Data File : BM033351.D

Acq On : 09 Dec 2021 10:16

Operator : CG/JU Sample : SSTD01012

Misc

ALS Vial : 4 Sample Multiplier: 1

Quant Time: Dec 09 13:10:24 2021

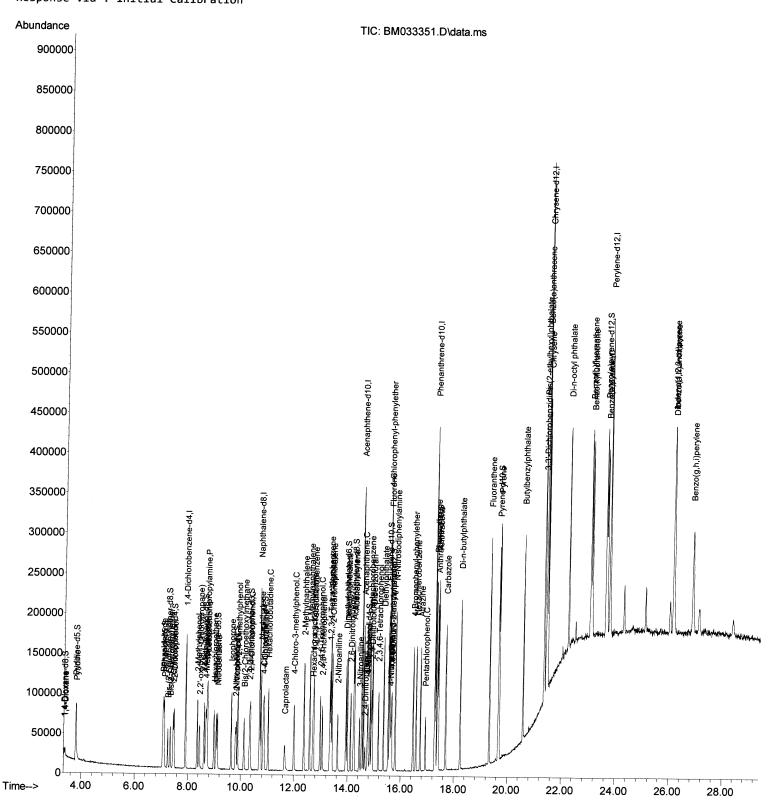
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 09 13:01:40 2021
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
SSTD010012

Manual IntegrationsAPPROVED

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



Data File: BM033351.D

Acq On : 09 Dec 2021 10:16

Operator : CG/JU Sample : SSTD01012

Misc

ALS Vial : 4 Sample Multiplier: 1

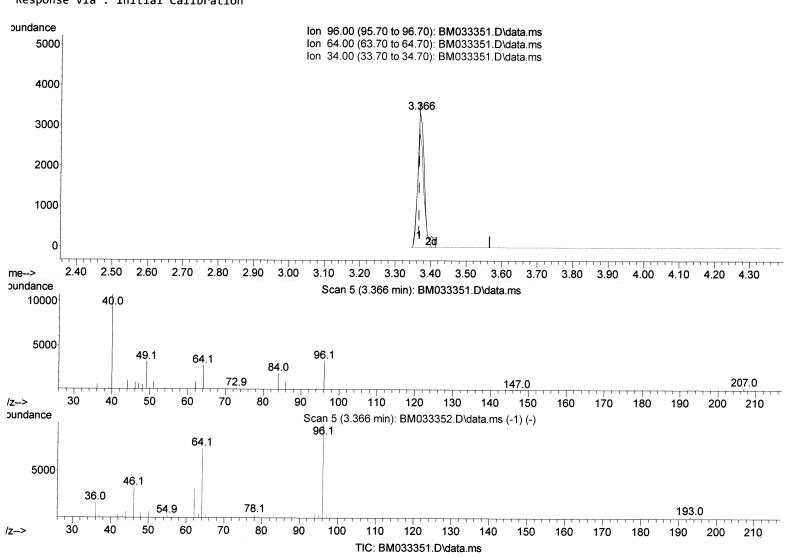
Quant Time: Dec 09 13:10:24 2021

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

3.366min (+ 0.000) 3.57 ng/uL

response	3655		
Ion	Ежр%	Act%	
96.00	100.00	100.00	
64.00	74.20	84.12	
34.00	0.00	0.00	
0.00	0.00	0.00	

