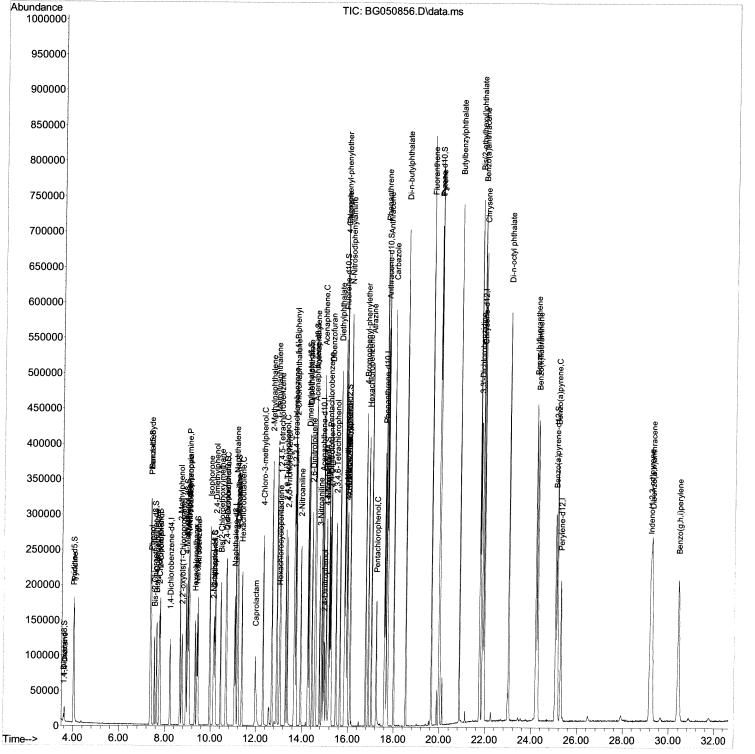
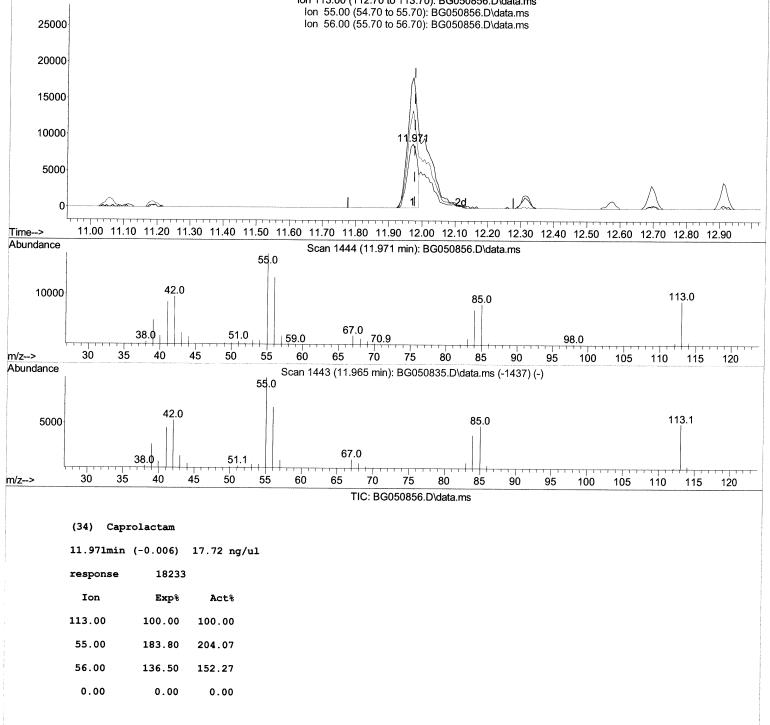
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\	
Data File : BG050856.D	Instrument :
Acq On : 3 Nov 2021 18:59	BNA_G
Operator : CG/JU	ClientSampleId :
Sample : PB140350BS	SLCS350
Misc :	
ALS Vial : 41 Sample Multiplier: 1	Manual IntegrationsAPPROVED
Quant Time: Nov 07 07:39:23 2021	Reviewed By :Jagrut Upadhyay 11/08/2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA	-BG110321.M Supervised By :mohammad ahmed 11/08/2021
Quant Title : SVOA CALIBRATION	
QLast Update : Tue Nov 02 14:49:05 2021	
Response via : Initial Calibration	

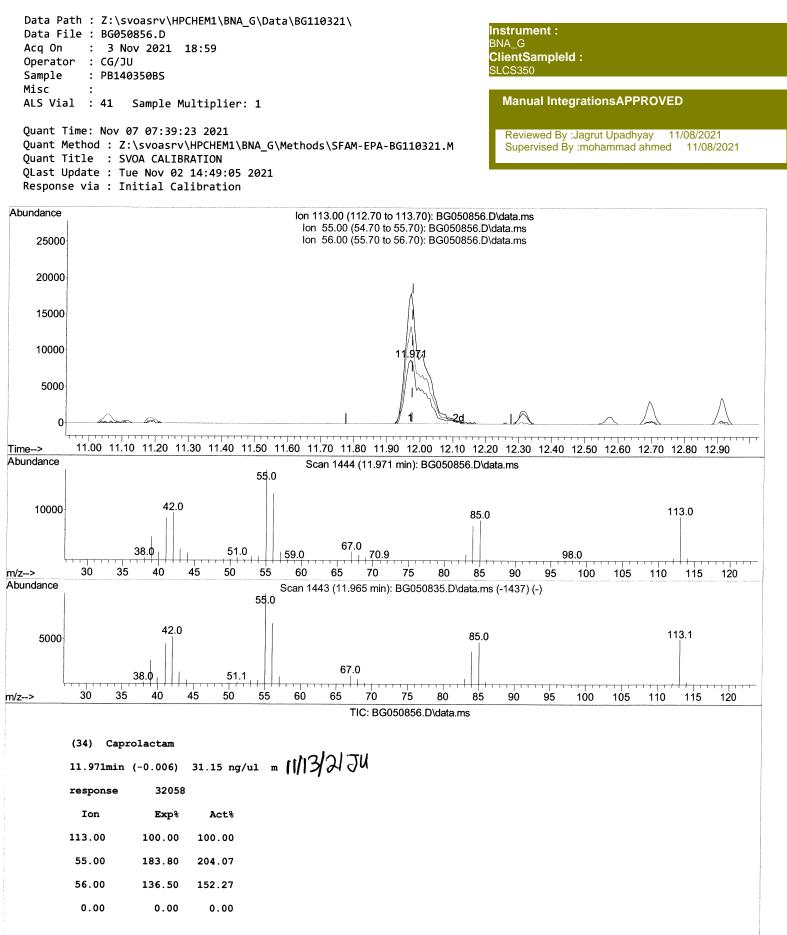


Quantitation Report (Qedit)





SFAM-EPA-BG110321.M Sun Nov 07 07:40:45 2021



Operator : CG/JU Sample : PB14035 Misc :	5.D 2021 18:59 DBS mple Multiplier: 1 07:39:23 2021 voasrv\HPCHEM1\BNA_ CALIBRATION Nov 02 14:49:05 202	G\Metho		PA-BG110321.M	Reviewed By	d: grationsAPPROVED :Jagrut Upadhyay 11/08/2021 / :mohammad ahmed 11/08/2021
Compound		-	•	Conc Units Dev	· /	
Internal Standards						
1) 1,4-Dichlorobe	enzene-d4 8.22	8 152	33967	20.000 ng/ul	0.00	
20) Naphthalene-d&		4 136	156125	20.000 ng/ul	0.00	
38) Acenaphthene-c			110676	20.000 ng/ul	0.00	
64) Phenanthrene-c			244017	20.000 ng/ul	0.00	
79) Chrysene-d12		5 240	207035	20.000 ng/ul	-0.01	
88) Perylene-d12		1 264	207033	20.000 ng/ul		
oby refyrene urz	23.29	1 204	203100	20.000 ng/u1	-0.02	
System Monitoring (ompounds					
3) 1,4-Dioxane-d8	•	7 06	5315		0.00	
4) Pyridine-d5			5215	4.955 ng/uL	0.00	
7) Phenol-d5	4.01		79770	25.337 ng/ul	0.00	
9) Bis-(2-Chloroe	7.37		100519	27.739 ng/ul	0.00	
11) 2-Chlorophenol			63372	27.073 ng/ul	0.00	
15) 4-Methylphenol			69990 78731	27.870 ng/ul	0.00	
21) Nitrobenzene-d			37857	27.598 ng/ul	0.00	
24) 2-Nitrophenol-				28.533 ng/ul	0.00	
