









Data Path : Z:\svoasrv\HPCHEM1 Data File : BG051350.D Acq On : 6 Dec 2021 14:48 Dperator : CG/JU Sample : PB141174BS Misc : ALS Vial : 7 Sample Multipl: Quant Time: Dec 06 15:22:06 202 Quant Method : Z:\svoasrv\HPCHE	ier: 1 21 EM1\BNA_G	Instrument : BNA_G ClientSampleId : SLCS174 Manual IntegrationsAPPROVED Reviewed By :Jagrut Upadhyay 12/06/2021			
Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23 Response via : Initial Calibrat	3:09 2021	Supervised By :mohammad ahmed 12/07/2021			
Compound	R.T.	QIon	Response	Conc Units Dev	/(Min)
Internal Standards					
 1,4-Dichlorobenzene-d4 	8.194		19937	20.000 ng/ul	0.00
20) Naphthalene-d8	11.020		85899	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.827		57337	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.577		135407	20.000 ng/ul	0.00
79) Chrysene-d12 88) Perylene-d12	21.878	240	128070	20.000 ng/ul	0.00
ob) Peryrene-urz	25.280	264	132097	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.529	96	3536	6.163 ng/uL	-0.02
4) Pyridine-d5	3.957	84	51099	30.353 ng/ul	-0.02
7) Phenol-d5	7.359	99	71747	36.412 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth		67	41685	33.684 ng/ul	0.00
11) 2-Chlorophenol-d4	7.724	132	51890	36.570 ng/ul	0.00
15) 4-Methylphenol-d8	8.910	113	55629	34.985 ng/ul	0.00
21) Nitrobenzene-d5	9.369	128	25564	35.255 ng/ul	0.00
24) 2-Nitrophenol-d4	10.097	143	29650	36.249 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.644	165	52413	37.767 ng/ul	0.00
31) 4-Chloroaniline-d4	11.161	131	101022	49.749 ng/ul	0.00
46) Dimethylphthalate-d6	14.216	166	155033	35.141 ng/ul	0.00
49) Acenaphthylene-d8	14.522	160	199471	35.856 ng/ul	0.00
54) 4-Nitrophenol-d4 60) Fluorene-d10	15.056 15.814	143 176	29471 144366	41.269 ng/ul	0.00
65) 4,6-Dinitro-2-methylph			16907	36.339 ng/ul 20.235 ng/ul	0.00 0.00
73) Anthracene-d10	17.677	188	238567	36.838 ng/ul	0.00
81) Pyrene-d10	19.956	212	267734	34.550 ng/ul	0.00
92) Benzo(a)pyrene-d12	25.050	264	249396	35.351 ng/ul	0.00
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Target Compounds	2 564	00	7604		alue
2) 1,4-Dioxane 5) Pyridine	3.564 3.981	88 79	7621	11.778 ng/uL#	
6) Benzaldehyde	7.324	79	53987 48127	30.818 ng/ul 38.353 ng/ul	99 98
8) Phenol	7.383	94	71407	34.982 ng/ul	100
10) Bis(2-Chloroethyl)ether	7.600	93	50044	32.406 ng/ul	94
12) 2-Chlorophenol	7.759	128	50817	35.145 ng/ul	96
<pre>13) 2-Methylphenol</pre>	8.640	108	52021	34.214 ng/ul	99
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.717	45	70171	31.488 ng/ul	95
16) Acetophenone	9.022	105	78782	32.032 ng/ul	96
17) N-Nitroso-di-n-propyla	8.993	70	44185	31.263 ng/ul	98
18) 4-Methylphenol	8.975	108	55385	34.065 ng/ul	96
19) Hexachloroethane	9.275	117	20621	33.764 ng/ul	96
22) Nitrobenzene 23) Isophorone	9.416	77 92	66510 110009	34.981 ng/ul	97
25) 2-Nitrophenol	9.933 10.133	82 139	119908 29431	32.461 ng/ul 34.738 ng/ul	99 90
26) 2,4-Dimethylphenol	10.133	107	62585	36.131 ng/ul	99
27) Bis(2-Chloroethoxy)met	10.409	93	66208	32.467 ng/ul	99
29) 2,4-Dichlorophenol	10.673	162	48832	35.745 ng/ul	97
30) Naphthalene	11.073	128	159576	3/11/2 ng/ul	07
32) 4-Chloroaniline	11.184	127	65061	31.914 ng/ul	98 1 222
33) Hexachlorobutadiene	11.331	225	31894	33.847 ng/ul	_95 R121011 -1
34) Caprolactam	11.948	113		>36.860 ng/ul	398 595 JU12/07/21
35) 4-Chloro-3-methylphenol	12.301	107	59274	36.119 ng/ul	96

