Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120921\

Data File : BM033367.D

Acq On : 09 Dec 2021 21:00

Operator : CG/JU Sample : PB141232BS

Misc

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 01:15:53 2021

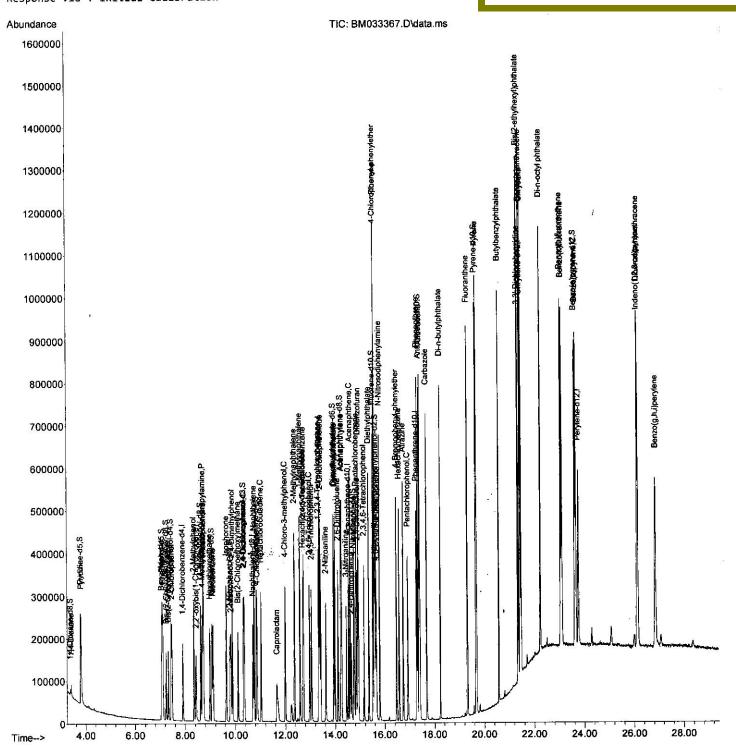
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

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

Manual IntegrationsAPPROVED

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



SFAM-EPA-BM120921.M Fri Dec 10 03:10:22 2021

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120921\

Data File : BM033367.D

: 09 Dec 2021 21:00 Acq On

Operator : CG/JU Sample : PB141232BS

Misc

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 01:15:53 2021

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

Quant Title : SVOA CALIBRATION

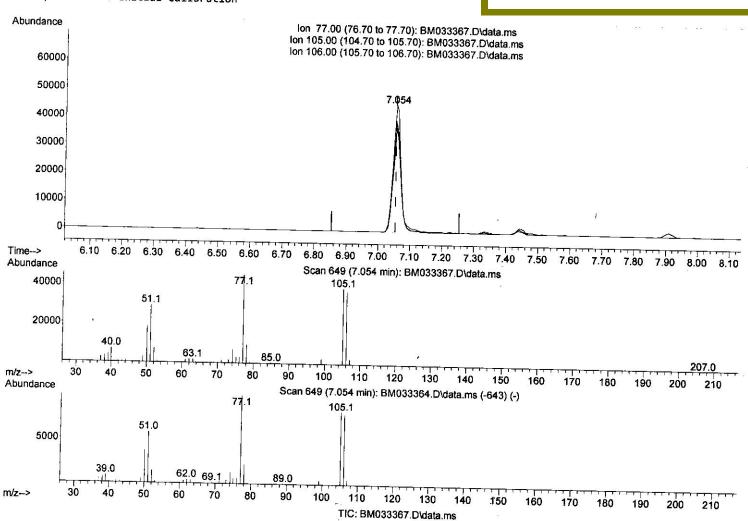
QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration

Instrument: BNA_M

ClientSampleId:

Manual IntegrationsAPPROVED

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



Benzaldehyde

36.74 ng/ul 7.054min (-0.000)

response	77747				
Ion	Ехр%	Act%			
77.00	100.00	100.00			
105.00	82.00	85.65			
106.00	75.70	82.04			
0.00	0.00	0.00			

