Data File : BM033356.D

Acq On : 09 Dec 2021 13:48

Operator : CG/JU Sample : SSTDICV020

Misc

ALS Vial : 9 Sample Multiplier: 1

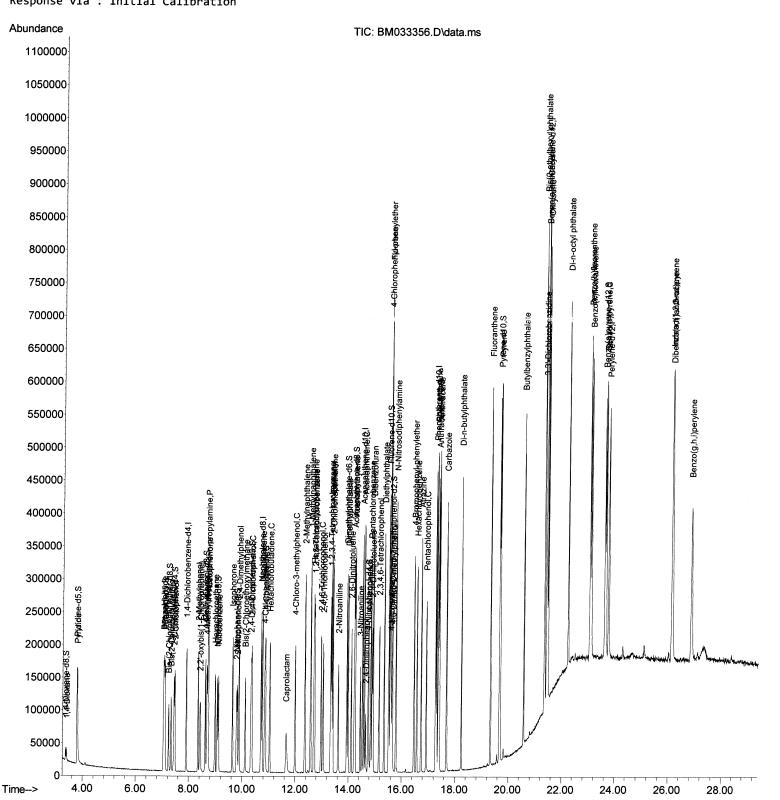
Quant Time: Dec 10 01:19:53 2021

 $\label{thm:local_power_power_local} Quant \ \ \mbox{Methods\sham-epa-bm120921.M}$

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument :
BNA_M
ClientSampleId :
SICV017

Manual IntegrationsAPPROVED

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



Data File: BM033356.D

Acq On : 09 Dec 2021 13:48

Operator : CG/JU Sample : SSTDICV020

Misc

ALS Vial : 9 Sample Multiplier: 1

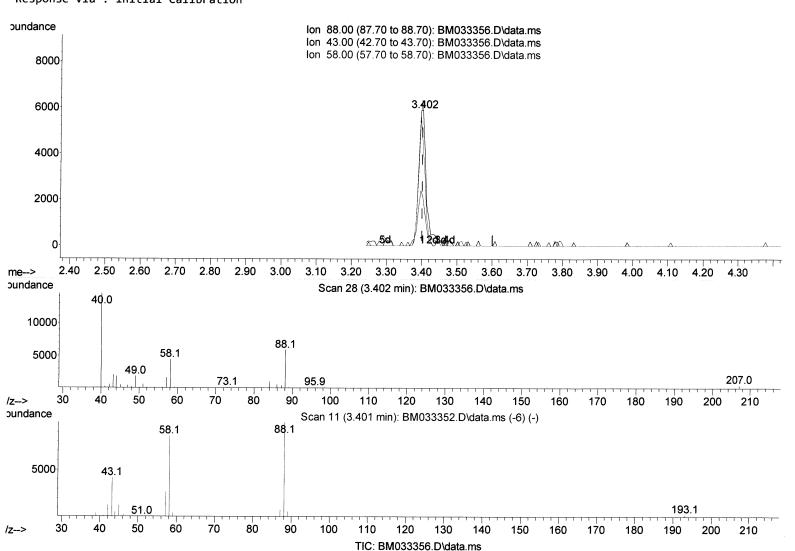
Quant Time: Dec 10 01:19:53 2021

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:
SICV017

Manual IntegrationsAPPROVED

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(2) 1,4-Dioxane

3.402min (+ 0.001) 6.66 ng/uL

response	8417	
Ion	Exp%	Act%
88.00	100.00	100.00
43.00	45.30	36.59
58.00	85.60	75.60
0.00	0.00	0.00

