#### (LSC Reviewed) Quantitation Report

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120621\

Data File : BG051399.D

8 Dec 2021 2:08 Acq On

: CG/JU Operator : M4942-03ME Sample

Misc

Sample Multiplier: 1 ALS Vial : 57

Quant Time: Dec 08 05:31:01 2021

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

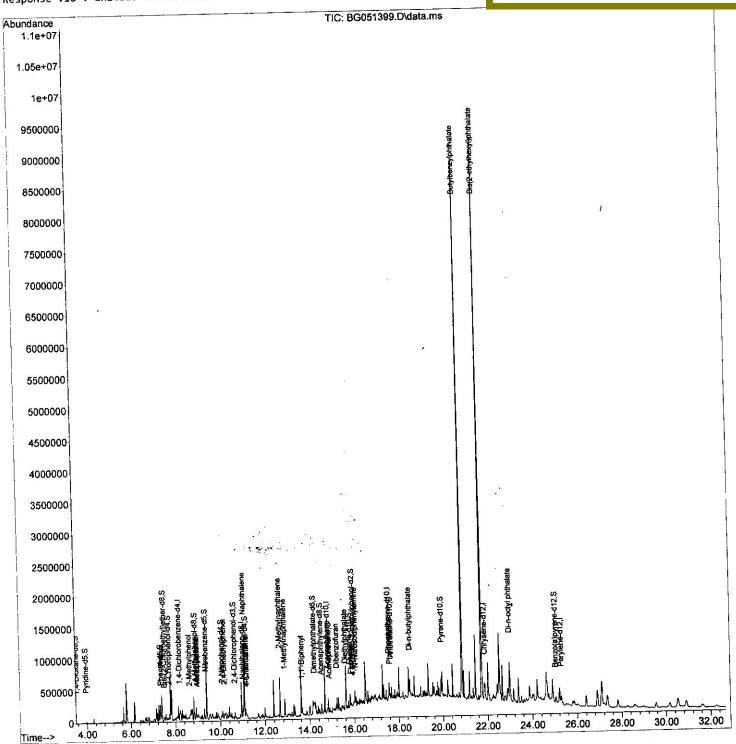
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration

Instrument: BNA\_G ClientSampleId:

# **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 12/08/2021 Supervised By :mohammad ahmed 12/15/2021



SFAM-EPA-BG112321.M Wed Dec 22 00:48:58 2021

# Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120621\

Data File : BG051399.D

: 8 Dec 2021 2:08 Acq On

CG/JU Operator : M4942-03ME Sample

Misc

Sample Multiplier: 1 : 57 ALS Vial

Quant Time: Dec 08 05:31:01 2021

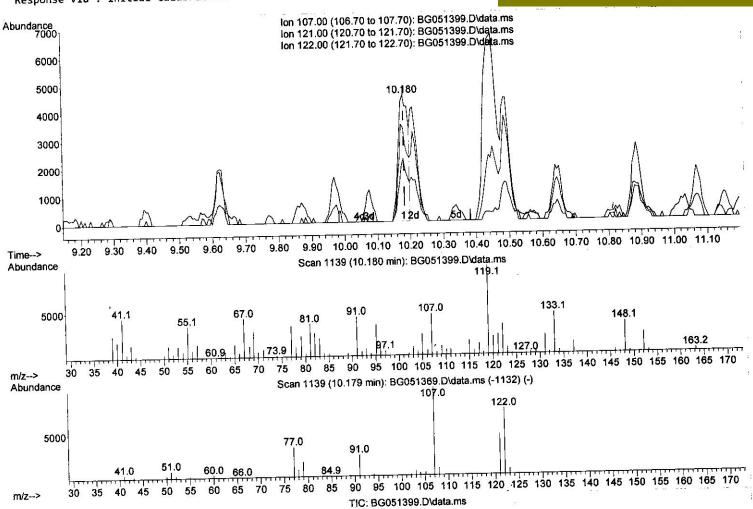
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration

Instrument: BNA\_G ClientSampleId: BGKQ1ME

### Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021 Supervised By:mohammad ahmed 12/15/2021



### (26) 2,4-Dimethylphenol

10.180min (-0.003) 3.58 ng/ul

response	9505		
Ion	Exp%	Act*	
107.00	100.00	100.00	
121.00	49.10	49.98	
122.00	79.60	74.10	
0.00	0.00	0.00	

## Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120621\

Data File : BG051399.D

8 Dec 2021 2:08 Acq On

: CG/JU Operator : M4942-03ME Sample

Misc

Sample Multiplier: 1 : 57 ALS Vial

Quant Time: Dec 08 05:31:01 2021

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

Quant Title : SVOA CALIBRATION

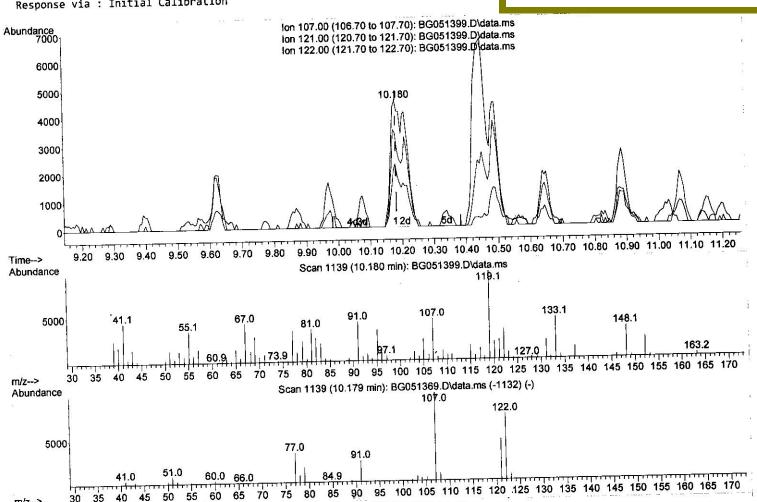
QLast Update : Fri Dec 03 15:23:09 2021

Response via : Initial Calibration



### Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021 Supervised By :mohammad ahmed 12/15/2021



