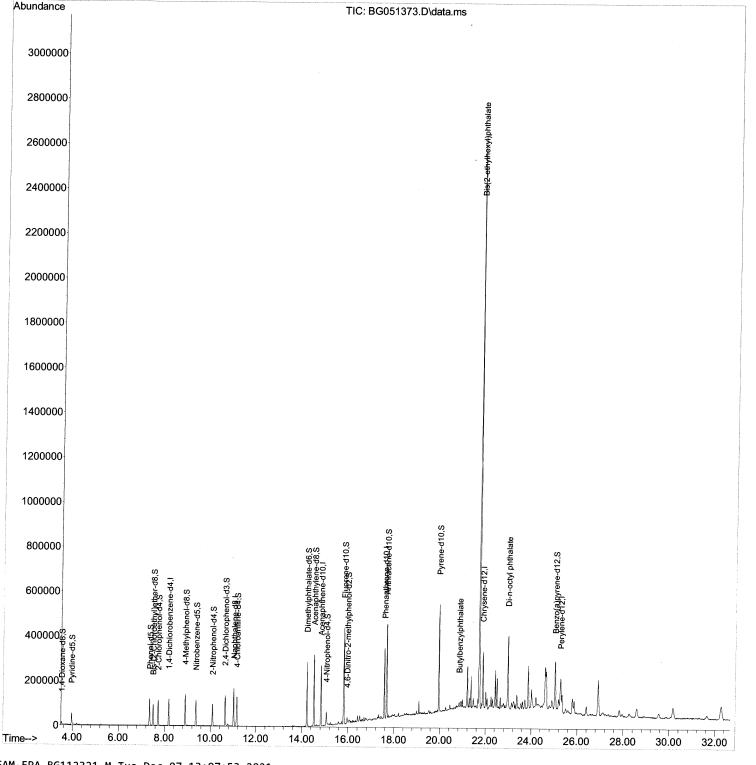
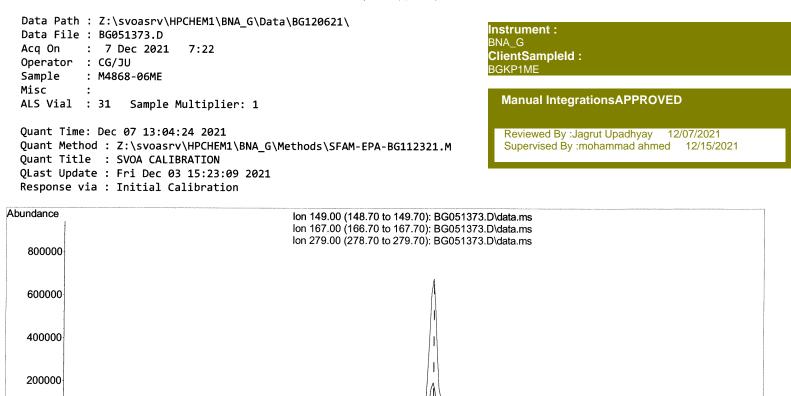
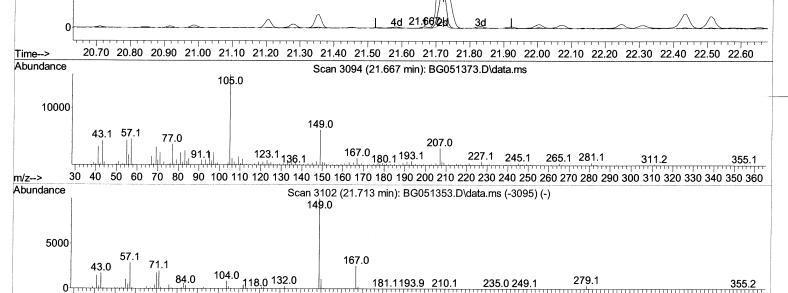
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\	
Data File : BG051373.D	Instrument :
Acq On : 7 Dec 2021 7:22	BNA G
Operator : CG/JU	ClientSampleId :
Sample : M4868-06ME	BGKP1ME
Misc :	
ALS Vial : 31 Sample Multiplier: 1	Manual IntegrationsAPPROVED
Quant Time: Dec 07 13:04:24 2021	
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M	Reviewed By :Jagrut Upadhyay 12/07/2021
Quant Title : SVOA CALIBRATION	Supervised By :mohammad ahmed 12/15/2021
QLast Update : Fri Dec 03 15:23:09 2021	
Response via : Initial Calibration	



SFAM-EPA-BG112321.M Tue Dec 07 13:07:53 2021





30 40 50 60 70 80 90 100 110 120 130 140 150 160 170 180 190 200 210 220 230 240 250 260 270 280 290 300 310 320 330 340 350 360