Data File: BM033356.D

Acq On : 09 Dec 2021 13:48

Operator : CG/JU Sample : SSTDICV020

Misc

ALS Vial : 9 Sample Multiplier: 1

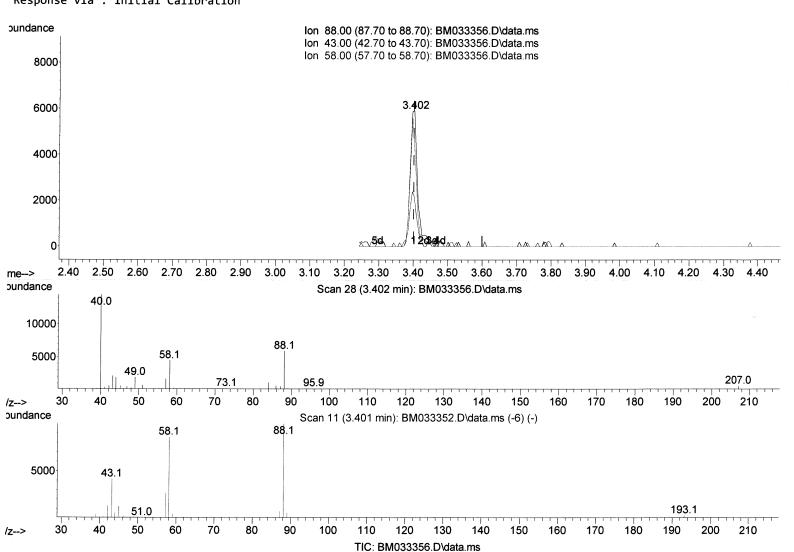
Quant Time: Dec 10 01:19:53 2021

 $\label{thm:local_model} Quant \ \ \mbox{Methods\sfam-epa-bm120921.M} \\ \ \ \mbox{Construction} \ \mbox$

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument:
BNA_M
ClientSampleId:
SICV017

Manual IntegrationsAPPROVED

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



(2) 1,4-Dioxane

3.402min	(+ 0.001)	7.18 ng/uL	T412/25/2
response	9069		3417
Ion	Ежр%	Act%	
88.00	100.00	100.00	
43.00	45.30	36.59	
58.00	85.60	75.60	
0.00	0.00	0.00	