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Misc

ALS Vial : 4 Sample Multiplier: 1

Quant Time: Dec 09 13:10:24 2021

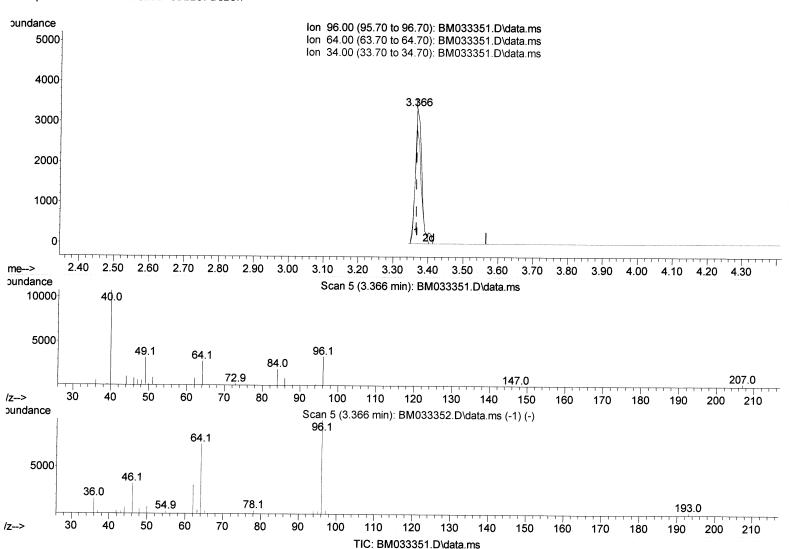
 $\label{lem:quant_method} \textbf{Quant Methods} : \textbf{Z:} \\ \textbf{SVOASRV} \\ \textbf{HPCHEM1} \\ \textbf{BNA_M} \\ \textbf{METHODS} \\ \textbf{SFAM-EPA-BM120921.M} \\ \textbf{METHODS} \\ \textbf{METH$

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:01:40 2021 Response via : Initial Calibration



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

3.366min	(+ 0.000)	4.24 ng/uL	m	12/24/21
response	4346		04	12/ - / .
Ion	Exp %	Act%		
96.00	100.00	100.00		
64.00	74.20	84.12		
34.00	0.00	0.00		
0.00	0.00	0.00		

Data File: BM033351.D

Acq On : 09 Dec 2021 10:16

Operator : CG/JU Sample : SSTD01012

Misc

ALS Vial : 4 Sample Multiplier: 1

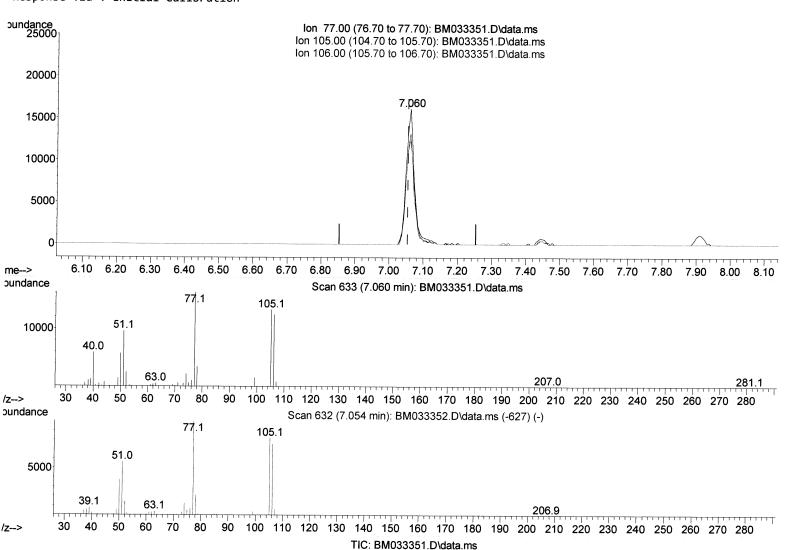
Quant Time: Dec 09 13:10:24 2021

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

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:01:40 2021 Response via : Initial Calibration Instrument:
BNA_M
ClientSampleId:
SSTD010012

