Data File: BM033376.D

Acq On : 10 Dec 2021 02:59

Operator : CG/JU Sample : PB141276BS

Misc

ALS Vial : 31 Sample Multiplier: 1

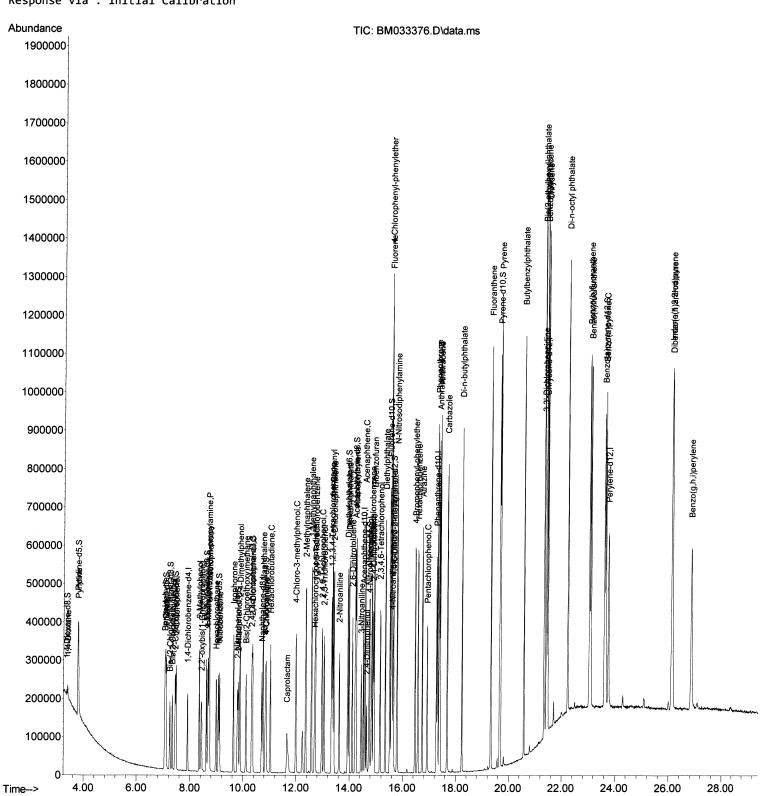
Quant Time: Dec 10 05:47:28 2021

 $\label{thm:linear_matter} Quant \ \ \mbox{Methods\span-epa-bm120921.M}$

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



Data File: BM033376.D

Acq On : 10 Dec 2021 02:59

Operator : CG/JU Sample : PB141276BS

Misc

ALS Vial : 31 Sample Multiplier: 1

Quant Time: Dec 10 05:47:28 2021

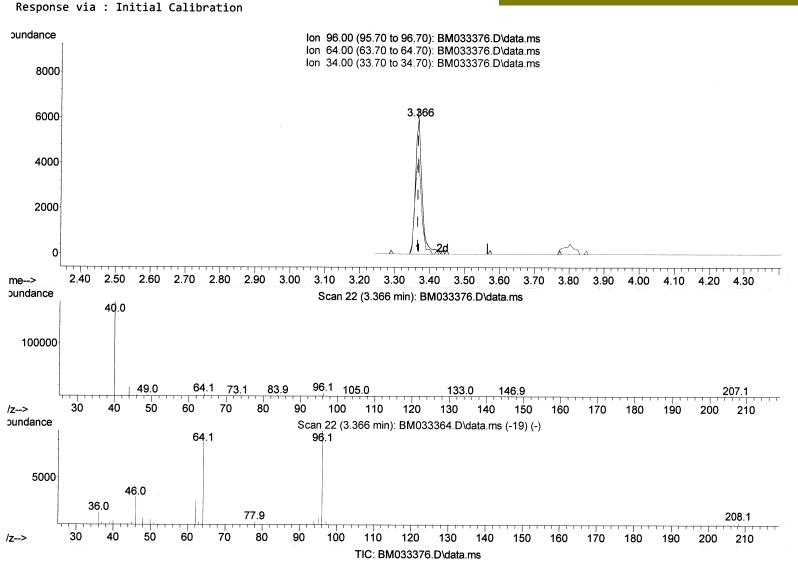
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