Data File : BM033356.D

Acq On : 09 Dec 2021 13:48

Operator : CG/JU Sample : SSTDICV020

Misc

ALS Vial : 9 Sample Multiplier: 1

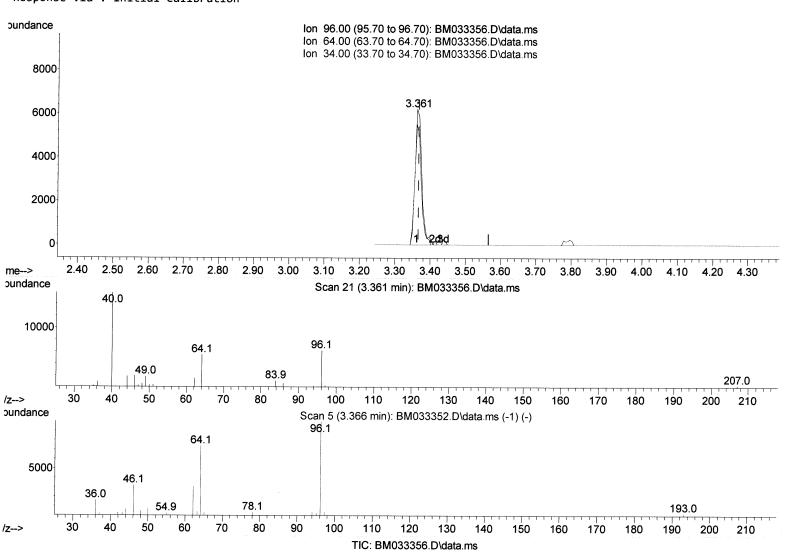
Quant Time: Dec 10 01:19:53 2021

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 : SICV017

Manual IntegrationsAPPROVED

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



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

3.361min (-0.005) 6.98 ng/uL

response	7916	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	88.70
34.00	0.00	0.00
0.00	0.00	0.00

Data File: BM033356.D

: 09 Dec 2021 13:48 Acq On

Operator | : CG/JU Sample : SSTDICV020

Misc

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Dec 10 01:19:53 2021

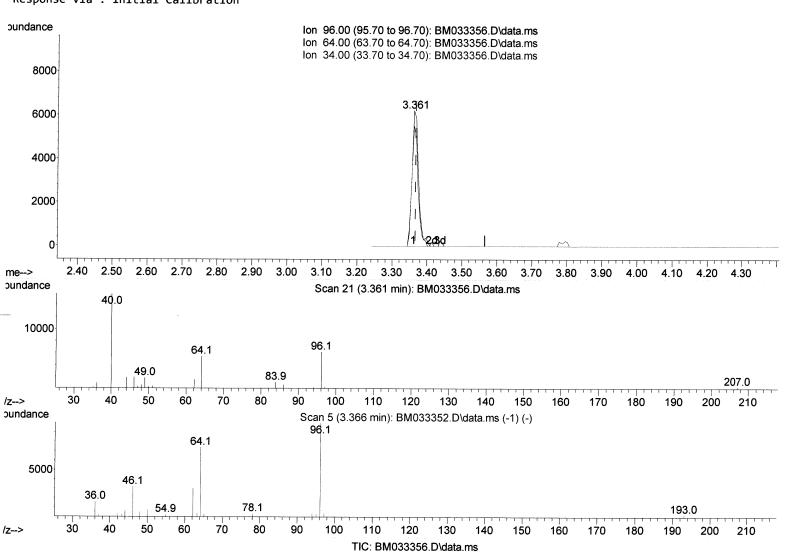
Quant Method: Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-BM120921.M

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

Instrument: BNA_M ClientSampleId : SICV017

Manual Integrations APPROVED

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(3)	Τ,	4-D1	oxan	e-as	(S)

m 12/28/21 3.361min (-0.005)7.16 ng/uL m response 8125 $\mathbf{Exp} \$$ Ion Act% 96.00 100.00 100.00 64.00 74.20 88.70 34.00 0.00 0.00

0.00

0.00

0.00

Data File: BM033356.D

Acq On : 09 Dec 2021 13:48

Operator : CG/JU Sample : SSTDICV020

Misc

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Dec 10 01:19:53 2021

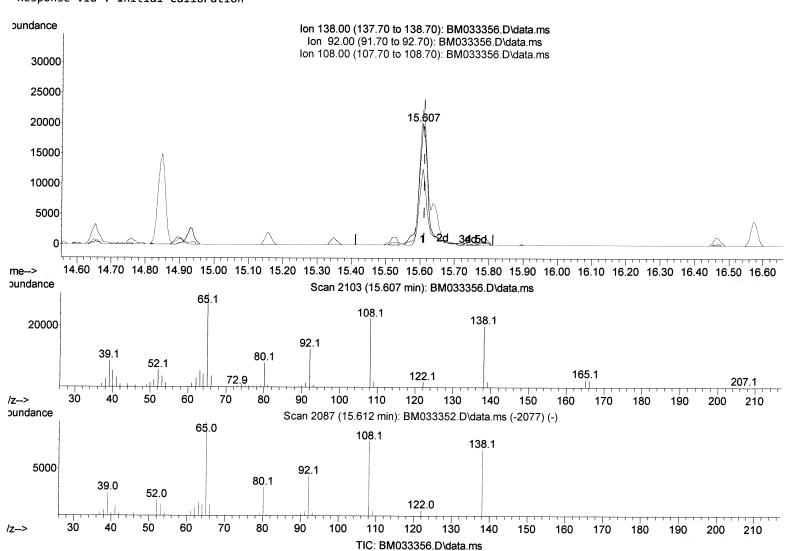
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

