Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\

Data File : BG050998.D

Acq On : 12 Nov 2021 13:01

Operator : CG/JU Sample : M4615-06

Misc

ALS Vial : 37 Sample Multiplier: 1

Quant Time: Nov 12 14:12:33 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M

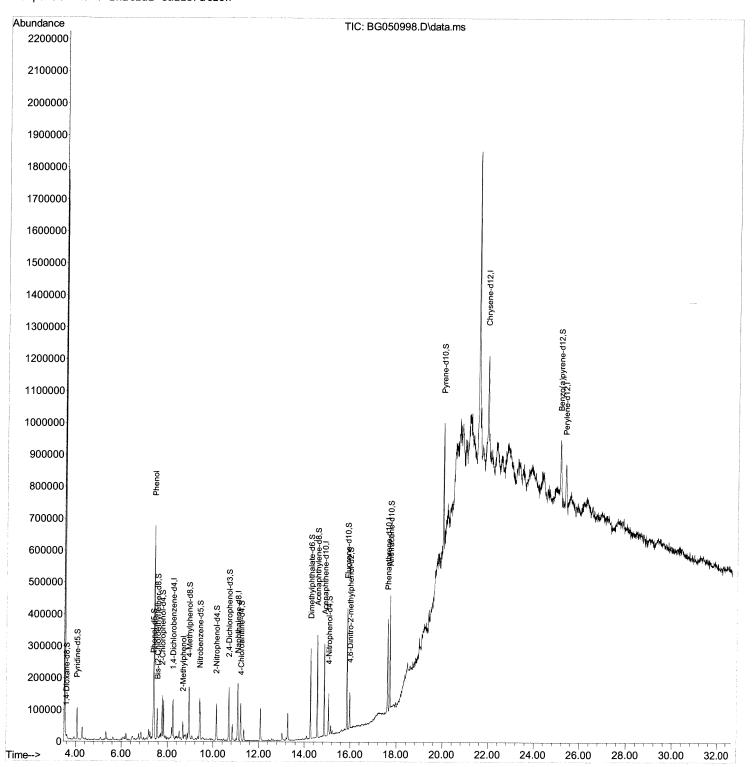
Quant Title : SVOA CALIBRATION QLast Update : Thu Nov 11 12:40:48 2021

Response via : Initial Calibration

Instrument : BNA_G ClientSampleId : C0V12

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/12/2021 Supervised By :mohammad ahmed 11/17/2021



Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\

Data File : BG050998.D

Acq On : 12 Nov 2021 13:01

Operator : CG/JU Sample : M4615-06

Misc

ALS Vial : 37 Sample Multiplier: 1

Quant Time: Nov 12 14:12:33 2021

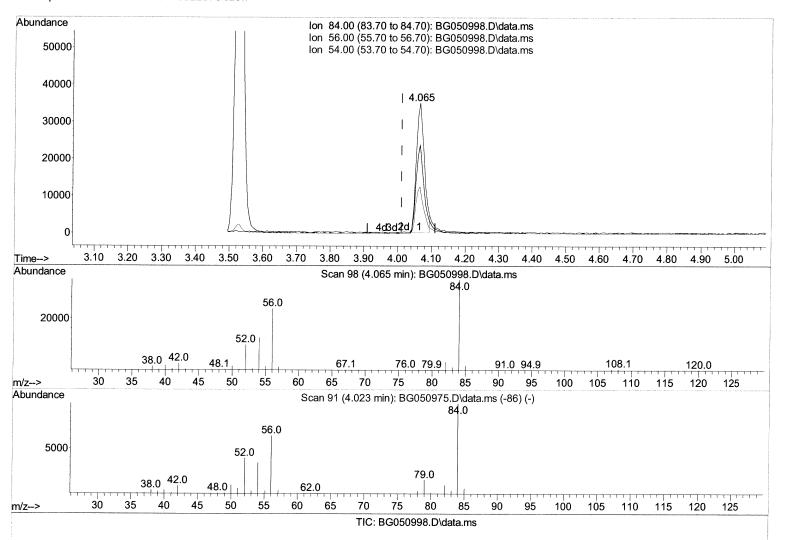
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Nov 11 12:40:48 2021 Response via : Initial Calibration Instrument:
BNA_G
ClientSampleId:
C0V12

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/12/2021 Supervised By :mohammad ahmed 11/17/2021



(4) Pyridine-d5 (S)

4.065min (+ 0.054) 19.14 ng/ul

response	58670	
Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	68.07
54.00	31.50	35.72
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\

