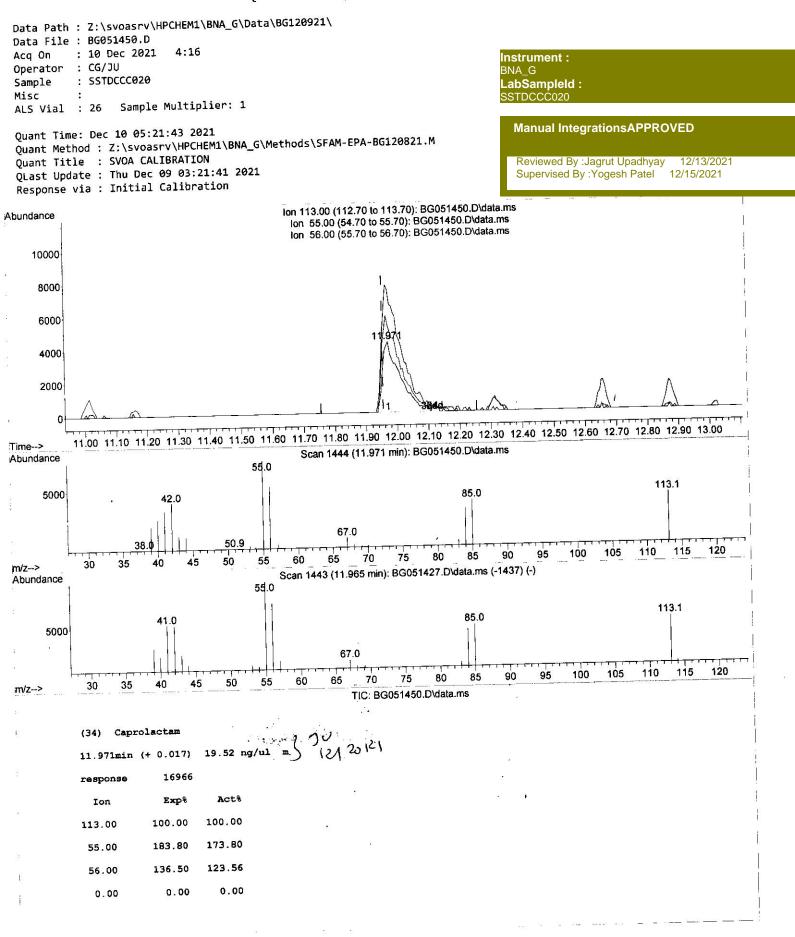
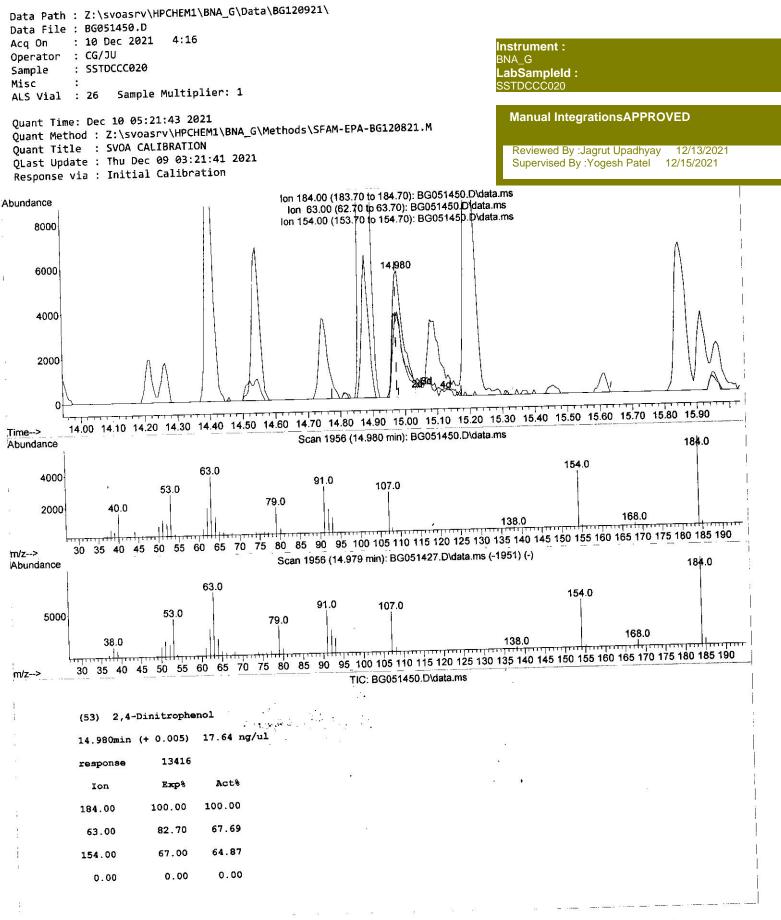
