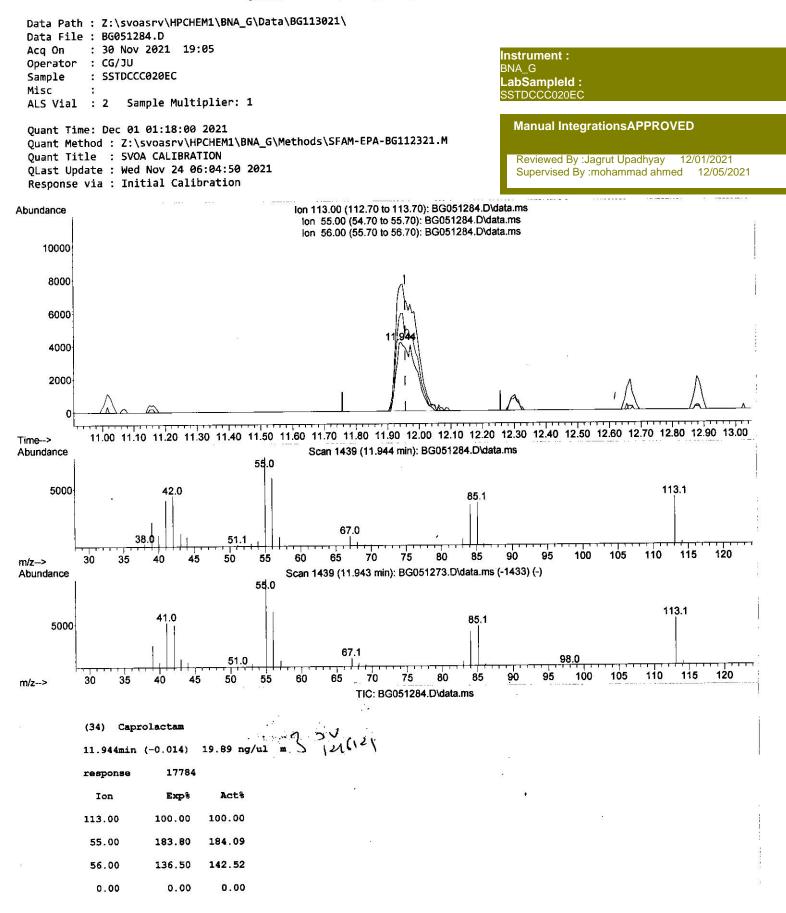


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Quantitation Report (Qedit)



