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

Data File : BG051413.D

Acq On : 8 Dec 2021 19:04

Operator : CG/JU Sample : SSTD02030

Misc

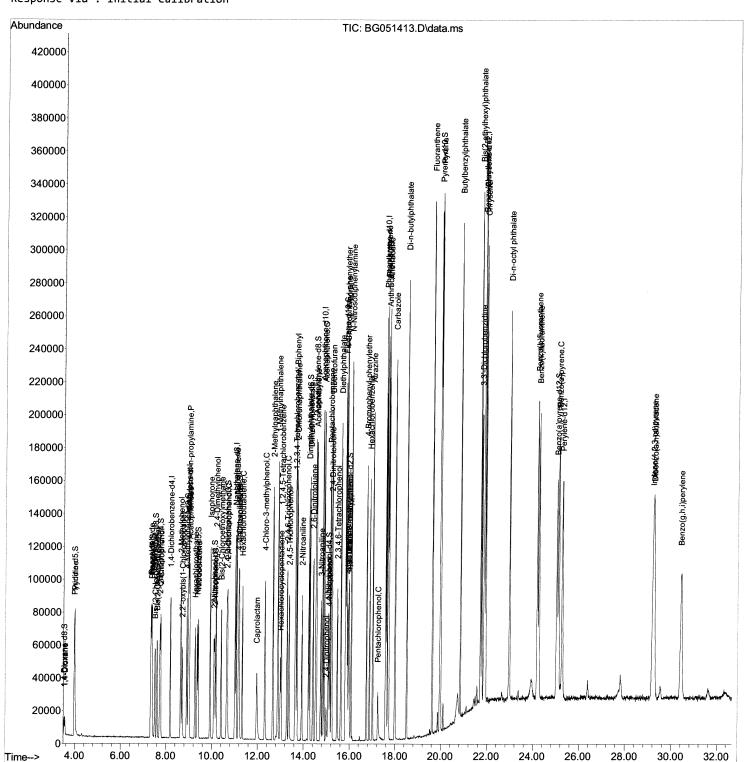
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 09 02:16:27 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 : SSTD020430

Manual IntegrationsAPPROVED



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

Data File : BG051413.D

Acq On : 8 Dec 2021 19:04

Operator : CG/JU Sample : SSTD02030

Misc

ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 09 02:16:27 2021

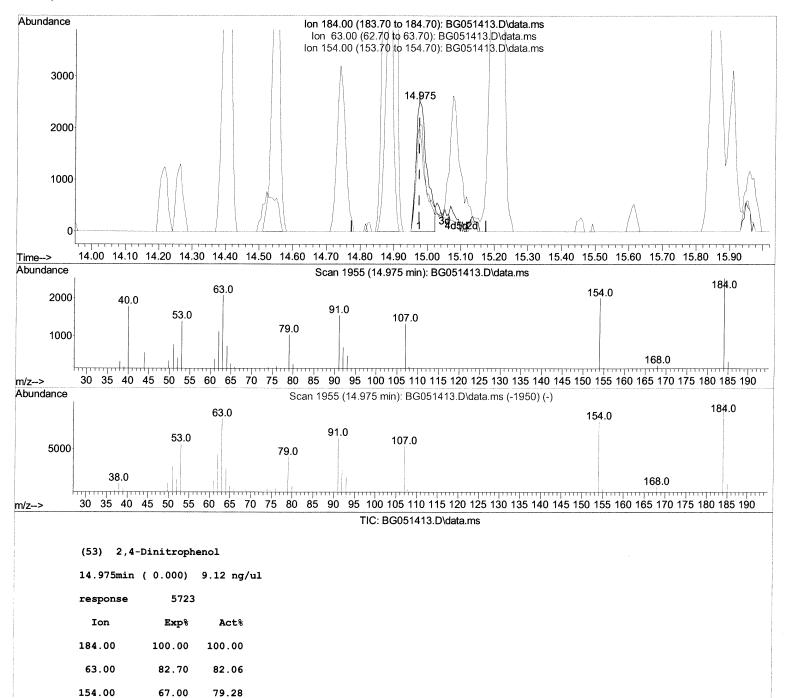
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M