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

Instrument :
BNA_M
ClientSampleId :
SLCS276

Manual IntegrationsAPPROVED

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

3.366min (+ 0.000) 6.15 ng/uL

response	7570			
Ion	Exp%	Act%		
96.00	100.00	100.00		
64.00	74.20	70.62		
34.00	0.00	0.00		
0.00	0.00	0.00		

Data File: BM033376.D

Acq On : 10 Dec 2021 02:59

Operator : CG/JU Sample : PB141276BS

Misc

ALS Vial : 31 Sample Multiplier: 1

Quant Time: Dec 10 05:47:28 2021

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

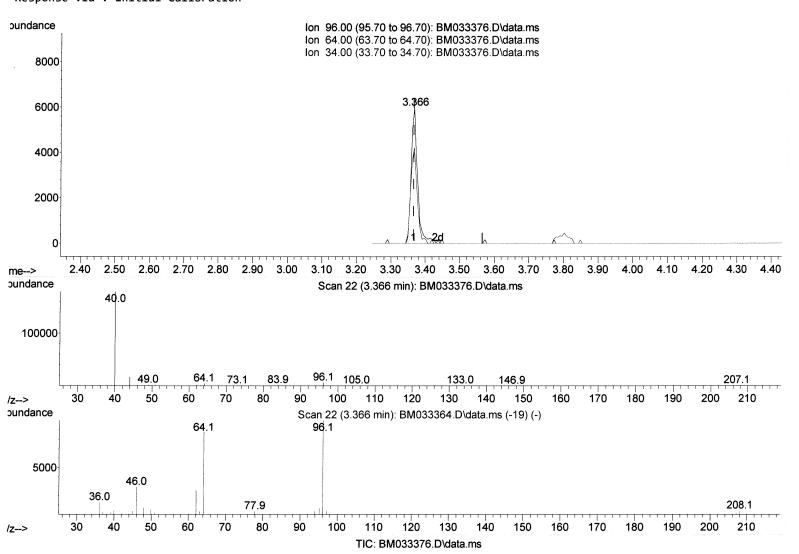
Quant Title : SVOA CALIBRATION

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



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

3.366min	(+ 0.000)	6.32 ng/	Jym m
response	7773		•
Ion	Exp %	Act%	
96.00	100.00	100.00	
64.00	74.20	70.62	
34.00	0.00	0.00	

0.00

0.00

0.00

Data File : BM033376.D

Acq On : 10 Dec 2021 02:59

Operator : CG/JU Sample : PB141276BS

Misc

ALS Vial : 31 Sample Multiplier: 1

Quant Time: Dec 10 05:47:28 2021

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

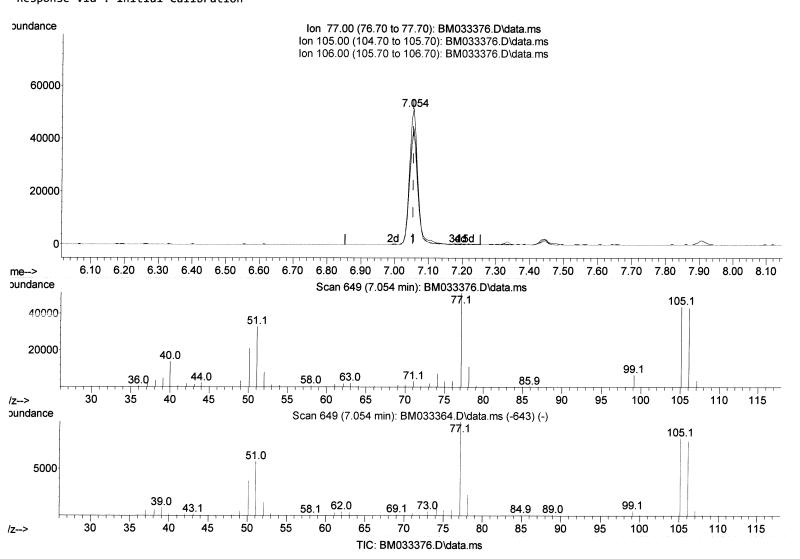
Quant Title : SVOA CALIBRATION

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



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(6) Benzaldehyde

7.054min (+ 0.000) 35.84 ng/ul

response	87454				
Ion	Exp%	Act%			
77.00	100.00	100.00			
105.00	82.00	85.56			
106.00	75.70	84.24			
0.00	0.00	0.00			

