Data File : BG051291.D

Acq On : 1 Dec 2021 18:48

Operator : CG/JU Sample : M4868-01

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Dec 01 20:23:18 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

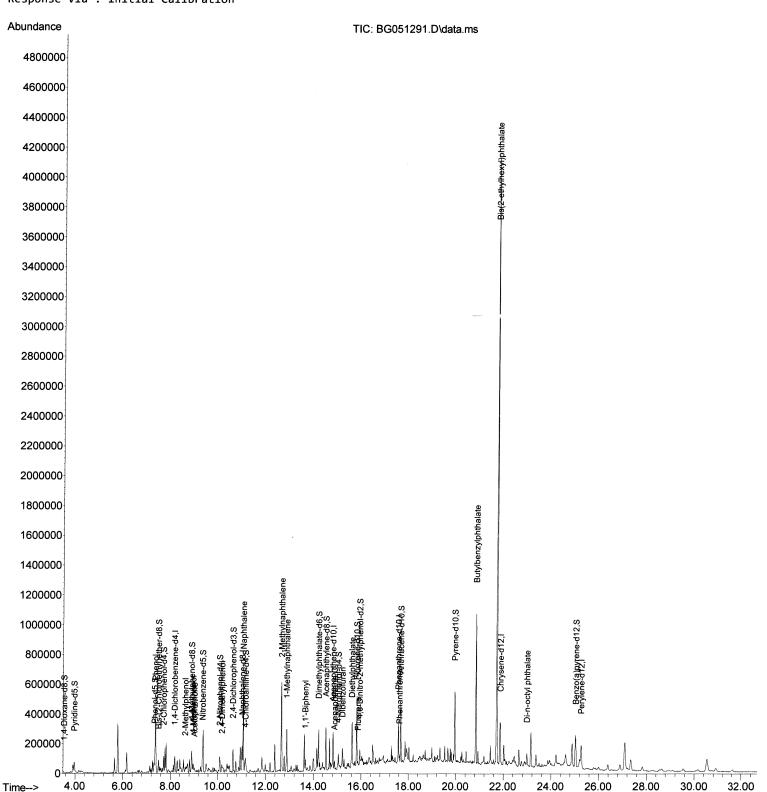


ClientSampleId:

BGKN8



Reviewed By: Jagrut Upadhyay 12/02/2021 Supervised By: mohammad ahmed 12/05/2021



Data File: BG051291.D

Acq On : 1 Dec 2021 18:48

Operator : CG/JU Sample : M4868-01

Misc

ALS Vial : 7 Sample Multiplier: 1

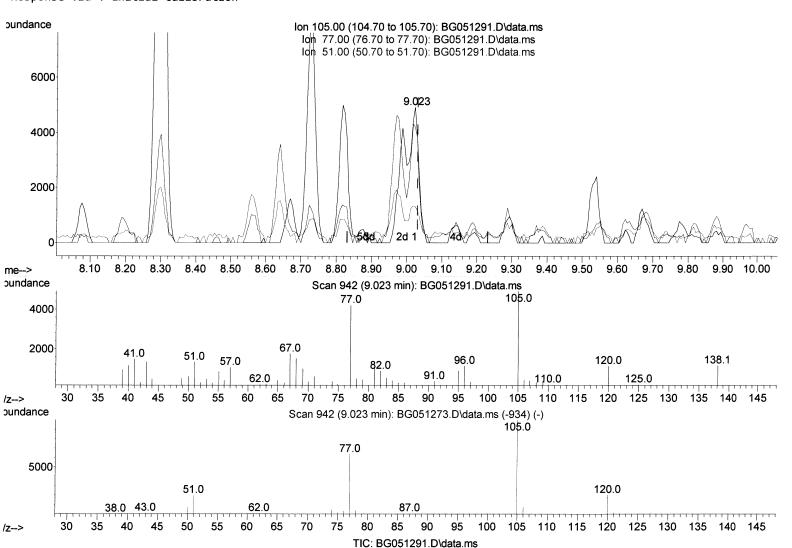
Quant Time: Dec 01 20:23:18 2021

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 Instrument:
BNA_G
ClientSampleId:
BGKN8

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/02/2021 Supervised By :mohammad ahmed 12/05/2021



(16) Acetophenone

9.023min (-0.009) 1.96 ng/ul

response	7488	
Ion	Ехр%	Act%
105.00	100.00	100.00
77.00	84.10	85.65
51.00	30.00	27.05
0.00	0.00	0.00

Data File : BG051291.D

Acq On : 1 Dec 2021 18:48

Operator : CG/JU Sample : M4868-01

Misc :