Manual IntegrationsAPPROVED

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(6) Benzaldehyde

7.060min (+ 0.006) 14.34 ng/ul

response	27294				
Ion	Exp%	Act%			
77.00	100.00	100.00			
105.00	82.00	81.79			
106.00	75.70	76.78			
0.00	0.00	0.00			

Data File: BM033351.D

Acq On : 09 Dec 2021 10:16

Operator : CG/JU Sample : SSTD01012

Misc

ALS Vial : 4 Sample Multiplier: 1

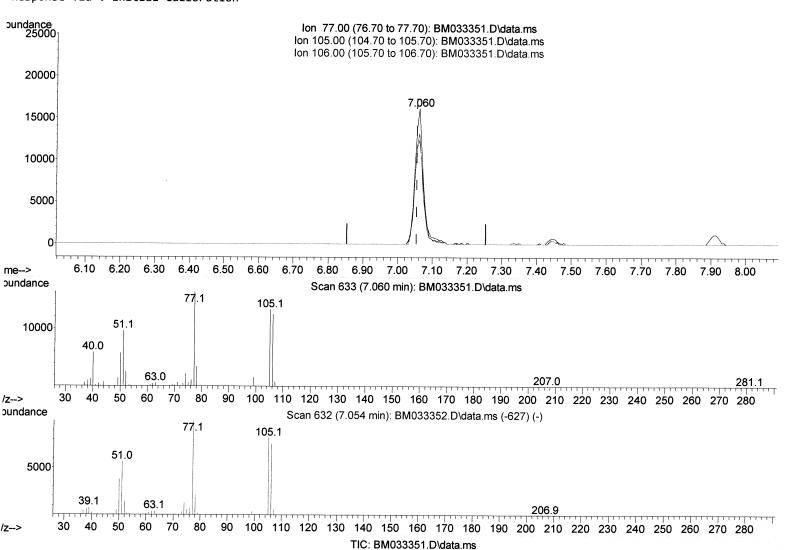
Quant Time: Dec 09 13:10:24 2021

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

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:01:40 2021 Response via : Initial Calibration Instrument : BNA_M ClientSampleId : SSTD010012

Manual IntegrationsAPPROVED

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6) Benzaldehyde

7.060min	(+ 0.006)	13.68 ng/ul	T412/23/27
response	26046		JUIN
Ion	Ехр%	Act%	
77.00	100.00	100.00	
105.00	82.00	81.79	
106.00	75.70	76.78	
0.00	0.00	0.00	

Data File : BM033351.D

Acq On : 09 Dec 2021 10:16

Operator : CG/JU Sample : SSTD01012

Misc

ALS Vial : 4 Sample Multiplier: 1

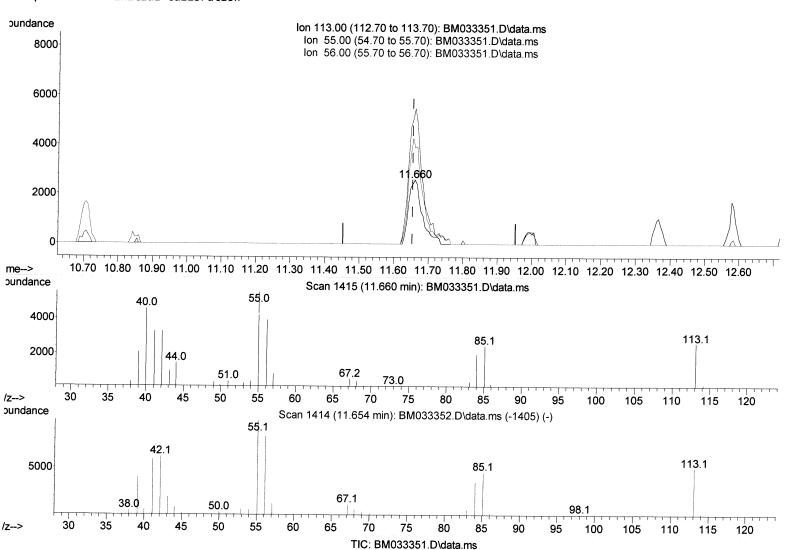
Quant Time: Dec 09 13:10:24 2021

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

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:01:40 2021 Response via : Initial Calibration Instrument : BNA_M ClientSampleId : SSTD010012

