Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\

Data File : BG051412.D

Acq On : 8 Dec 2021 18:23

Operator : CG/JU Sample : SSTD01029

Misc

ALS Vial : 4 Sample Multiplier: 1

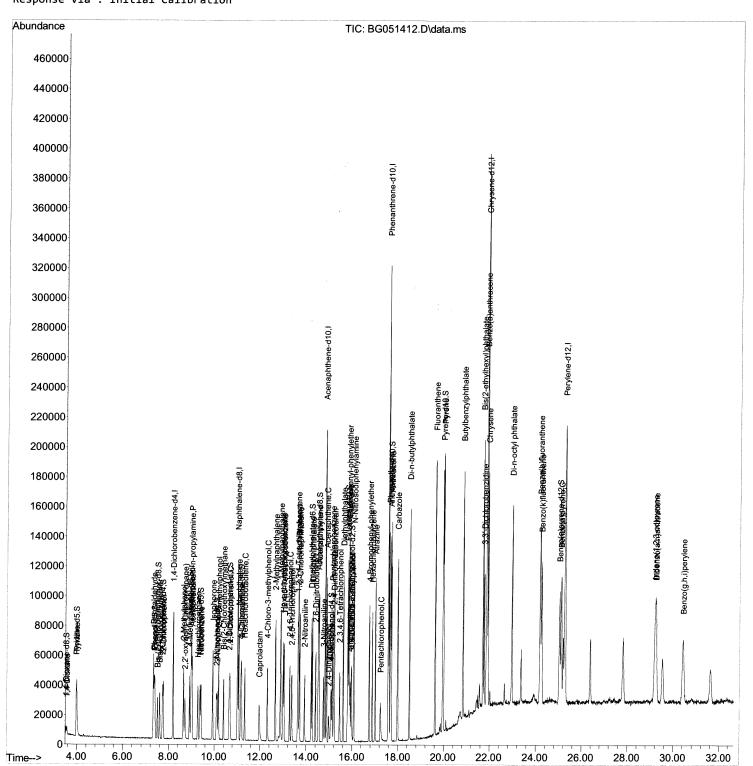
Quant Time: Dec 09 02:29:03 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 09 02:14:28 2021
Response via : Initial Calibration

Instrument : BNA_G ClientSampleId : SSTD010429

Manual IntegrationsAPPROVED



Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG120821\

Data File : BG051412.D

Acq On : 8 Dec 2021 18:23

Operator : CG/JU Sample : SSTD01029

Misc

ALS Vial : 4 Sample Multiplier: 1

Quant Time: Dec 09 02:29:03 2021

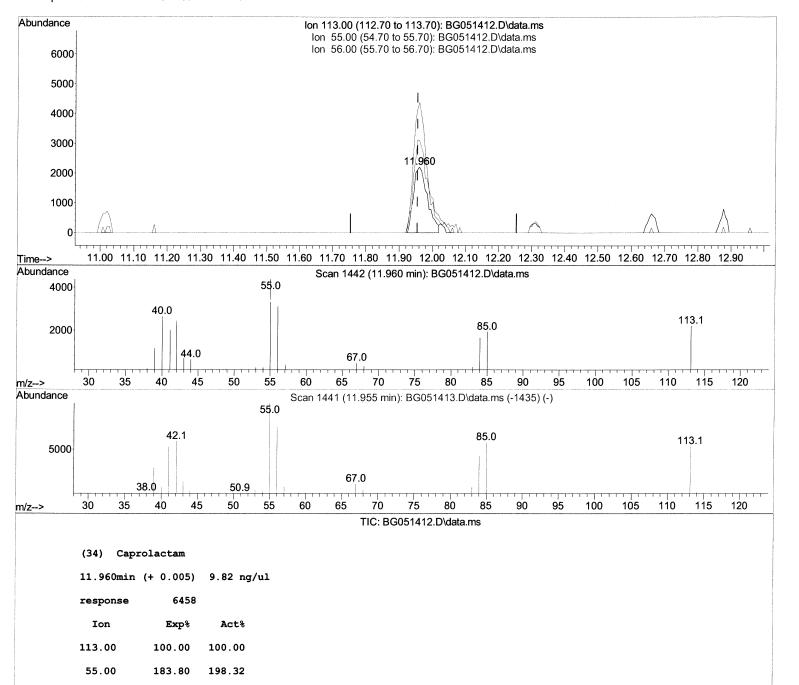
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Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 02:14:28 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/09/2021 Supervised By :Yogesh Patel 12/16/2021



