Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\

Data File : BM033130.D

Acq On : 17 Nov 2021 21:36

Operator : CG/JU Sample : PB140812BS

Misc :

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 18 00:36:28 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M

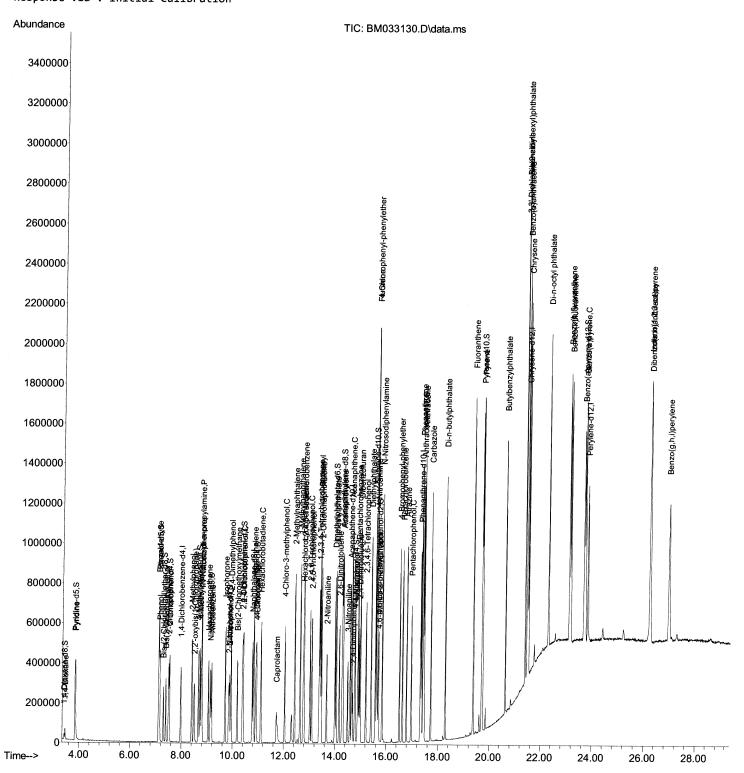
Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 17 14:14:11 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/18/2021 Supervised By :mohammad ahmed 11/26/2021



SFAM-EPA-BM111721.M Thu Nov 18 01:07:22 2021

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\

Data File : BM033130.D

Acq On : 17 Nov 2021 21:36

Operator : CG/JU Sample : PB140812BS

Misc

ALS Vial : 7 Sample Multiplier: 1

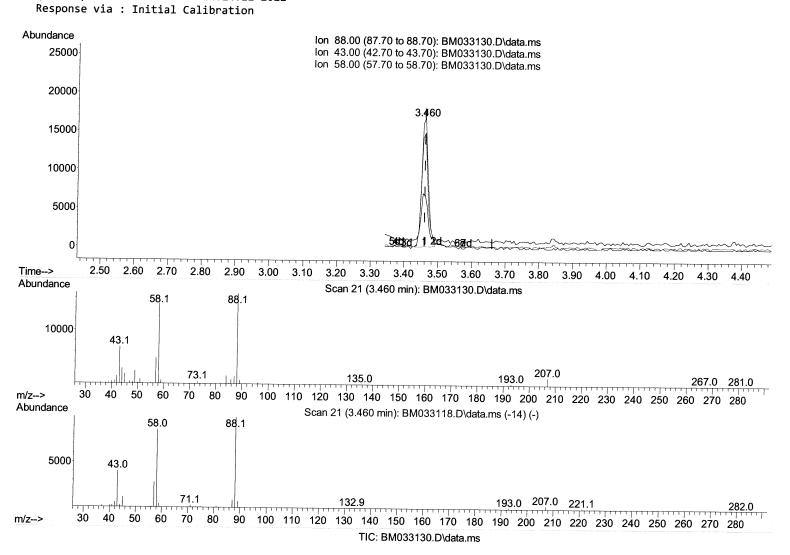
Quant Time: Nov 18 00:36:28 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M

Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 17 14:14:11 2021 Instrument:
BNA_M
ClientSampleId:
SLCS812

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/18/2021 Supervised By: mohammad ahmed 11/26/2021



(2) 1,4-Dioxane

3.460min (+ 0.000) 9.17 ng/uL

response	22673			
Ion	Ехр%	Act%		
88.00	100.00	100.00		
43.00	43.60	40.97		
58.00	87.50	88.85		
0.00	0.00	0.00		

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\

Data File : BM033130.D

Acq On : 17 Nov 2021 21:36

Operator : CG/JU Sample : PB140812BS

Misc

ALS Vial : 7 Sample Multiplier: 1

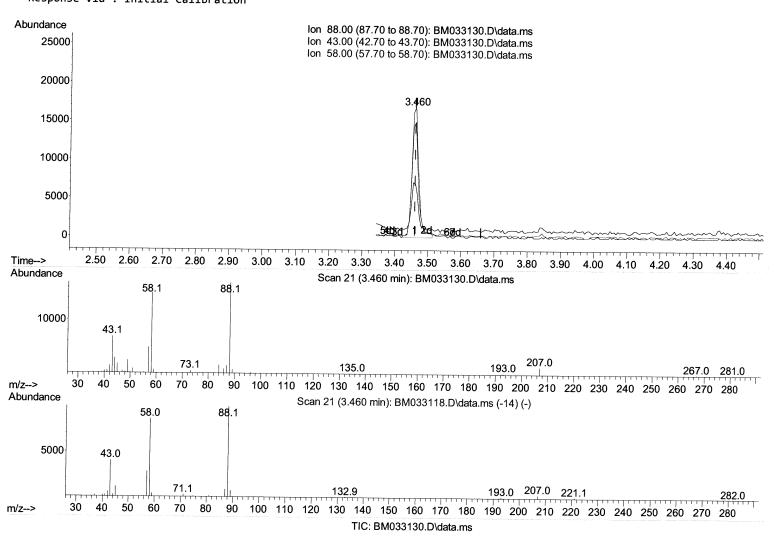
Quant Time: Nov 18 00:36:28 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M

Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 17 14:14:11 2021 Response via : Initial Calibration Instrument:
BNA_M
ClientSampleId:
SLCS812

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/18/2021 Supervised By :mohammad ahmed 11/26/2021



(2) 1,4-Dioxane

3.460min (+ 0.000) 9.84 ng/uL m 11/29/2174

response	24324	
Ion	Ехр%	Act%
88.00	100.00	100.00
43.00	43.60	40.97
58.00	87.50	88.85
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\

Data File : BM033130.D

Acq On : 17 Nov 2021 21:36

Operator : CG/JU Sample : PB140812BS

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 18 00:36:28 2021

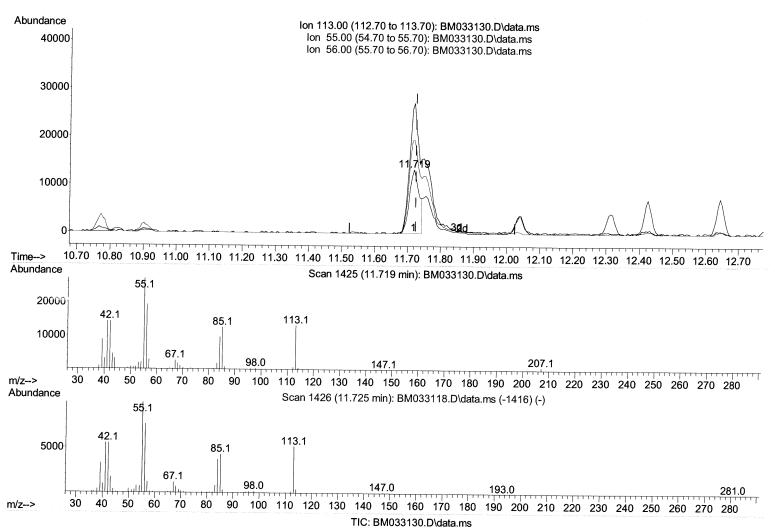
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M

Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 17 14:14:11 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/18/2021 Supervised By :mohammad ahmed 11/26/2021



(34) Caprolactam

11.719min (-0.006) 15.62 ng/ul

response	27964			
Ion	Ежр%	Act%		
113.00	100.00	100.00		
55.00	196.60	203.90		
56.00	147.80	146.44		
0.00	0.00	0.00		

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\

Data File : BM033130.D

Acq On : 17 Nov 2021 21:36

Operator : CG/JU Sample : PB140812BS

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 18 00:36:28 2021

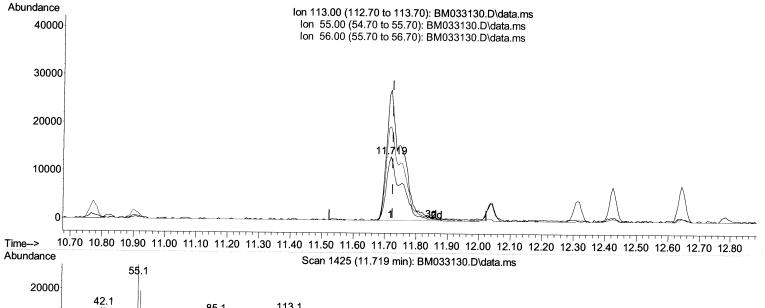
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M

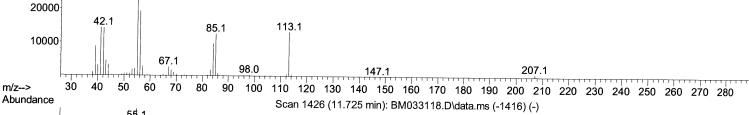
Quant Title : SVOA CALIBRATION

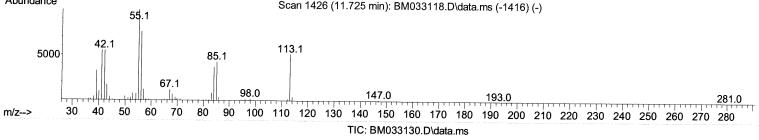
QLast Update : Wed Nov 17 14:14:11 2021 Response via : Initial Calibration Instrument :
BNA_M
ClientSampleId :
SLCS812

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/18/2021 Supervised By: mohammad ahmed 11/26/2021







(34) Caprolactam

11.719min (-0.006) 25.47 ng/ul m 1/29/2/10

response	45603		
Ion	Ехр%	Act%	
113.00	100.00	100.00	
55.00	196.60	203.90	
56.00	147.80	146.44	
0.00	0.00	0.00	

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\

Data File : BM033130.D Acq On : 17 Nov 2021 21:36

Operator : CG/JU Sample : PB140812BS

Misc :

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 18 00:36:28 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 17 14:14:11 2021 Response via : Initial Calibration Instrument: BNA_M ClientSampleId: SLCS812

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/18/2021 Supervised By :mohammad ahmed 11/26/2021