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120921\

Data File : BM033367.D

: 09 Dec 2021 21:00 Acq On

Operator : CG/JU Sample : PB141232BS

Misc

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 01:15:53 2021

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

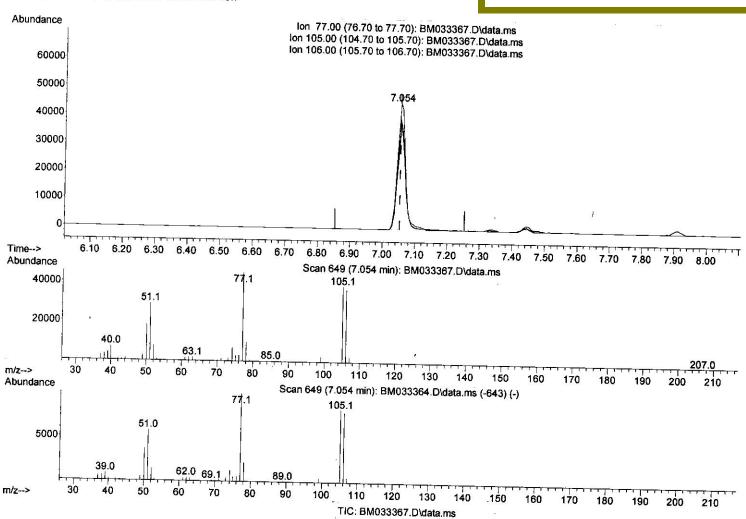
Quant Title : SVOA CALIBRATION

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

Instrument: BNA_M ClientSampleId:

Manual IntegrationsAPPROVED

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



Benzaldehyde

7.054min (-0.000) 36.17 ng/ul m

response	76550				
Ion	Exp%	Act%			
77.00	100.00	100.00			
105.00	82.00	85.65			
106.00	75.70	82.04			
0.00	0.00	0.00			

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120921\

Data File : BM033367.D

Acq On : 09 Dec 2021 21:00

Operator : CG/JU Sample : PB141232BS

Misc

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 01:15:53 2021

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

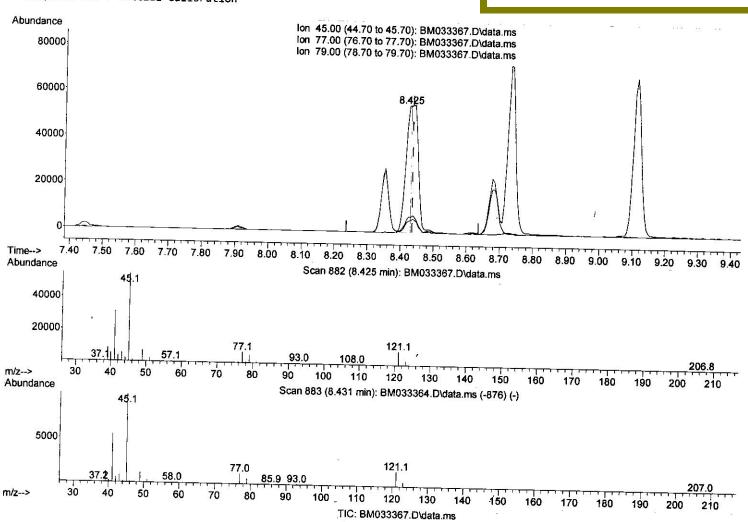
Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:25:37 2021

Response via : Initial Calibration

Instrument : BNA_M ClientSampleId :

Manual IntegrationsAPPROVED

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



(14) 2,2'-oxybis (1-Chloropropane)

8.425min (-0.012) 14.27 ng/ul

response	72453	
Ion	Exp%	Act*
45.00	100.00	100.00
77.00	12.40	12.44
79.00	10.40	9.09
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120921\