136.50

0.00

140.63

0.00

56.00

0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\

Data File : BG051412.D

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Operator : CG/JU Sample : SSTD01029

Misc

ALS Vial : 4 Sample Multiplier: 1

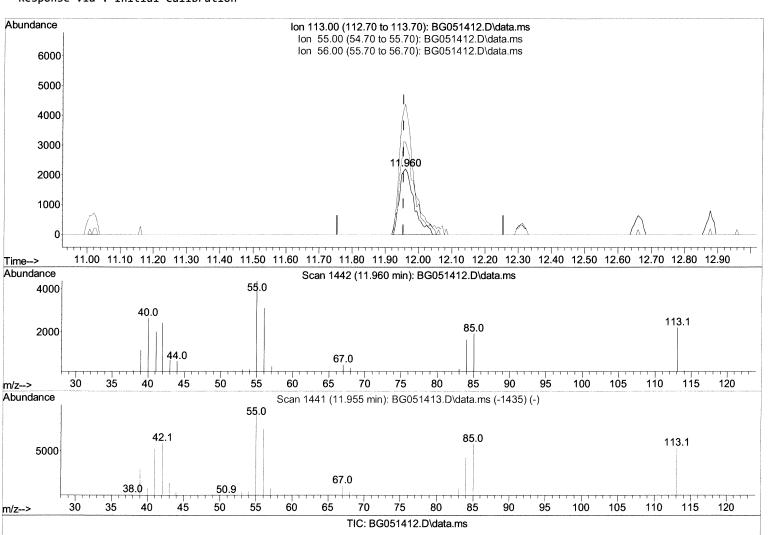
Quant Time: Dec 09 02:29:03 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 02:14:28 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleId : SSTD010429

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/09/2021 Supervised By :Yogesh Patel 12/16/2021



(34) Caprolactam

11.960min (+ 0.005) 10.21 ng/ul m 12/16/21 JU

response	6709	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	183.80	198.32
56.00	136.50	140.63
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\

Data File : BG051412.D

Acq On : 8 Dec 2021 18:23

Operator : CG/JU Sample : SSTD01029

Misc :

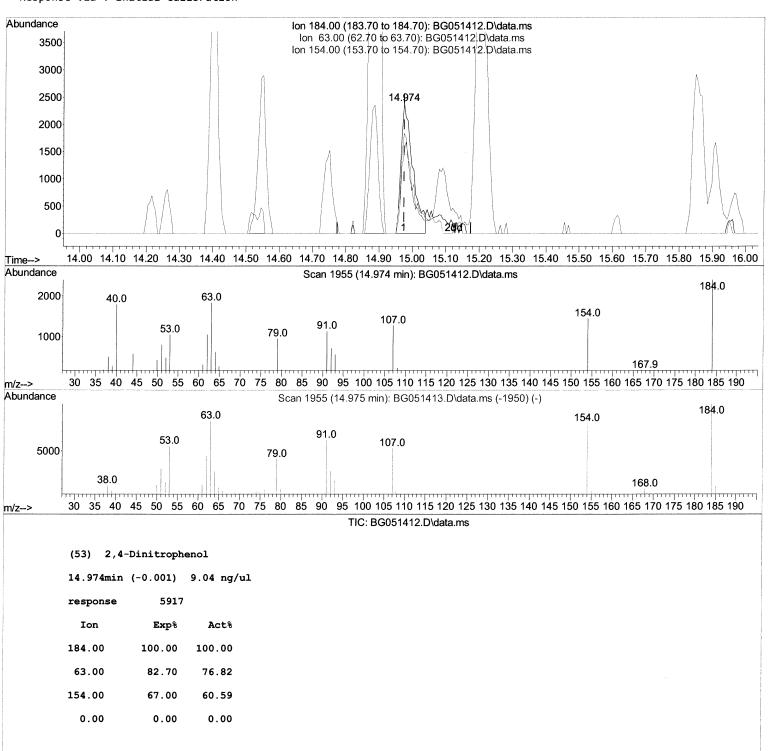
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Dec 09 02:29:03 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 02:14:28 2021 Response via : Initial Calibration Instrument:
BNA_G
ClientSampleId:
SSTD010429

