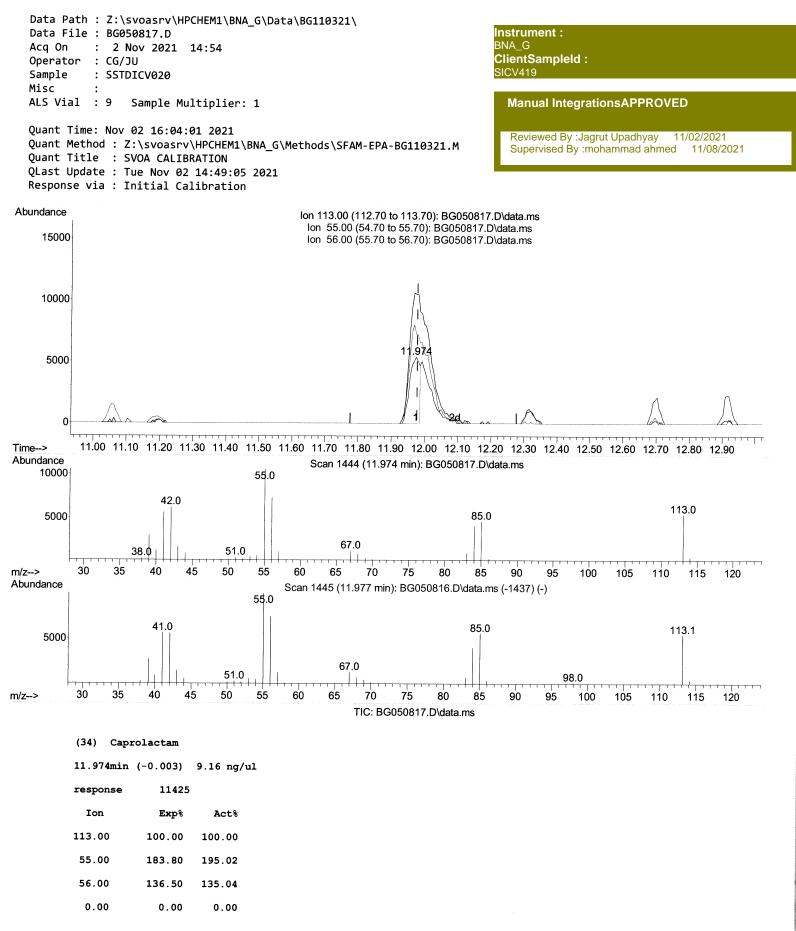
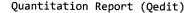
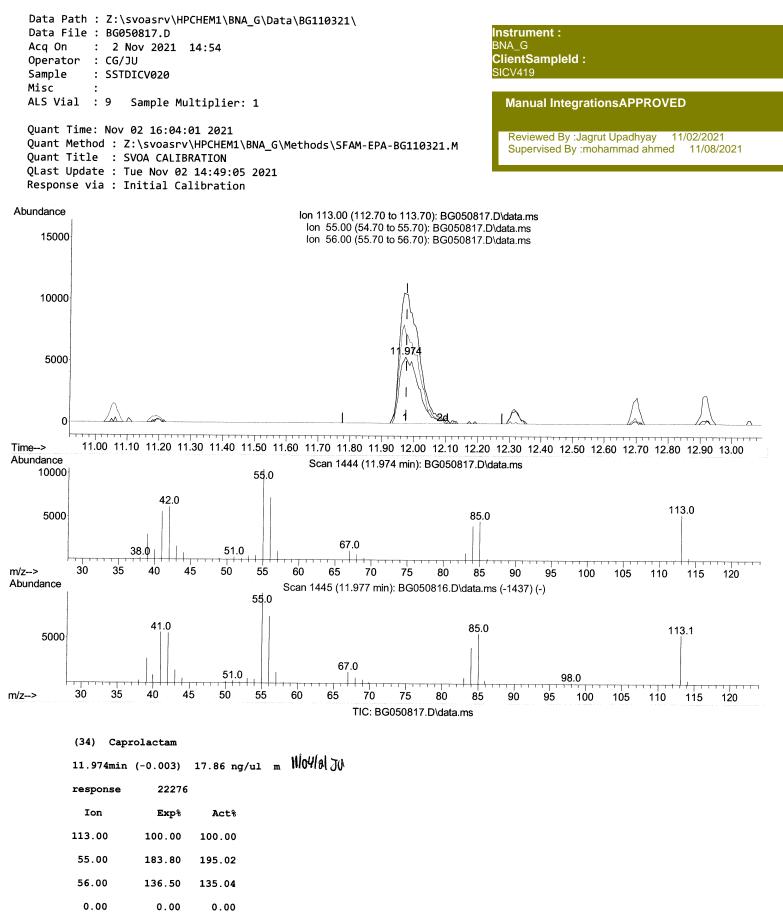
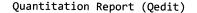
Data File : Acq On : Operator :	2 Nov 2021 1		5∖Data∖BG1	10321\			BNA_	itSample	eld :				
ALS Vial :									Manual IntegrationsAPPROVED				
Quant Method Quant Title QLast Update	Nov 02 16:04:0 : Z:\svoasrv\ : SVOA CALIBF : Tue Nov 02 : Initial Cal	HPCHEM1\BN ATION 14:49:05 2		ds\SFAM-EP	A-BG1103	21.M	ReSu	eviewed B ipervised	y :Jagrut L By :mohar	Jpadhyay nmad ahn	11/02/20 ned 11/08		
Abundance				TIC:	BG050817	.D\data.ms	6						
650000													
600000					le-d10,1		ate palate						
550000				Howyphratesets.S. Accentacementaries.Kenaphthene-Acenaphthene-d10,I Accentacementaries. Accentacementaries. Achieved a construction of the second action of	—————————————————————————————————————	Fluoranthene Pyrene d:10,S	Butylbenzylphthalate						
500000				10,1 	a魏mm御地略	- Linor Linor	But	thalate					
450000				enaphthene-d' liate Votene-d10 S	Anth Carbazole		ine	Di-n-octyl phthalate					
400000			er Ivnehaid	Diethylphthe	phenylether Atrazine		3.3 - Dichlorobenzidine	Ben <mark>2ิจุตรุสน์</mark> ไปใช้เหลี่กลิกษ์nene					
350000			ប្រុសនៃជាអង់គេមេនាំ សំអីអីរីបិនក្រុងកំណើរទទេក អំអីតអំអីរីសំអារីសាម័ណនៃតែ A - Bithenvi	Dimetterity in the second seco	adnBr0BBR2envl-phenylether Atrazine		<del>8</del>	Ben Bonk	ଅଖିକୀ)୪% କିne,C				