AM-EPA-BG112321.M Mon Dec 06 15:24:04 2021

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Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051350.D
Acq On : 6 Dec 2021 14:48
Dperator : CG/JU
Sample : PB141174BS
Misc :
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Dec 06 15:22:06 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 03 15:23:09 2021
Response via : Initial Calibration

Instrument: BNA_G ClientSampleId: SLCS174

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/06/2021 Supervised By :mohammad ahmed 12/07/2021

	Compound	R.T.	QIon	Response	Conc Un	its Dev	(Min)	
36`) 2-Methylnaphthalene	12.665	142	105537	33.197	ng/ul	97	
) 1-Methylnaphthalene	12.882	142	110367	33.744		99	
) 1,2,4,5-Tetrachloroben			62879	34.932		97	
) Hexachlorocyclopentadiene		237	5279		ng/ul	92	
	2,4,6-Trichlorophenol	13.270	196	42389	37.526		99	
	2,4,5-Trichlorophenol	13.358	196	47270	39.961		97	
	1,1'-Biphenyl	13.658	154	143047	33.403	-	98	
	2-Chloronaphthalene	13.711	162	115846	34.006	-	100	
	2-Nitroaniline	13.916	65	45096	38.249		95	
	Dimethylphthalate	14.269		148286	33.206	-	100	
	2,6-Dinitrotoluene	14.404	165	33576	35.795		93	
	Acenaphthylene	14.551	152	191457	34.834		97	
	3-Nitroaniline	14.739						
		14.739		36451	39.313		91 07	
	Acenaphthene 2,4-Dinitrophenol		153	122852	33.892	-	97	
	4-Nitrophenol	14.968	184	4606		ng/ul	91 02	
	Dibenzofuran	15.068 15.227	109	25589		-	92	
			168	178975	34.232	-	99	
	2,4-Dinitrotoluene	15.197	165	49315	36.809	-	97	
	2,3,4,6-Tetrachlorophenol		232	36523	39.319		99	
	Diethylphthalate	15.620	149	163354	34.850		99	
	Fluorene	15.873	166	147782	35.288		99	
	4-Chlorophenyl-phenyle	15.855	204	76953	34.096		96	
	4-Nitroaniline	15.908	138	40484	44.868		95	
	4,6-Dinitro-2-methylph	15.973	198	14903	18.494	-	96	
	N-Nitrosodiphenylamine	16.073	169	138548	35.741		96	
	4-Bromophenyl-phenylether		248	49361	34.013	-	95	
	Hexachlorobenzene	16.878	284	52816	35.691		96	
	Atrazine	17.013	200	58737	36.054		97	
	Pentachlorophenol	17.236	266	19154	29.211		96	
•	Phenanthrene	17.618	178	261213	34.938		99	
	Anthracene	17.712	178	264161	35.577		99	
	1,2,3,4-Tetrachloroben	13.634	216	66073	33.454		99	
	Pentachlorobenzene	15.144	250	59008	32.065	ng/uL	98	
	Carbazole	17.982	167	249652	38.305		99	
	Di-n-butylphthalate	18.505	149	304554	36.240		99	
	Fluoranthene	19.621	202	316655	33.270		96	
	Pyrene	19.986	202	316204	33.963		94	
	Butylbenzylphthalate	20.844	149	126924	32.792		94	
	3,3'-Dichlorobenzidine	21.760	252	101260	33.959		98	
	Benzo(a)anthracene	21.860	228	300429	34.586		99	
•	Bis(2-ethylhexyl)phtha	21.713	149	188565	33.855		99	
	Chrysene	21.931	228	289809	34.729	ng/ul	100	
	Di-n-octyl phthalate	22.976	149	317063	33.131	•	100	
	Benzo(b)fluoranthene	24.193	252	305054	34.219		99	
	Benzo(k)fluoranthene	24.257	252	279340	33.391	ng/ul	99	
	Benzo(a)pyrene	25.121	252	295437	34.737	ng/ul	98	
94)	Indeno(1,2,3-cd)pyrene	29.198	276	337192	35.430	ng/ul	96	
95)	Dibenzo(a,h)anthracene	29.251	278	287943	35.663	ng/ul	> > 7412/	
	<pre>Benzo(g,h,i)perylene</pre>	30.444	276	277237m	34.623	ng/117 >		۶,

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