Data File: BM033376.D

Acq On : 10 Dec 2021 02:59

Operator : CG/JU Sample : PB141276BS

Misc

ALS Vial : 31 Sample Multiplier: 1

Quant Time: Dec 10 05:47:28 2021

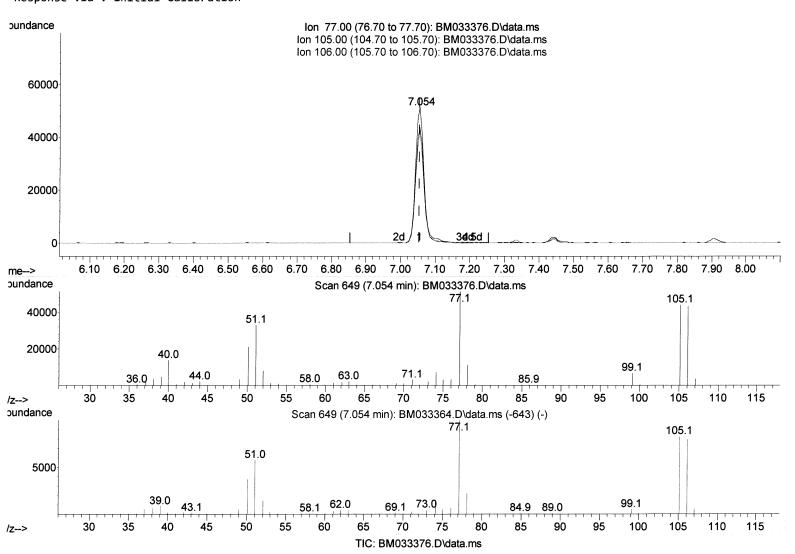
Quant Method: Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument :
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ClientSampleId :
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Benzaldehyde 35.02 ng/ul m 7.054min (+ 0.000) 85468 response Act% Ion Ехр% 77.00 100.00 100.00 105.00 82.00 85.56 106.00 75.70 84.24

0.00

0.00

0.00

Data File : BM033376.D

Acq On : 10 Dec 2021 02:59

Operator : CG/JU Sample : PB141276BS

Misc

ALS Vial : 31 Sample Multiplier: 1

Quant Time: Dec 10 05:47:28 2021

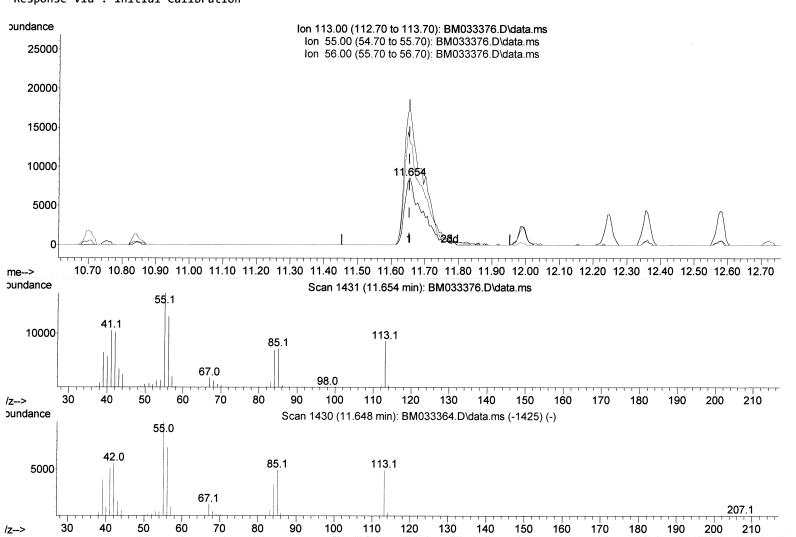
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument:
BNA_M
ClientSampleId:
SLCS276