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	Qua	murua	cion nepor	c (g. ne	/	
Data Path : Z:\svoasrv\HPCHEM1\BM	IA_G\Data	\BG11	3021\			
Data File : BG051284.D						
Acq On : 30 Nov 2021 19:05					In	strument :
Operator : CG/JU Sample : SSTDCCC020EC						IA_G
Sample : SSTDCCC020EC Misc :						bSampleId :
ALS Vial : 2 Sample Multiplier	r: 1				SS	TDCCC020EC
Quant Time: Dec 01 01:18:00 2021				DC112221 M		Manual Integrati
Quant Method : Z:\svoasrv\HPCHEM	1\BNA_G\N	Method	IS\SFAM-EPA	-B0112321.M		
Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:	6 2021					Reviewed By :Jagr
Response via : Initial Calibratio						Supervised By :mo
.3						
	R.T. (			Conc Units Dev	(Min)	
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.189	152		20.000 ng/ul	-0.01	
20) Naphthalene-d8	11.021		142994		0.00	
	14.823		97443	20.000 ng/ul	0.00	
	17.578			20.000 ng/ul		
· · · · · · · · · · · · · · · · · · ·	21.873		198699	20.000 ng/ul 20.000 ng/ul	0.00	
88) Perylene-d12	25.275	264	195571	20.000 ng/ui	0.00	
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.530	96	7305	8.160 ng/uL		4
4) Pyridine-d5	3.959	84	49560	18.865 ng/ul	-0.02	
7) Phenol-d5	7.349	99	59705	19.417 ng/ul	0.00	
9) Bis-(2-Chloroethyl)eth	7.508	67	37864	19.607 ng/ul	0.00	•
11) 2-Chlorophenol-d4	7.725		42550	19.217 ng/ul	0.00	
15) 4-Methylphenol-d8	8.906		47425		0.00 0.00	
21) Nitrobenzene-d5	9.370		23347	19.342 ng/ul	0.00	
24) 2-Nitrophenol-d4	10.099		26319 44818	19.329 ng/ul 19.400 ng/ul	0.00	
28) 2,4-Dichlorophenol-d3	10.645		44818 64781	19.164 ng/ul	0.00	
31) 4-Chloroaniline-d4	11.156 14.217		147172	19.629 ng/ul	0.00	
46) Dimethylphthalate-d6	14.523		185262	and a second	0.00	
49) Acenaphthylene-d8 54) 4-Nitrophenol-d4	15.046		20354		0.00	
60) Fluorene-d10	15.816		129469	19.176 ng/ul	0.00	
65) 4,6-Dinitro-2-methylph			23257	17.104 ng/ul	0.00	
73) Anthracene-d10	17.672	188	206142	19.561 ng/ul	0.00	
81) Pyrene-d10	19.952	212	238557		0.00	
92) Benzo(a)pyrene-d12	25.034	264	201931	19.333 ng/ul	0.00	
Target Compounds					alue	
2) 1.4-Dioxane	3.565	88	8055	7.978 ng/uL	91	
5) Pyridine	3.982	79	53427 n	19.544 ng/ul	94	
6) Benzaldehyde	7.325	77	44428m	) 22.689 ng/ul	07	
8) Phenol	7.378	94	61348 (	19.260 ng/ul	97 97	
10) Bis(2-Chloroethyl)ether	7.602	93	48251	20.023 ng/ul	97 97	•
12) 2-Chlorophenol	7.754	128	43902	19.457 ng/ul	98	*
13) 2-Methylphenol	8.642	108	45756 68697	19.755 ng/ul	97	
14) 2,2'-oxybis(1-Chloropr	8.718 9.023		75855	19.764 ng/ul	99	
16) Acetophenone 17) N-Nitroso-di-n-propyla	8.994	2	45040	20.422 ng/ul	98	
18) 4-Methylphenol	8.971		49189	19.388 ng/ul	94·	0.1
19) Hexachloroethane	9.276		18344	19.248 ng/ul	95	1121
22) Nitrobenzene	9.411		62924	19.881 ng/ul	97	306121
23) Isophorone	9,928		124859	20.305 ng/ul	100	,
25) 2-Nitrophenol	10.128		26824	19.019 ng/ul	97	
26) 2,4-Dimethylphenol	10.181		57680	20.003 ng/ul	97	
<pre>27) Bis(2-Chloroethoxy)met</pre>	10.410		67467	19.874 ng/ul	99	
29) 2,4-Dichlorophenol	10.669		43780	19.251 ng/ul	98 97	
30) Naphthalene	11.068		149749	19.246 ng/ul 19.258 ng/ul	100	
32) 4-Chloroaniline	11.180			19.258 ng/ul	98	
33) Hexachlorobutadiene	11.333 11.944		· · · · · · · · · · · · · · · · · · ·	<b>\</b>	20	
34) Caprolactam 35) 4-Chloro-3-methylphenol	12.302				99	
22) 4-CUTOLO-2-WeinArbuenor		/				

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## al IntegrationsAPPROVED

wed By :Jagrut Upadhyay 12/01/2021 rvised By :mohammad ahmed 12/05/2021

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Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG113021\ Data File : BG051284.D Acq On : 30 Nov 2021 19:05 Operator : CG/JU Sample : SSTDCCC020EC Misc : ALS Vial : 2 Sample Multiplier: 1 Quant Time: Dec 01 01:18:00 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

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Instrument :

BNA\_G LabSampleId :