Data File : BM033367.D

Acq On : 09 Dec 2021 21:00

Operator : CG/JU Sample : PB141232BS

Misc

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 01:15:53 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

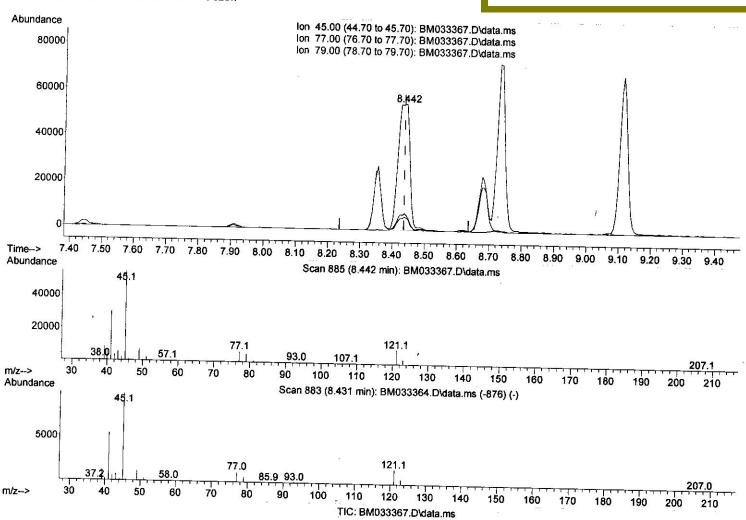
Instrument : BNA_M

ClientSampleId :

SLCS232

Manual IntegrationsAPPROVED

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



(14) 2,2'-oxybis(1-Chloropropane) 8.442min (+ 0.006) 27.53 ng/ul m

response	139819	
Ion	Екр%	Act%
45.00	100.00	100.00
77.00	12.40	11.81
79.00	10.40	9.46
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120921\

Data File : BM033367.D

: 09 Dec 2021 21:00 Acq On

Operator : CG/JU Sample : PB141232BS

Misc

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 01:15:53 2021

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

Quant Title : SVOA CALIBRATION

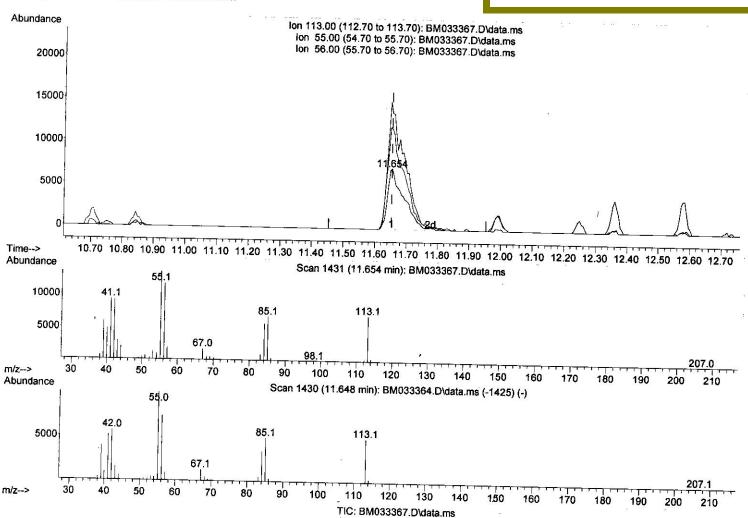
QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration

Instrument: BNA_M

ClientSampleId:

Manual IntegrationsAPPROVED

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



(34) Caprolactam

11.654min (-0.000) 27.56 ng/ul

response	25243					
Ion	Ежр%	Act*				
113.00	100.00	100.00				
55.00	197.40	192.86				
56.00	164.70	166.86				
0.00	0.00	0.00				

