Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051368.D

3:19 Acq On : 7 Dec 2021

Operator : CG/JU : M4942-05 Sample

Misc

ALS Vial : 25 Sample Multiplier: 1

Quant Time: Dec 07 03:53:46 2021

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

Quant Title : SVOA CALIBRATION

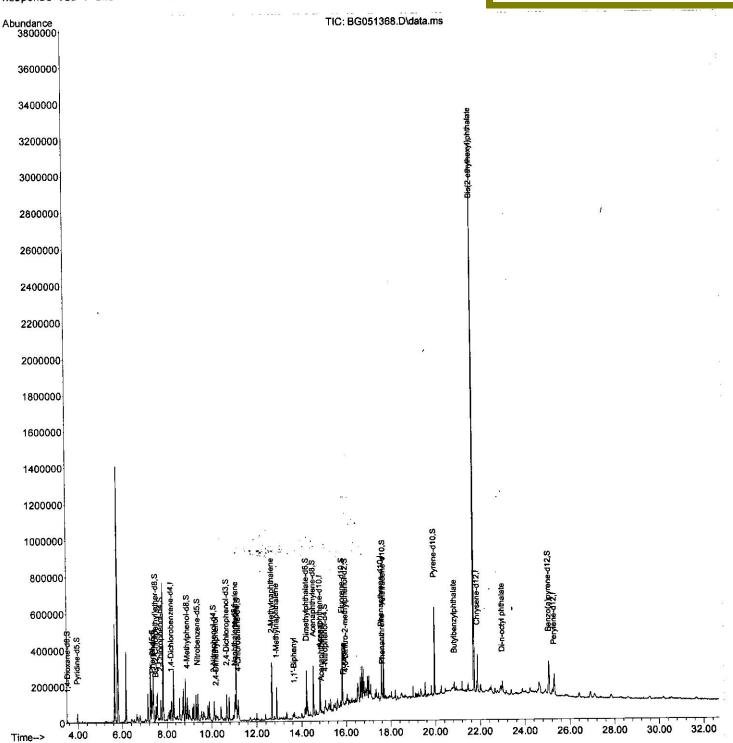
QLast Update : Fri Dec 03 15:23:09 2021

Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/07/2021 Supervised By :mohammad ahmed 12/07/2021



SFAM-EPA-BG112321.M Wed Dec 08 09:51:23 2021

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051368.D

Acq On : 7 Dec 2021 3:19

Operator : CG/JU Sample : M4942-05

Misc

ALS Vial : 25 Sample Multiplier: 1

Quant Time: Dec 07 03:53:46 2021

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

Quant Title : SVOA CALIBRATION

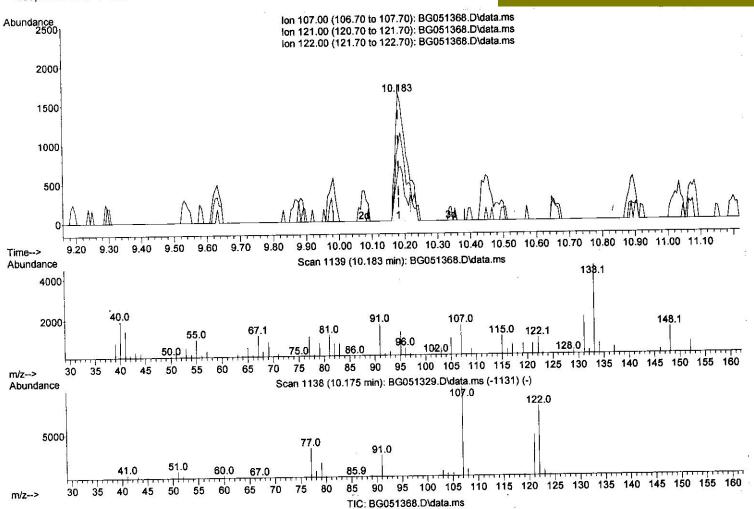
QLast Update : Fri Dec 03 15:23:09 2021

Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/07/2021 Supervised By :mohammad ahmed 12/07/2021



. .