Manual Integrations APPROVED



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Operator : CG/JU Sample : SSTD01029

Misc

ALS Vial : 4 Sample Multiplier: 1

Quant Time: Dec 09 02:29:03 2021

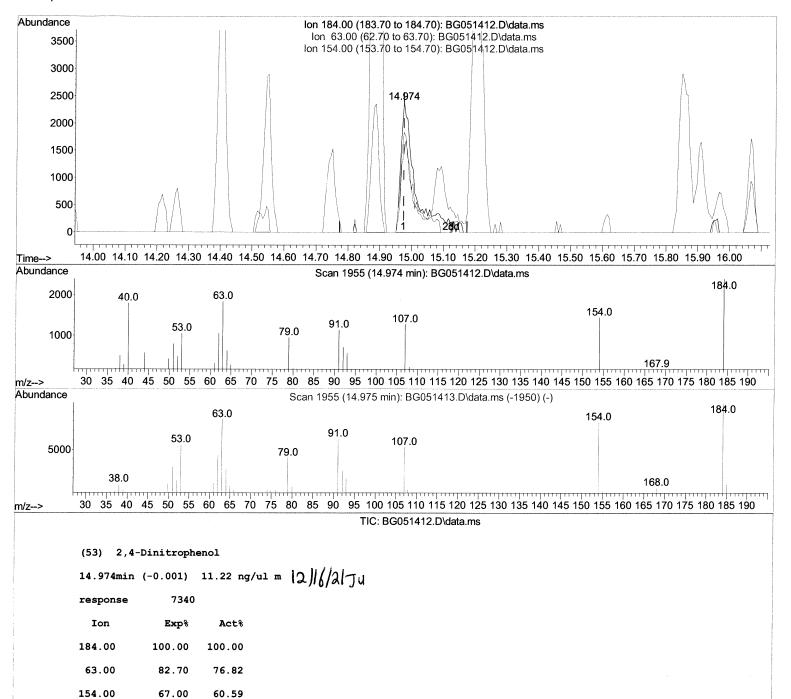
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 02:14:28 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/09/2021 Supervised By :Yogesh Patel 12/16/2021



0.00

0.00

0.00

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Data File : BG051412.D

Acq On : 8 Dec 2021 18:23

Operator : CG/JU Sample : SSTD01029

Misc :

ALS Vial : 4 Sample Multiplier: 1

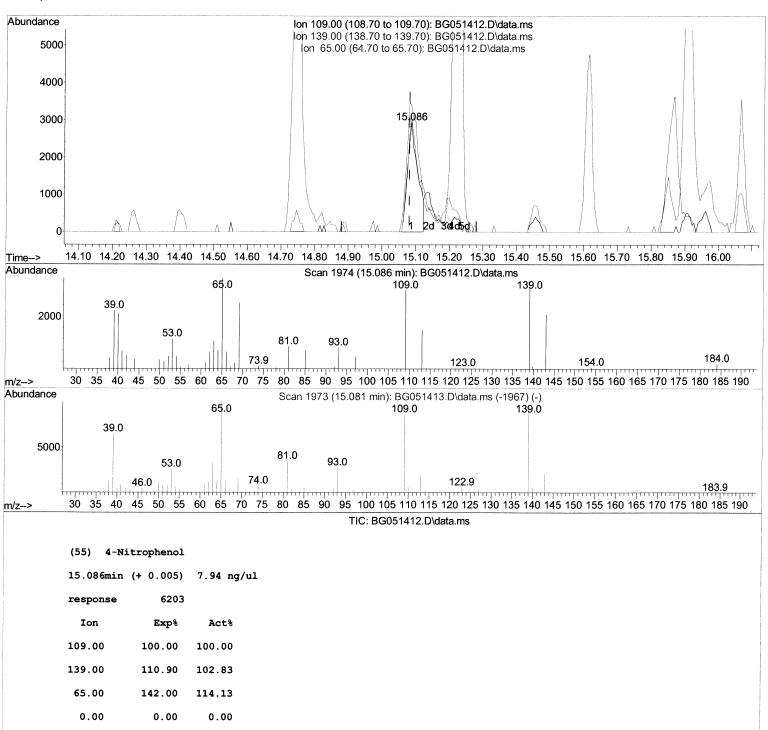
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION
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Data File : BG051412.D