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120921\

Data File : BM033367.D

Acq On : 09 Dec 2021 21:00

Operator : CG/JU Sample : PB141232BS

Misc

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 01:15:53 2021

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

Quant Title : SVOA CALIBRATION

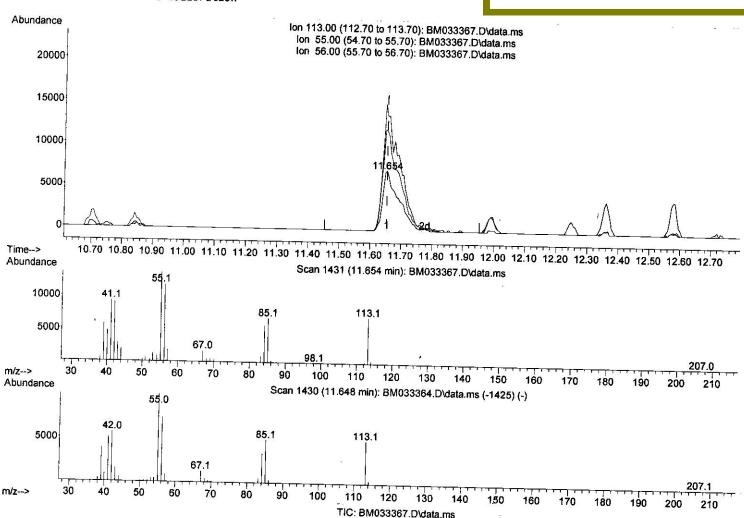
QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration

Instrument: BNA_M

ClientSampleId:

Manual IntegrationsAPPROVED

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



(34) Caprolactam

11.654min (-0.000) 27.98 ng/ul

response	25623					
Ion	Ежр%	Act%				
113.00	100.00	100.00				
55.00	197.40	192.86				
56,00	164.70	166.86				
0.00	0.00	0.00				

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120921\

Data File : BM033367.D

Acq On : 09 Dec 2021 21:00

Operator : CG/JU Sample : PB141232BS

Misc

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 01:15:53 2021

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

Quant Title : SVOA CALIBRATION

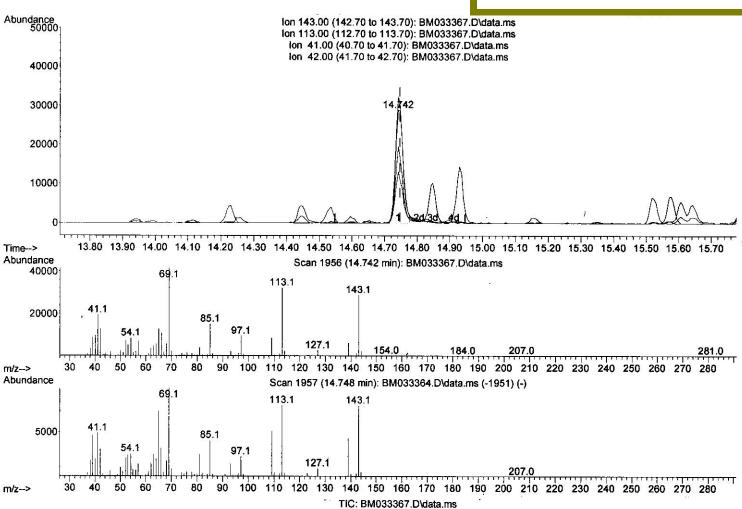
QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument : BNA_M

ClientSampleId:

SLCS232

Manual IntegrationsAPPROVED

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



(54) 4-Nitrophenol-d4 (S)

14.742min (-0.006) 28.86 ng/ŭl

response	47740	
Ion	Exp&	Act%
143.00	100.00	100.00
113.00	105.00	112.14
41.00	57.20	67.45
42.00	39.50	45.25