Manual IntegrationsAPPROVED

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



(34) Caprolactam

11.660min (+ 0.006) 8.99 ng/ul

response	6911	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	197.40	211.99
56.00	164.70	150.50
0.00	0.00	0.00

Data File : BM033351.D

Acq On : 09 Dec 2021 10:16

Operator : CG/JU Sample : SSTD01012

Misc

ALS Vial : 4 Sample Multiplier: 1

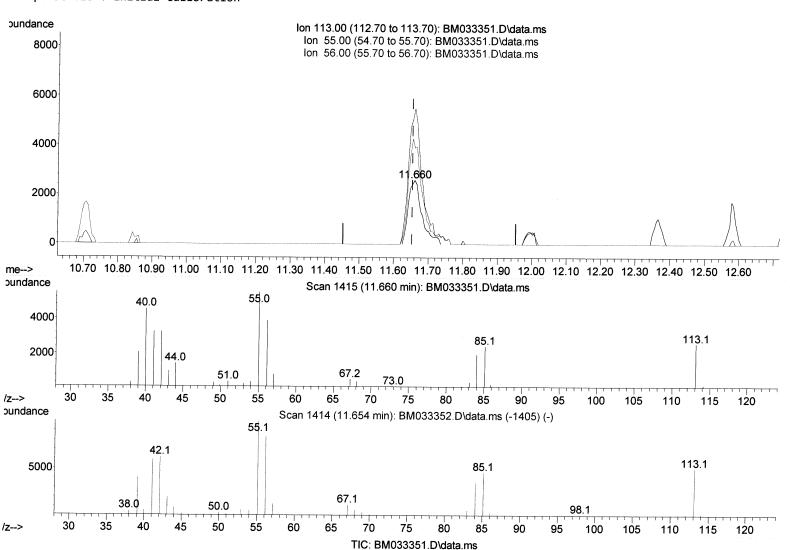
Quant Time: Dec 09 13:10:24 2021

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

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:01:40 2021 Response via : Initial Calibration Instrument:
BNA_M
ClientSampleId:
SSTD010012

Manual IntegrationsAPPROVED

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(34) Caprolactam

J(2/23/21 11.660min (+ 0.006) 9.24 ng/ul 7099 response Ion Ехр% Act% 113.00 100.00 100.00 55.00 197.40 211.99 56.00 164.70 150.50 0.00 0.00 0.00

Data File : BM033351.D

Acq On : 09 Dec 2021 10:16

Operator : CG/JU Sample : SSTD01012

4isc

ALS Vial : 4 Sample Multiplier: 1

Quant Time: Dec 09 13:10:24 2021

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

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:01:40 2021 Response via : Initial Calibration

