Data File : BM033364.D

Acq On : 09 Dec 2021 19:12

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 18 Sample Multiplier: 1

Quant Time: Dec 10 02:47:03 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

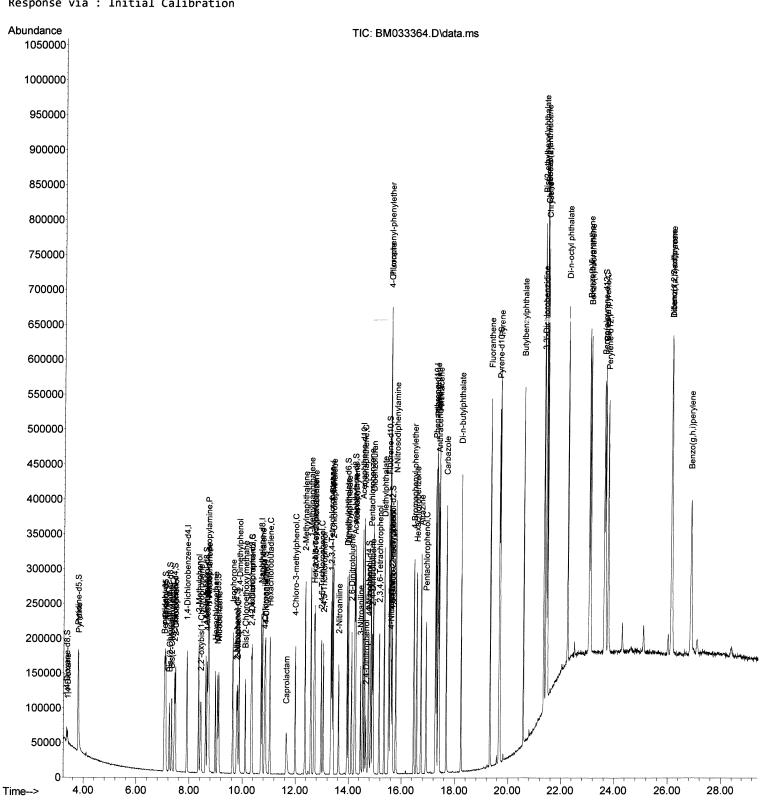
QLast Update : Thu Dec 09 13:25:37 2021

Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



Data File : BM033364.D

Acq On : 09 Dec 2021 19:12

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 18 Sample Multiplier: 1

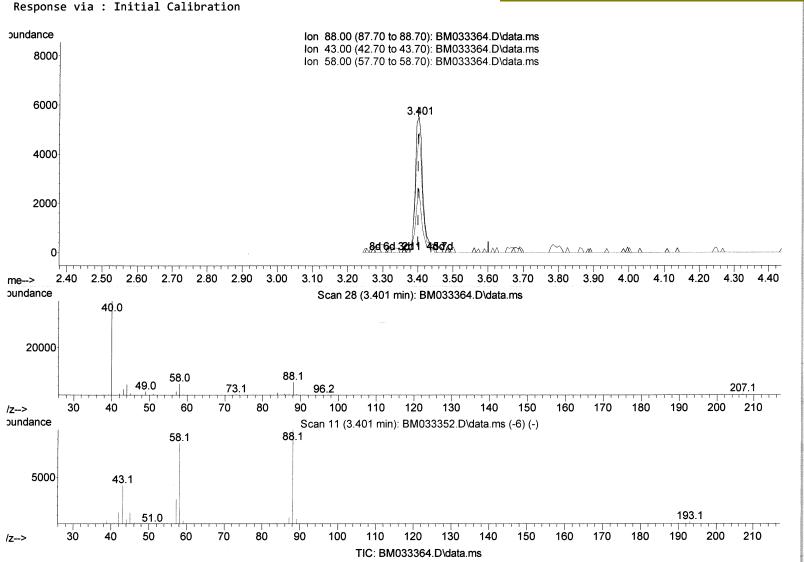
Quant Time: Dec 10 02:47:03 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:25:37 2021 Instrument :
BNA_M
LabSampleId :
SSTDCCC020