SSTDCCC020EC

Manual IntegrationsAPPROVED

1

Reviewed By :Jagrut Upadhyay 12/01/2021 Supervised By :mohammad ahmed 12/05/2021

sponse via : Initial Calibratio					
Compound	R.T.	QIon	Response	Conc Units Dev(Mi	
36) 2-Methylnaphthalene	12.666	142	102274	19.325 ng/ul	99 99
36) 2-Methylnaphthalene	12.884	142	104390	19.173 ng/ul	98
39) 1,2,4,5-Tetrachloroben	13.025	216	58529	19.133 ng/ul	95
40) Hexachlorocyclopentadiene	12.989	237	32251	26.083 ng/ul	99
40) hexacilito ocyclophenol 41) 2,4,6-Trichlorophenol	13.271	196	35191	18.331 ng/ul	98
41) 2,4,5-Trichlorophenol	13.354	196	37110	18.460 ng/ul 19.501 ng/ul	97
42) 2,4,9 11 10120 07 43) 1,1'-Biphenyl	13.659		141930	19.063 ng/ul	97
44) 2-Chloronaphthalene	13.706		110363	20.383 ng/ul	92
45) 2-Nitroaniline	13.918	-	40841	19.304 ng/ul	100
47) Dimethylphthalate	14.264		146500	19.796 ng/ul	91
48) 2,6-Dinitrotoluene	14.405		31558	19.619 ng/ul	99
50) Acenaphthylene	14.552		183259	21.059 ng/ul	94
51) 3-Nitroaniline	14.740		33183	19.289 ng/ul	97
52) Acenaphthene	14.887		118825	23.707 ng/ul	90
53) 2,4-Dinitrophenol	14.964		20889	22.435 ng/ul	92
55) 4-Nitrophenol	15.058		23620	19.181 ng/ul	98
56) Dibenzofuran	15.222		170428 45421	19.949 ng/ul	93
57) 2 4-Dinitrotoluene	15.199	Carl Changers		17.129 ng/ul	95
58) 2,3,4,6-Tetrachlorophenol	15.45		27041	19.626 ng/ul	99
59) Diethylphthalate	15.62		156340	19.383 ng/ul	100
61) Fluorene	15.87		137957 72412	18.879 ng/ul	98
62) 4-Chlorophenyl-phenyle	15.85			22.945 ng/ul	98
63) 4-Nitroaniline	15.90			17.029 ng/ul#	97
66) 4.6-Dinitro-2-methylph	15.96			19.411 ng/ul	98
67) N-Nitrosodiphenylamine	10.0/			18.706 ng/ul	92
68) 4-Bromophenyl-phenyletner	16.75			19,292 ng/ul	97
69) Hexachlorobenzene	10.01			19,381 ng/ul	98
70) Atrazine	17.00			19.641 ng/ul	98
71) Pentachlorophenol	17.23	Courses and the second second		19.367 ng/ul	99
72) Phenanthrene	17.6			19.653 ng/ul	97
74) Anthracene	17.7			19.236 ng/uL	98
75) 1.2.3.4-Tetrachloroben	. 13.6			19.145 ng/uL	98
76) Pentachlorobenzene	T.2 • T.			20.319 ng/ul	99
77) Carbazole	17.9			20.028 ng/ul	100
78) Di-n-butylphthalate	18.5			19.857 ng/ul	96
80) Fluoranthene	19.6 19.9			20.021 ng/ul	99
82) Pyrene	20.8				96
83) Butylbenzylphthalate	20.0			19.245 ng/ul	98
84) 3,3'-Dichlorobenzidine	21.8			19.381 ng/ul	99
85) Benzo(a)anthracene				19.699 ng/ul	100
86) Bis(2-ethylhexyl)phtha	21.9			19.251 ng/ul	. 99
87) Chrysene	22.9			20.319 ng/ul	100
89) Di-n-octyl phthalate	24.		52 251911	19.086 ng/ul	98
90) Benzo(b)fluoranthene	24.		52 24244	3 19.575 ng/ul	98
91) Benzo(k)fluoranthene	24.		52 24312	3 19.308 ng/ul	97
93) Benzo(a)pyrene	29.		76 27299	2 19.374 ng/ul	98
94) Indeno(1,2,3-cd)pyrene			78 22922	2 19.176 ng/ul	98
95) Dibenzo(a, h)anthracene			76 22750		98
96) Benzo(g,h,i)perylene					
	22 19 19 19 19 19 19 19 19 19 19 19 19 19				

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

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