Data File : BG050998.D

Acq On : 12 Nov 2021 13:01

Operator : CG/JU Sample : M4615-06

Misc

ALS Vial : 37 Sample Multiplier: 1

Quant Time: Nov 12 14:12:33 2021

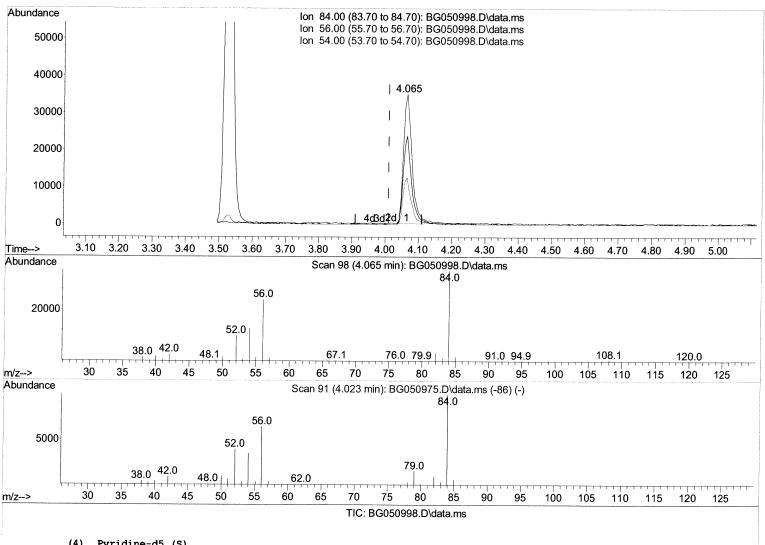
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Nov 11 12:40:48 2021 Response via : Initial Calibration Instrument: BNA_G ClientSampleId: C0V12

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/12/2021 Supervised By :mohammad ahmed 11/17/2021



(4) Pyridine-d5 (S)

response	61676		
Ion	Ехр%	Act%	
84.00	100.00	100.00	
56.00	68.00	68.07	
54.00	31.50	35.72	
0.00	0.00	0.00	

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\

Data File : BG050998.D

Acq On : 12 Nov 2021 13:01

Operator : CG/JU Sample : M4615-06

Misc :

ALS Vial : 37 Sample Multiplier: 1

Quant Time: Nov 12 14:12:33 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Nov 11 12:40:48 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleId :

C0V12

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/12/2021 Supervised By: mohammad ahmed 11/17/2021

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)	
Internal Standards						
 1,4-Dichlorobenzene-d4 	8.242	152	33073	20.000 ng/ul	0.02	
20) Naphthalene-d8	11.074	136	153371	20.000 ng/ul	0.02	
38) Acenaphthene-d10	14.870	164	99073	20.000 ng/ul	0.02	
64) Phenanthrene-d10	17.614	188	183374	20.000 ng/ul	0.02	
79) Chrysene-d12	21.921	240	155872	20.000 ng/ul	0.03	
88) Perylene-d12	25.340	264	162036	20.000 ng/ul	0.06	
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.601	96	4013	3.916 ng/uL	0.01	
4) Pyridine-d5	4.065	84	61676m>	20.119 ng/ul>	· 0.05 11/17/21 JU	
7) Phenol-d5	7.391	99	80389	22.784 ng/ul		
9) Bis-(2-Chloroethyl)eth	7.561	67	47022	20.631 ng/ul	0.02	
<pre>11) 2-Chlorophenol-d4</pre>	7.772	132	58611	23.970 ng/ul	0.02	
<pre>15) 4-Methylphenol-d8</pre>	8.942	113	65171	23.463 ng/ul	0.02	
21) Nitrobenzene-d5	9.418	128	31422	24.109 ng/ul	0.02	
24) 2-Nitrophenol-d4		143	35610	24.572 ng/ul	0.01	
28) 2,4-Dichlorophenol-d3	10.687	165	60581	24.816 ng/ul	0.02	
31) 4-Chloroaniline-d4	11.210	131	65572	17.737 ng/ul	0.02	
46) Dimethylphthalate-d6	14.265	166	187309	24.712 ng/ul	0.02	
<pre>49) Acenaphthylene-d8</pre>	14.565	160	237423	25.142 ng/ul	0.02	
54) 4-Nitrophenol-d4	15.058	143	30646	22.299 ng/ul	0.02	
60) Fluorene-d10	15.857	176	152371	22.693 ng/ul	0.02	
65) 4,6-Dinitro-2-methylph	15.975	200	23587	21.216 ng/ul	0.02	
73) Anthracene-d10	17.714	188	214271	24.715 ng/ul	0.02	
81) Pyrene-d10	19.993	212	222570	22.108 ng/ul	0.02	
92) Benzo(a)pyrene-d12	25.105	264	204848	22.869 ng/ul	0.05	
Target Compounds				Ovalue		
8) Phenol	7.420	94	384174	105.257 ng/ul	96	
13) 2-Methylphenol	8.677	108		6.485 ng/ul	93	

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