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120921\

Data File : BM033367.D

Acq On : 09 Dec 2021 21:00

Operator : CG/JU Sample : PB141232BS

Misc

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 01:15:53 2021

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

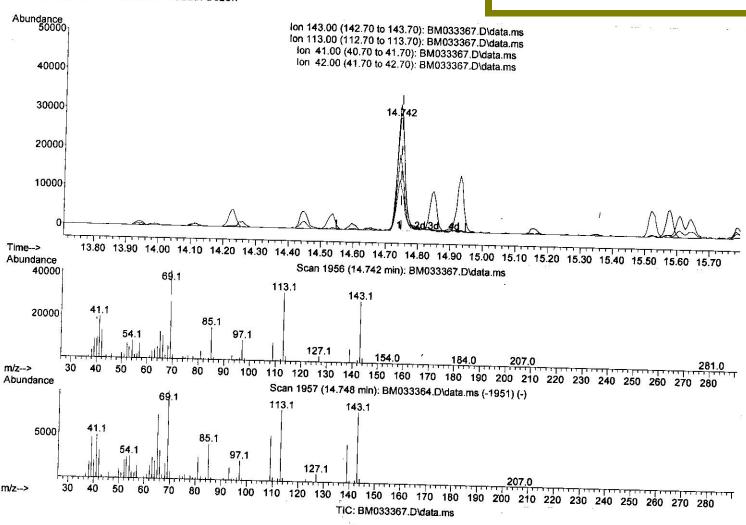
Quant Title : SVOA CALIBRATION

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

Instrument: BNA_M ClientSampleId:

Manual IntegrationsAPPROVED

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



(54) 4-Nitrophenol-d4 (S) 14.742min (-0.006) 29.59 ng/ul m

response	48951					
Ion	Exp%	Act*				
143.00	100.00	100.00				
113.00	105.00	112.14				
41.00	57.20	67.45				
42.00	39.50	45.25				

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120921\

Data File : BM033367.D

Acq On : 09 Dec 2021 21:00 Operator : CG/JU

Sample : PB141232BS

Misc

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 01:15:53 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 ClientSampleld: SLCS232