300000	S, 3D-lor	lamine,P nol Bisserane-d8,I	ohenol,C 2-Methyl 1-Tetrachlorob6 henol,C 3-4-Tetrachleri	uene Dimettat Mutan Spentach Chlorophenol	Hexad				Perylene-a 2,4		hracene		
250000	Bisa 2.6 នាមសាស់ស្តាំងត្រូវថ្មានទាំងទាល់-d5.S 14-Dichlorobenzene-d4,I	Optimization         State           Proversition         State	Hexacinoroutaments,C 4-Chloro-3-methylphenol,C 2.4.5746161616616461610 2.4.574461161616466161001 2.4.57446116161646161001 2.4.57446116161646161001 2.4.574461161616	z-vurceanline Litroanline Litroanline - <u>4.Nitrontrotourceanline</u> - 2.3.4.6. Tetrachiorophenc - 4.4 Miscrittione onternegorgen							Indeno <b>(ປີເຊີຣຳຂຢູ່(ສ_ກັງຢູາຕ</b> ີກາສcene (g.h.i)perylene		
200000 ويتركي المراجع		<u>-oxybis(1-Chiogonal)()()()()()</u> Hewanikashanga, s 2000 -	нехаслю 4-С 16-С 2,4,52	د-۲۰۹۵ کی 16-2 3-Nitroaniline 4-Nitro 3-23	ophenol,C						Indeno <b>();ଌିର୍ଜ୍ୟୁ</b> Benzo(g,h,i)perylene		
150000	Bital	2.2 <sup>-</sup> oxybis(1-( Hergerals			Pentachlorophenol, C								
100000 <sup>%</sup> 87 <b>9</b> <b>9</b> <b>9</b> <b>9</b> <b>9</b> <b>1</b> <b>1</b> <b>1</b> <b>1</b> <b>1</b> <b>1</b> <b>1</b> <b>1</b> <b>1</b> <b>1</b>			Caprolactam										
50000							Jury Human					<b></b>	
0 Time> 4.00	6.00 8.00	<u>ייי קירקי אין אין אין אין אין אין אין אין אין אי</u>	12.00 14.	00 16.00	18.00 م) المالية 18.00	مستعابة الليب 20.00	22.00	24.00	26.00	28.00	30.00	32.00	

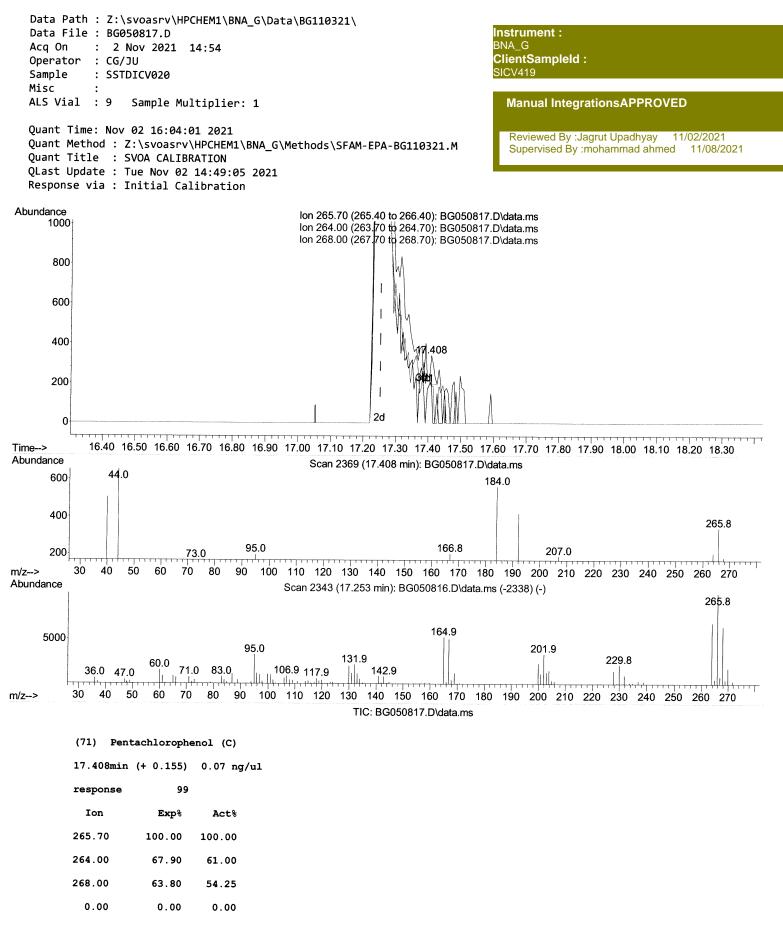


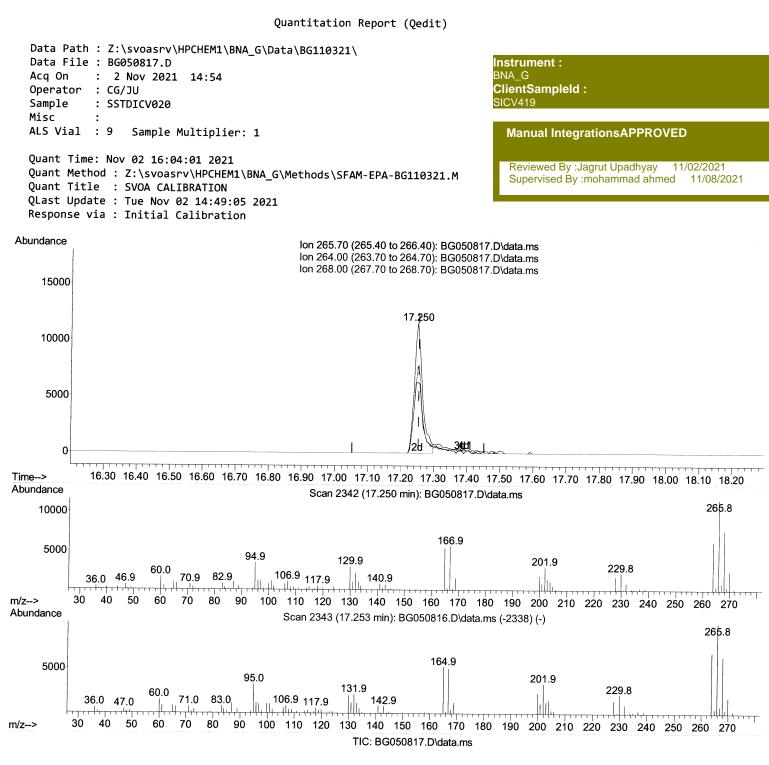








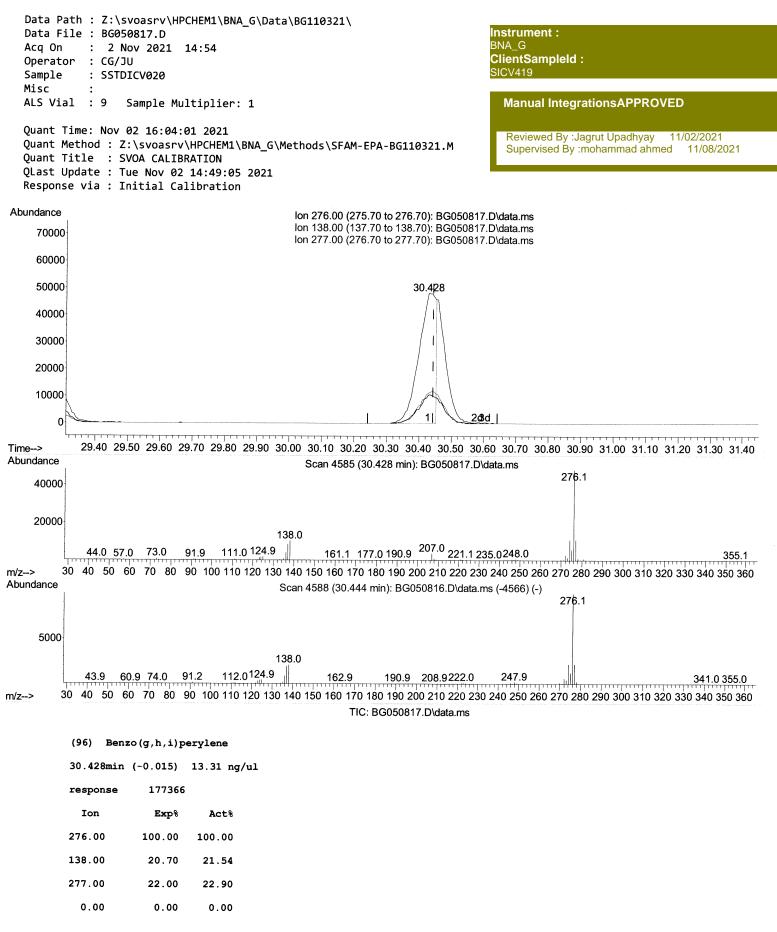


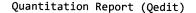