Instrument : BNA_M ClientSampleId : SSTD010012

Manual IntegrationsAPPROVED

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

Compound	R.T.	QIon	Response	Conc Units De	ev(Min)	
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.913	152	39493	20.000 ng/u]	0.00	
20) Naphthalene-d8	10.707		167421	20.000 ng/u]		
38) Acenaphthene-d10	14.536		110791	20.000 ng/u]		
64) Phenanthrene-d10	17.271		245987	20.000 ng/u]		
79) Chrysene-d12	21.436		266649	20.000 ng/u]		
88) Perylene-d12	23.759		276811	20.000 ng/u]		
55, 16. 31 00 411	23.733	204		20.000 lig/ul	. 0.00	401
System Monitoring Compounds						7412/23/21
3) 1,4-Dioxane-d8	3.366	96	4346m~	> 4.242 ng/ul	>0.00>	-T414
4) Pyridine-d5	3.790	84	26868	9.522 ng/ul	0.00	0 1
7) Phenol-d5	7.084	99	33045	9.892 ng/ul	0.00	
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.243	67	23411	11.049 ng/ul		
11) 2-Chlorophenol-d4	7.448	132	24886	9.836 ng/ul		
15) 4-Methylphenol-d8	8.619	113	25312	9.763 ng/ul		
21) Nitrobenzene-d5	9.072	128	12417	10.330 ng/ul		
24) 2-Nitrophenol-d4	9.795	143	12104	10.055 ng/ul		
28) 2,4-Dichlorophenol-d3	10.331	165	23814	8.638 ng/ul		
31) 4-Chloroaniline-d4	10.848	131	33438	9.121 ng/ul		
46) Dimethylphthalate-d6	13.942	166	79627	9.799 ng/ul		
49) Acenaphthylene-d8	14.230	160	96962	9.262 ng/ul		
54) 4-Nitrophenol-d4	14.754	143	11244	8.494 ng/ul		
60) Fluorene-d10	15.524	176	70568	9.680 ng/ul		
65) 4,6-Dinitro-2-methylph	15.642	200	10660	9.221 ng/ul		
73) Anthracene-d10	17.371	188	114126	9.625 ng/ul		
81) Pyrene-d10	19.653	212	138406	8.783 ng/ul		
92) Benzo(a)pyrene-d12	23.612	264	142648	9.602 ng/ul		
Target Compounds				0	value	
2) 1,4-Dioxane	3.402	88	4969	بر 4.777 ng/uL	Value 01	1
5) Pyridine	3.813	79	28420	9.870 ng/ul	91	1412/28/21
6) Benzaldehyde	7.060	77	26046m		<u> </u>	7419 m
8) Phenol	7.113	94	34659	10.419 ng/ul	88	
10) Bis(2-Chloroethyl)ether	7.337	93	27008	10.220 ng/ul	98	
12) 2-Chlorophenol	7.478	128	26223	10.220 ng/ul	91	
13) 2-Methylphenol	8.360	108	24565	9.640 ng/ul	99	
14) 2,2'-oxybis(1-Chloropr	8.437	45	48428	11.866 ng/ul	98	
16) Acetophenone	8.737	105	43888	10.725 ng/ul	96	
17) N-Nitroso-di-n-propyla	8.713	70	25855	11.727 ng/ul	95	
18) 4-Methylphenol	8.684	108	27191	10.178 ng/ul	96	
19) Hexachloroethane	8.989	117	13989	11.758 ng/ul	92	
22) Nitrobenzene	9.119	77	38200	11.388 ng/ul	99	
23) Isophorone	9.636	82	63073	10.287 ng/ul	99	
25) 2-Nitrophenol	9.825	139	13550	10.619 ng/ul	95	
26) 2,4-Dimethylphenol	9.884	107	33135	9.883 ng/ul	97	
<pre>27) Bis(2-Chloroethoxy)met</pre>	10.119	93	36326	9.956 ng/ul	96	
29) 2,4-Dichlorophenol	10.360	162	24697	9.239 ng/ul	99	
30) Naphthalene	10.754	128	90384	10.152 ng/ul	99	
32) 4-Chloroaniline	10.872	127	34078	9.201 ng/ul		•
33) Hexachlorobutadiene	11.031	225	19146	9.195 ng/ul	93	412/28/21
34) Caprolactam	11.660	113	7099m ~	> 9.237 ng/ul	$\supset \mathcal{I}$	4121-01
35) 4-Chloro-3-methylphenol	11.995	107	27428	9.422 ng/ul#	: 89	'
				-		

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Sample : SSTD01012

Misc :

ALS Vial : 4 Sample Multiplier: 1

Quant Time: Dec 09 13:10:24 2021

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

R.T. QIon Response Conc Units Dev(Min)