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



Manual IntegrationsAPPROVED

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



(63) 4-Nitroaniline

15.607min (-0.006) 19.88 ng/ul

response	35084	
Ion	Exp %	Act%
138.00	100.00	100.00
92.00	60.00	62.64
108.00	111.90	111.50
0.00	0.00	0.00

Data File : BM033356.D

Acq On : 09 Dec 2021 13:48

Operator : CG/JU Sample : SSTDICV020

Misc

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Dec 10 01:19:53 2021

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

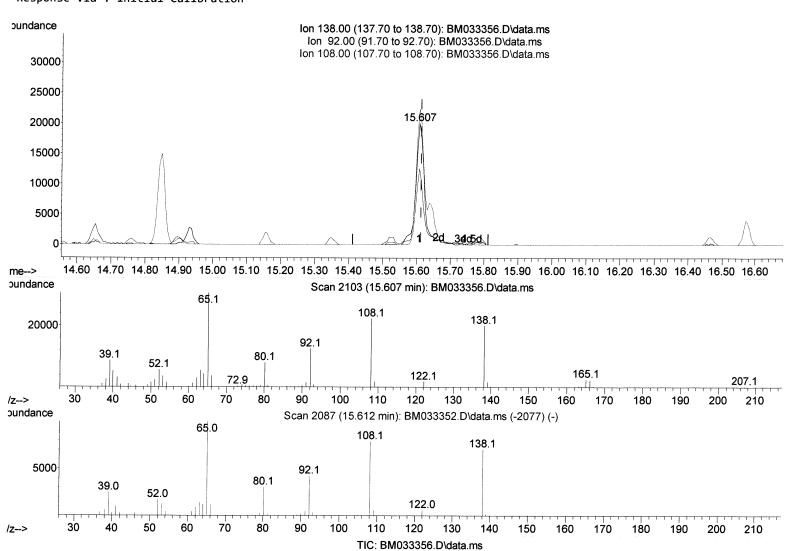
Quant Title : SVOA CALIBRATION

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



Manual IntegrationsAPPROVED

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



(63) 4-Nitroaniline

20.29 ng/ul m 1 2 2 2 1 15.607min (-0.006) response 35805 Ion Ехр% Act% 138.00 100.00 100.00 92.00 60.00 62.64 108.00 111.90 111.50 0.00 0.00 0.00

Data File : BM033356.D

Acq On : 09 Dec 2021 13:48

Jperator : CG/JU
Sample : SSTDICV020

Compound

Misc :