28) 2,4-Dichloroph			42776 73432	28.996 ng/ul	0.00	
31) 4-Chloroanilin			100831	29.549 ng/ul	0.00	
46) Dimethylphthal				26.793 ng/ul	0.00	
49) Acenaphthylene			248425 301714	29.339 ng/ul	0.00	
54) 4-Nitrophenol-				28.600 ng/ul	0.00	
60) Fluorene-d10	15.843		43477 213460	28.319 ng/ul	0.00	
	methylph 15.960		41670	28.458 ng/ul	0.00	
73) Anthracene-d10			324288	28.166 ng/ul	0.00	
81) Pyrene-d10	19.973		379571	28.109 ng/ul	0.00	
92) Benzo(a)pyrene			321762	28.385 ng/ul	0.00	
	25.007	204	521702	28.378 ng/ul	0.00	
Target Compounds				0.42	lue	
2) 1,4-Dioxane	3.622	88	12990	11.238 ng/uL#	90	
5) Pyridine	4.033		88217	27.069 ng/ul	97	
6) Benzaldehyde	7.364		73485	32.152 ng/ul	96	
8) Phenol	7.400		111751	29.812 ng/ul	98	
10) Bis(2-Chloroet			83669	29.820 ng/ul	97	
12) 2-Chlorophenol	7.787		76562	30.027 ng/ul	94	
13) 2-Methylphenol	8.663		82920	29.927 ng/ul	99	
14) 2,2'-oxybis(1-0			130565	29.543 ng/ul	98	
16) Acetophenone	. 9.057		131077	29.578 ng/ul	95	
17) N-Nitroso-di-n			80621	30.152 ng/ul	98	
18) 4-Methylphenol	8.992	108	89639	30.385 ng/ul	99	
19) Hexachloroetham	ne 9.321	117	31546	29.591 ng/ul	97	
22) Nitrobenzene	9.444	77	114115	30.839 ng/ul	99	
23) Isophorone	9.967	82	226662	31.562 ng/ul	99	
25) 2-Nitrophenol	10.161	139	46726	31.571 ng/ul	96	
26) 2,4-Dimethylphe		107	97603	29.965 ng/ul	98	
27) Bis(2-Chloroeth	oxy)met 10.443	93	121313	31.352 ng/ul	98	
29) 2,4-Dichlorophe	nol 10.696	162	77717	32.062 ng/ul	95	
30) Naphthalene	11.107	128	260793	30.548 ng/ul	98	
32) 4-Chloroaniline		127	110211	29.493 ng/ul	99	
33) Hexachlorobutad	liene 11.377	225	48489	30.478 ng/ul	96	
34) Caprolactam	11.971	113	32058m 🗲	31.152 ng/ul >		
35) 4-Chloro-3-meth	ylphenol 12.312	107	100969	32.607 ng/ul	97	

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Data Path : Z:\svoasrv\HPCHEM1\	BNA_G\Da	ta\BG1	L10321\		Instrument			