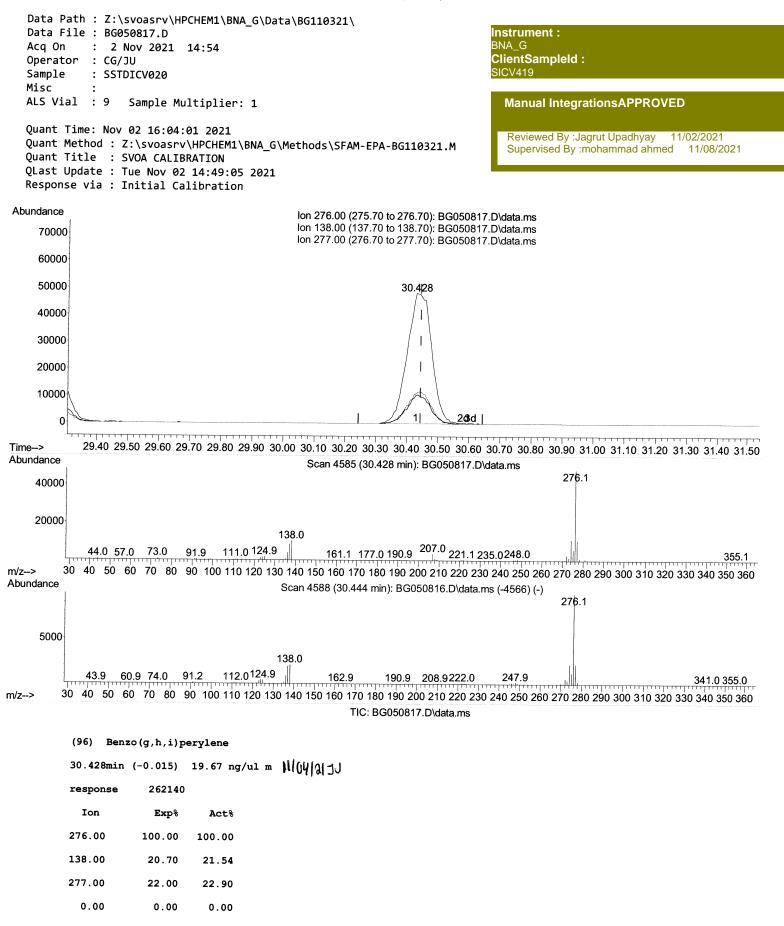
(71) Pentachlorophenol (C)

17.250min (-0.003) 14.26 ng/ul m 1/64/8/JU response 20233 Ion Exp\* Act% 265.70 100.00 100.00 264.00 67.90 54.31# 268.00 63.80 66.61 0.00 0.00 0.00









Data Path : Z:\svoasrv\HPCHEM1 Data File : BG050817.D Acq On : 2 Nov 2021 14:54 Operator : CG/JU Sample : SSTDICV020 Misc : ALS Vial : 9 Sample Multipl Quant Time: Nov 02 16:04:01 20	ier: 1	Instrument : BNA_G ClientSampleId : SICV419 Manual IntegrationsAPPROVED				
Quant Method : Z:\svoasrv\HPCH Quant Title : SVOA CALIBRATIO QLast Update : Tue Nov 02 14:49 Response via : Initial Calibra	EM1\BNA_0 N 9:05 2021		ods\SFAM-E	PA-BG110321.M	Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021	
Compound	R.T.	QIon	Response	Conc Units Dev	(Min)	
Internal Standards						
<ol> <li>1,4-Dichlorobenzene-d4</li> </ol>	8.231	152	39092	20.000 ng/ul	0.00	
20) Naphthalene-d8	11.057		189259	20.000 ng/ul	0.00	
38) Acenaphthene-d10	14.858		132408	20.000 ng/ul	0.00	
64) Phenanthrene-d10	17.602		302274	20.000 ng/ul	0.00	
79) Chrysene-d12	21.903		247939	20.000 ng/ul	0.00	
88) Perylene-d12	25.305		210791	20.000 ng/ul	0.00	
, , , , , , , , , , , , , , , , , , , ,	201000	204	210/51	20.000 mg/ui	0.00	
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.589	96	8998	7.429 ng/uL	0.00	
4) Pyridine-d5	4.012		64839	17.894 ng/ul	0.00	
7) Phenol-d5	7.373		72972	17.497 ng/ul	0.00	
9) Bis-(2-Chloroethyl)eth	7.543	67	49288	18.296 ng/ul	0.00	