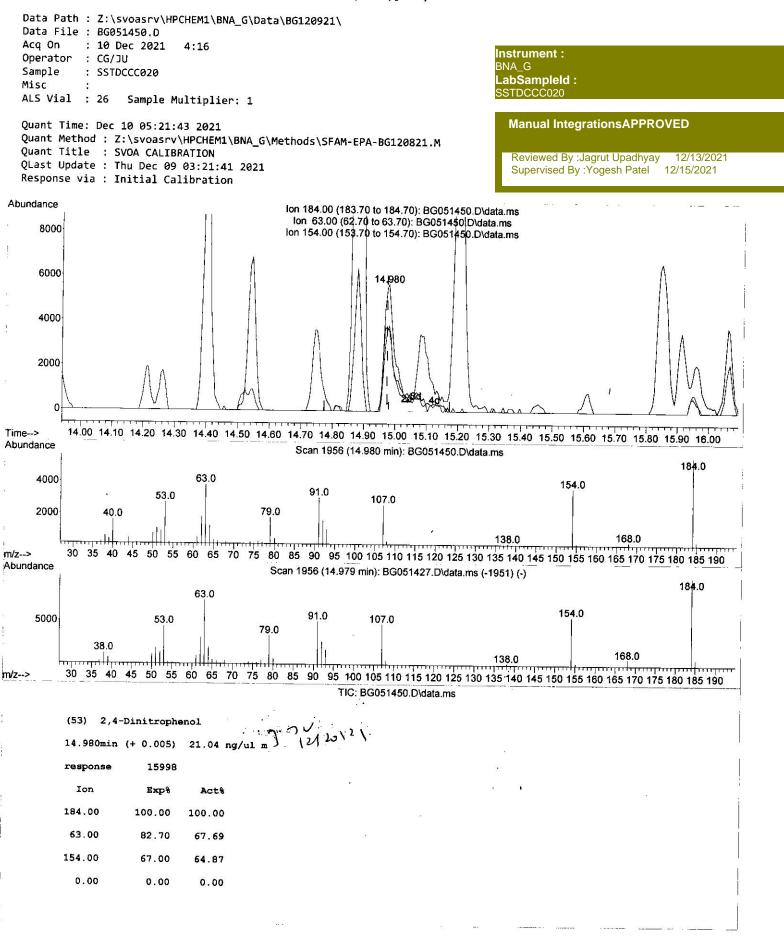