TIC: BG051373.D\data.ms

(86) Bis (2-ethylhexyl) phthalate

m/z-->

 21.667min (-0.056)
 1.03 ng/ul

 response
 7246

 Ion
 Exp%
 Act%

 149.00
 100.00
 100.00

 167.00
 26.40
 21.91

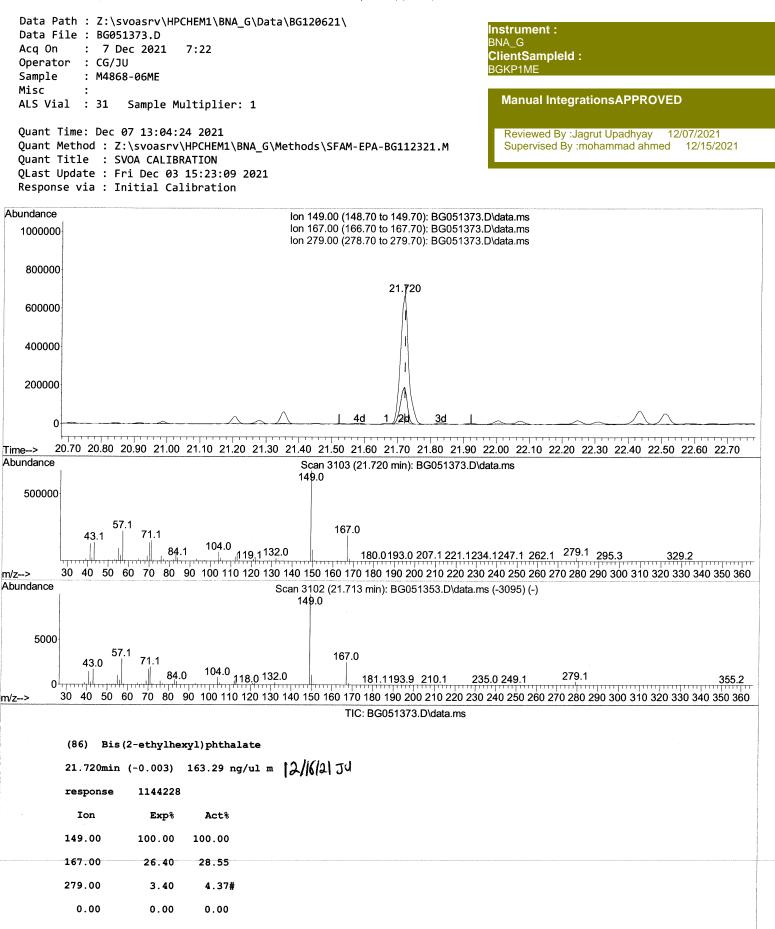
 279.00
 3.40
 3.13

0.00

0.00

0.00





Data Path : Z:\svoasrv\HPCHEM1\ Data File : BG051373.D Acq On : 7 Dec 2021 7:22 Operator : CG/JU Sample : M4868-06ME Misc :	_	ta∖BG	120621\		Instrument : BNA_G ClientSampleId : BGKP1ME
ALS Vial : 31 Sample Multipl	ier: 1				Manual IntegrationsAPPROVED
Quant Time: Dec 07 13:04:24 202 Quant Method : Z:\svoasrv\HPCHE Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23 Response via : Initial Calibrat	Reviewed By :Jagrut Upadhyay 12/07/2021 Supervised By :mohammad ahmed 12/15/2021				
Compound				Conc Units Dev	
Internal Standards					
1) 1,4-Dichlorobenzene-d4	8 188	152	33097	20.000 ng/ul	-0.01
20) Naphthalene-d8	11.020		143449	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.822		97748	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.577		193387	20.000 ng/ul	0.00
79) Chrysene-d12	21.878		161121	20.000 ng/ul	0.00
88) Perylene-d12	25.286		162791	20.000 ng/ul	0.00
				-	
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.523	96	5916	6.212 ng/uL	-0.02
4) Pyridine-d5	3.958	84	34253	12.256 ng/ul	-0.02
7) Phenol-d5	7.354	99	72173	22.064 ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.507	67	46083	22.431 ng/ul	0.00
11) 2-Chlorophenol-d4	7.724		53393	22.667 ng/ul	0.00
15) 4-Methylphenol-d8	8.905	113	57200	21.669 ng/ul	0.00
21) Nitrobenzene-d5	9.369		28404	23.457 ng/ul	0.00
24) 2-Nitrophenol-d4	10.092	143	31730	23.229 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.644		53722	23.180 ng/ul	0.00
-31) 4-Chloroaniline-d4 —	11.155		75327	22.213 ng/ul	0.00
46) Dimethylphthalate-d6	14.217		181307	24.106 ng/ul	0.00
49) Acenaphthylene-d8	14.522		234624	24.739 ng/ul	0.00
54) 4-Nitrophenol-d4	15.057		20576	16.901 ng/ul	0.00
60) Fluorene-d10	15.815		165764	24.475 ng/ul	0.00
65) 4,6-Dinitro-2-methylph 73) Anthracene-d10			6072	5.088 ng/ul	0.00
81) Pyrene-d10	17.677		250393	27.072 ng/ul	0.00
92) Benzo(a)pyrene-d12	19.957		268915	27.584 ng/ul	0.00
52) Benzo(a)pyrene-uiz	25.057	204	229627	26.412 ng/ul	0.01
Target Compounds				Qva	lue
83) Butylbenzylphthalate	20.844	149	6460	1.327 ng/ul#	84
83) Butylbenzylphthalate 86) Bis(2-ethylhexyl)phtha	21.720	149	1144228m 🖒	>163.295 ng/ul >	12/16/21
89) Di-n-octyl phthalate	22.989	149	340501	28.872 ng/ul	100

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