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TIC: BM033376.D\data.ms

(34) Caprolactam

11.654min (+ 0.000) 29.15 ng/ul

response	30078				
Ion	Ежр%	Act%			
113.00	100.00	100.00			
55.00	197.40	202.69			
56.00	164.70	153.24			
0.00	0.00	0.00			

Data File : BM033376.D

Acq On : 10 Dec 2021 02:59

Operator : CG/JU Sample : PB141276BS

Misc

ALS Vial : 31 Sample Multiplier: 1

Quant Time: Dec 10 05:47:28 2021

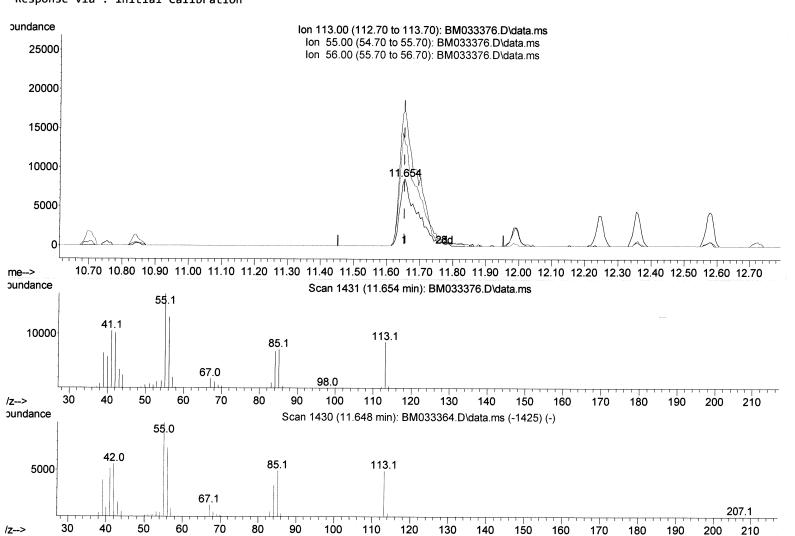
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument:
BNA_M
ClientSampleId:
SLCS276

Manual IntegrationsAPPROVED

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



TIC: BM033376.D\data.ms

(34) Caprolactam
11.654min (+ 0.000) 29.64 ng/ul m

response 30586 Ion Ехр% Act% 113.00 100.00 100.00 55.00 197.40 202.69 56.00 164.70 153.24 0.00 0.00 0.00

Data File : BM033376.D

Acq On : 10 Dec 2021 02:59

Operator : CG/JU Sample : PB141276BS

Misc

ALS Vial : 31 Sample Multiplier: 1

Quant Time: Dec 10 05:47:28 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 : SLCS276