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

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

Data File : BG051413.D

Acq On : 8 Dec 2021 19:04

Operator : CG/JU Sample : SSTD02030

Misc :

ALS Vial : 5 Sample Multiplier: 1

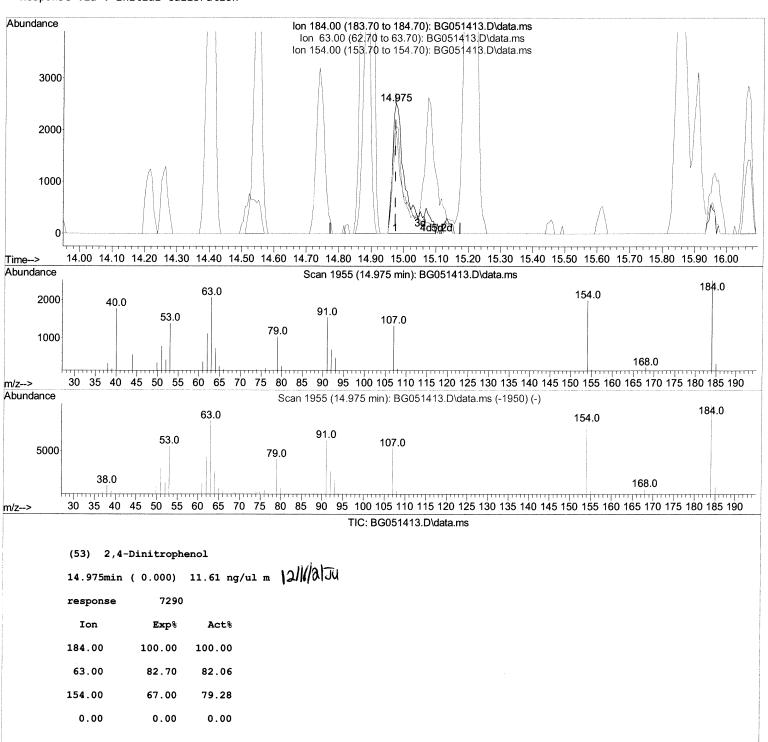
Quant Time: Dec 09 02:16:27 2021

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



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Misc

ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 09 02:16:27 2021

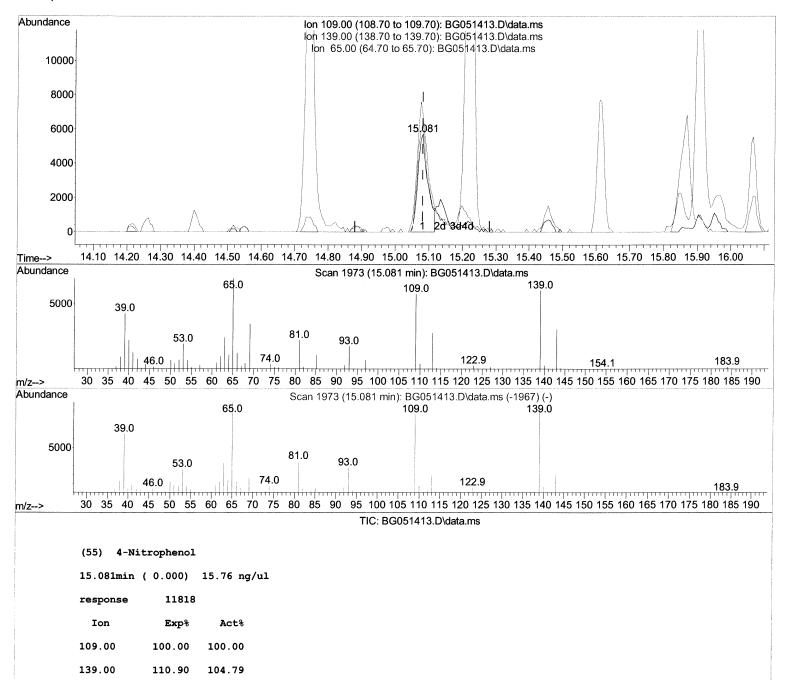
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Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 09 02:14:28 2021
Response via : Initial Calibration