Data File : BG050856.D					Instrument : BNA_G			
Acq On : 3 Nov 2021 18:59					ClientSampleId :			
Operator : CG/JU					SLCS350			
Sample : PB140350BS Misc :								
ALS Vial : 41 Sample Multipl	ier·1				Manual IntegrationsAPPROVED			
, io viai i ii Sampie Haitipi	101.1							
Quant Time: Nov 07 07:39:23 202	1				Reviewed By :Jagrut Upadhyay 11/08/2021			
Quant Method : Z:\svoasrv\HPCHE	M1\BNA_G	\Methc	ds\SFAM-EF	A-BG110321.M	Supervised By :mohammad ahmed 11/08/2021			
Quant Title : SVOA CALIBRATION								
QLast Update : Tue Nov 02 14:49								
Response via : Initial Calibrat	ion							
Compound	вт	OTon	Response	Conc Units Dev	(Min)			
36) 2-Methylnaphthalene	12.699	142	181329	31.177 ng/ul	96			
37) 1-Methylnaphthalene	12.911		184413	31.292 ng/ul	97			
39) 1,2,4,5-Tetrachloroben	13.058	216	95784	29.702 ng/ul	99			
40) Hexachlorocyclopentadiene	13.028	237	38700	24.990 ng/ul	99			
<pre>41) 2,4,6-Trichlorophenol</pre>	13.293	196	66337	31.437 ng/ul	99			
42) 2,4,5-Trichlorophenol	13.369		71353	31.496 ng/ul	98			
43) 1,1'-Biphenyl	13.692		246339	30.451 ng/ul	98			
44) 2-Chloronaphthalene	13.739		190768	30.092 ng/ul	98			
45) 2-Nitroaniline 47) Dimethylphthalate	13.939 14.298	65 163	79461 263447	31.549 ng/ul	94			
48) 2,6-Dinitrotoluene	14.298		56701	31.116 ng/ul 31.998 ng/ul	100 95			
50) Acenaphthylene	14.580	152	324894	30.729 ng/ul	99			
51) 3-Nitroaniline	14.756	138	58246	31.781 ng/ul	92			
52) Acenaphthene	14.920	153	214655	30.876 ng/ul	98			
53) 2,4-Dinitrophenol	14.967	184	27177	27.801 ng/ul	96			
55) 4-Nitrophenol	15.055	109	43012	30.548 ng/ul	94			
56) Dibenzofuran	15.249	168	304605	30.610 ng/ul	98			
57) 2,4-Dinitrotoluene	15.208	165	81692	32.303 ng/ul	99			
· · · ·	15.473	232	57105	32.075 ng/ul	99			