11) 2-Chlorophenol-d4	7.755	132	53400	18.476 ng/ul	0.00	
15) 4-Methylphenol-d8	8.930	113	58586	17.844 ng/ul	0.00	
21) Nitrobenzene-d5	9.406	128	28435	17.680 ng/ul	0.00	
24) 2-Nitrophenol-d4	10.129	143	31851	17.810 ng/ul	0.00	
<pre>28) 2,4-Dichlorophenol-d3</pre>	10.675	165	53272	17.684 ng/ul	0.00	
31) 4-Chloroaniline-d4	11.192	131	79993	17.535 ng/ul	0.00	
46) Dimethylphthalate-d6	14.253	166	174985	17.274 ng/ul	0.00	
<pre>49) Acenaphthylene-d8</pre>	14.553		216547	17.158 ng/ul	0.00	
54) 4-Nitrophenol-d4	15.046	143	31652	17.233 ng/ul	0.00	
60) Fluorene-d10	15.845		151890	16.926 ng/ul	0.00	
65) 4,6-Dinitro-2-methylph			30028	16.385 ng/ul	0.00	
73) Anthracene-d10	17.702		242908	16.997 ng/ul	0.00	
81) Pyrene-d10	19.976		286818	17.910 ng/ul	0.00	
92) Benzo(a)pyrene-d12	25.070	264	232363	19.940 ng/ul	0.00	
Target Compounds Qvalue						
2) 1,4-Dioxane	3.624	88	10357	-		
5) Pyridine	4.030	79	69556	7.785 ng/uL# 18.545 ng/ul	93 96	
6) Benzaldehyde	7.361	77	48566	18.463 ng/ul	97	
8) Phenol	7.402	94	77068	17.864 ng/ul	98	
<pre>10) Bis(2-Chloroethyl)ether</pre>	7.637	93	59565	18.446 ng/ul	99	
12) 2-Chlorophenol	7.790	128	53864	18.355 ng/ul	99	
13) 2-Methylphenol	8.666	108	57119	17.913 ng/ul	100	
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.748	45	92917	18.268 ng/ul	100	
16) Acetophenone	9.059	105	93054	18.245 ng/ul	98	
17) N-Nitroso-di-n-propyla	9.036	70	55726	18.109 ng/ul#	96	
18) 4-Methylphenol	8.995	108	61760	18.190 ng/ul	99	
19) Hexachloroethane	9.324	117	23111	18.836 ng/ul	95	