Acq On : 8 Dec 2021 18:23

Operator : CG/JU Sample : SSTD01029

Misc

ALS Vial : 4 Sample Multiplier: 1

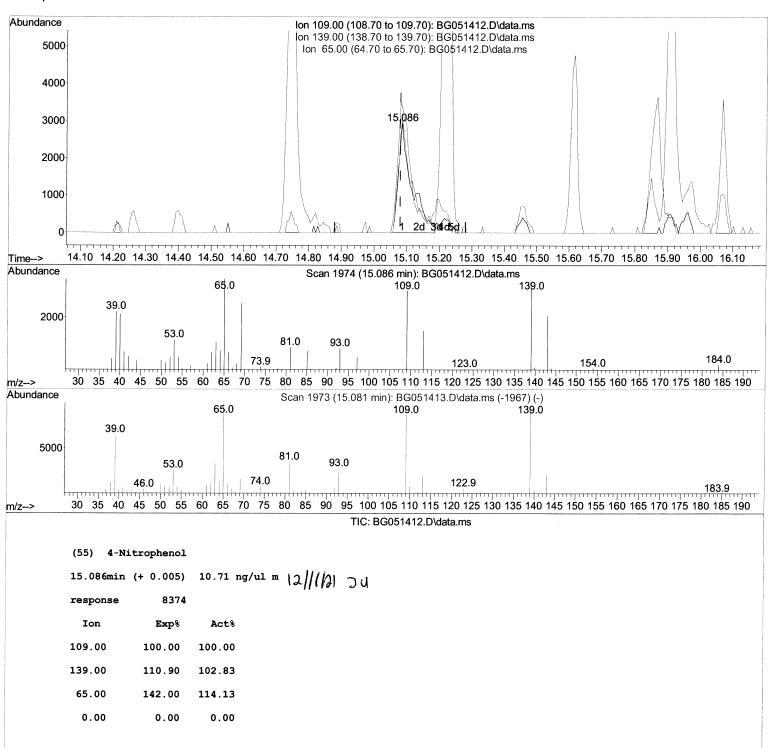
Quant Time: Dec 09 02:29:03 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION
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Misc