Instrument:
BNA_G
ClientSampleId:
SSTD020430

Manual IntegrationsAPPROVED

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



142.00

0.00

120.05

0.00

65.00

0.00

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Operator : CG/JU Sample : SSTD02030

Misc

ALS Vial : 5 Sample Multiplier: 1

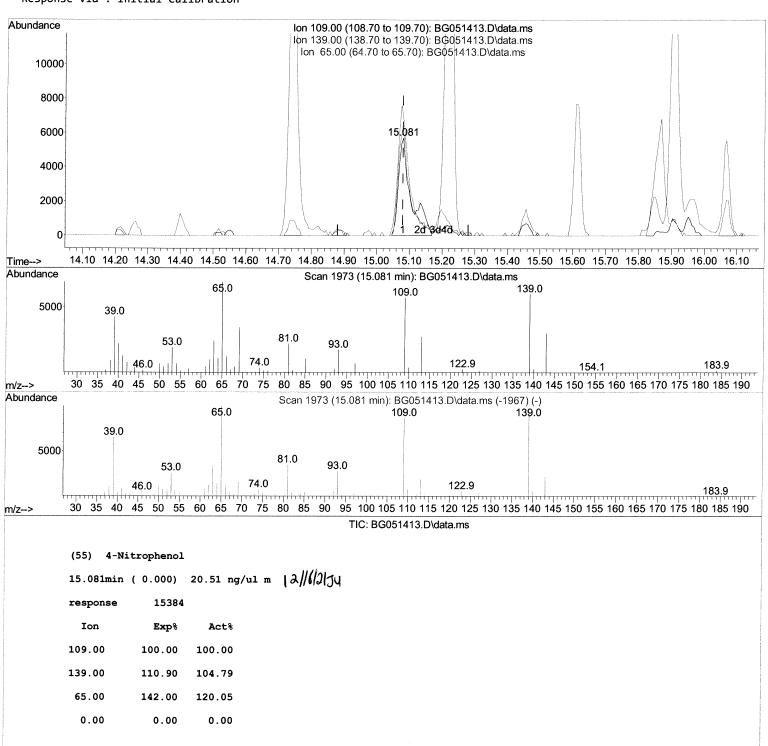
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Manual Integrations APPROVED



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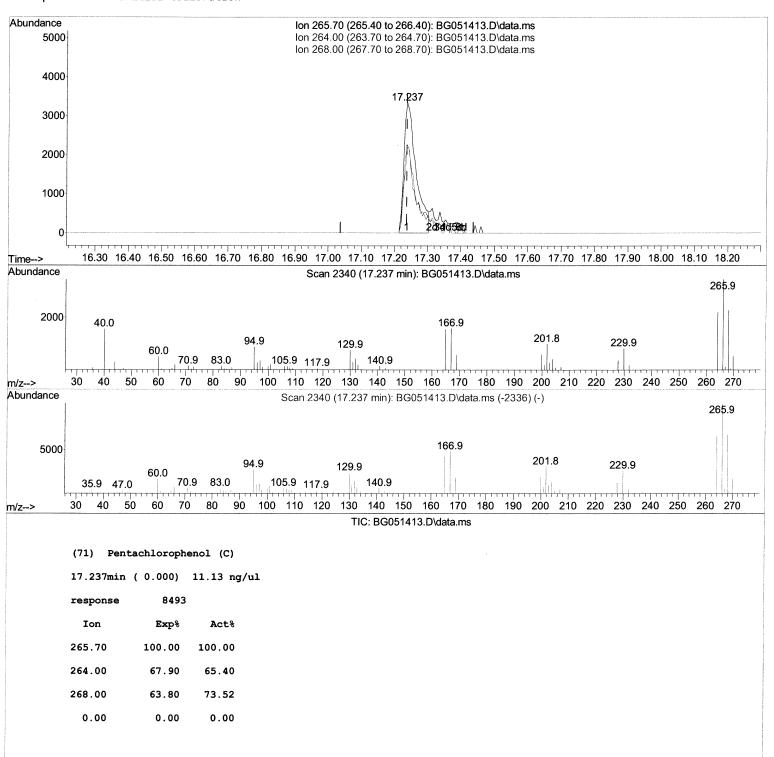
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ALS Vial : 5 Sample Multiplier: 1