Manual IntegrationsAPPROVED

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

Compound	R.T.	QIon	Response	Conc Unit	ts Dev(M	lin)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.907	152	46260	20.000 r	าศ/นไ	0.00
20) Naphthalene-d8	10.701	136	192100	20.000 r	_	0.00
38) Acenaphthene-d10	14.530		137385	20.000 r	_	0.00
64) Phenanthrene-d10	17.271	188	309773	20.000 r	_	0.00
79) Chrysene-d12	21.436	240	322210	20.000 r	_	0.00
88) Perylene-d12	23.753	264	308272	20.000 r		0.00
System Monitoring Compounds						0.00 \ TY(1) 2) 0.00 0.00
3) 1,4-Dioxane-d8	3.366	96	7779mプ	6.319 n	ng/uL 🤊	0.00 >TY(Y)
4) Pyridine-d5	3.784	84	101699	28.493 n	g/ul	0.00
7) Phenol-d5	7.078	99	122993	28.078 n	g/ul	0.00
9) Bis-(2-Chloroethyl)eth	7.237	67	82323	28.706 n	g/ul	0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.437	132	89430	29.159 n	g/ul -	0.01
<pre>15) 4-Methylphenol-d8</pre>	8.613	113	92698	27.036 n	g/ul (0.00
21) Nitrobenzene-d5	9.066	128	44675	28.661 n	g/ul (0.00
24) 2-Nitrophenol-d4	9.790	143	48435	30.287 n	g/ul (0.00
28) 2,4-Dichlorophenol-d3	10.325	165	92006	30.443 n	g/ul (0.00
31) 4-Chloroaniline-d4	10.842	131	108175	24.141 n	g/ul (9.00
46) Dimethylphthalate-d6	13.942	166	305518	29.763 n		9.00
49) Acenaphthylene-d8	14.225	160	377576	29.672 n		9.00
54) 4-Nitrophenol-d4	14.742	143	55464	29.759 n		9.00
60) Fluorene-d10	15.525	176	272876	29.719 n	-	9.00
65) 4,6-Dinitro-2-methylph	15.642	200	51592	27.594 n	_	0.00
73) Anthracene-d10	17.371	188	464622	30.349 n		9.00
81) Pyrene-d10	19.654	212	552774	30.696 n		0.00
92) Benzo(a)pyrene-d12	23.606	264	516521	30.912 n	g/ul 0	0.00
Target Compounds					Qvalu	ie
2) 1,4-Dioxane	3.402	88	15622	11.401 ng	g/uL	88 88 96 97
5) Pyridine	3.802	79	104803	28.510 ng		88
6) Benzaldehyde	7.054	77	85468m_D			s soully of
8) Phenol	7.102	94	124125	27.510 ng		96
10) Bis(2-Chloroethyl)ether	7.331	93	93758	27.621 ng		= :
12) 2-Chlorophenol	7.472	128	90160	28.429 ng	-	97
13) 2-Methylphenol	8.349	108	86347	26.396 ոք		96
14) 2,2'-oxybis(1-Chloropr	8.437	45	163535	27.921 ng		97
16) Acetophenone	8.731	105	152453	26.841 ng	-	99
17) N-Nitroso-di-n-propyla	8.713	70	88314	28.446 ng		97
18) 4-Methylphenol	8.678		97205	27.159 ng		98
19) Hexachloroethane	8.978	117	45050	28.117 ng		90
22) Nitrobenzene	9.113	77	135942	29.789 ng		99
23) Isophorone	9.637	82	230306	29.472 ng		99
25) 2-Nitrophenol	9.819	139	50137	29.572 ng		96
26) 2,4-Dimethylphenol27) Bis(2-Chloroethoxy)met	9.878	107	119291	29.019 ng		98
27) Bis(2-Chioroethoxy)met 29) 2,4-Dichlorophenol	10.113	93	127942	29.111 ng		98
30) Naphthalene	10.354	162	90022	29.425 ng		95
32) 4-Chloroaniline	10.754	128	302464	28.254 ng		99
33) Hexachlorobutadiene	10.866 11.025	127 225	104750	23.260 ng		98
34) Caprolactam	11.625	113	65090 30586m)	29.066 ng		92 7412/2/2/
35) 4-Chloro-3-methylphenol	11.034	107	111049	29.645 ng 30.923 ng		92 7412/23/21
, . chizoro s meenyiphenoi	11.909	10/	111047	30.323 fig	;/ u1	3 2

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Operator : CG/JU Sample : PB141276BS

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ALS Vial : 31 Sample Multiplier: 1

Quant Time: Dec 10 05:47:28 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 : SLCS276