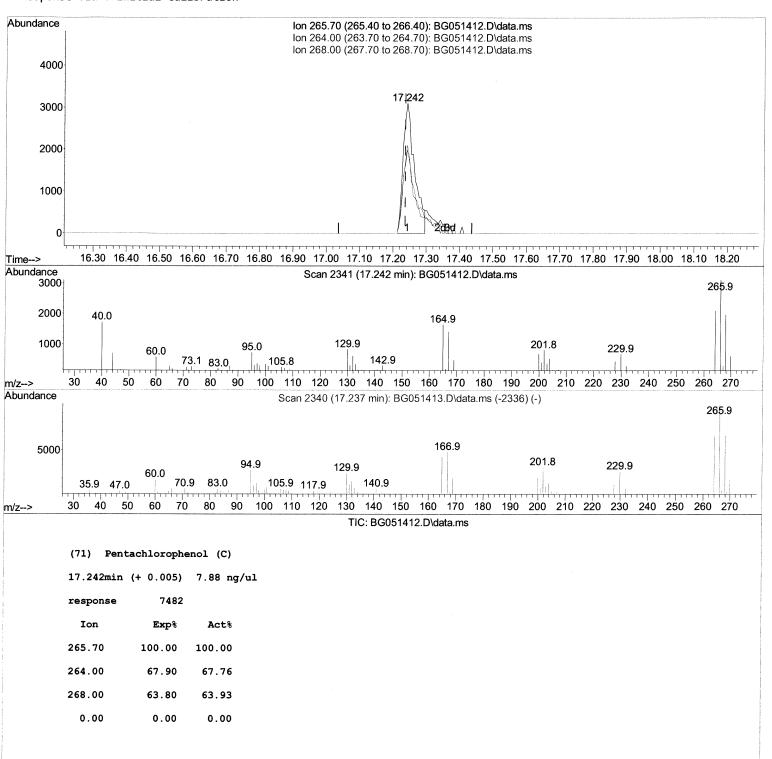
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Dec 09 02:29:03 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 02:14:28 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleId : SSTD010429

Manual IntegrationsAPPROVED



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\

Data File : BG051412.D

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Operator : CG/JU Sample : SSTD01029

Misc

ALS Vial : 4 Sample Multiplier: 1

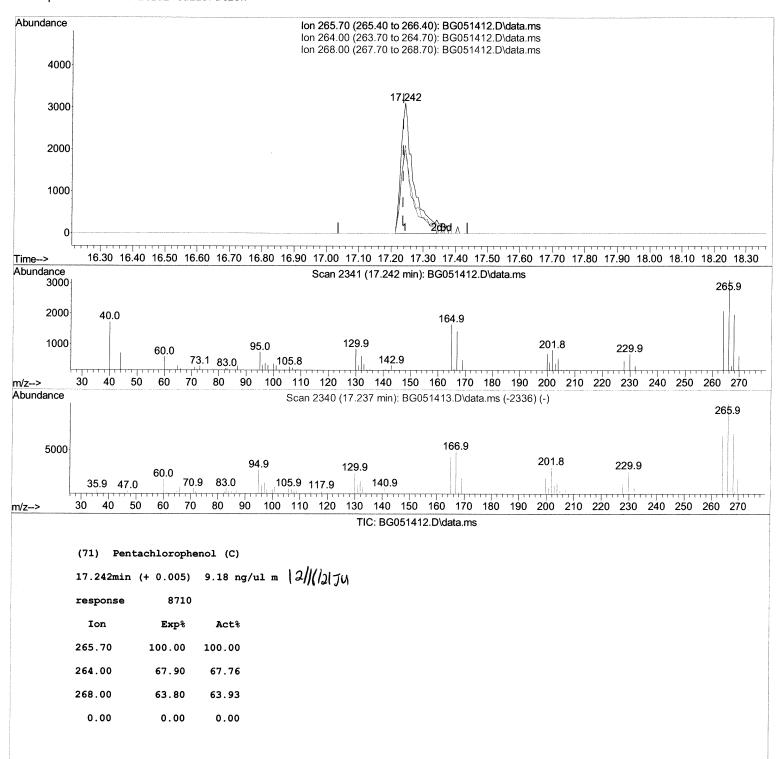
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Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 02:14:28 2021 Response via : Initial Calibration