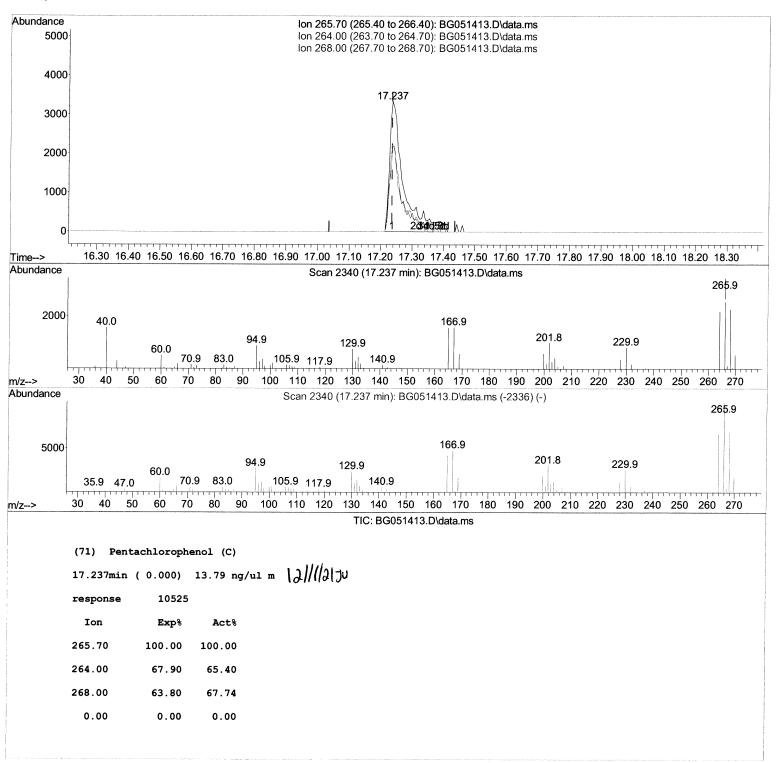
Quant Time: Dec 09 02:16:27 2021

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



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Data File : BG051413.D

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Operator : CG/JU Sample : SSTD02030