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



(2) 1,4-Dioxane

3.401min (+ 0.000) 6.62 ng/uL

response	8186	
Ion	Ежр%	Act%
88.00	100.00	100.00
43.00	45.30	47.40
58.00	85.60	88.53
0.00	0.00	0.00

Data File : BM033364.D

Acq On : 09 Dec 2021 19:12

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 18 Sample Multiplier: 1

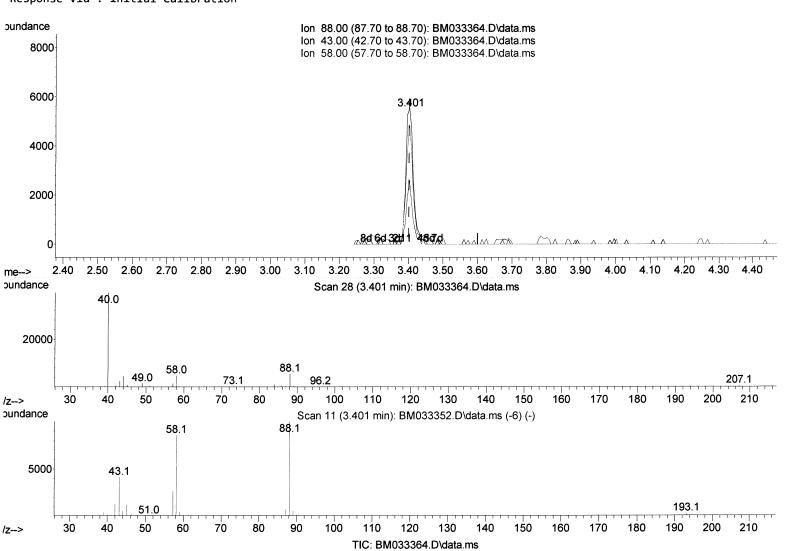
Quant Time: Dec 10 02:47:03 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument:
BNA_M
LabSampleId:
SSTDCCC020

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 12/10/2021 Supervised By: mohammad ahmed 12/15/2021



(2) 1,4-Dioxane

3.401min	(+ 0.000)	6.99 ng/u	"Ju 2/23/27
response	8640		garring
Ion	Exp %	Act%	
88.00	100.00	100.00	
43.00	45.30	47.40	
58.00	85.60	88.53	
0 00	0.00	0.00	

Data File: BM033364.D

Acq On : 09 Dec 2021 19:12

Operator : CG/JU Sample : SSTDCCC020

Misc :

ALS Vial : 18 Sample Multiplier: 1

Quant Time: Dec 10 02:47:03 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

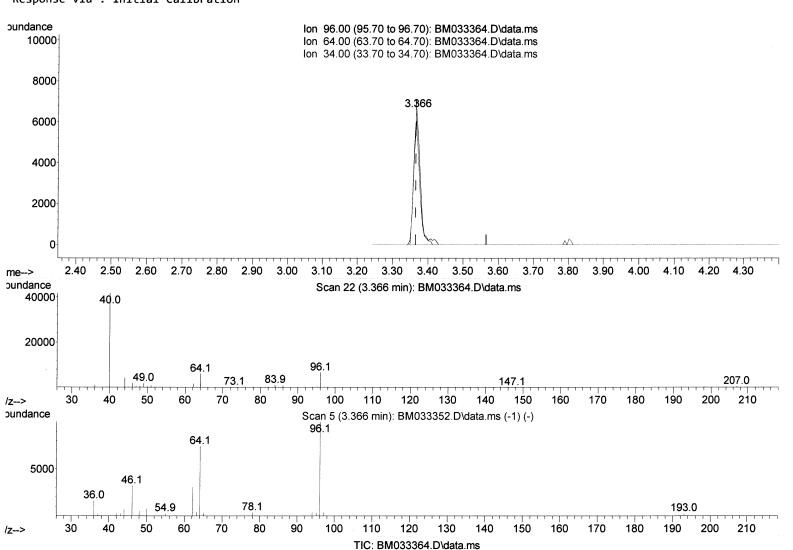
Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration

Instrument:
BNA_M
LabSampleId:
SSTDCCC020

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 12/10/2021 Supervised By: mohammad ahmed 12/15/2021



(3) 1,4-Dioxane-d8 (S)

3.366min (+ 0.000) 7.37 ng/uL

response	8188				
Ion	Exp%	Act%			
96.00	100.00	100.00			
64.00	74.20	91.25#			
34.00	0.00	0.00			
0.00	0.00	0.00			

Data File: BM033364.D

Acq On : 09 Dec 2021 19:12

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 18 Sample Multiplier: 1

Quant Time: Dec 10 02:47:03 2021

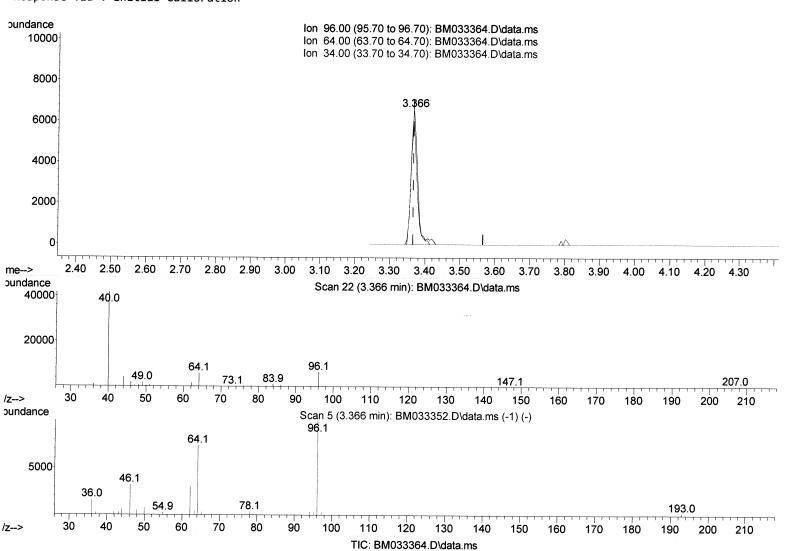
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument:
BNA_M
LabSampleId:
SSTDCCC020

Manual Integrations APPROVED

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(3) 1,4-Dioxane-d8 (S)

3.366min	(+ 0.000)	7.53 ng/uL	"Ju 2 20 2)
response	8368		0-(1-1
Ion	Ехр%	Act%	
96.00	100.00	100.00	
64.00	74.20	91.25#	
34.00	0.00	0.00	
0.00	0.00	0.00	

