

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122624\
 Data File : BG063891.D
 Acq On : 28 Dec 2024 5:57
 Operator : RC/JU
 Sample : SSTDCC020
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTD020587

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/30/2024
 Supervised By :mohammad ahmed 12/30/2024

Quant Time: Dec 29 23:38:35 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 24 00:44:39 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	Qvalue
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.794	152	108494	20.000	ng/u1	0.00	
20) Naphthalene-d8	10.585	136	438192	20.000	ng/u1	0.00	
38) Acenaphthene-d10	14.439	164	301488	20.000	ng/u1	0.00	
64) Phenanthrene-d10	17.194	188	705236	20.000	ng/u1	0.00	
79) Chrysene-d12	21.448	240	812649	20.000	ng/u1	0.01	
88) Perylene-d12	24.468	264	828738	20.000	ng/u1	0.02	
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.264	96	27465	9.920	ng/uL	0.00	
4) Pyridine-d5	3.669	84	190384	21.799	ng/u1	0.00	
7) Phenol-d5	6.983	99	202715	20.287	ng/u1	0.00	
9) Bis-(2-Chloroethyl)eth...	7.130	67	135302	19.605	ng/u1	0.00	
11) 2-Chlorophenol-d4	7.335	132	133996	20.714	ng/u1	0.00	
15) 4-Methylphenol-d8	8.516	113	152079	19.603	ng/u1	0.00	
21) Nitrobenzene-d5	8.951	128	72288	23.105	ng/u1	0.00	
24) 2-Nitrophenol-d4	9.674	143	80541	22.965	ng/u1	0.00	
28) 2,4-Dichlorophenol-d3	10.220	165	153081	22.545	ng/u1	0.00	
31) 4-Chloroaniline-d4	10.726	131	195165	22.637	ng/u1	0.00	
46) Dimethylphthalate-d6	13.845	166	453463	22.104	ng/u1	0.00	
49) Acenaphthylene-d8	14.133	160	541524	23.156	ng/u1	0.00	
54) 4-Nitrophenol-d4	14.674	143	59978	20.635	ng/u1	0.00	
60) Fluorene-d10	15.438	176	405041	22.331	ng/u1	0.00	
65) 4,6-Dinitro-2-methylph...	15.573	200	66876	18.264	ng/u1	0.00	
73) Anthracene-d10	17.294	188	705851	23.280	ng/u1	0.00	
81) Pyrene-d10	19.598	212	873779	20.662	ng/u1	0.01	
92) Benzo(a)pyrene-d12	24.263	264	884286	22.722	ng/u1	0.02	
Target Compounds							
2) 1,4-Dioxane	3.293	88	29800	9.095	ng/uL	95	
5) Pyridine	3.687	79	200522	21.435	ng/u1	98	
6) Benzaldehyde	6.942	77	144397	23.559	ng/u1	92	
8) Phenol	7.012	94	217361	20.411	ng/u1	93	
10) Bis(2-Chloroethyl)ether	7.224	93	164480	20.187	ng/u1	94	
12) 2-Chlorophenol	7.371	128	137379	20.980	ng/u1	97	
13) 2-Methylphenol	8.252	108	155713	20.231	ng/u1	96	
14) 2,2'-oxybis(1-Chloropr...	8.317	45	248434	20.539	ng/u1	99	
16) Acetophenone	8.616	105	261838	19.899	ng/u1	96	
17) N-Nitroso-di-n-propyla...	8.599	70	150241	19.136	ng/u1#	87	
18) 4-Methylphenol	8.581	108	175007	20.775	ng/u1	89	
19) Hexachloroethane	8.869	117	65027	20.210	ng/u1	97	
22) Nitrobenzene	8.998	77	225625	21.595	ng/u1	97	
23) Isophorone	9.515	82	399355	20.230	ng/u1	98	
25) 2-Nitrophenol	9.703	139	78160	21.685	ng/u1	97	
26) 2,4-Dimethylphenol	9.774	107	169928	20.777	ng/u1	98	
27) Bis(2-Chloroethoxy)met...	9.997	93	230445	21.583	ng/u1	96	
29) 2,4-Dichlorophenol	10.250	162	152932	22.667	ng/u1	96	
30) Naphthalene	10.637	128	477453	21.793	ng/u1	97	
32) 4-Chloroaniline	10.749	127	187943	22.911	ng/u1	100	
33) Hexachlorobutadiene	10.919	225	110553	20.169	ng/u1	95	
34) Caprolactam	11.501	113	52285m	22.418	ng/u1		
35) 4-Chloro-3-methylphenol	11.895	107	163847	20.685	ng/u1	97	
36) 2-Methylnaphthalene	12.253	142	336034	21.341	ng/u1	100	

