

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121720\
 Data File : BG047764.D
 Acq On : 17 Dec 2020 13:11
 Operator : CG/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

mohammad
 12/18/2020 10:29:16 AM

Quant Time: Dec 17 14:04:06 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG121720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 17 13:16:53 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.204	152	29590	20.00	ng	# 0.00
21) Naphthalene-d8	11.030	136	107494	20.00	ng	# 0.00
39) Acenaphthene-d10	14.826	164	84491	20.00	ng	0.00
64) Phenanthrene-d10	17.564	188	206481	20.00	ng	# 0.00
76) Chrysene-d12	21.841	240	207803	20.00	ng	# 0.00
86) Perylene-d12	25.190	264	223056	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.731	112	180499	98.88	ng	0.00
7) Phenol-d6	7.347	99	261982	99.14	ng	0.00
23) Nitrobenzene-d5	9.374	82	300960	120.15	ng	0.00
42) 2,4,6-Tribromophenol	16.307	330	150040	106.05	ng	0.00
45) 2-Fluorobiphenyl	13.451	172	644588	97.85	ng	0.00
79) Terphenyl-d14	20.149	244	1290980	109.25	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.563	88	37010	40.94	ng	# 80
3) Pyridine	3.980	79	105511	43.64	ng	# 85
4) n-Nitrosodimethylamine	3.886	42	79284	65.39	ng	# 45
6) Aniline	7.517	93	146411	46.93	ng	94
8) 2-Chlorophenol	7.764	128	92840	47.90	ng	98
9) Benzaldehyde	7.329	77	75402	42.42	ng	# 81
10) Phenol	7.370	94	120872	48.22	ng	# 51
11) bis(2-Chloroethyl)ether	7.611	93	84808	42.65	ng	97
12) 1,3-Dichlorobenzene	8.093	146	120668	47.84	ng	# 84
13) 1,4-Dichlorobenzene	8.240	146	120099m	47.41	ng	
14) 1,2-Dichlorobenzene	8.563	146	113403	47.20	ng	92
15) Benzyl Alcohol	8.433	79	120038	59.11	ng	# 79
16) 2,2'-oxybis(1-Chloropr...	8.733	45	81228	26.47	ng	79
17) 2-Methylphenol	8.633	107	80988	47.74	ng	# 82
18) Hexachloroethane	9.303	117	47642	51.37	ng	92
19) n-Nitroso-di-n-propyla...	9.009	70	88805	51.14	ng	90
20) 3+4-Methylphenols	8.968	107	115762	50.94	ng	# 83
22) Acetophenone	9.033	105	158109	51.01	ng	# 86
24) Nitrobenzene	9.415	77	142157	55.78	ng	# 81
25) Isophorone	9.938	82	237169	55.26	ng	# 92
26) 2-Nitrophenol	10.131	139	57989	52.76	ng	# 54
27) 2,4-Dimethylphenol	10.178	122	78560	53.23	ng	87
28) bis(2-Chloroethoxy)met...	10.419	93	118652	45.68	ng	99
29) 2,4-Dichlorophenol	10.666	162	106435	52.19	ng	97
30) 1,2,4-Trichlorobenzene	10.889	180	135533	53.91	ng	97
31) Naphthalene	11.083	128	304764	49.52	ng	99
32) Benzoic acid	10.320	122	30478	29.66	ng	# 84
33) 4-Chloroaniline	11.177	127	133510	51.57	ng	# 89
34) Hexachlorobutadiene	11.359	225	121172	63.70	ng	97
35) Caprolactam	11.953	113	34379	51.59	ng	# 70
36) 4-Chloro-3-methylphenol	12.282	107	117477	56.52	ng	# 88
37) 2-Methylnaphthalene	12.670	142	241028	52.11	ng	# 86
38) 1-Methylnaphthalene	12.887	142	216853	49.34	ng	# 96
40) 1,2,4,5-Tetrachloroben...	13.034	216	172138	51.60	ng	# 96
41) Hexachlorocyclopentadiene	13.010	237	102846	59.48	ng	95
43) 2,4,6-Trichlorophenol	13.263	196	115420	53.22	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.334	196	111867	50.80	ng	#	90
46) 1,1'-Biphenyl	13.663	154	315729	45.64	ng		97
47) 2-Chloronaphthalene	13.710	162	261282	46.31	ng		95
48) 2-Nitroaniline	13.904	65	100970	54.00	ng	#	77
49) Acenaphthylene	14.550	152	388202	47.04	ng		98
50) Dimethylphthalate	14.268	163	366879	49.17	ng		98
51) 2,6-Dinitrotoluene	14.391	165	75221	48.60	ng	#	45
52) Acenaphthene	14.891	154	283733	53.78	ng		98
53) 3-Nitroaniline	14.720	138	75747	49.17	ng	#	78
54) 2,4-Dinitrophenol	14.926	184	50497	50.86	ng	#	34
55) Dibenzofuran	15.220	168	398966	45.99	ng		94
56) 4-Nitrophenol	15.014	139	59703	50.03	ng	#	28
57) 2,4-Dinitrotoluene	15.173	165	110798	49.04	ng	#	63
58) Fluorene	15.866	166	334582	47.99	ng		99
59) 2,3,4,6-Tetrachlorophenol	15.443	232	116287	50.30	ng		93
60) Diethylphthalate	15.619	149	359626	48.79	ng		96
61) 4-Chlorophenyl-phenyle...	15.854	204	214893	52.04	ng		96
62) 4-Nitroaniline	15.878	138	82244	49.71	ng	#	13
63) Azobenzene	16.148	77	336202	50.90	ng		85
65) 4,6-Dinitro-2-methylph...	15.936	198	73360	51.10	ng		82
66) n-Nitrosodiphenylamine	16.066	169	305513	48.10	ng		92
67) 4-Bromophenyl-phenylether	16.747	248	142533	49.43	ng	#	90
68) Hexachlorobenzene	16.871	284	154444	49.29	ng	#	91
69) Atrazine	17.000	200	129856	47.80	ng		98
70) Pentachlorophenol	17.211	266	91778	53.75	ng		94
71) Phenanthrene	17.605	178	577567	48.20	ng		95
72) Anthracene	17.693	178	557678	47.91	ng		97
73) Carbazole	17.958	167	495973	48.26	ng		97
74) Di-n-butylphthalate	18.498	149	592812	47.17	ng	#	96
75) Fluoranthene	19.597	202	771815	50.57	ng		93
77) Benzidine	19.767	184	275926	39.59	ng		96
78) Pyrene	19.961	202	769090	53.15	ng		95
80) Butylbenzylphthalate	20.825	149	272599	49.93	ng	#	86
81) Benzo(a)anthracene	21.818	228	762414	51.76	ng		97
82) 3,3'-Dichlorobenzidine	21.724	252	266105	52.21	ng	#	93
83) Chrysene	21.888	228	709375	50.03	ng		98
84) Bis(2-ethylhexyl)phtha...	21.700	149	364065	47.67	ng		96
85) Di-n-octyl phthalate	22.952	149	620536	48.35	ng	#	95
87) Indeno(1,2,3-cd)pyrene	29.050	276	889950	52.56	ng	#	86
88) Benzo(b)fluoranthene	24.127	252	811032	53.26	ng	#	94
89) Benzo(k)fluoranthene	24.197	252	776928	52.23	ng	#	94
90) Benzo(a)pyrene	25.032	252	752713	52.96	ng	#	93
91) Dibenzo(a,h)anthracene	29.115	278	723588	52.51	ng	#	91
92) Benzo(g,h,i)perylene	30.255	276	710790	52.01	ng	#	89

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