Instrument : BNA_G ClientSampleId : SSTD010429

Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc U	nits De	ev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.188	152	23339	20.00	ng/ul	0.00
20) Naphthalene-d8	11.014		105134		ng/ul	0.00
38) Acenaphthene-d10	14.821	164	72349		ng/ul	0.00
64) Phenanthrene-d10	17.571	188	195949		ng/ul	0.00
79) Chrysene-d12	21.872		203026		ng/ul	0.00
88) Perylene-d12	25.274		195762		ng/ul	0.00
,, ,,,,,,,,,,						
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.529	96	2726	4.06	ng/uL	0.00
4) Pyridine-d5	3.970	84	18974	9.63	ng/ul	0.00
7) Phenol-d5	7.360	99	22545	9.77	ng/ul	0.00
Bis-(2-Chloroethyl)eth	7.507	67	14752	10.18	ng/ul	0.00
11) 2-Chlorophenol-d4	7.724	132	16631		ng/ul	0.00
15) 4-Methylphenol-d8	8.911	113	17839		ng/ul	0.00
21) Nitrobenzene-d5	9.369	128	8788		ng/ul	0.00
24) 2-Nitrophenol-d4	10.092	143	9959		ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.650	165	16130		ng/ul	0.00
31) 4-Chloroaniline-d4	11.161	131	23297		ng/ul	0.00
46) Dimethylphthalate-d6	14.216	166	55977		ng/ul	0.00
49) Acenaphthylene-d8	14.522	160	69896		ng/ul	0.00
54) 4-Nitrophenol-d4	15.074	143	6634		ng/ul	0.01
60) Fluorene-d10	15.809	176	49603		ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.955	200	8620		ng/ul	0.00
73) Anthracene-d10	17.671	188	81552		ng/ul	0.00
81) Pyrene-d10	19.951	212	96627		ng/ul	0.00
92) Benzo(a)pyrene-d12	25.039	264	77720		ng/ul	0.00
					_	-
Target Compounds	2 570	-00	2426		-	value
2) 1,4-Dioxane	3.570	88	3136		ng/uL#	
5) Pyridine	3.987	79	20284		ng/ul	97
6) Benzaldehyde	7.324	77	24329	16.56		96
8) Phenol	7.389	94	23131		ng/ul	98
10) Bis(2-Chloroethyl)ether	7.601	93	18432	10.20	_	97
12) 2-Chlorophenol	7.753	128	17623	10.41	_	99
13) 2-Methylphenol	8.640	108	17026		ng/ul	97
14) 2,2'-oxybis(1-Chloropr	8.705	45	28222	10.82		98
16) Acetophenone	9.022	105	29063	10.09		96
17) N-Nitroso-di-n-propyla	8.993	70	17601	10.64		98
18) 4-Methylphenol	8.975	108	18782		ng/ul	100
19) Hexachloroethane	9.269	117	7334	10.26		95
22) Nitrobenzene	9.410	77	24363		ng/ul#	
23) Isophorone	9.927	82	47594	10.53		96
25) 2-Nitrophenol	10.127	139	10050		ng/ul	91
26) 2,4-Dimethylphenol	10.180	107	21415	10.10		96
27) Bis(2-Chloroethoxy)met	10.403	93	25147	10.08	-	96
29) 2,4-Dichlorophenol	10.679	162	16266		ng/ul	95
30) Naphthalene	11.067	128	58483	10.22	-	100
32) 4-Chloroaniline	11.185	127	24152		ng/ul	100
33) Hexachlorobutadiene	11.326	225	11183		ng/ul	5 12/11/12/ Ju
34) Caprolactam	11.960	113	6709m >			
35) 4-Chloro-3-methylphenol	12.313	107	20407	10.16	ng/u1	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\

Data File : BG051412.D

Acq On : 8 Dec 2021 18:23

Operator : CG/JU Sample : SSTD01029

Misc

ALS Vial : 4 Sample Multiplier: 1

Quant Time: Dec 09 02:29:03 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 02:14:28 2021 Response via : Initial Calibration Instrument: BNA_G ClientSampleId: SSTD010429

Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc Units Dev(Min)
36) 2-Methylnaphthalene	12.659	142	38587	9.92 ng/ul	97
37) 1-Methylnaphthalene	12.877	142	39908	9.97 ng/ul	98
39) 1,2,4,5-Tetrachloroben	13.024	216	22034	9.70 ng/ul	98
40) Hexachlorocyclopentadiene	12.988	237	18215	19.84 ng/ul	99
41) 2,4,6-Trichlorophenol	13.270	196	13708	9.62 ng/ul	96
42) 2,4,5-Trichlorophenol	13.364	196	15499	10.38 ng/ul	94
43) 1,1'-Biphenyl	13.652	154	53408	9.88 ng/ul	99
44) 2-Chloronaphthalene	13.705	162	42356	9.85 ng/ul	98
45) 2-Nitroaniline	13.917	65	15471	10.40 ng/ul	92
47) Dimethylphthalate	14.263	163	56653	10.05 ng/ul	99
48) 2,6-Dinitrotoluene	14.404	165	11659	9.85 ng/ul	98
50) Acenaphthylene	14.545	152	71382	10.29 ng/ul	98
51) 3-Nitroaniline	14.745	138	11822	10.10 ng/ul	100
52) Acenaphthene	14.886	153	45996	10.06 ng/ul	99
53) 2,4-Dinitrophenol	14.974	184	7340m 🥄	11.22 ng/ul 🔪	12/16/21 JU
55) 4-Nitrophenol	15.086	109	8374m 🖊	> _{10.71 ng/ul} >	
56) Dibenzofuran	15.221	168	65783	9.97 ng/ul	99
57) 2,4-Dinitrotoluene	15.198	165	17392	10.29 ng/ul#	97
58) 2,3,4,6-Tetrachlorophenol	15.456	232	11218	9.57 ng/ul#	92
59) Diethylphthalate	15.615	149	61219	10.35 ng/ul	98
61) Fluorene	15.867	166	52748	9.98 ng/ul	97
62) 4-Chlorophenyl-phenyle	15.850	204	28070	9.86 ng/ul	96
63) 4-Nitroaniline	15.908	138	11654	10.24 ng/ul	93
66) 4,6-Dinitro-2-methylph	15.973	198	8442	7.24 ng/ul#	90
67) N-Nitrosodiphenylamine	16.067	169	46700	8.32 ng/ul	98
68) 4-Bromophenyl-phenylether	16.749	248	16834	8.02 ng/ul	94
69) Hexachlorobenzene	16.872	284	16970	7.92 ng/ul	97
70) Atrazine	17.013	200	20944	8.88 ng/ul	982/16/21JU
71) Pentachlorophenol	17.242	266	8710m >		•
72) Phenanthrene	17.618	178	91738	8.48 ng/ul	100
74) Anthracene	17.706	178	94687	8.81 ng/ul	98
75) 1,2,3,4-Tetrachloroben	13.629	216	22854	8.00 ng/uL	97
76) Pentachlorobenzene	15.139	250	20782	7.80 ng/uL	96
77) Carbazole	17.982	167	85387	9.05 ng/ul	100
78) Di-n-butylphthalate	18.500	149	113770	9.36 ng/ul	99
80) Fluoranthene	19.622	202	117438	7.78 ng/ul	97
82) Pyrene	19.980	202	117554	7.96 ng/ul	99
83) Butylbenzylphthalate	20.838	149	49895	8.13 ng/ul	99
84) 3,3'-Dichlorobenzidine	21.760	252	32893	6.96 ng/ul	99
85) Benzo(a)anthracene	21.854	228	105798 70519	7.68 ng/ul	99 98
86) Bis(2-ethylhexyl)phtha87) Chrysene	21.708	149		7.99 ng/ul 7.55 ng/ul	98 97
89) Di-n-octyl phthalate	21.925 22.971	228 149	99818 115389	8.14 ng/ul	100
90) Benzo(b)fluoranthene	24.187	252	102547	7.76 ng/ul	97
91) Benzo(k)fluoranthene	24.167	252	94230	7.76 ng/ul 7.60 ng/ul	99
93) Benzo(a)pyrene	25.109	252	95645	7.59 ng/ul	99
94) Indeno(1,2,3-cd)pyrene	29.199	276	106507	7.55 ng/ul 7.55 ng/ul	96
95) Dibenzo(a,h)anthracene	29.246	278	89070	7.44 ng/ul	96
96) Benzo(g,h,i)perylene	30.438	276	88093	7.42 ng/ul	98

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