

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101921\
 Data File : BG050628.D
 Acq On : 19 Oct 2021 9:56
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Quant Time: Oct 19 11:42:39 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 14 10:56:22 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.250	152	40965	20.00	ng	0.00
21) Naphthalene-d8	11.076	136	170340	20.00	ng	# 0.00
39) Acenaphthene-d10	14.871	164	110744	20.00	ng	0.00
64) Phenanthrene-d10	17.615	188	252150	20.00	ng	# 0.00
76) Chrysene-d12	21.916	240	234046	20.00	ng	# 0.00
86) Perylene-d12	25.330	264	240579	20.00	ng	0.00

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System Monitoring Compounds						
5) 2-Fluorophenol	5.788	112	212719	77.66	ng	0.00
7) Phenol-d6	7.392	99	309194	77.96	ng	0.00
23) Nitrobenzene-d5	9.425	82	309072	75.11	ng	0.00
42) 2,4,6-Tribromophenol	16.358	330	81070	76.75	ng	0.00
45) 2-Fluorobiphenyl	13.496	172	550738	74.85	ng	0.00
79) Terphenyl-d14	20.206	244	915915	76.27	ng	0.00

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Target Compounds						Qvalue
2) 1,4-Dioxane	3.655	88	52085	36.71	ng	91
3) Pyridine	4.066	79	147327	38.02	ng	93
4) n-Nitrosodimethylamine	3.978	42	71725	39.27	ng	# 52
6) Aniline	7.574	93	175385	38.06	ng	# 93
8) 2-Chlorophenol	7.815	128	104913	39.78	ng	86
9) Benzaldehyde	7.386	77	102654	41.21	ng	88
10) Phenol	7.421	94	149838	38.86	ng	87
11) bis(2-Chloroethyl)ether	7.662	93	114212	37.95	ng	83
12) 1,3-Dichlorobenzene	8.144	146	115841	38.05	ng	95
13) 1,4-Dichlorobenzene	8.291	146	117679	38.34	ng	96
14) 1,2-Dichlorobenzene	8.614	146	111055	37.88	ng	95
15) Benzyl Alcohol	8.485	79	121985	38.92	ng	91
16) 2,2'-oxybis(1-Chloropr...	8.773	45	181724	37.02	ng	86
17) 2-Methylphenol	8.684	107	99158	39.08	ng	94
18) Hexachloroethane	9.348	117	44115	37.98	ng	95
19) n-Nitroso-di-n-propyla...	9.055	70	107753	37.28	ng	# 87
20) 3+4-Methylphenols	9.013	107	139702	39.12	ng	95
22) Acetophenone	9.078	105	177651	36.68	ng	# 97
24) Nitrobenzene	9.466	77	152433	37.26	ng	95
25) Isophorone	9.989	82	266915	36.57	ng	# 95
26) 2-Nitrophenol	10.183	139	60407	40.21	ng	# 92
27) 2,4-Dimethylphenol	10.224	122	97504	37.56	ng	92
28) bis(2-Chloroethoxy)met...	10.465	93	153313	36.74	ng	93
29) 2,4-Dichlorophenol	10.717	162	101317	38.30	ng	94
30) 1,2,4-Trichlorobenzene	10.935	180	113481	37.76	ng	97
31) Naphthalene	11.129	128	340535	37.35	ng	99
32) Benzoic acid	10.353	122	52751	27.38	ng	# 72
33) 4-Chloroaniline	11.234	127	146237	38.23	ng	97
34) Hexachlorobutadiene	11.399	225	69871	37.03	ng	93
35) Caprolactam	12.004	113	38560	38.08	ng	# 74
36) 4-Chloro-3-methylphenol	12.327	107	124346	37.60	ng	# 89
37) 2-Methylnaphthalene	12.715	142	233307	37.09	ng	92
38) 1-Methylnaphthalene	12.932	142	222968	36.55	ng	93
40) 1,2,4,5-Tetrachloroben...	13.073	216	123039	37.62	ng	98
41) Hexachlorocyclopentadiene	13.050	237	61370	32.47	ng	92
43) 2,4,6-Trichlorophenol	13.308	196	87906	38.15	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.379	196	93612	38.85	ng	90
46) 1,1'-Biphenyl	13.708	154	301349	37.31	ng	92
47) 2-Chloronaphthalene	13.755	162	238285	38.42	ng	98
48) 2-Nitroaniline	13.955	65	98706	40.00	ng	94
49) Acenaphthylene	14.601	152	382026	37.59	ng	96
50) Dimethylphthalate	14.313	163	318676	38.07	ng	99
51) 2,6-Dinitrotoluene	14.442	165	64783	38.64	ng	91
52) Acenaphthene	14.936	154	239602m	36.69	ng	
53) 3-Nitroaniline	14.771	138	77812	39.24	ng	92
54) 2,4-Dinitrophenol	14.977	184	34987	33.64	ng	86
55) Dibenzofuran	15.271	168	377443	37.37	ng	97
56) 4-Nitrophenol	15.065	139	60575	37.97	ng	# 82
57) 2,4-Dinitrotoluene	15.224	165	93955	39.76	ng	93
58) Fluorene	15.917	166	288919	37.24	ng	95
59) 2,3,4,6-Tetrachlorophenol	15.488	232	78146	37.23	ng	96
60) Diethylphthalate	15.670	149	333214	37.89	ng	97
61) 4-Chlorophenyl-phenyle...	15.900	204	159095	37.48	ng	97
62) 4-Nitroaniline	15.935	138	81735	39.61	ng	# 66
63) Azobenzene	16.193	77	376436	36.97	ng	83
65) 4,6-Dinitro-2-methylph...	15.988	198	57588	38.02	ng	84
66) n-Nitrosodiphenylamine	16.117	169	261673	36.52	ng	97
67) 4-Bromophenyl-phenylether	16.798	248	95158	37.45	ng	97
68) Hexachlorobenzene	16.916	284	94579	37.45	ng	96
69) Atrazine	17.051	200	105332	37.32	ng	97
70) Pentachlorophenol	17.257	266	54909	31.94	ng	96
71) Phenanthrene	17.662	178	487695	36.70	ng	97
72) Anthracene	17.750	178	491189	37.20	ng	98
73) Carbazole	18.015	167	500414	38.03	ng	99
74) Di-n-butylphthalate	18.555	149	604055	39.11	ng	# 97
75) Fluoranthene	19.654	202	609063	37.29	ng	96
77) Benzidine	19.830	184	273871	36.10	ng	97
78) Pyrene	20.018	202	616208	37.11	ng	98
80) Butylbenzylphthalate	20.888	149	281308	40.05	ng	98
81) Benzo(a)anthracene	21.898	228	599231	38.69	ng	100
82) 3,3'-Dichlorobenzidine	21.804	252	195347	38.06	ng	# 96
83) Chrysene	21.969	228	554185	38.04	ng	95
84) Bis(2-ethylhexyl)phtha...	21.769	149	397139	39.79	ng	# 95
85) Di-n-octyl phthalate	23.050	149	676816	40.66	ng	97
87) Indeno(1,2,3-cd)pyrene	29.249	276	647175	38.62	ng	# 89
88) Benzo(b)fluoranthene	24.237	252	566067	38.42	ng	# 92
89) Benzo(k)fluoranthene	24.307	252	567022	39.12	ng	# 97
90) Benzo(a)pyrene	25.165	252	485364	38.75	ng	# 94
91) Dibenzo(a,h)anthracene	29.319	278	529703	38.22	ng	# 94
92) Benzo(g,h,i)perylene	30.488	276	519959	38.30	ng	# 91

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