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
37) 1-Methylnaphthalene	12.471	142	343453	21.414	ng/ul	99
39) 1,2,4,5-Tetrachloroben...	12.629	216	207654	21.515	ng/ul	96
40) Hexachlorocyclopentadiene	12.600	237	68209	17.877	ng/ul	96
41) 2,4,6-Trichlorophenol	12.876	196	124282	22.783	ng/ul	97
42) 2,4,5-Trichlorophenol	12.958	196	128105	22.503	ng/ul	96
43) 1,1'-Biphenyl	13.270	154	453345	22.801	ng/ul	96
44) 2-Chloronaphthalene	13.311	162	372165	23.310	ng/ul	96
45) 2-Nitroaniline	13.522	65	124651	23.308	ng/ul	96
47) Dimethylphthalate	13.892	163	452193	21.993	ng/ul	96
48) 2,6-Dinitrotoluene	14.022	165	90706	23.755	ng/ul#	92
50) Acenaphthylene	14.163	152	569706	22.720	ng/ul	99
51) 3-Nitroaniline	14.351	138	78017	22.099	ng/ul	91
52) Acenaphthene	14.504	153	403520	22.704	ng/ul	98
53) 2,4-Dinitrophenol	14.586	184	24865	12.742	ng/ul#	90
55) 4-Nitrophenol	14.692	109	56069	18.577	ng/ul	89
56) Dibenzofuran	14.844	168	572875	22.863	ng/ul	97
57) 2,4-Dinitrotoluene	14.821	165	137420	24.220	ng/ul	98
58) 2,3,4,6-Tetrachlorophenol	15.079	232	108362	22.082	ng/ul#	90
59) Diethylphthalate	15.261	149	447900	22.534	ng/ul	99
61) Fluorene	15.496	166	460759	23.029	ng/ul	98
62) 4-Chlorophenyl-phenyle...	15.485	204	235902	21.163	ng/ul	98
63) 4-Nitroaniline	15.520	138	65876m	18.639	ng/ul	
66) 4,6-Dinitro-2-methylph...	15.590	198	67394	17.866	ng/ul#	92
67) N-Nitrosodiphenylamine	15.702	169	388609	22.755	ng/ul	97
68) 4-Bromophenyl-phenylether	16.384	248	158962	22.117	ng/ul	96
69) Hexachlorobenzene	16.507	284	177035	21.985	ng/ul	96
70) Atrazine	16.654	200	161965	22.436	ng/ul	96
71) Pentachlorophenol	16.860	266	59828m	17.621	ng/ul	
72) Phenanthrene	17.236	178	800666	23.559	ng/ul	100
74) Anthracene	17.330	178	824948	23.983	ng/ul	98
75) 1,2,3,4-Tetrachloroben...	13.234	216	206734	21.971	ng/ul	96
76) Pentachlorobenzene	14.768	250	207186	22.226	ng/ul	97
77) Carbazole	17.606	167	731952	26.288	ng/ul	97
78) Di-n-butylphthalate	18.164	149	835572	24.849	ng/ul	99
80) Fluoranthene	19.263	202	1009113	20.995	ng/ul	99
82) Pyrene	19.627	202	1094248	21.398	ng/ul	95
83) Butylbenzylphthalate	20.520	149	395491	25.194	ng/ul	96
84) 3,3'-Dichlorobenzidine	21.354	252	321291	21.901	ng/ul	93
85) Benzo(a)anthracene	21.431	228	1111638	21.827	ng/ul	99
86) Bis(2-ethylhexyl)phtha...	21.343	149	602745	26.396	ng/ul	96
87) Chrysene	21.495	228	1076469	22.177	ng/ul	99
89) Di-n-octyl phthalate	22.477	149	993639	27.779	ng/ul	100
90) Benzo(b)fluoranthene	23.516	252	1069757	22.712	ng/ul	99
91) Benzo(k)fluoranthene	23.575	252	1082529	23.040	ng/ul	97
93) Benzo(a)pyrene	24.327	252	999061	22.598	ng/ul	97
94) Indeno(1,2,3-cd)pyrene	27.876	276	1226299	22.712	ng/ul	96
95) Dibenzo(a,h)anthracene	27.917	278	1014201m	23.476	ng/ul	
96) Benzo(g,h,i)perylene	28.940	276	951144	22.200	ng/ul	98

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

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