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Dec 01 20:23:18 2021

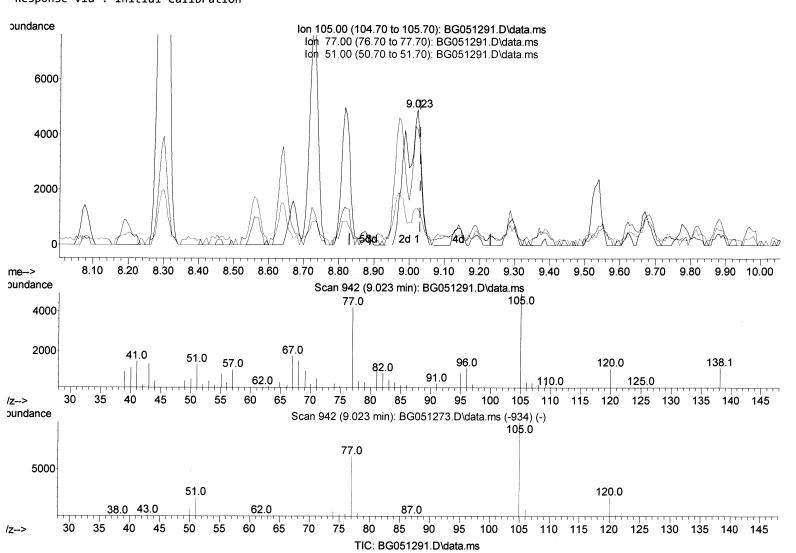
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration Instrument:
BNA_G
ClientSampleId:
BGKN8

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/02/2021 Supervised By :mohammad ahmed 12/05/2021



(16) Acetophenone

9.023min (-0.009) 3.95 ng/ul m \2/20/2/JU

response	15065		
Ion	Exp%	Act%	
105.00	100.00	100.00	
77.00	84.10	85.65	
51.00	30.00	27.05	
0.00	0.00	0.00	

Data File : BG051291.D

Acq On : 1 Dec 2021 18:48

Operator : CG/JU Sample : M4868-01

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Dec 01 20:23:18 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

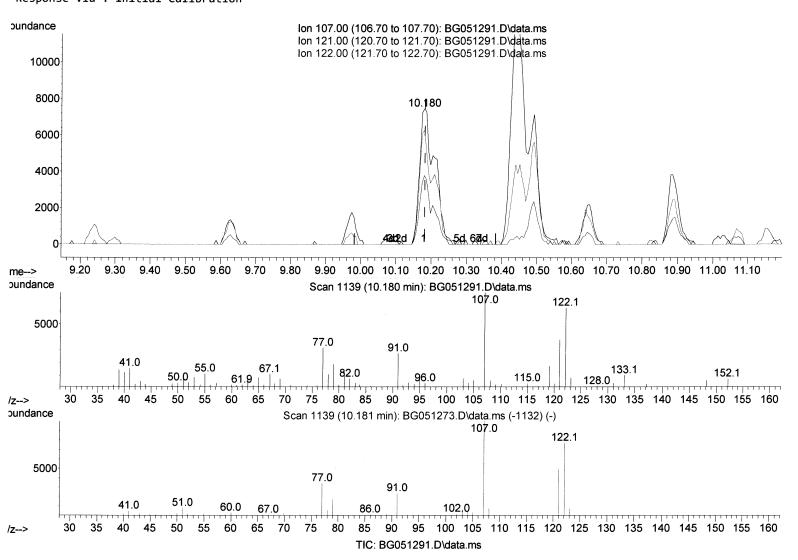
Quant Title : SVOA CALIBRATION

QLast Update: Wed Nov 24 06:04:50 2021 Response via: Initial Calibration

Instrument: BNA_G ClientSampleId: BGKN8

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/02/2021 Supervised By :mohammad ahmed 12/05/2021



(26) 2,4-Dimethylphenol

10.180min (-0.003) 5.39 ng/ul

response	14361	
Ion	Exp%	Act%
107.00	100.00	100.00
121.00	49.10	50.83
122.00	79.60	84.16
0.00	0.00	0.00