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)
36) 2-Methylnaphthalene	12.360	142	209389	28.842 ng/ul	97
37) 1-Methylnaphthalene	12.578	142	219125	29.044 ng/ul	98
39) 1,2,4,5-Tetrachloroben	12.725	216	114225	27.595 ng/ul	99
40) Hexachlorocyclopentadiene	12.695	237	70519	26.103 ng/ul	95
41) 2,4,6-Trichlorophenol	12.966	196	72524	29.765 ng/ul	98
42) 2,4,5-Trichlorophenol	13.042	196	79786	30.227 ng/ul	91
43) 1,1'-Biphenyl	13.366	154	296257	28.187 ng/ul	99
44) 2-Chloronaphthalene	13.413	162	227058	28.079 ng/ul	100
45) 2-Nitroaniline	13.619	65	89749	31.121 ng/ul	98
47) Dimethylphthalate	13.989	163	296444	29.085 ng/ul	98
48) 2,6-Dinitrotoluene	14.113	165	59507	30.203 ng/ul	94
50) Acenaphthylene	14.254	152	388124	29.294 ng/ul	99
51) 3-Nitroaniline	14.448	138	53950	27.819 ng/ul	92
52) Acenaphthene	14.595	153	251822	28.638 ng/ul	98
53) 2,4-Dinitrophenol	14.648	184	28046	24.223 ng/ul	95
55) 4-Nitrophenol	14.754	109	61290	30.425 ng/ul	92
56) Dibenzofuran	14.930	168	374593	29.377 ng/ul	99
57) 2,4-Dinitrotoluene	14.895	165	92057	31.806 ng/ul#	99
58) 2,3,4,6-Tetrachlorophenol	15.154	232	69306	30.887 ng/ul	98
59) Diethylphthalate	15.348	149	322291	30.467 ng/ul	98
61) Fluorene	15.577	166	316226	30.161 ng/ul	100
62) 4-Chlorophenyl-phenyle	15.572	204	154619	29.563 ng/ul	98
63) 4-Nitroaniline	15.607	138	69242	34.769 ng/ul	99
66) 4,6-Dinitro-2-methylph	15.654	198	52829	28.366 ng/ul	91
67) N-Nitrosodiphenylamine	15.783	169	270908	29.744 ng/ul	99
68) 4-Bromophenyl-phenylether	16.466	248	93003	29.808 ng/ul	99
69) Hexachlorobenzene	16.572	284	102710	28.571 ng/ul	98
70) Atrazine	16.736	200	101507	28.087 ng/ul	99
71) Pentachlorophenol	16.924	266	57505	28.526 ng/ul	96
72) Phenanthrene	17.313	178	524659	29.359 ng/ul	99
74) Anthracene	17.407	178	538930	29.745 ng/ul	99
75) 1,2,3,4-Tetrachloroben	13.331	216	118922	26.618 ng/uL	100
76) Pentachlorobenzene	14.848	250	116603	26.429 ng/uL	98
77) Carbazole	17.677	167	489891	29.916 ng/ul	99
78) Di-n-butylphthalate	18.230	149	577992	31.456 ng/ul	99
80) Fluoranthene	19.318	202	639932	30.148 ng/ul	98
82) Pyrene	19.683	202	667062	30.018 ng/ul	99
83) Butylbenzylphthalate	20.571	149	268213	31.509 ng/ul	95
84) 3,3'-Dichlorobenzidine	21.354	252	185116	25.277 ng/ul	97
85) Benzo(a)anthracene	21.418	228	624935	29.520 ng/ul	99
86) Bis(2-ethylhexyl)phtha	21.336	149	386930	31.628 ng/ul	99
87) Chrysene	21.471	228	621906	29.890 ng/ul	98
89) Di-n-octyl phthalate	22.236	149	668910	29.595 ng/ul	100
90) Benzo(b)fluoranthene	23.053	252	658347	31.170 ng/ul	99
91) Benzo(k)fluoranthene	23.101	252	581562	29.738 ng/ul	99
93) Benzo(a)pyrene	23.653	252	617299	30.460 ng/ul	99
94) Indeno(1,2,3-cd)pyrene	26.130	276	641894	29.124 ng/ul	98
95) Dibenzo(a,h)anthracene	26.141	278	560025	29.211 ng/ul	99
96) Benzo(g,h,i)perylene	26.859	276	545655	29.026 ng/ul	98

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