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Dec 10 01:19:53 2021

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

Quant Title : SVOA CALIBRATION

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

Instrument : BNA_M ClientSampleld : SICV017

Manual IntegrationsAPPROVED

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

	Compound	к. г.	Q1011	response	COILC OIL	rrs per	(MTII)		
 Tota	annol Chandanda								
	ernal Standards	7 007	150	42649	20.000	n ~ /1	0.00		
-	1,4-Dichlorobenzene-d4	7.907	152	42648	20.000		0.00		
	Naphthalene-d8 Acenaphthene-d10	10.701	136	181035 121755	20.000		0.00		
	Phenanthrene-d10	14.536 17.277	164		20.000		0.00		
	Chrysene-d12	21.436		265175 269909	20.000	-	0.00		
	Perylene-d12	23.765	240		20.000		0.00		
88)	rei yielle-uiz	23.703	264	263114	20.000	-	0.00		
Svst	em Monitoring Compounds						>0.00 0.00	1	2/21
	1,4-Dioxane-d8	3.361	96	8125m>	7.159	nø/ul	Da 00 >	770121	20(-1
-	Pyridine-d5	3.784	84	55603	16.898	ng/ul	0.00	J9 1 1	
	Phenol-d5	7.078	99	70055	17.347	ng/ul	0.00		
	Bis-(2-Chloroethyl)eth	7.237	67	46856	17.722		0.00		
	2-Chlorophenol-d4	7.443	132	51812	18.325		0.00		
	4-Methylphenol-d8	8.613	113	56540	17.887		0.00		
	Nitrobenzene-d5	9.066	128	27257	18.556		0.00		
	2-Nitrophenol-d4	9.790	143	28688	19.035	-	0.00		
	2,4-Dichlorophenol-d3	10.325	165	53556	18.804		0.00		
	4-Chloroaniline-d4	10.843		73385	17.378	_	0.00		
	Dimethylphthalate-d6	13.942	166	173130	19.031	_	0.00		
	Acenaphthylene-d8	14.231	160	216674	19.213		0.00		
	4-Nitrophenol-d4	14.742	143	28352	17.165		0.00		
•	Fluorene-d10	15.525	176	153980	18.923	_	0.00		
	4,6-Dinitro-2-methylph	15.642	200	27041	16.895		0.00		
	Anthracene-d10	17.372	188	249489	19.037		0.00		
-	Pyrene-d10	19.660	212	288873	19.149	_	0.00		
	Benzo(a)pyrene-d12	23.612	264	262211	18.386		0.00		
						O.		4	
Targ	et Compounds						alue	12/28	2)
2)	1,4-Dioxane	3.402	88	9069m 🤊	7.179	ng/uL_	> >II	190	,
5)	Pyridine	3.802	79	58142	17.156		93		
	Benzaldehyde	7.055	77	52025	23.123		89		
	Phenol	7.102	94	72970	17.542		95		
	Bis(2-Chloroethyl)ether	7.331	93	56821	18.157	ng/ul	96		
	2-Chlorophenol	7.472	128	55328	18.923	ng/ul	95		
	2-Methylphenol	8.349	108	52759	17.494	ng/ul	92		
	2,2'-oxybis(1-Chloropr	8.425	45	97783	18.109		98		
	Acetophenone	8.731	105	92371	17.640	ng/ul	98		
	N-Nitroso-di-n-propyla	8.713	70	53131	18.563		98		
	4-Methylphenol	8.678	108	57959	17.565	ng/ul	93		
	Hexachloroethane	8.984	117	27776	18.804		85		
	Nitrobenzene	9.113	77	79454	18.475	_	98		
	Isophorone	9.631	82	136327	18.512	_	99		
	2-Nitrophenol	9.819	139	30269	18.945	_	98		
	2,4-Dimethylphenol	9.878	107	71891	18.557		97		
	Bis(2-Chloroethoxy)met	10.113	93	77206	18.641		99		
	2,4-Dichlorophenol	10.354	162	55255	19.165		94		
	Naphthalene	10.754	128	188486	18.683		100		
	4-Chloroaniline	10.866	127	76754	18.085		99		
	Hexachlorobutadiene	11.031	225	39223	18.585	_	98		
	Caprolactam	11.654	113	16085	16.543		98		
35)	4-Chloro-3-methylphenol	11.990	107	62682	18.521	ng/ul	91		

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

Data File : BM033356.D

Acq On : 09 Dec 2021 13:48

Dperator : CG/JU
Sample : SSTDICV020

Misc :

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Dec 10 01:19:53 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 : SICV017