22) Nitrobenzene	9.447	77	78535	17.508 ng/ul	97	
23) Isophorone 25) 2-Nitrophenol	9.964 10.164	82	154375	17.733 ng/ul	100	
		139	32478	18.102 ng/ul	98	
<pre>26) 2,4-Dimethylphenol 27) Bis(2-Chloroethoxy)met</pre>	10.211 10.446	107	70010	17.730 ng/ul	95	
27) Bis(2-Chlorophenol 29) 2,4-Dichlorophenol	10.446	93 162	83167	17.730 ng/ul	97	
30) Naphthalene		162	52353	17.817 ng/ul	97	
32) 4-Chloroaniline	11.110 11.216	128	180310	17.423 ng/ul	97	
33) Hexachlorobutadiene	11.216	127 225	79910 34448	17.641 ng/ul	100	
34) Caprolactam	11.974	113	22276m >	17.862 ng/ul 17.857 ng/ul >	97  1  64  21  Ju	
35) 4-Chloro-3-methylphenol	12.314	107	65580	17.471 ng/ul >	96	
· · · · · · · · · · · · · · · · · · ·					20	

Data Path : Z:\svoasrv\HPCHEM1	BNA_G\Da	ata\BG	110321\					
Data File : BG050817.D					Instrument :			
Acq On : 2 Nov 2021 14:54		BNA_G						
Operator : CG/JU					ClientSampleId: SICV419			
Sample : SSTDICV020 Misc :					000413			
ALS Vial : 9 Sample Multipli	er·1				Manual IntegrationsAPPROVED			
itat i s sampte harcipi								
Quant Time: Nov 02 16:04:01 202	1				Decisional Decisional Line disease 44/00/0004			
Quant Method : Z:\svoasrv\HPCHE	M1\BNA_G	i\Metho	ds\SFAM-EF	PA-BG110321.M	Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021			
Quant Title : SVOA CALIBRATION								
QLast Update : Tue Nov 02 14:49								
Response via : Initial Calibrat	ion							
Compound	ВТ	OTon	Response	Conc Units Dev(	(Min)			
					·····			
36) 2-Methylnaphthalene	12.702	142	123263	17.483 ng/ul	99			