SFAM-EPA-BG120821.M Fri Dec 10 05:22:07 2021

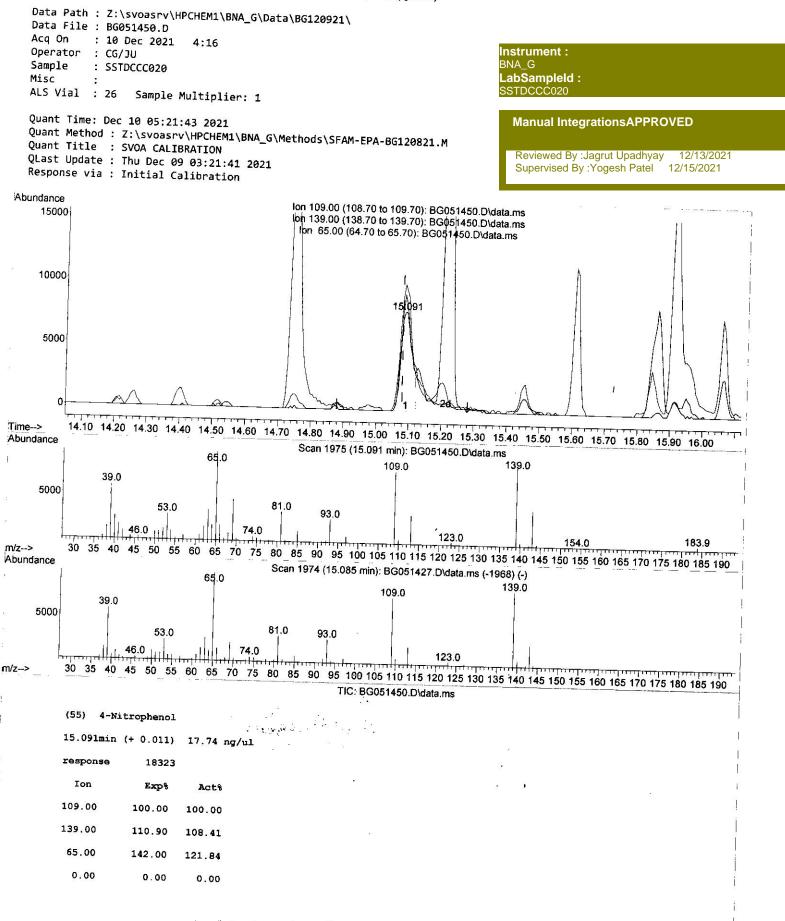




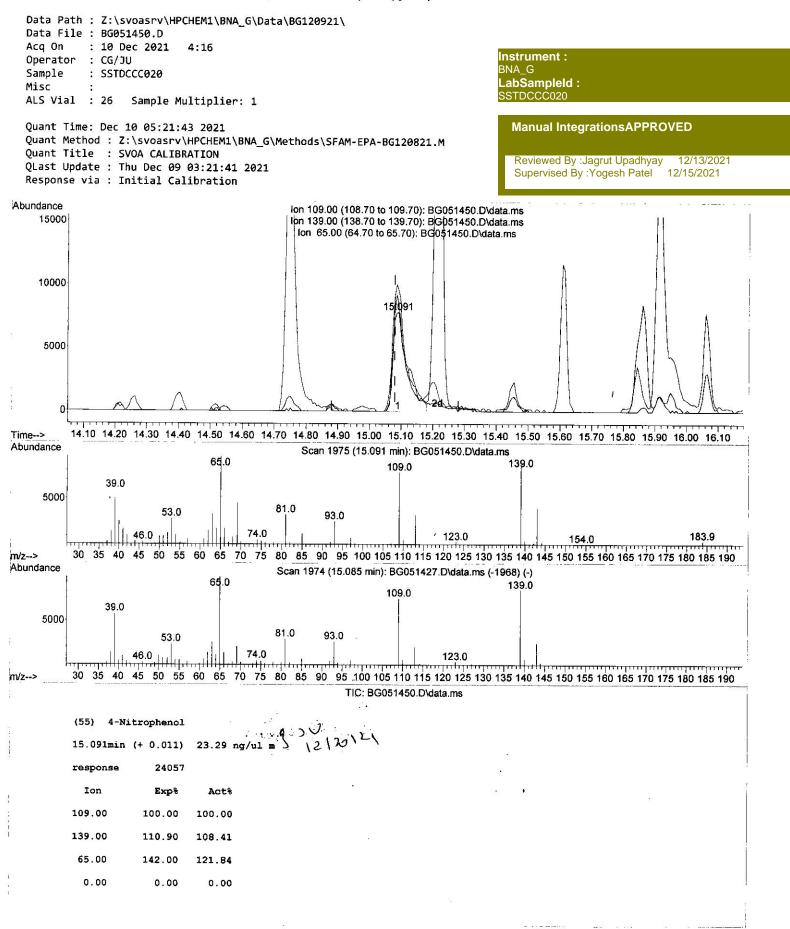




