

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061121\
 Data File : BG048909.D
 Acq On : 11 Jun 2021 13:54
 Operator : CG/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

mohammad
 6/14/2021 5:48:01 PM

Quant Time: Jun 11 14:29:15 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 11 13:47:28 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.129	152	40709	20.000	ng	0.00
21) Naphthalene-d8	10.949	136	162122	20.000	ng	# 0.00
39) Acenaphthene-d10	14.769	164	118208	20.000	ng	0.00
64) Phenanthrene-d10	17.518	188	266738	20.000	ng	# 0.00
76) Chrysene-d12	21.819	240	251389	20.000	ng	# 0.00
86) Perylene-d12	25.150	264	264872	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.667	112	310973	127.762	ng	0.00
7) Phenol-d6	7.266	99	466459	135.486	ng	0.00
23) Nitrobenzene-d5	9.298	82	468070	129.258	ng	0.00
42) 2,4,6-Tribromophenol	16.255	330	219023	138.192	ng	0.00
45) 2-Fluorobiphenyl	13.394	172	976559	119.304	ng	0.00
79) Terphenyl-d14	20.127	244	1588942	108.786	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.546	88	74295	77.903	ng	# 81
3) Pyridine	3.958	79	212672m	85.552	ng	
4) n-Nitrosodimethylamine	3.870	42	108995	57.429	ng	# 70
6) Aniline	7.448	93	293427	75.936	ng	# 91
8) 2-Chlorophenol	7.689	128	169213	65.532	ng	85
9) Benzaldehyde	7.260	77	115661	50.700	ng	92
10) Phenol	7.295	94	238456	70.437	ng	92
11) bis(2-Chloroethyl)ether	7.542	93	174832	76.628	ng	83
12) 1,3-Dichlorobenzene	8.018	146	186487	62.814	ng	95
13) 1,4-Dichlorobenzene	8.165	146	187486	61.969	ng	97
14) 1,2-Dichlorobenzene	8.482	146	182476	65.367	ng	94
15) Benzyl Alcohol	8.364	79	216098	71.231	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.652	45	338193	136.739	ng	80
17) 2-Methylphenol	8.558	107	156094	72.223	ng	96
18) Hexachloroethane	9.222	117	79114	66.572	ng	95
19) n-Nitroso-di-n-propyla...	8.940	70	181255	78.034	ng	# 95
20) 3+4-Methylphenols	8.893	107	219600	73.642	ng	98
22) Acetophenone	8.958	105	295821	70.104	ng	# 98
24) Nitrobenzene	9.340	77	260815	72.208	ng	98
25) Isophorone	9.874	82	490116	76.538	ng	# 97
26) 2-Nitrophenol	10.051	139	94635	60.533	ng	90
27) 2,4-Dimethylphenol	10.103	122	157728	65.628	ng	94
28) bis(2-Chloroethoxy)met...	10.350	93	227841	81.666	ng	93
29) 2,4-Dichlorophenol	10.585	162	178379	65.812	ng	94
30) 1,2,4-Trichlorobenzene	10.808	180	200211	63.699	ng	99
31) Naphthalene	11.002	128	572858	68.863	ng	99
32) Benzoic acid	10.250	122	119786	85.067	ng	82
33) 4-Chloroaniline	11.102	127	247978	71.640	ng	97
34) Hexachlorobutadiene	11.290	225	142368	61.016	ng	92
35) Caprolactam	11.901	113	51467	55.880	ng	# 42
36) 4-Chloro-3-methylphenol	12.213	107	220018	71.675	ng	# 89
37) 2-Methylnaphthalene	12.600	142	435248	65.473	ng	98
38) 1-Methylnaphthalene	12.824	142	409295	66.174	ng	93
40) 1,2,4,5-Tetrachloroben...	12.965	216	224001	58.220	ng	98
41) Hexachlorocyclopentadiene	12.947	237	149865	83.317	ng	90
43) 2,4,6-Trichlorophenol	13.200	196	153877	59.925	ng	94

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.270	196	158298	58.052	ng	91
46) 1,1'-Biphenyl	13.605	154	521820	59.763	ng	98
47) 2-Chloronaphthalene	13.646	162	414520	62.149	ng	98
48) 2-Nitroaniline	13.840	65	164318	68.560	ng	95
49) Acenaphthylene	14.492	152	688940	66.208	ng	97
50) Dimethylphthalate	14.222	163	549348	61.954	ng	100
51) 2,6-Dinitrotoluene	14.334	165	122709	59.985	ng	96
52) Acenaphthene	14.833	154	457463	69.230	ng	94
53) 3-Nitroaniline	14.669	138	123551	62.912	ng	82
54) 2,4-Dinitrophenol	14.868	184	61170	57.350	ng	# 83
55) Dibenzofuran	15.168	168	694957	63.533	ng	98
56) 4-Nitrophenol	14.957	139	103524	71.153	ng	# 88
58) Fluorene	15.820	166	558422	62.161	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.385	232	164307	65.123	ng	100
60) Diethylphthalate	15.585	149	593340	64.682	ng	97
61) 4-Chlorophenyl-phenyle...	15.808	204	301894	61.109	ng	93
62) 4-Nitroaniline	15.832	138	126760	60.372	ng	# 80
63) Azobenzene	16.102	77	671826	72.725	ng	87
66) n-Nitrosodiphenylamine	16.020	169	499688	66.280	ng	97
67) 4-Bromophenyl-phenylether	16.707	248	204467	66.601	ng	97
68) Hexachlorobenzene	16.819	284	220320	70.851	ng	97
69) Atrazine	16.972	200	156743	49.998	ng	97
70) Pentachlorophenol	17.160	266	141111	85.158	ng	98
71) Phenanthrene	17.565	178	894931	64.643	ng	97
72) Anthracene	17.653	178	885800	64.662	ng	97
73) Carbazole	17.918	167	853900	65.075	ng	97
74) Di-n-butylphthalate	18.482	149	966428	65.697	ng	98
75) Fluoranthene	19.569	202	1054492	60.229	ng	98
77) Benzidine	19.745	184	313526	45.926	ng	98
78) Pyrene	19.933	202	1066096	61.297	ng	96
80) Butylbenzylphthalate	20.820	149	421917	62.038	ng	99
81) Benzo(a)anthracene	21.796	228	1066044	62.056	ng	98
82) 3,3'-Dichlorobenzidine	21.707	252	362851	61.594	ng	98
83) Chrysene	21.866	228	1033096	63.133	ng	96
84) Bis(2-ethylhexyl)phtha...	21.707	149	589467	61.912	ng	97
85) Di-n-octyl phthalate	22.977	149	1007713	62.172	ng	# 95
87) Indeno(1,2,3-cd)pyrene	28.999	276	1219851	62.706	ng	# 96
88) Benzo(b)fluoranthene	24.099	252	1131125	62.948	ng	# 95
89) Benzo(k)fluoranthene	24.169	252	1053307	62.332	ng	# 97
90) Benzo(a)pyrene	25.004	252	1022747	61.782	ng	# 97
91) Dibenzo(a,h)anthracene	29.075	278	1027394	61.576	ng	# 95
92) Benzo(g,h,i)perylene	30.209	276	1026432	62.314	ng	# 95

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