Misc

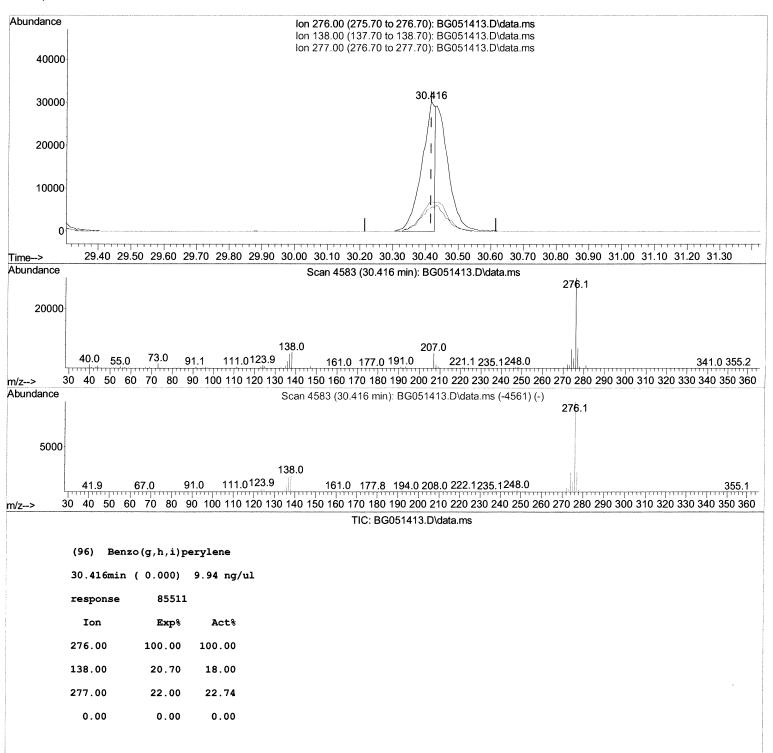
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 09 02:16:27 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:
SSTD020430

Manual Integrations APPROVED



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Operator : CG/JU Sample : SSTD02030

Misc

ALS Vial : 5 Sample Multiplier: 1

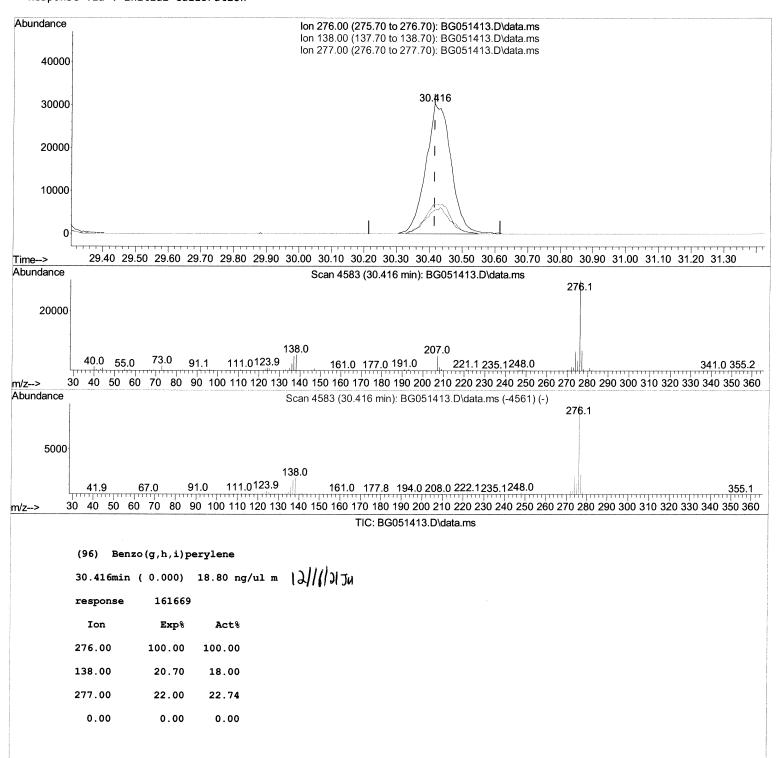
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Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 02:14:28 2021 Response via : Initial Calibration Instrument: BNA_G ClientSampleId: SSTD020430