SFAM-EPA-BG120821.M Fri Dec 10 05:22:56 2021

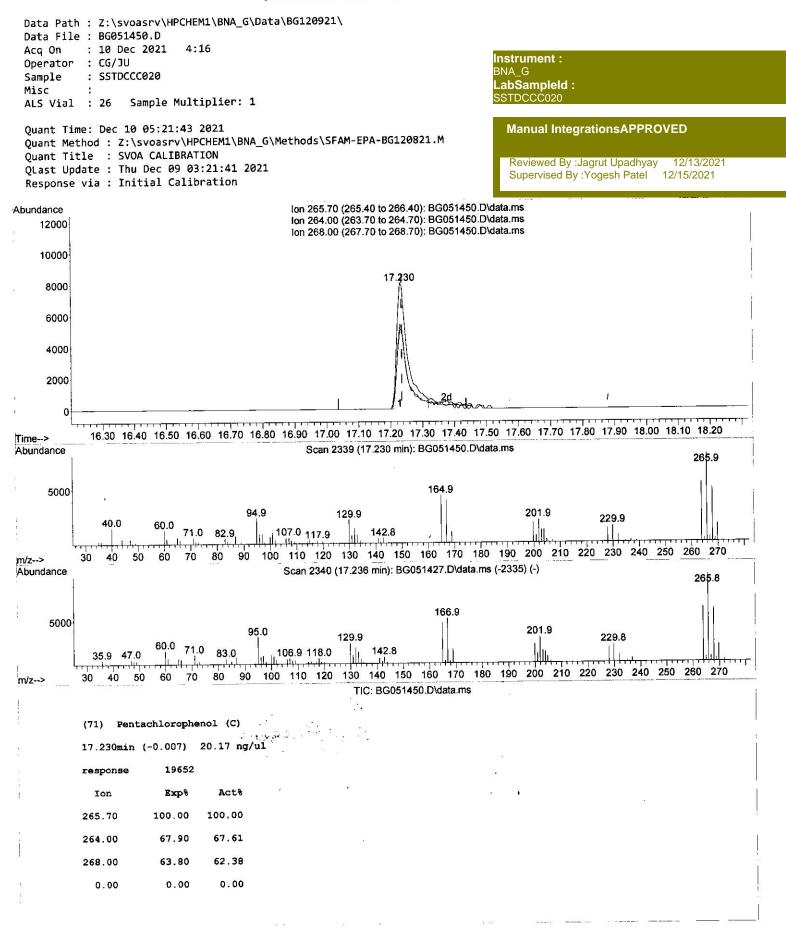


FAM-EPA-BG120821.M Fri Dec 10 05:23:10 2021