Data File: BM033364.D

Acq On : 09 Dec 2021 19:12

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 18 Sample Multiplier: 1

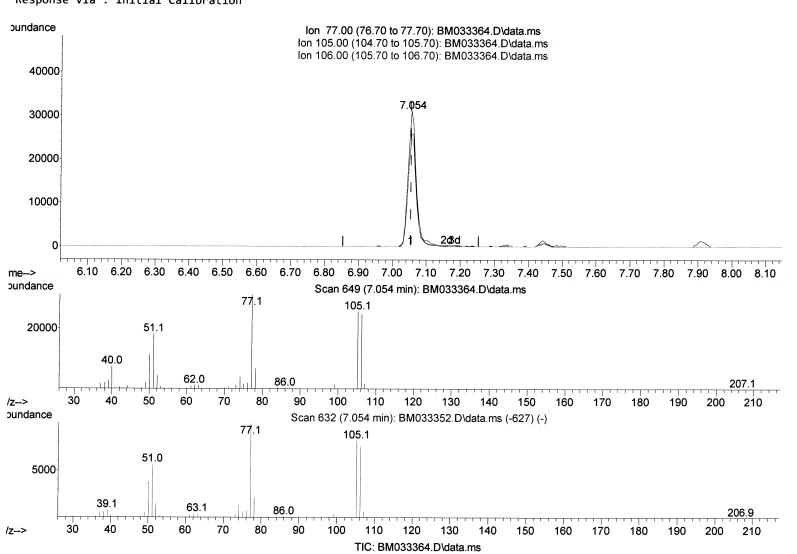
Quant Time: Dec 10 02:47:03 2021

Quant Method: Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument:
BNA_M
LabSampleId:
SSTDCCC020

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



(6) Benzaldehyde

7.054min (0.000) 23.92 ng/ul

response	52682	
Ion	Exp %	Act%
77.00	100.00	100.00
105.00	82.00	82.42
106.00	75.70	79.52
0.00	0.00	0.00

Data File: BM033364.D

Acq On : 09 Dec 2021 19:12

Operator : CG/JU Sample : SSTDCCC020

Misc :

ALS Vial : 18 Sample Multiplier: 1

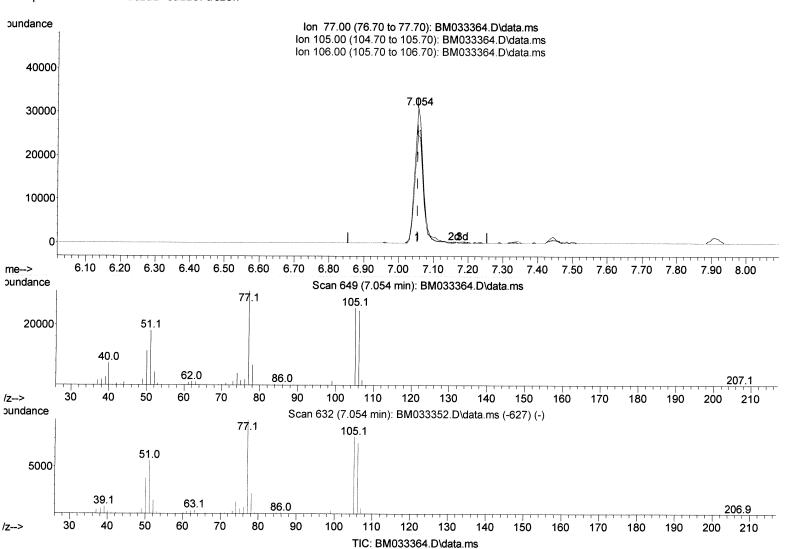
Quant Time: Dec 10 02:47:03 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument:
BNA_M
LabSampleId:
SSTDCCC020

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



(6) Benzaldehyde

23.08 ng/ul m 7.054min (0.000) response 50825 Ion Exp% Act% 77.00 100.00 100.00 105.00 82.00 82.42 106.00 75.70 79.52 0.00 0.00 0.00