Manual IntegrationsAPPROVED

Response via . Initial cultivat.	1011				
Compound	R.T.	QIon	Response	Conc Units De	v(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.189	152	23514	20.00 ng/ul	0.00
20) Naphthalene-d8	11.015		104901	20.00 ng/ul	0.00
38) Acenaphthene-d10	14.822	164	69412	20.00 ng/ul	0.00
64) Phenanthrene-d10	17.572	188	157622	20.00 ng/ul	0.00
79) Chrysene-d12	21.873	240	143719	20.00 ng/ul	0.00
88) Perylene-d12	25.274	264	141854	20.00 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.529	96	5559	8.22 ng/uL	0.00
4) Pyridine-d5	3.964	84	38396	19.34 ng/ul	0.00
7) Phenol-d5	7.354	99	43825	18.86 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth	7.501	67	28625	19.61 ng/ul	0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.724	132	31434	18.78 ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.911	113	34495	18.39 ng/ul	0.00
21) Nitrobenzene-d5	9.370	128	16789	18.96 ng/ul	0.00
24) 2-Nitrophenol-d4	10.098	143	19584	19.61 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.650	165	32220	19.01 ng/ul	0.00
31) 4-Chloroaniline-d4	11.162	131	45833	18.48 ng/ul	0.00
46) Dimethylphthalate-d6	14.217	166	101551	19.01 ng/ul	0.00
49) Acenaphthylene-d8	14.516	160	131582	19.54 ng/ul	0.00
54) 4-Nitrophenol-d4	15.063	143	13995	16.19 ng/ul	0.00
60) Fluorene-d10	15.809	176	91579	19.04 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.950	200	16809	17.28 ng/ul	0.00
73) Anthracene-d10	17.672	188	143933	19.09 ng/ul	0.00
81) Pyrene-d10	19.951	212	171554	19.73 ng/ul	0.00
92) Benzo(a)pyrene-d12	25.034	264	142153	18.76 ng/ul	0.00
Target Compounds				Qv	/alue
2) 1,4-Dioxane	3.565	88	6194	8.12 ng/uL	92
5) Pyridine	3.988	79	40196	19.46 ng/ul	94
Benzaldehyde	7.325	77	32702	22. 1 0 ng/ul	88
8) Phenol	7.384	94	45625	18.95 ng/ul	97
<pre>10) Bis(2-Chloroethyl)ether</pre>	7.595	93	34815	19.11 ng/ul	96
<pre>12) 2-Chlorophenol</pre>	7.754	128	32662	19.15 ng/ul	97
<pre>13) 2-Methylphenol</pre>	8.641	108	33069	18.44 ng/ul	99
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.712	45	53790	20.47 ng/ul	97
16) Acetophenone	9.023	105	55995	19.30 ng/ul	98
17) N-Nitroso-di-n-propyla	8.994	70	33520	20.11 ng/ul	95
18) 4-Methylphenol	8.976	108	34695	18.09 ng/ul	98
19) Hexachloroethane	9.270	117	14547	20.20 ng/ul	97
22) Nitrobenzene	9.411	77	47007	20.24 ng/ul	98
23) Isophorone	9.928	82	90560	20.07 ng/ul	99
<pre>25) 2-Nitrophenol</pre>	10.128	139	19266	18.62 ng/ul	95
26) 2,4-Dimethylphenol	10.180	107	41302	19.52 ng/ul	97
27) Bis(2-Chloroethoxy)met	10.404	93	49363	19.82 ng/ul	98
29) 2,4-Dichlorophenol	10.674	162	30911	18.5 3 ng/ul	97
30) Naphthalene	11.068	128	109454	19.18 ng/ul	97
32) 4-Chloroaniline	11.179	127	45882	18.43 ng/ul	99
33) Hexachlorobutadiene	11.326	225	20727	18.01 ng/ul	99
34) Caprolactam	11.955	113	12441	18.97 ng/ul	95
35) 4-Chloro-3-methylphenol	12.307	107	37915	18.92 ng/ul	94

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Data File : BG051413.D

Acq On : 8 Dec 2021 19:04

Operator : CG/JU Sample : SSTD02030

Misc

ALS Vial : 5 Sample Multiplier: 1

Quant Time: Dec 09 02:16:27 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 : SSTD020430

Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc U	nits Dev(Min)
36) 2-Methylnaphthalene	12.660	142	72380	18 64	ng/ul	97
37) 1-Methylnaphthalene	12.877	142	75147		ng/ul#	100
39) 1,2,4,5-Tetrachloroben	13.024	216	41050	18.84		99
40) Hexachlorocyclopentadiene	12.989	237	14262		ng/ul#	96
41) 2,4,6-Trichlorophenol	13.271	196	26976	19.73	•	98
42) 2,4,5-Trichlorophenol	13.359	196	27180	18.98	_	96
43) 1,1'-Biphenyl	13.653	154	100173	19.32	-	98
44) 2-Chloronaphthalene	13.706	162	78424	19.02		99
45) 2-Nitroaniline	13.917	65	29301		ng/ul	96
47) Dimethylphthalate	14.264	163	102938	19.04		98
48) 2,6-Dinitrotoluene	14.405	165	21289	18.75	ng/ul	91
50) Acenaphthylene	14.552	152	128498	19.31	ng/ul	97
51) 3-Nitroaniline	14.746	138	22237	19.81	ng/ul	99
52) Acenaphthene	14.887	153	84320	19.22	ng/ul	97
53) 2,4-Dinitrophenol	14.975	184	7290m 🥆	11.61	ng/ul 🔪	1011/101-1
55) 4-Nitrophenol	15.081	109	15384m/	20.51	ng/ul >	12/16/2/54
56) Dibenzofuran	15.222	168	119511	18.88	ng/ul	100
57) 2,4-Dinitrotoluene	15.198	165	30837	19.01	ng/ul	99
58) 2,3,4,6-Tetrachlorophenol	15.457	232	21410	19.04	ng/ul	98
59) Diethylphthalate	15.615	149	109686	19.33	ng/ul	99
61) Fluorene	15.868	166	96027	18.94	ng/ul	98
62) 4-Chlorophenyl-phenyle	15.850	204	49794	18.22	ng/ul	96
63) 4-Nitroaniline	15.909	138	19683	18.02	ng/ul	93
66) 4,6-Dinitro-2-methylph	15.968	198	16727	17.83	ng/ul	96
67) N-Nitrosodiphenylamine	16.068	169	85183	18.88	ng/ul	97
68) 4-Bromophenyl-phenylether	16.743	248	30682	18.16	ng/ul	97
69) Hexachlorobenzene	16.873	284	30921	17.95	-	96
70) Atrazine	17.008	200	37185	19.61	-	99
71) Pentachlorophenol	17.237	266	10525m >		ng/ul >	ialllal Ju
72) Phenanthrene	17.613	178	165378	19.00	-	100
74) Anthracene	17.707	178	166653	19.28	•	98
75) 1,2,3,4-Tetrachloroben	13.629	216	42827	18.63	-	99
76) Pentachlorobenzene	15.139	250	38874	18.15	•	98
77) Carbazole	17.983	167	151131	19.92	-	99
78) Di-n-butylphthalate	18.500	149	197286	20.17	_	99
80) Fluoranthene	19.616	202	207617	19.44	-	99
82) Pyrene	19.981	202	206243	19.74	_	98
83) Butylbenzylphthalate	20.838	149	88853	20.46	-	96
<pre>84) 3,3'-Dichlorobenzidine 85) Benzo(a)anthracene</pre>	21.761	252	64474	19.27	-	99
	21.855	228	186483	19.13		99
86) Bis(2-ethylhexyl)phtha87) Chrysene	21.708		122963	19.67		98
89) Di-n-octyl phthalate	21.925 22.965	228 149	177877 209728	18.99 20.41		99 1 00
90) Benzo(b)fluoranthene	24.182	252	182931	19.11	-	99
91) Benzo(k)fluoranthene	24.258	252	169078	18.82		99
93) Benzo(a)pyrene	25.116	252	174105	19.06		98
94) Indeno(1,2,3-cd)pyrene	29.199	276	190097	18.60	-	99
95) Dibenzo(a,h)anthracene	29.246	278	162630	18.76		99
96) Benzo(g,h,i)perylene	30.416	276	161669m >		ng/ul >	12/16/21 JU

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