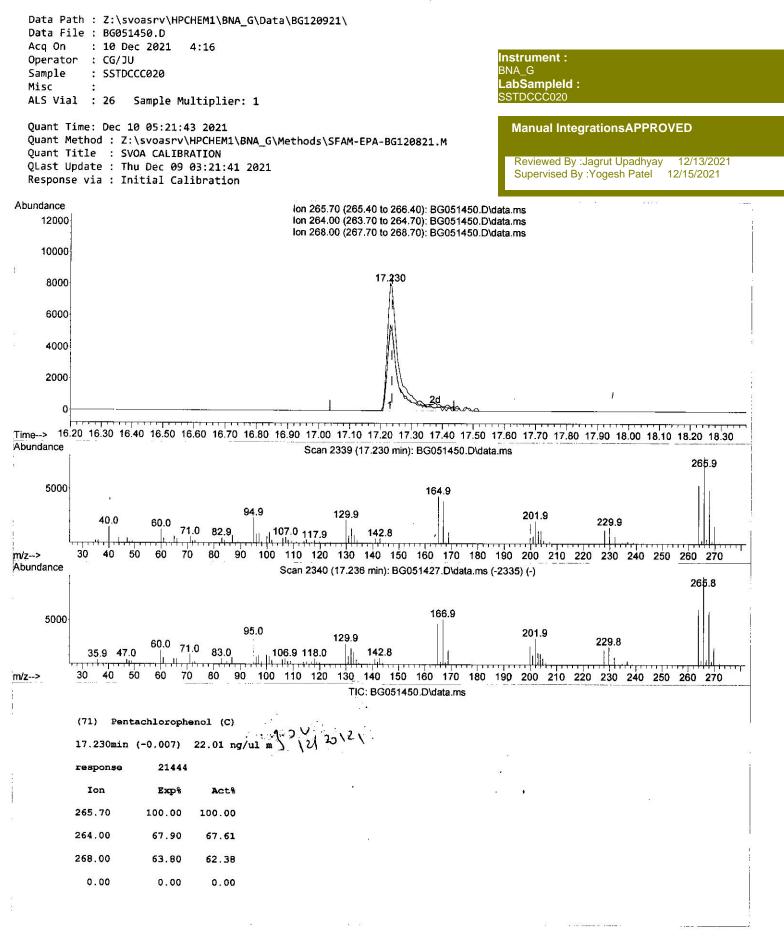


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SFAM-EPA-BG120821.M Fri Dec 10 05:23:43 2021