Data File : BM033364.D

Acq On : 09 Dec 2021 19:12

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 18 Sample Multiplier: 1

Quant Time: Dec 10 02:47:03 2021

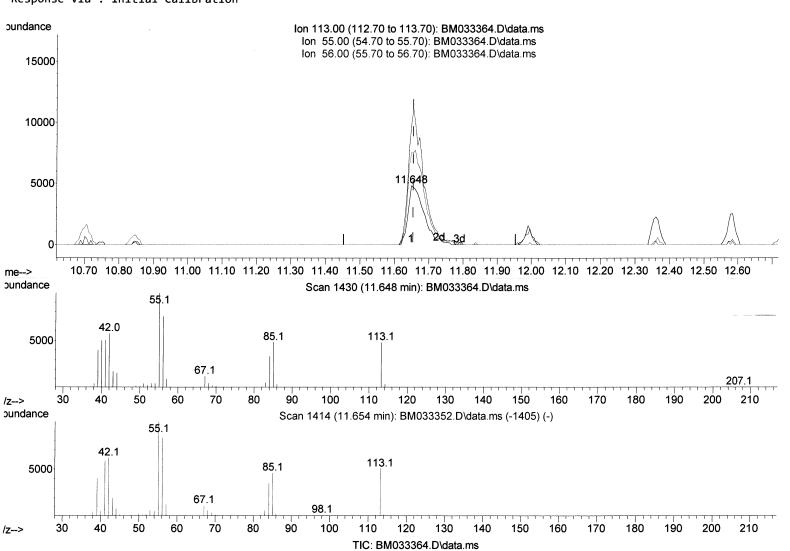
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument :
BNA_M
LabSampleId :
SSTDCCC020

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



(34) Caprolactam

11.648min (-0.006) 16.27 ng/ul

response	14886	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	197.40	206.67
56.00	164.70	156.49
0.00	0.00	0.00

Data File : BM033364.D

Acq On : 09 Dec 2021 19:12

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 18 Sample Multiplier: 1

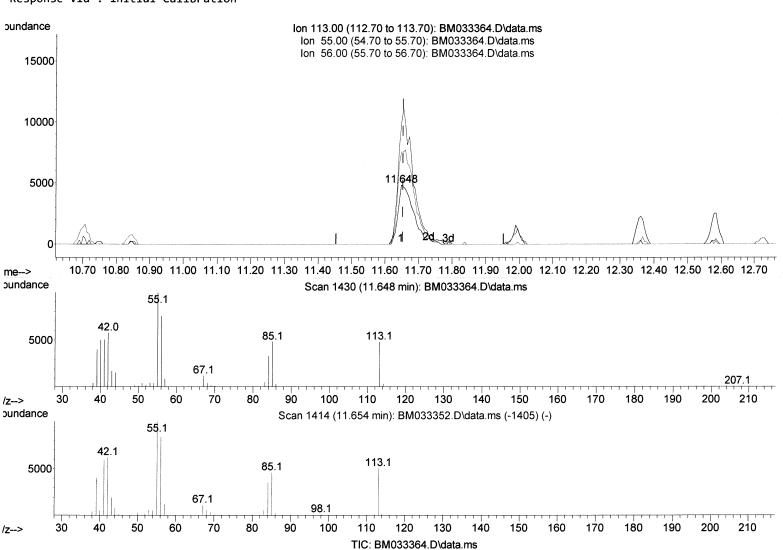
Quant Time: Dec 10 02:47:03 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument:
BNA_M
LabSampleId:
SSTDCCC020

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



(34) Caprolactam

11.648min	(-0.006)	17.03 ng/ul	Te(2/25/21
response	15580		1912/ 51
Ion	Exp%	Act%	
113.00	100.00	100.00	
55.00	197.40	206.67	
56.00	164.70	156.49	
0.00	0.00	0.00	