Data File: BG051291.D

Acq On : 1 Dec 2021 18:48

Operator : CG/JU Sample : M4868-01

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Dec 01 20:23:18 2021

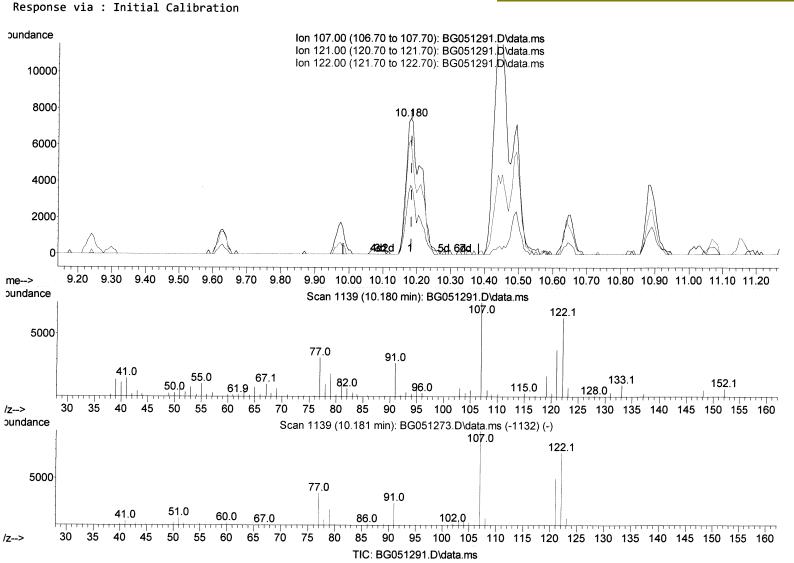
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 24 06:04:50 2021

Instrument:
BNA_G
ClientSampleId:
BGKN8

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/02/2021 Supervised By :mohammad ahmed 12/05/2021



(26) 2,4-Dimethylphenol

10.180min (-0.003) 8.35 ng/ul m 12/20/21JU

response	22256			
Ion	Ежр%	Act%		
107.00	100.00	100.00		
121.00	49.10	50.83		
122.00	79.60	84.16		
0.00	0.00	0.00		

Data File : BG051291.D

Acq On : 1 Dec 2021 18:48

Dperator : CG/JU
Sample : M4868-01

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Dec 01 20:23:18 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

Instrument : BNA_G ClientSampleId : BGKN8

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/02/2021 Supervised By :mohammad ahmed 12/05/2021

Compound	R.T.	QIon	Response	Conc Ur	its Dev((Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	0 100	150	20022	20.000		0.01
20) Naphthalene-d8	8.188 11.020	152 136	30932		ng/ul ng/ul	-0.01
38) Acenaphthene-d10	14.822	164	132153 86642		ng/ul ng/ul	0.00
64) Phenanthrene-d10	17.571	188				0.00
79) Chrysene-d12	21.872	240	176979 159 0 12		ng/ul ng/ul	0.00
88) Perylene-d12	25.268	264	164246		ng/ul ng/ul	0.00
ob) Perylene-ulz	23.208	204	104240	20.000	ng/ui	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.535	96	3257	3.659	ng/uL	0.00
4) Pyridine-d5	3.964	84	44302		ng/ul	-0.01
7) Phenol-d5	7.348	99	67372		ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.507	67	44170		ng/ul	0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.724	132	47550		ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.905	113	55523		ng/ul	0.00
21) Nitrobenzene-d5	9.369	128	26549		ng/ul	0.00
24) 2-Nitrophenol-d4	10.092	143	30576		ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.644	165	51224		ng/ul	0.00
31) 4-Chloroaniline-d4	11.161	131	59453	19.030		0.00
46) Dimethylphthalate-d6	14.217	166	163535	24.531		0.00
49) Acenaphthylene-d8	14.522	160	208090		ng/ul	0.00
54) 4-Nitrophenol-d4	15.051	143	25847	23.952		0.00
60) Fluorene-d10	15.815	176	144495	24.069		0.00
65) 4,6-Dinitro-2-methylph	15.944	200	22421	20.531		0.00
73) Anthracene-d10	17.671	188	230523	27.235		0.00
81) Pyrene-d10	19.951	212	259558	26.977	_	0.00
92) Benzo(a)pyrene-d12	25.033	264	229521	26.166	_	0.00
Target Compounds					Qva	
8) Phenol	7.378	94	234756	74.126		100
<pre>13) 2-Methylphenol</pre>	8.641	108	12176		ng/ul	99
16) Acetophenone	9.023	105	15065m>		ng/ul >	
18) 4-Methylphenol	8.976	108	20867		ng/ul	94
26) 2,4-Dimethylphenol	10.180	107	22256m ≤			uzlaolal ju
30) Naphthalene	11.073	128	515936	71.750		97
36) 2-Methylnaphthalene	12.665	142	237554	48.569	_	100
37) 1-Methylnaphthalene	12.877	142	136963	27.219		96
43) 1,1'-Biphenyl	13.658	154	33570	5.188		96
52) Acenaphthene	14.886	153	20492	3.741		97
56) Dibenzofuran	15.221	168	23057	2.918		99
59) Diethylphthalate	15.615	149	15709	2.218		98
61) Fluorene	15.873	166	14433	2.281		98
72) Phenanthrene	17.618	178	29118	2.980	-	97
83) Butylbenzylphthalate	20.844	149	309601	64.423	_	96
86) Bis(2-ethylhexyl)phtha	21.720	149	1636497	236.645	-	92
89) Di-n-octyl phthalate	22.983	149	107452	9.030	ng/ul	100

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