37) 1-Methylnaphthalene	12.919	142	125617	17.584 ng/ul	98			
<pre>39) 1,2,4,5-Tetrachloroben</pre>			65740	17.040 ng/ul	96			
<ol><li>40) Hexachlorocyclopentadiene</li></ol>			28905	15.602 ng/ul	97			
41) 2,4,6-Trichlorophenol	13.296		43407	17.194 ng/ul	99			
42) 2,4,5-Trichlorophenol	13.372		46611	17.198 ng/ul	97			
43) 1,1'-Biphenyl	13.695		166897	17.245 ng/ul	96			
44) 2-Chloronaphthalene 45) 2-Nitroaniline	13.742 13.942		129195 51948	17.035 ng/ul	96			
47) Dimethylphthalate	13.942		175853	17.240 ng/ul 17.361 ng/ul	97 98			
48) 2,6-Dinitrotoluene	14.424		37129	17.514 ng/ul	94			
50) Acenaphthylene	14.582		217033	17.158 ng/ul	98			
51) 3-Nitroaniline	14.759		36839	16.802 ng/ul	84			
52) Acenaphthene	14.923	153	140318	16.871 ng/ul	94			
53) 2,4-Dinitrophenol	14.970	184	17430	14.904 ng/ul	95			
55) 4-Nitrophenol	15.058		29495	17.510 ng/ul	90			
56) Dibenzofuran	15.252		203021	17.053 ng/ul	99			
57) 2,4-Dinitrotoluene	15.211	165	52972	17.509 ng/ul	97			
58) 2,3,4,6-Tetrachlorophenol 59) Diethylphthalate		232	36963	17.354 ng/ul	97			
61) Fluorene	15.652 15.904	149 166	189295 159619	17.459 ng/ul 16.939 ng/ul	99			
62) 4-Chlorophenyl-phenyle	15.887	204	84613	17.244 ng/ul	96 97			
63) 4-Nitroaniline	15.922	138	37550	17.263 ng/ul	97			
66) 4,6-Dinitro-2-methylph	15.975	198	29744	16.643 ng/ul	94			
67) N-Nitrosodiphenylamine	16.098	169	143962	17.038 ng/ul	99			
68) 4-Bromophenyl-phenylether	16.780	248	51165	17.016 ng/ul	97			