59) Diethylphthalate	15.649	149	283123	31.241 ng/ul	99			
<pre>61) Fluorene 62) 4-Chlorophenyl-phenyle</pre>	15.902 15.884	166 204	241552 126495	30.668 ng/ul	100 98			
63) 4-Nitroaniline	15.919	138	58201	30.842 ng/ul 32.011 ng/ul	96			
66) 4,6-Dinitro-2-methylph	15.972	198	43947	30.461 ng/ul	99			
67) N-Nitrosodiphenylamine	16.095	169	217546	31.894 ng/ul	97			
68) 4-Bromophenyl-phenylether	16.777	248	76971	31.711 ng/ul	98			
69) Hexachlorobenzene	16.900	284	78667	31.525 ng/ul	96			
70) Atrazine	17.035	200	88193	30.491 ng/ul	99			
71) Pentachlorophenol	17.241	266	36307	31.691 ng/ul	95			
72) Phenanthrene 74) Anthracene	17.641	178	409386	31.429 ng/ul	99			
75) 1,2,3,4-Tetrachloroben	17.735 13.657	178 216	397850 100750	30.442 ng/ul	99			
76) Pentachlorobenzene	15.167	250	93756	30.323 ng/uL 30.455 ng/uL	99 98			
77) Carbazole	17.999	167	379261	32.377 ng/ul	98			
78) Di-n-butylphthalate	18.534	149	486068	31.579 ng/ul	100			
80) Fluoranthene	19.638	202	494801	30.832 ng/ul	99			
82) Pyrene	20.003	202	479213	30.561 ng/ul	99			
83) Butylbenzylphthalate	20.866	149	213027	31.593 ng/ul	98			
84) 3,3'-Dichlorobenzidine	21.783	252	150006	29.751 ng/ul	98			
85) Benzo(a)anthracene	21.877	228	449074	31.332 ng/ul	100			
86) Bis(2-ethylhexyl)phtha	21.748	149	307209	31.740 ng/ul	98			
87) Chrysene 89) Di-n-octyl phthalate	21.942 23.017	228 149	420554 518879	30.716 ng/ul	98 100			
90) Benzo(b)fluoranthene	23.017	252	453646	31.075 ng/ul 31.048 ng/ul	100 99			
91) Benzo(k)fluoranthene	24.280	252	411272	29.996 ng/ul	99			
93) Benzo(a)pyrene	25.138	252	423704	30.447 ng/ul	98			
94) Indeno(1,2,3-cd)pyrene	29.198	276	478325	30.877 ng/ul	98			
95) Dibenzo(a,h)anthracene	29.268	278	403695	30.799 ng/ul	98			
96) Benzo(g,h,i)perylene	30.432		397037	30.618 ng/ul	98			

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