TIC: BG051399.D\data.ms

(26) 2,4-Dimethylphenol

m/z-->

5.87 ng/ul m 10.180min (-0.003)

response	15605		
Ion	Exp%	Act%	
107.00	100.00	100.00	
121.00	49.10	49.98	
122.00	79.60	74.10	
0.00	0.00	0.00	

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120621\

Data File : BG051399.D

Acq On : 8 Dec 2021 2:08

Operator : CG/JU Sample : M4942-03ME

Misc

ALS Vial : 57 Sample Multiplier: 1

Quant Time: Dec 08 05:31:01 2021

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

Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units Dev(	Min)
	. <b></b>	<b>-</b> -			
Internal Standards			24303	20.000 ng/ul	-0.01
<ol> <li>1,4-Dichlorobenzene-d4</li> </ol>	8.189	152	31393	20.000 ng/ul	-0.01
20) Naphthalene-d8	11.015	136	131723	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.822	164	82479	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.572	188	155573	20.000 ng/ul	0.00
79) Chrysene-d12	21.879	240	139337	20.000 ng/ul	0.00
88) Perylene-d12	25.286	264	142106	20.000 Hg/ di	0.00
System Monitoring Compounds			70 202002		0.01
3) 1,4-Dioxane-d8	3.529	96	1930	2.136 ng/uL	-0.01
4) Pyridine-d5	3.964	84	14530	5.481 ng/ul	-0.01
7) Phenol-d5	7.354	99	43766	14.106 ng/ul	0.00
<li>9) Bis-(2-Chloroethyl)eth</li>	7.501	67	29610	15.195 ng/ul	-0.01
11) 2-Chlorophenol-d4	7.719	132	32398	14.501 ng/ul	-0.01
15) 4-Methylphenol-d8	8.906	113	36084	14.412 ng/ul	0.00
21) Nitrobenzene-d5	9.370	128	18083	16.263 ng/ul	0.00
24) 2-Nitrophenol-d4	10.092	143	19888	15.856 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.645	165	31889	14.984 ng/ul	0.00
31) 4-Chloroaniline-d4	11,156	131	40675	13.062 ng/ul	0.00
46) Dimethylphthalate-d6	14.211	166	95145	14.992 ng/ul	-0.01
49) Acenaphthylene-d8	14.517	160	125721	15.710 ng/ul	-0.01
54) 4-Nitrophenol-d4	15.028	143	601	0.585 ng/ul	-0.02
60) Fluorene-d10	15.815	176	81999	14.348 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.956	200	4090	4.260 ng/ul	0.00
73) Anthracene-d10	17.672	188	119803	16.101 ng/ul	0.00
81) Pyrene-d10	19.951	212	130016	15.421 ng/ul	0.00
92) Benzo(a)pyrene-d12	25.045	264	116698	15.376 ng/ul	0.00
Target Compounds Qvalue					
8) Phenol	7.378	94	167206	52.021 ng/ul	97
13) 2-Methylphenol	8.635		24147	10.086 ng/ul	98
16) Acetophenone	8.982	105	31361	8.098 ng/ul#	
18) 4-Methylphenol	8.970		18983	ን 7.415 ng/ul	92
26) 2,4-Dimethylphenol	10.180	er reconstruction	15605n		
30) Naphthalene	11.068		977235		98
32) 4-Chloroaniline	11.179		6606	2.113 ng/ul#	
36) 2-Methylnaphthalene	12.660		292150	59.927 ng/ul	97
37) 1-Methylnaphthalene	12.877		135436	27.003 ng/ul	99
43) 1,1'-Biphenyl	13.653			7.679 ng/ul	96
52) Acenaphthene	14.887			4.523 ng/ul	95
56) Dibenzofuran	15.222			5.416 ng/ul	98
59) Diethylphthalate	15.609			3.215 ng/ul	100
61) Fluorene	15.868		23190	. 3.849 ng/ul:	# 91
67) N-Nitrosodiphenylamine	16.06			4.294 ng/ul	# 71
72) Phenanthrene	17.61				# 96
78) Di-n-butylphthalate	18.50				99
83) Butylbenzylphthalate	20.87			758.048 ng/ul	
86) Bis(2-ethylhexyl)phtha					
89) Di-n-octyl phthalate	22.98				100
			<b></b>		

Instrument: BNA\_G ClientSampleId: BGKQ1ME

### **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 12/08/2021 Supervised By :mohammad ahmed 12/15/2021

11/22/21

<sup>(#)</sup> = qualifier out of range (m) = manual integration (+) = signals summed