SFAM-EPA-BG120821.M Fri Dec 10 05:24:43 2021

	Quantitudeta	2						
Data Path : Z:\svoasrv\HPCHEM1\BNA	_G\Data\BG120921\							
Data File : BG051450.D	- *							
Aco On : 10 Dec 2021 4:16								
Acq On : 10 Dec 2021 4.10 Operator : CG/JU			Instrument :					
Sample : SSTDCCC020			BNA_G					
2 4.0 F = -			LabSampleId :					
Misc ALS Vial : 26 Sample Multiplie	r: 1		SSTDCCC020					
Quant Time: Dec 10 05:21:43 2021	\BNA_G\Methods\SFAM-EPA-F	BG120821.M	Manual IntegrationsAPPROVED					
Quant litle : Svon Children 09 03:21:4	1 2021		Reviewed By :Jagrut Upadhyay 12/13/2021 Supervised By :Yogesh Patel 12/15/2021					
QLast Opdate . Initial Calibration Response via : Initial Calibration	R.T. QIon Response C	one Units Dev(Min)						
Compound	R.T. QION Response C							
Internal Standards	o 188 152 30255	20.000 ng/ul 0.0						
1) 1.4-Dichlorobenzene-04	8.188 152 30255 11.014 136 133966	20.000 ng/ul 0.0						
20) Naphthalene-d8	14.815 164 89112	20.000 ng/ul 0.0						
38) Acenaphthene-d10	17.571 188 218656	20.000 ng/ul 0.0						
64) Phenanthrene-d10	21.871 240 205859	20.000 ng/ul 0.0						
79) Chrysene-d12	25.262 264 201772	20.000 ng/ul -0.0	1					
88) Perylene-d12	231202							
System Monitoring Compounds	1700	7.359 ng/uL 0.0	9					
3) 1,4-Dioxane-d8	3.528 96 6780	16.712 ng/u1 0.0	0					
4) Pyridine-d5	3.969 84 44214	18.107 ng/ul 0.0	0					
7) Phenol-d5	7.359 99 55771	18.126 ng/ul 0.0	0					
9) Bis-(2-Chloroethyl)eth	7.500 67 35802 7.723 132 40917	18.672 ng/ul 0.0	0					
11) 2-Chlorophenol-d4	11/20 47551	17.999 ng/ul 0.0	0					
15) 4-Methylphenol-d8	0.910 111	18.295 ng/ul 0.0	00					
21) Nitrobenzene-d5	5.500	18.803 ng/ul 0.0						
adv 2-Nitrophenol-04	10.031 1.0	18.835 ng/ul 0.6						
28) 2.4-Dichlorophenol-us	10.045	18.016 ng/ul 0.0						
21) A-Chloroanlline-04	11.160 131 56372 14.216 166 140536	20.381 ng/ul 0.0						
46) Dimethylphthalate-00	14.515 160 168349	19.277 ng/ul 0.						
49) Acenaphthylene-08	15.068 143 18548	17.866 ng/ul 0.						
54) 4-Nitrophenol-04	15.808 176 120969	19.700 18/ 4-	00					
60) Fluorene-d10		19.302 115/ 44	00					
65) 4,6-Dinitro-2-methylph	17.671 188 204174	T2'22' (19'	00					
73) Anthracene-d10	19.950 212 242098		00					
81) Pyrene-d10 92) Benzo(a)pyrene-d12	25.032 264 205849	19.780 MB/ 44						
92) Benzo(a)pyrene ===	Qvalue							
Target Compounds	2 563 88 7237	7.040 ng/uL	96					
2) 1,4-Dioxane	5.505 00 17506	17.235 ng/ul	96					
5) Pyridine	3.307 75 40170	20.515 ng/ul	95					
6) Benzaldehyde	7.330 77 40170 7.388 94 58522	18.563 ng/ul	98					
e) Phonol	7.300 57 44079	18.254 ng/ul	98 94					
10) Bis(2-Chloroethyl)ether	7.753 128 41391	18.440 ng/ul	94 96					
12) 2-Chlorophenol	8,646 108 43413		96					
13) 2-Methylphenol		10 007 00/11	97					
14) 2,2'-oxybis(1-Chloropr			99					
16) Acetophenone 17) N-Nitroso-di-n-propyla	8.987 70 41695		97					
17) N-Nitroso-ul i p.ep. 18) 4-Methylphenol	01.01.0	40 305 0g/11]	- 98					
19) Hexachloroethane	9.263 117 17844		, 99					
22) Nitrobenzene	9.410 77 59818		99					
23) Isophorone	9.927 82 111687		99					
25) 2-Nitrophenol	10.120		97					
ach a A-Dimethylphenol	10.1/0	8 18.707 ng/ul	99					
27) Bis(2-Chloroetnoxy)mer	101-102 4053	4 19.321 ng/ul	97					
29) 2,4-Dichlorophenoi	12012	9 18.913 ng/ul	98 0V					
30) Nanhthalene	11.001	6 18.674 ng/ul	35 DI 21.11					
22) A-Chloroaniline	11.184 127 58/7 11.319 225 2637	1 a 18.433 ng/ul	91 101					
33) Hexachlorobutadiene	11, 212	56m \ 19.521 ng/ul	06					
ad concolactam	11.3/1 444	- +0 001 ng/11	96					
35) 4-Chloro-3-methylphen			1					
	0 05-25-18 2021							

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SFAM-EPA-BG120821.M Fri Dec 10 05:25:18 2021

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120921\ Data File : BG051450.D Acq On : 10 Dec 2021 4:16 Operator : CG/JU Sample : SSTDCCC020 Misc : ALS Vial : 26 Sample Multiplier: 1 Quant Time: Dec 10 05:21:43 2021 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG120821.M Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 03:21:41 2021 Response via : Initial Calibration

Instrument :

BNA\_G LabSampleId :

SSTDCCC020

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/13/2021 Supervised By :Yogesh Patel 12/15/2021