(26) 2,4-Dimethylphenol

10.183min (-0.000) 1.50 ng/ul

response	3487	
Ion	Ежр%	Act*
107.00	100.00	100.00
121.00	49.10	42.14
122.00	79.60	60.64#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051368.D

; 7 Dec 2021 3:19 Acq On

Operator : CG/JU Sample : M4942-05

Misc

Sample Multiplier: 1 : 25 ALS Vial

Quant Time: Dec 07 03:53:46 2021

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

Quant Title : SVOA CALIBRATION

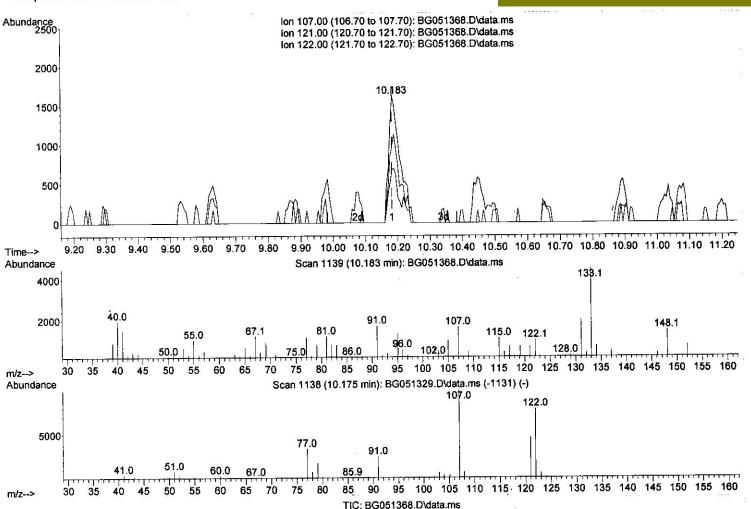
QLast Update : Fri Dec 03 15:23:09 2021

Response via: Initial Calibration

Instrument: BNA_G ClientSampleId:

Manual IntegrationsAPPROVED

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(26) 2,4-Dimethylphenol

10.183min (-0.000) 1.71 ng/ul m 3

response	3967	
Ion	Exp%	Act*
107.00	100.00	100.00
121.00	49.10	42.14
122.00	79.60	60.64#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

BG051368.D Data File :

7 Dec 2021 3:19 Acq On

: CG/JU Operator Sample M4942-05

Misc

Sample Multiplier: 1 : 25 ALS Vial

Quant Time: Dec 07 03:53:46 2021

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

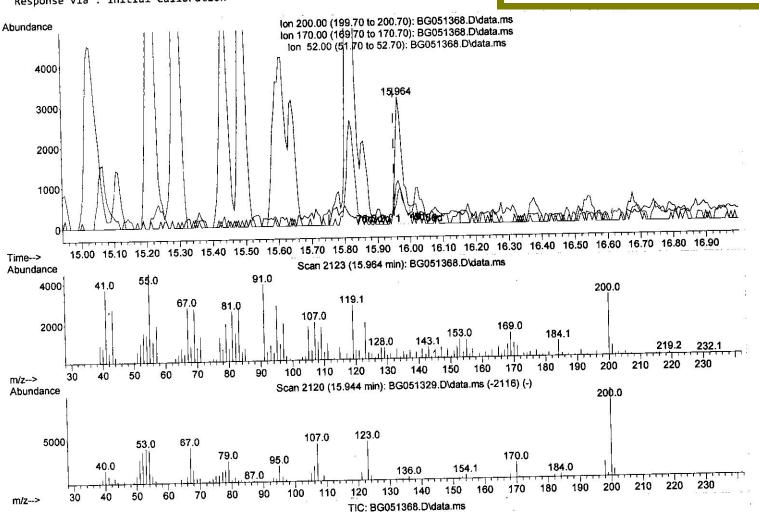
Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:09 2021

Response via : Initial Calibration



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4,6-Dinitro-2-methylphenol-d2 (S)

5.46 ng/ul 15.964min (+ 0.011)

response	4892			
Ion	Ежр%	Acts		
200.00	100.00	100.00		
170.00	19.80	26.11#		
52.00	47.40	34.56#		
0.00	0.00	0.00		