Compound	R.T.	QIon	Response	Conc Units Dev	/(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.972	152	93834	20 000 ng/ul	0.00
20) Naphthalene-d8	10.772		389989	20.000 ng/ul 20.000 ng/ul	0.00 0.00
38) Acenaphthene-d10	14.589		269208	20.000 ng/ul	
64) Phenanthrene-d10	17.330		580622	20.000 ng/ul	0.00
79) Chrysene-d12	21.483		568750		0.00
88) Perylene-d12	23.830		561689	20.000 ng/ul 20.000 ng/ul	0.00 0.00
• • •		204	301003	20.000 Hg/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.425	96	11309	4.646 ng/uL	0.00
4) Pyridine-d5	3.843	84	161513	24.090 ng/ul	0.00
7) Phenol-d5	7.125	99	201346	25.368 ng/ul	0.00
Bis-(2-Chloroethyl)eth	7.301	67	127316	25.290 ng/ul	0.00
11) 2-Chlorophenol-d4	7.507	132	153798	25.583 ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.672	113	160571	26.068 ng/ul	0.00
21) Nitrobenzene-d5	9.136	128	73077	26.099 ng/ul	0.00
24) 2-Nitrophenol-d4	9.854	143	76112	27.142 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.389		166977	26.002 ng/ul	0.00
31) 4-Chloroaniline-d4	10.907	131	196929	23.060 ng/ul	0.00
46) Dimethylphthalate-d6	14.001	166	513845	26.023 ng/ul	0.00
49) Acenaphthylene-d8	14.289		641875	25.232 ng/ul	0.00
54) 4-Nitrophenol-d4	14.771	143	83355	25.915 ng/ul	0.00
60) Fluorene-d10	15.577	176	465358	26.271 ng/ul	0.00
65) 4,6-Dinitro-2-methylph		200	71669	26.265 ng/ul	0.00
73) Anthracene-d10	17.424	188	726099	25.943 ng/ul	0.00
81) Pyrene-d10	19.706	212	859193	25.564 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.677	264	796202	26.412 ng/ul	0.00
Target Compounds				0.45	ılue
2) 1,4-Dioxane	3.460	88	24324m>		
5) Pyridine	3.860	79	165423	9.843 ng/uL > 24.179 ng/ul	' '' ''
6) Benzaldehyde	7.119	77	126220	27.903 ng/ul	99
8) Phenol	7.154	94	204716	25.900 ng/ul	99
10) Bis(2-Chloroethyl)ether	7.395	93	161959	25.795 ng/ul	99
12) 2-Chlorophenol	7.537	128	161509	26.029 ng/ul	98
13) 2-Methylphenol	8.407	108	156368	25.825 ng/ul	99
14) 2,2'-oxybis(1-Chloropr	8.501	45	263163	27.138 ng/ul	99 98
16) Acetophenone	8.795	105	257568	_	
17) N-Nitroso-di-n-propyla	8.784	70	141736	26.492 ng/ul 27.057 ng/ul	98
18) 4-Methylphenol	8.731	108	167056		98
19) Hexachloroethane	9.054	117	74159	26.319 ng/ul	99
22) Nitrobenzene	9.178	77	206392	26.235 ng/ul	96
23) Isophorone	9.707	82	378095	26.413 ng/ul	98
25) 2-Nitrophenol	9.883	139	83843	26.473 ng/ul	98
26) 2,4-Dimethylphenol	9.936	107	207013	28.208 ng/ul	96
27) Bis(2-Chloroethoxy)met	10.178	93	219046	26.507 ng/ul	97
29) 2,4-Dichlorophenol	10.178	162	164278	25.772 ng/ul	97 06
30) Naphthalene	10.413	128	531584	26.382 ng/ul	96
32) 4-Chloroaniline	10.930	127		25.633 ng/ul	99
33) Hexachlorobutadiene	11.101	225	197301	22.870 ng/ul	97
34) Caprolactam	11.719	113	124061	25.577 ng/ul	99
35) 4-Chloro-3-methylphenol	12.036	107	45603m > 183887	G,,	11/54/5/29
,	-2.000	10/	10300/	27.119 ng/ul	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\

Data File : BM033130.D

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Misc :

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 18 00:36:28 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 17 14:14:11 2021 Response via : Initial Calibration Instrument: BNA_M ClientSampleId: SLCS812

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/18/2021 Supervised By: mohammad ahmed 11/26/2021