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12/20121

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Response via : inicial calls acto	••				
Compound	R.T. Q	Ion R		Conc Units Dev(Mi	n)  99
36) 2-Methylnaphthalene		142	93408	19.040 ng/ul 19.235 ng/ul	94
27) 1-Methylnaphthalene		142	97130	19.153 ng/ul	97
39) 1,2,4,5-Tetrachloroben	13.017	216	53145	18.683 ng/ul	96
40) Hexachlorocyclopentadiene	12.982	237	27451	19.576 ng/ul	97
41) 2,4,6-Trichlorophenol	13.270	196	35089	18.566 ng/ul	95
42) 2,4,5-Trichlorophenol	13.364	196	35628	19.299 ng/ul	95
43) 1,1'-Biphenyl	13.652	154	128571	19.504 ng/ul	98
44) 2-Chloronaphthalene	13.705	162	101949	19.661 ng/ul	93
45) 2-Nitroaniline	13.922	65	38845	20.338 ng/ul	100
47) Dimethylphthalate	14.263	163	141334 30039	20.427 ng/ul	97
48) 2,6-Dinitrotoluene	14.404	165	163238	18.943 ng/ul	99
50) Acenaphthylene	14.545	152	31070	21.940 ng/ul	98
51) 3-Nitroaniline	14.745	138	108575		98
52) Acenaphthene	14.880	153	15998m	21.036 ng/ul	
53) 2,4-Dinitrophenol	14.980	184	24057m	23.294 ng/ul	
55) 4-Nitrophenol	15.091	109	157178	19.607 ng/ul	100
56) Dibenzofuran	15.215	168 165	44096	20.978 ng/ul	95
57) 2 A-Dinitrotoluene	15.197	232	31162	21.432 ng/ul#	97
58) 2.3.4,6-Tetrachlorophenol	15.455 15.614	149	151910	20.254 ng/ul	99
59) Diethylphthalate		166	129609	19.964 ng/ul	99
61) Fluorene	15.861 15.843	204	67020	19.664 ng/ul	98
62) 4-Chlorophenyl-phenyle	15.845	138	28964	23.041 ng/ul	92
63) 4-Nitroaniline		198	25174	19.928 ng/ul	94
66) 4,6-Dinitro-2-methylph	16.067	169	119519	19.608 ng/ul	99
67) N-Nitrosodiphenylamine			41748	18.909 ng/ul	94
68) 4-Bromophenyl-phenylether	16.871		44325	19.696 ng/ul	98
69) Hexachlorobenzene	17.007		49543	18.837 ng/ul	98
70) Atrazine	17.230		21444m	22.011 ng/ul	
71) Pentachlorophenol	17.612		232132	19.704 ng/ul	99
72) Phenanthrene	17.706	1	233669	19.814 ng/ul	98
74) Anthracene			55432	18.129 ng/uL	97
75) 1,2,3,4-Tetrachloroben	15.132		51839	18.726 ng/uL	97 99
76) Pentachlorobenzene	17.982		215565	20.534 ng/ul	100
77) Carbazole 78) Di-n-butylphthalate	18.499		278120	19.760 ng/ul	98
80) Fluoranthene	19.61	5 202		19.375 ng/ul	97
80) Fluoranchene	19.98	0 202		19.679 ng/ul	96
82) Pyrene 83) Butylbenzylphthalate	20.83	1, 149			98
84) 3,3'-Dichlorobenzidine	21.75				99
85) Benzo(a)anthracene	21.84				99
86) Bis(2-ethylhexyl)phtha.	. 21.70			10 705	.99
87) Chrysene	21.91		3 255884		100
89) Di-n-octyl phthalate	22.95				99
90) Benzo(b)fluoranthene	24.18	Second Second			99
91) Benzo(k)fluoranthene	24.24				98
93) Benzo(a)pyrene	25.10			1 / ]	96
94) Indeno(1,2,3-cd)pyrene	29.18		5		95
95) Dibenzo(a,h)anthracene	29.23			10 202 20/11	97
96) Benzo(g,h,i)perylene	30.42	20 27	0 21445		
					cummed

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

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