Manual Integrations APPROVED

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

Compound	R.T.	QIon	Response	Conc Units De	v(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4					
20) Naphthalene-d8	7,907		40115	20.000 ng/ul	0.00
38) Acenaphthene-d10	10.701		170506	20.000 ng/ul	0.00
64) Phenanthrene-d10	14.536	3500000000	121943	20.000 ng/ul	0.00
79) Chrysene-d12	17.271		276505	20.000 ng/ul	0.00
88) Perylene-d12	21.436		295870	20.000 ng/ul	0.00
ob) Ferylene-ul2	23.759	264	279100	20.000 ng/ul	0.00
System Monitoring Compounds				æa.	
3) 1,4-Dioxane-d8	3 366		202120		
4) Pyridine-d5	3.366	96	5975	5.597 ng/uL	0.00
7) Phenol-d5	3.784		82827	26.761 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth	7.078	99	107143	28.206 ng/ul	0.00
11) 2-Chlorophenol-d4		67	69851	28.088 ng/ul	0.00 .
15) 4-Methylphenol-d8	7.442	132	78137	29.380 ng/ul	0.00
21) Nitrobenzene-d5	8.619	113	81733	27.489 ng/ul	0.00
24) 2-Nitrophenol-d4	9.072	128	41085	29.696 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.789	143	43598	30.715 ng/ul	0.00
31) 4-Chloroaniline-d4	10.330	165	82352	30.700 ng/ul	0.00
46) Dimethylphthalate-d6	10.842	131	119776	30.116 ng/ul	0.00
49) Acenaphthylene-d8	13,942	166	271741	29.825 ng/ul	0.00 AU
54) 4-Nitrophenol-d4	14.230	160	336711	29.811 ng/ul	0.00 12/2X121
60) Fluorene-d10	14.742	143	48951m	29.590 ng/ul	0.00
65) 4 6 Dinitha 2	15.524	176	246211 _	30.210 ng/ul	0.00
65) 4,6-Dinitro-2-methylph		200	47437	28.424 ng/ul	0.00
73) Anthracene-d10	17.371	188	409157	29.942 ng/ul	0.00
81) Pyrene-d10	19.653	212	496881	30.048 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.612	264	467726	30.918 ng/ul	0.00
Target Compounds					
2) 1,4-Dioxane	3.401		42422		lue
5) Pyridine		88	12633	10.632 ng/uL	94
6) Benzaldehyde	3.801	79	84392 ^	26.475 ng/ul	96
8) Phenol	7.054	77	76550m	36.172 ng/ul	
<pre>10) Bis(2-Chloroethyl)ether</pre>	7.107	94	103543	26.463 ng/ul	98
12) 2-Chlorophenol	7.331	93	78625	26.711 ng/ul	99 .
13) 2-Methylphenol	7.478	128	75910	27.602 ng/ul	98
14) 2,2'-oxybis(1-Chloropr	8.354	108	74999	26.439 ng/ul	97
16) Acetophenone	10 <u>-</u> 10 1222222 1040	45	139819m	27.529 ng/ul	
17) N-Nitroso-di-n-propyla		105	134577	27.324 ng/ul	99
18) 4-Methylphenol	8.713		77568	28.812 ng/ul	96 ^ V
19) Hexachloroethane	8.683		82215	26.490 ng/ul	96 JU 99 12/2X12 \
22) Nitrobenzene		117	39405	28.361 ng/ul	87 \2\371
23) Isophorone	9.113	77	118141	29.167 ng/ul	97
25) 2-Nitrophenol	9.636	82	201308	29.024 ng/ul	100
26) 2,4-Dimethylphenol		139	42828	28.460 ng/ul	97
27) Bis(2-Chloroethoxy)met		107	99076	27.154 ng/ul	96
29) 2,4-Dichlorophenol	10.113		110710	28.381 ng/ul	98
30) Naphthalene		162	78312	28.839 ng/ul	97
32) 4-Chloroaniline			263855	27.769 ng/ul	99
33) Hexachlorobutadiene			101585	25.414 ng/ul	99
34) Caprolactam		225	58100	29.230 ng/ul	96
35) 4-Chloro-3-methylphenol		L13	25623m	27.980 ng/ul	
22) 4 Chitor 0-3-metnyiphenol	11.989 1	107	95917	30.092 ng/ul	92

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120921\

Data File : BM033367.D

Acq On : 09 Dec 2021 21:00

Operator : CG/JU : PB141232BS Sample

Misc

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 01:15:53 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 ClientSampleId : SLCS232