Data File: BM033364.D

Acq On : 09 Dec 2021 19:12

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 18 Sample Multiplier: 1

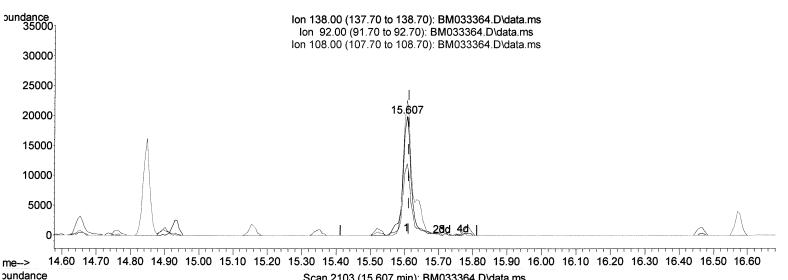
Quant Time: Dec 10 02:47:03 2021

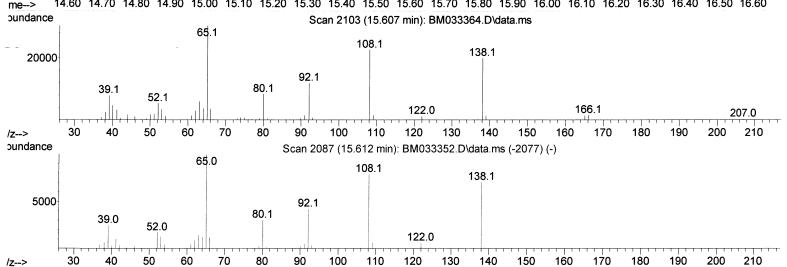
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument :
BNA_M
LabSampleId :
SSTDCCC020

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021





TIC: BM033364.D\data.ms

(63) 4-Nitroaniline

15.607min (-0.006) 19.11 ng/ul

response	31268	
Ion	Ехр%	Act%
138.00	100.00	100.00
92.00	60.00	60.01
108.00	111.90	113.56
0.00	0.00	0.00

Data File: BM033364.D

Acq On : 09 Dec 2021 19:12

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 18 Sample Multiplier: 1

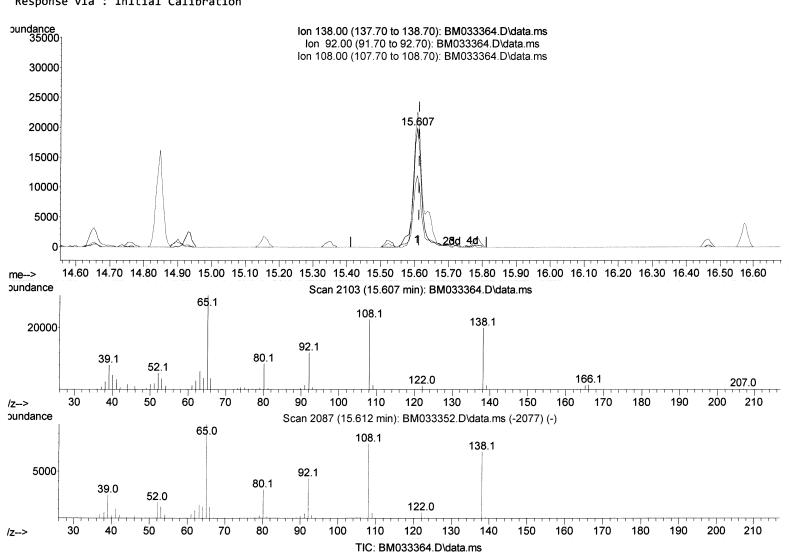
Quant Time: Dec 10 02:47:03 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument: BNA_M LabSampleId: SSTDCCC020

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



(63) 4-Nitroaniline

21.16 ng/ul m J4(2/25/2) 15.607min (-0.006) response 34632 Ion Ехр% Act% 138.00 100.00 100.00 92.00 60.00 60.01 108.00 111.90 113.56 0.00 0.00 0.00

Data File : BM033364.D

Acq On : 09 Dec 2021 19:12

Dperator : CG/JU
Sample : SSTDCCC020

Misc :

ALS Vial : 18 Sample Multiplier: 1

Quant Time: Dec 10 02:47:03 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021
Response via : Initial Calibration