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
36) 2-Methylnaphthalene	12.425	142	381308	26.537 ng/ul	99
37) 1-Methylnaphthalene	12.642	142	384204	26.193 ng/ul	100
39) 1,2,4,5-Tetrachloroben	12.783		222999	24.898 ng/ul	99
40) Hexachlorocyclopentadiene	12.766	237	145439	23.077 ng/ul	96
41) 2,4,6-Trichlorophenol	13.024	196	139521	26.450 ng/ul	100
42) 2,4,5-Trichlorophenol	13.095	196	149101	26.361 ng/ul	97
43) 1,1'-Biphenyl	13.430	154	527456	25.147 ng/ul	98
44) 2-Chloronaphthalene	13.472	162	405391	25.062 ng/ul	99
45) 2-Nitroaniline	13.671	65	122426	28.715 ng/ul	95
47) Dimethylphthalate	14.048	163	506845	26.336 ng/ul	99
48) 2,6-Dinitrotoluene	14.166	165	93863	28.505 ng/ul	98
50) Acenaphthylene	14.319	152	660429	25.584 ng/ul	100
51) 3-Nitroaniline	14.495	138	83900	25.288 ng/ul	92
52) Acenaphthene	14.654	153	435500	25.861 ng/ul	99
53) 2,4-Dinitrophenol	14.695	184	43368	25.804 ng/ul	93
55) 4-Nitrophenol	14.789	109	87539	26.647 ng/ul	96
56) Dibenzofuran	14.989	168	645528	26.155 ng/ul	100
57) 2,4-Dinitrotoluene	14.948	165	138062	30.487 ng/ul#	99
58) 2,3,4,6-Tetrachlorophenol	15.207	232	126845	27.474 ng/ul	95
59) Diethylphthalate	15.407	149	534591	27.733 ng/ul	99
61) Fluorene	15.636	166	520141	26.514 ng/ul	99
62) 4-Chlorophenyl-phenyle	15.630	204	272844	26.647 ng/ul	98
63) 4-Nitroaniline	15.648	138	94100	28.844 ng/ul	99
66) 4,6-Dinitro-2-methylph	15.707	198	72152	26.382 ng/ul	100
67) N-Nitrosodiphenylamine	15.842	169	453589	26.561 ng/ul	99
68) 4-Bromophenyl-phenylether	16.518	248	172347	26.692 ng/ul	96
69) Hexachlorobenzene	16.630	284	190569	25.763 ng/ul	97
70) Atrazine	16.789	200	170107	25.779 ng/ul	99
71) Pentachlorophenol	16.971	266	109650	25.645 ng/ul	99
72) Phenanthrene	17.371	178	850473	26.443 ng/ul	98
74) Anthracene	17.459	178	858072	26.604 ng/ul	99
75) 1,2,3,4-Tetrachloroben	13.395	216	228040	23.911 ng/uL	97
76) Pentachlorobenzene	14.907	250	226867	24.265 ng/uL	99
77) Carbazole	17.724	167	765352	26.711 ng/ul	99
78) Di-n-butylphthalate	18.289	149	901921	28.787 ng/ul	100
80) Fluoranthene	19.371	202	1018722	25.872 ng/ul	99
82) Pyrene	19.736	202	1040501	26.026 ng/ul	99
83) Butylbenzylphthalate	20.624	149	383042	28.043 ng/ul	96
84) 3,3'-Dichlorobenzidine	21.395	252	307656	24.011 ng/ul	99
85) Benzo(a)anthracene86) Bis(2-ethylhexyl)phtha	21.465	228	984603	26.766 ng/ul	99
87) Chrysene	21.389	149	550331	28.425 ng/ul	99
89) Di-n-octyl phthalate	21.518	228	957115	26.643 ng/ul	99
90) Benzo(b)fluoranthene	22.300	149	929500	27.159 ng/ul	100
91) Benzo(k)fluoranthene	23.118	252	1026976	27.148 ng/ul	98
93) Benzo(a)pyrene	23.165	252	912577	26.334 ng/ul	99
94) Indeno(1,2,3-cd)pyrene	23.730	252	953024	26.657 ng/ul	100
95) Dibenzo(a,h)anthracene	26.241	276	1050728	26.392 ng/ul	97
96) Benzo(g,h,i)perylene	26.253 26.977	278 276	902094 914759	26.485 ng/ul 26.473 ng/ul	98 97
			J14/JJ	iig/ul	97

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