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051368.D

Acq On : 7 Dec 2021 3:19

Operator : CG/JU Sample : M4942-05

Misc

ALS Vial : 25 Sample Multiplier: 1

Quant Time: Dec 07 03:53:46 2021

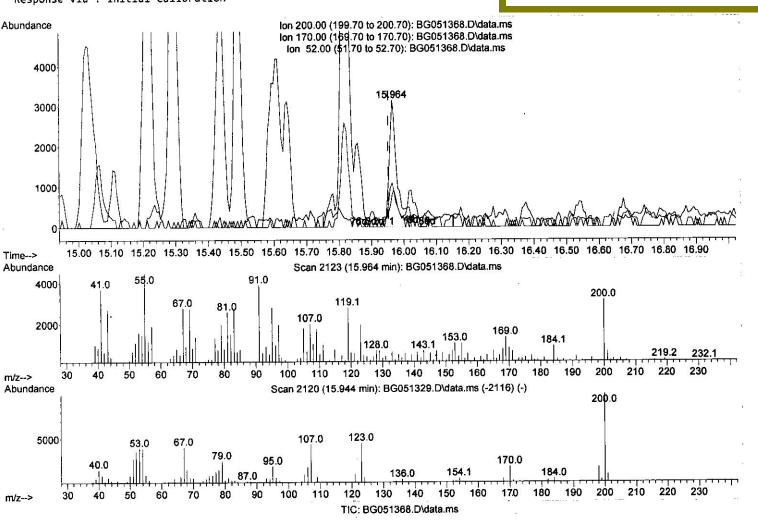
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION
OLast Update : Fri Dec 03 15:23:09 :

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

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/07/2021 Supervised By :mohammad ahmed 12/07/2021



response	5463	
Ion	Ехр%	Act%
200.00	100.00	100.00
170.00	19.80	26.11#
52.00	47.40	34.56#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051368.D

: 7 Dec 2021 3:19 Acq On

Operator : CG/JU : M4942-05 Sample

Misc

ALS Vial : 25 Sample Multiplier: 1

Quant Time: Dec 07 03:53:46 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

nstrument :
BNA_G
ClientSampleId:
BGKQ5

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/07/2021 Supervised By :mohammad ahmed 12/07/2021

Compound	R.T.	QIon	Response	Conc Unit	ts Dev(Min)	
Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.197	152	26127	20.000 1	- 10 Th	0.00	
20) Naphthalene-d8	11.023	136	115036	20.000		0.00	
38) Acenaphthene-d10	14.830	164	75850	20.000		0.00	
64) Phenanthrene-d10	17.580	188	145229	20.000		0.00	
79) Chrysene-d12	21.881	240	126733	20.000		0.00	
88) Perylene-d12	25.294	264	129028	20.000	ng/ul	0.01	
System Monitoring Compounds					200	70 012	
3) 1,4-Dioxane-d8	3.532	96	3168	4.214	0.	-0.01	
4) Pyridine-d5	3.966	84	33371	15.126	•	-0.01	
7) Phenol-d5	7.357	99	63085	24.430		0.00	
9) Bis-(2-Chloroethyl)eth	7.503	67	42615	26.277		-0.01	X
11) 2-Chlorophenol-d4	7.727	132	49088	26.399		0.00	
15) 4-Methylphenol-d8	8.908	113	51919	24.916		0.00	
21) Nitrobenzene-d5	9.372	128	26930	27.732		0.00	
24) 2-Nitrophenol-d4	10.100	143	30156	27.529		0.00	
28) 2,4-Dichlorophenol-d3	10.653	165	491.02	26.419		0.00	21
31) 4-Chloroaniline-d4	11.164	131	66427	24.427		0.00	
46) Dimethylphthalate-d6	14.219		163343	27.988		0.00	
49) Acenaphthylene-d8	14.525	160	205923	27.981		0.00	
54) 4-Nitrophenol-d4	15.059	143	18888	19.994		0.01	24 12121
60) Fluorene-d10	15.823	176	143947	27.390		0.00	2/2/20181
65) 4,6-Dinitro-2-methylph	15.964	200	5463m '			0.01	
73) Anthracene-d10	17.680	188	213653 ~	30.760		0.00	
81) Pyrene-d10	19.959	212	223887	29,196		0.00	
92) Benzo(a)pyrene-d12	25.059	264	197363	28.641	ng/ul	0.02	
Target Compounds						alue	27 -111
8) Phenol	7.386	94	2869 7		ng/ul#	1	12/22/11
26) 2,4-Dimethylphenol	10.183	107	3967m(ng/ul		1
30) Naphthalene	11.076	128	293910 ~			98	
36) 2-Methylnaphthalene	12.668	142	147773	34.709		97	
37) 1-Methylnaphthalene	12.885	142	83427			99	
43) 1,1'-Biphenyl	13.661	. 154	17244		ng/ul	93	*
52) Acenaphthene	14.889	153		1.290		95	
61) Fluorene	15.876		6025		ng/ul#		
72) Phenanthrene	17.621	178		2,410			
83) Butylbenzylphthalate	20.847	149			ng/ul	98	
86) Bis(2-ethylhexyl)phtha	21.722	149		215.828			
89) Di-n-octyl phthalate	22.989	149	81759	8.746	ng/ul	100	9.7
		-					. 1

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