Instrument : BNA_M **LabSampleld :** SSTDCCC020

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021

Compound	R.T.	QIon	Response	Conc Units Dev((Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.907	152	41747	20.000 ng/ul	0.00
20) Naphthalene-d8	10.707	136	170359	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.530	164	112890	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.271	188	245940	20.000 ng/ul	0.00
79) Chrysene-d12	21.436	240	257007	20.000 ng/ul	0.00
88) Perylene-d12	23.759	264	250202	20.000 ng/ul	0.00
System Monitoring Compounds					0.00 541425 21
3) 1,4-Dioxane-d8	3.366	96	8368ŋ ≫	7.532 ng/uL\(\sigma\)	0.0053419
4) Pyridine-d5	3.784	84	55483	17.225 ng/ul	0.00
7) Phenol-d5	7.078	99	68052	17.215 ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.242	67	45720	17.666 ng/ul	0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.442	132	50794	18.352 ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.619	113	54203	17.518 ng/ul	0.00
21) Nitrobenzene-d5	9.072	128	26145	18.914 ng/ul	0.00
24) 2-Nitrophenol-d4	9.789	143	26883	18.955 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.330	165	51530	19.226 ng/ul	0.00
31) 4-Chloroaniline-d4	10.842	131	68248	17.175 ng/ul	0.00
46) Dimethylphthalate-d6	13.942	166	157274	18.646 ng/ul	0.00
<pre>49) Acenaphthylene-d8</pre>	14.230	160	198801	19. 01 3 ng/ul	0.00
54) 4-Nitrophenol-d4	14.748	143	25389	16.578 ng/ul	0.00
60) Fluorene-d10	15.524	176	141569	18.764 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.642	200	22340	15.050 ng/ul	0.00
73) Anthracene-d10	17.371	188	225739	18.572 ng/ul	0.00
81) Pyrene-d10	19.653	212	265183	18.462 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.606	264	248405	18.317 ng/ul	0.00
Target Compounds				Qva	1ue 54129 21 91 54(2/28/2) 95 54(2/28/2)
2) 1,4-Dioxane	3.401	88	8640m	6.987 ng/uL	5 3919
5) Pyridine	3.807	79	57761	17.412 ng/ul	91 1 1 1 1 1 1 1 1 1
Benzaldehyde	7.054	77	50825m	23.077 ng/ul	> Ju(2/20/2)
8) Phenol	7.107	94	69903	17.167 ng/ul	95
<pre>10) Bis(2-Chloroethyl)ether</pre>	7.331	93	54212	17.697 ng/ul	96
<pre>12) 2-Chlorophenol</pre>	7.478	128	52905	18.485 ng/ul	96
<pre>13) 2-Methylphenol</pre>	8.354	108	50174	16.996 ng/ul	93
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.431	45	93800	17.746 ng/ul#	96
16) Acetophenone	8.736	105	89185	17.400 ng/ul	96
17) N-Nitroso-di-n-propyla	8.713	70	50773	18.122 ng/ul	99
<pre>18) 4-Methylphenol</pre>	8.678	108	55619	17.220 ng/ul	93
19) Hexachloroethane	8.983	117	26075	18.033 ng/ul	92
22) Nitrobenzene	9.113	77	76194	18.827 ng/ul	98
23) Isophorone	9.636	82	129570	18.697 ng/ul	99
25) 2-Nitrophenol	9.825	139	28371	18.869 ng/ul	99
26) 2,4-Dimethylphenol	9.878	107	69355	19.024 ng/ul	99
27) Bis(2-Chloroethoxy)met	10.113	93	73779	18.930 ng/ul	99
29) 2,4-Dichlorophenol	10.354	162	51796	19.091 ng/ul	98
30) Naphthalene	10.754	128	174862	18.419 ng/ul	99
32) 4-Chloroaniline	10.872	127	72667	18.195 ng/ul	96
33) Hexachlorobutadiene	11.030	225	37576	18.921 ng/ul	- 97 5414201 -1
34) Caprolactam	11.648	113	1558@m	17.028 ng/ul	> 96 541428 2.)
35) 4-Chloro-3-methylphenol	11.989	107	60125	18.879 ng/ul	93

Data File : BM033364.D

Acq On : 09 Dec 2021 19:12

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Sample : SSTDCCC020

Misc :

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 09 13:25:37 2021
Response via : Initial Calibration