Quant Title : SVOA CALIBRATION

Compound

QLast Update : Thu Dec 09 13:01:40 2021 Response via : Initial Calibration

Instrument: BNA_M ClientSampleId: SSTD010012

Manual IntegrationsAPPROVED

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

Compound	K. I.	ÓTO!!	kesponse	conc units bev	MIN)
36) 2-Methylnaphthalene	12.366	142	59378	9.626 ng/ul	96
37) 1-Methylnaphthalene	12.583	142	62485	9.923 ng/ul	95
39) 1,2,4,5-Tetrachloroben	12.730	216	32155	8.724 ng/ul	97
40) Hexachlorocyclopentadiene	12.701	237	18444	7.111 ng/ul	99
41) 2,4,6-Trichlorophenol	12.972	196	17698	8.153 ng/ul	92
42) 2,4,5-Trichlorophenol	13.048	196	19559	8.402 ng/ul	92
43) 1,1'-Biphenyl	13.372	154			
44) 2-Chloronaphthalene			81401	9.430 ng/ul	99
45) 2-Nitroaniline	13.413	162	63730	9.573 ng/ul	99
•	13.624	65	20988	11.962 ng/ul	98
47) Dimethylphthalate	13.989	163	79849	10.081 ng/ul	99
48) 2,6-Dinitrotoluene	14.113	165	14197	10.476 ng/ul	88
50) Acenaphthylene	14.260	152	102793	9.676 ng/ul	98
51) 3-Nitroaniline	14.448	138	13325	9.759 ng/ul	93
52) Acenaphthene	14.601	153	69145	9.977 ng/ul	96
53) 2,4-Dinitrophenol	14.654	184	5542	8.012 ng/ul	95
55) 4-Nitrophenol	14.760	109	13305	9.841 ng/ul	86
56) Dibenzofuran	14.936	168	99394	9.785 ng/ul	100
57) 2,4-Dinitrotoluene	14.901	165	21005	11.271 ng/ul	95
58) 2,3,4,6-Tetrachlorophenol	15.160	232	16425	3.645 ng/ul#	97
59) Diethylphthalate	15.348	149	82068	10.345 ng/ul	99
61) Fluorene	15.583	166	82844	10.261 ng/ul	96
62) 4-Chlorophenyl-phenyle	15.571	204	40885	9.702 ng/ul	97
63) 4-Nitroaniline	15.607	138	14668	10.925 ng/ul	98
66) 4,6-Dinitro-2-methylph	15.654	198	11072	9.556 ng/ul#	99
67) N-Nitrosodiphenylamine	15.789	169	68473	9.464 ng/ul	97
68) 4-Bromophenyl-phenylether	16.465	248	23437	8.568 ng/ul	96
69) Hexachlorobenzene	16.577	284	27502	8.776 ng/ul	97
70) Atrazine	16.736	200	24998	8.942 ng/ul	99
71) Pentachlorophenol	16.924	266	11039	6.094 ng/ul	93
72) Phenanthrene	17.318	178	137672	10.104 ng/ul	99
74) Anthracene	17.407	178	137368	10.053 ng/ul	99
75) 1,2,3,4-Tetrachloroben	13.336	216	33291	8.239 ng/uL	94
76) Pentachlorobenzene	14.848	250	33675	8.501 ng/uL	94
77) Carbazole	17.683	167	117710	9.697 ng/ul	99
78) Di-n-butylphthalate	18.230	149	132535	9.985 ng/ul	98
80) Fluoranthene	19.324	202	162515	8.804 ng/ul	96
82) Pyrene	19.683	202	175381	9.357 ng/ul	98
83) Butylbenzylphthalate	20.571	149	61419	9.591 ng/ul	100
84) 3,3'-Dichlorobenzidine	21.353	252	55878	9.302 ng/ul	96
<pre>85) Benzo(a)anthracene</pre>	21.418	228	171331	9.934 ng/ul	99
<pre>86) Bis(2-ethylhexyl)phtha</pre>	21.342	149	89117	9.818 ng/ul	99
87) Chrysene	21.471	228	169336	10.054 ng/ul	98
89) Di-n-octyl phthalate	22.242	149	158923	9.422 ng/ul	100
90) Benzo(b)fluoranthene	23.053	252	177553	9.524 ng/ul	97
91) Benzo(k)fluoranthene	23.100	252	166197	9.732 ng/ul	99
93) Benzo(a)pyrene	23.653	252	171942	9.759 ng/ul	100
94) Indeno(1,2,3-cd)pyrene	26.130	276	191858	9.779 ng/ul	96
95) Dibenzo(a,h)anthracene	26.147	278	167936	10.005 ng/ul	97
96) Benzo(g,h,i)perylene	26.859	276	161448	9.481 ng/ul	100

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