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	126524	18.493 ng/ul	98	
37) 1-Methylnaphthalene	12.578		134158	18.869 ng/ul	100	
39) 1,2,4,5-Tetrachloroben	12.725		70201	19.137 ng/ul	95	
40) Hexachlorocyclopentadiene	12.701	237	64879	27.098 ng/ul	98	
41) 2,4,6-Trichlorophenol	12.972	196	42236	19.560 ng/ul	92	
42) 2,4,5-Trichlorophenol	13.048	196	45609	19.497 ng/ul	97	
43) 1,1'-Biphenyl	13.372	154	178774	19.193 ng/ul	96	
44) 2-Chloronaphthalene	13.413	162	138617	19.342 ng/ul	100	
45) 2-Nitroaniline	13.625	65	48648	19.035 ng/ul	95	
47) Dimethylphthalate	13.989	163	170912	18.921 ng/ul	99	
48) 2,6-Dinitrotoluene	14.113	165	32944	18.868 ng/ul	93	
50) Acenaphthylene	14.254	152	223902	19.069 ng/ul	99	
51) 3-Nitroaniline	14.448	138	32246	18.762 ng/ul#	99	
52) Acenaphthene	14.595	153	146807	18.839 ng/ul	94	
53) 2,4-Dinitrophenol	14.654	184	15288	14.899 ng/ul	97	
55) 4-Nitrophenol	14.760	109	31016	17.373 ng/ul	91	
56) Dibenzofuran	14.930	168	213328	18.877 ng/ul	98	
57) 2,4-Dinitrotoluene	14.901	165	49465	19.284 ng/ul	98	
58) 2,3,4,6-Tetrachlorophenol	15.160	232	37503	18.859 ng/ul#	98	
59) Diethylphthalate	15.348	149	175887	18.762 ng/ul	99	
61) Fluorene	15.578	166	176583	19.004 ng/ul	99	121
62) 4-Chlorophenyl-phenyle	15.572	204	88027	18.992 ng/ul	- 96 TO12	25/21
63) 4-Nitroaniline	15.607	138	35805m		> .791 1	28/21
66) 4,6-Dinitro-2-methylph	15.660	198	27670	17.356 ng/ul#		
67) N-Nitrosodiphenylamine	15.789	169	146706	18.816 ng/ul	98	
68) 4-Bromophenyl-phenylether	16.466	248	51031	19.106 ng/ul	98	
69) Hexachlorobenzene	16.577	284	56413	18.332 ng/ul	94	
70) Atrazine	16.736	200	54533	17.627 ng/ul	98	
71) Pentachlorophenol72) Phenanthrene	16.924 17.319	266 178	39921	23.134 ng/ul	97	
74) Anthracene	17.319	178	283140 286246	18.509 ng/ul	99	
75) 1,2,3,4-Tetrachloroben	13.331	216	72132	18.456 ng/ul 18.861 ng/uL	99 97	
76) Pentachlorobenzene	14.848	250	72132	18.951 ng/uL	97 95	
77) Carbazole	17.677	167	253880	18.111 ng/ul	99	
78) Di-n-butylphthalate	18.230	149	291278	18.518 ng/ul	99	
80) Fluoranthene	19.324	202	334198	18.795 ng/ul	99	
82) Pyrene	19.689	202	353085	18.968 ng/ul	98	
83) Butylbenzylphthalate	20.577	149	134704	18.891 ng/ul	94	
84) 3,3'-Dichlorobenzidine	21.359	252	115599	18.843 ng/ul	96	
85) Benzo(a)anthracene	21.424	228	331092	18.670 ng/ul	99	
86) Bis(2-ethylhexyl)phtha	21.342	149	190191	18.559 ng/ul	99	
87) Chrysene	21.477	228	324213	18.602 ng/ul	98	
89) Di-n-octyl phthalate	22.242	149	326642	16.932 ng/ul	100	
90) Benzo(b)fluoranthene	23.059	252	337678	18.732 ng/ul	98	
<pre>91) Benzo(k)fluoranthene</pre>	23.106	252	309479	18.541 ng/ul	99	
93) Benzo(a)pyrene	23.659	252	317977	18.383 ng/ul	99	
94) Indeno(1,2,3-cd)pyrene	26.136	276	339241	18.034 ng/ul	99	
95) Dibenzo(a,h)anthracene	26.153	278	292107	17.852 ng/ul	97	
96) Benzo(g,h,i)perylene	26.865	276	289475	18.042 ng/ul	98	

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