Instrument :
BNA_M
LabSampleId :
SSTDCCC020

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021

	Compound	R.T.	QIon	Response	Conc Un	its Dev	(Min)	
36)	2-Methylnaphthalene	12.360	142	118199	18.359	ng/ul	 99	
	1-Methylnaphthalene	12.577	142	123477		ng/ul	96	
	1,2,4,5-Tetrachloroben	12.724	216	64275		ng/ul	95	
	Hexachlorocyclopentadiene	12.701	237	50578		ng/ul	95	
	2,4,6-Trichlorophenol	12.971	196	39368		ng/ul	95	
	2,4,5-Trichlorophenol	13.048	196	42207	19.459		93	
	1,1'-Biphenyl	13.366	154	165436	19.155		96	
	2-Chloronaphthalene	13.413	162	126840	19.089		99	
	2-Nitroaniline	13.624	65	46901	19.792		99	
	Dimethylphthalate	13.989	163	157813	18.843	_	99	
	2,6-Dinitrotoluene	14.113	165	30287	18.708	_	91	
-	Acenaphthylene	14.260	152	210197	19.307	_	98	
	3-Nitroaniline	14.448	138	30539	19.164	-	95	
	Acenaphthene	14.595	153	135520	18.756		96	
	2,4-Dinitrophenol	14.654	184	14238	14.965	-	94	
	4-Nitrophenol	14.760	109	29800	18.003	_	89	
	Dibenzofuran	14.930	168	197896	18.887		96	
-	2,4-Dinitrotoluene	14.901	165	45803	19.259		98	
	2,3,4,6-Tetrachlorophenol	15.160	232	35429		ng/ul#	98	
	Diethylphthalate	15.348	149	162412	18.685		99	
	Fluorene	15.577	166	164080	19.045		aa	4
•	4-Chlorophenyl-phenyle	15.571	204	80344	18.695		95	100
•	4-Nitroaniline	15.607	138	34632m		ng/ul_	ر د	recy 201
	4,6-Dinitro-2-methylph	15.654	198	23136		ng/ul#	96	surfrol
	N-Nitrosodiphenylamine	15.789	169	137221	18.976		97	
	4-Bromophenyl-phenylether	16.465	248	46589	18.807		96	
	Hexachlorobenzene	16.571	284	52256	18.309		95	
-	Atrazine	16.736	200	50193	17.493	-	98	
	Pentachlorophenol	16.924	266	34136	21.329		93	
	Phenanthrene	17.312	178	266813	18.805	-	99	
•	Anthracene	17.406	178	267490	18.596	_	99	
	1,2,3,4-Tetrachloroben	13.330	216	66320	18.697	_	96	
	Pentachlorobenzene	14.848	250	67082	19.151		96	
	Carbazole	17.677	167	233363	17.949	_	99	
•	Di-n-butylphthalate	18.230	149	272629	18.688		99	
	Fluoranthene	19.324	202	308863	18.242	_	99	
	Pyrene	19.683	202	327910	18.500	-	98	
	Butylbenzylphthalate	20.571	149	125196	18.439		99	
	3,3'-Dichlorobenzidine	21.353	252	99805	17.085		95	
	Benzo(a)anthracene	21.418	228	309246	18.314	_	99	
	Bis(2-ethylhexyl)phtha	21.342	149	179513	18.396	-	100	
	Chrysene	21.471	228	307318	18.518		98	
	Di-n-octyl phthalate	22.241	149	312156	17.016		100	
	Benzo(b)fluoranthene	23.053	252	315161	18.385	_	99	
	Benzo(k)fluoranthene	23.100	252	294749	18.570		100	
	Benzo(a)pyrene	23.659	252	305952	18.601	-	98	
	Indeno(1,2,3-cd)pyrene	26.129	276	322179	18.011		100	
	Dibenzo(a,h)anthracene	26.141	278	278056	17.870		99	
	Benzo(g,h,i)perylene	26.859	276	275775	18.075		98	

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