Manual IntegrationsAPPROVED

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

Response via : Initial Calibration	•						Dov/Min'	١.	Superv
	R.T.	QIon	Resp	onse C	onc Un	172	Dev(Min)	<u>′</u>	
Compound			. -		28.490	nø/	ul '	97	
36) 2-Methylnaphthalene	12.360	142		3584	28.506	ng/	ul	97	
36) 2-Methylnaphthalene	12.577	142		0891	27.132	ng	/ul	98	
37) 1-Methylnaphthalene 39) 1,2,4,5-Tetrachloroben	12.724	216		9684	31.59	ng.	/ul	99	
40) Hexachlorocyclopentadiene	12.701	237		5762	29.53	ng.	/ul	93	
41) 2,4,6-Trichlorophenol	12.966		2	3864	29.37	1 ng	/ul	98	
41) 2,4,6-111chlorophenol 42) 2,4,5-Trichlorophenol	13.048			8814	27.74	0 ng	/ul	99	
42) 2,4,5-11 Ichian opini	13.365	154		8791	27.91	3 ng	/ul	98	
43) 1,1'-Biphenyl 44) 2-Chloronaphthalene	13.413	162		0345	31.36	2 ng	/ul	99	
45) 2-Nitroaniline	13,624			30278	28.51	5 ng	/ul	99	
47) Dimethylphthalate	13.989			57962	29.38	6 ng	/ul#	86	
48) 2,6-Dinitrotoluene	14.11		2 card	51389	28.78	34 ne	z/ul	98	
50) Acenaphthylene	14.26			38496	29.4	33 ns	z/ul	99	
51) 3-Nitroaniline	14.44		7° www	50665	28.5	30 ni	g/ul	97	
51) 3-NILI Odli III.	14.59		_	22440	29.0	68 n	g/ul	93	
52) Acenaphthene	14.64		100	29873	31.6	77 n	g/ul	90)
53) 2,4-Dinitrophenol	14.76	0 10		56638	20.0	75 n	g/ul	98	;
55) 4-Nitrophenol	14.93	0 16	8 3	329073	29.0	02 n	g/ul	93	3
56) Dibenzofuran	14.98	1 16	55	79363	30.0 30.7	96 n	g/ul	99	€
57) 2,4-Dinitrotoluene	15.15	9 2	32	61156	20.7	7/1 r	ig/ul	99	€.
58) 2,3,4,6-Tetrachlorophenol	15.34	18 14		281437	29.5	00 r	ng/ul	99	9
59) Diethylphthalate	15.5	77 1		276384	29.0	157 i	ng/ul	9	8
61) Fluoriene			04	136744	29.4	+3/ 1 377	ng/ul	9	2
61) Fluoriene 62) 4-Chlorophenyl-phenyle	15.6	07 1	38	60227	34.1	275	ng/ul	9	6
63) 4-Nitroaniline		59 1	98	45276	2/.	4.AE	ng/ul		8
66) 4,6-Dinitro-2-methylph	15.7		.69	236947	29.	147 747	ng/ul		8
67) N-Nitrosodiphenylamine			48	80061	28.	74/	ng/ul		96
68) 4-Bromophenyl-phenylethe	16.5		284	90950	28.	053	ng/ul		99
69) Hexachlorobenzene	16.7	736	200	86946	26.	222	ng/ul		92
70) Atrazine	16.9	24	266	56865	31.	C 4 6	ng/ul		98
71) Pentachlorophenol	17.		178	456943	28	424	ng/ul		99
72) Phenanthrene	17.	406	178	459685	28	020	ng/uL		96
74) Anthracene			216	103437		.930	ng/uL		98
75) 1,2,3,4-Tetrachloroben		848	250	104619		.500	ng/ul		99
76) Pentachlorobenzene	_	677	167	416538		.49/	ng/ul		99
77) Carbazole		230	149	500787		.554	ng/ul		98
78) Di-n-butylphthalate		324	202	557254		100	ng/ul	- 65	100
80) Fluoranthene		683	202	575274		0.194	ng/ul		96
82) Pyrene	20	571	149	234236			ng/ul		96
83) Butylbenzylphthalate	21	353	252	172818		707	2 ng/ul		99
84) 3,3'-Dichlorobenzidine		418	228	557948		3.70	2 ng/ul		99
85) Benzo(a)anthracene	. 21	. 336	149	33546	3 4	9.00	7 ng/ul	*	98
86) Bis(2-ethylhexyl)phtha.	21	.471	228	54751		8.00	o ng/ul	34	100
g7) Chrysene		.241	149	59195		8.92	8 ng/ul		99
89) Di-n-octyl phthalate		.053	252	57257	1 2	9.94	2 ng/ul	8	99
90) Benzo(b)fluoranthene	23	.100	252	53573		0.25	8 ng/ul		99
91) Benzo(k)fluoranthene	23	3.653	252	54211		9.54	6 ng/ul		98
on ponzo(a)nvrene		5.129	276			8.6	[3 ng/ul		97
94) Indeno(1,2,3-cd)pyrene		6.141	278	49862		28./	27 ng/ul		99
95) Dibenzo(a,h)anthracene		6.865	276				33 ng/ul		
96) Benzo(g,h,i)perylene	_ 							pp : 74 (20) (47)	
				17400		\ -	signals	SUM	ımed

(*) = qualifier out of range (m) = manual integration (+) = signals summed