69) Hexachlorobenzene	16.903	284	52427	16.961 ng/ul	94			
70) Atrazine	17.038	200	61714	17.224 ng/ul	99			
71) Pentachlorophenol 72) Phenanthrene	17.250	266	20233m>		noamija			
74) Anthracene	17.643 17.737	178 178	277691 277250	17.210 ng/ul	99			
75) 1,2,3,4-Tetrachloroben	13.660	216	69865	17.125 ng/ul 16.975 ng/uL	99 99			
76) Pentachlorobenzene	15.170	250	64701	16.966 ng/uL	97			
77) Carbazole	18.002	167	259328	17.872 ng/ul	99			
78) Di-n-butylphthalate	18.536	149	332907	17.460 ng/ul	99			
80) Fluoranthene	19.647	202	342961	17.845 ng/ul	98			
82) Pyrene	20.005	202	336076	17.897 ng/ul	99			
83) Butylbenzylphthalate	20.869	149	144573	17.903 ng/ul	99			
84) 3,3'-Dichlorobenzidine	21.786	252	108246	17.927 ng/ul	99			
85) Benzo(a)anthracene	21.880	228	303650	17.691 ng/ul	99			
86) Bis(2-ethylhexyl)phtha 87) Chrysene	21.750 21.950	149 228	205882 288151	17.762 ng/ul 17.573 ng/ul	100 98			
89) Di-n-octyl phthalate	23.025	228 149	349752	20.381 ng/ul	100			
90) Benzo(b)fluoranthene	24.212	252	293676	19.557 ng/ul	99			
91) Benzo(k)fluoranthene	24.288	252	285144	20.236 ng/ul	100			
93) Benzo(a)pyrene	25.140	252	288021	20.138 ng/ul	99			
94) Indeno(1,2,3-cd)pyrene	29.206	276	316284	19.866 ng/ul	96			
95) Dibenzo(a,h)anthracene	29.277		267111	19.828 ng/ul	97			
96) Benzo(g,h,i)perylene	30.428			. 19.670 ng/ul≻	11/04/21 Ju			

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

## SFAM-EPA-BG110321.M Tue Nov 02 16:06:22 2021

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