

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102121\
 Data File : BG050665.D
 Acq On : 21 Oct 2021 13:05
 Operator : CG/JU
 Sample : PB140074BSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB140074BSD

Manual Integrations
 APPROVED

mohammad
 10/22/2021 2:54:11 PM

Quant Time: Oct 21 13:45:36 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	8.253	152	39394	20.00	ng	0.00
21) Naphthalene-d8	11.079	136	176934	20.00	ng	# 0.00
39) Acenaphthene-d10	14.875	164	114523	20.00	ng	0.00
64) Phenanthrene-d10	17.619	188	251767	20.00	ng	# 0.00
76) Chrysene-d12	21.920	240	195124	20.00	ng	# 0.00
86) Perylene-d12	25.333	264	189927	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.792	112	340155	129.13	ng	0.00
7) Phenol-d6	7.401	99	457195	119.88	ng	0.00
23) Nitrobenzene-d5	9.428	82	325062	76.05	ng	0.00
42) 2,4,6-Tribromophenol	16.361	330	131438	120.33	ng	0.00
45) 2-Fluorobiphenyl	13.500	172	598862	78.70	ng	0.00
79) Terphenyl-d14	20.204	244	881214	88.01	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.647	88	36096	26.46	ng	90
3) Pyridine	4.058	79	96774	25.97	ng	91
4) n-Nitrosodimethylamine	3.970	42	72863	41.49	ng	# 52
6) Aniline	7.572	93	202812	45.77	ng	93
8) 2-Chlorophenol	7.813	128	123796	48.81	ng	89
9) Benzaldehyde	7.384	77	101259	42.41	ng	90
10) Phenol	7.425	94	171834	46.34	ng	86
11) bis(2-Chloroethyl)ether	7.660	93	126023	43.54	ng	85
12) 1,3-Dichlorobenzene	8.142	146	125889	43.00	ng	99
13) 1,4-Dichlorobenzene	8.289	146	128558	43.56	ng	95
14) 1,2-Dichlorobenzene	8.612	146	125229	44.42	ng	95
15) Benzyl Alcohol	8.488	79	141668	47.01	ng	92
16) 2,2'-oxybis(1-Chloropr...	8.770	45	208132	44.10	ng	85
17) 2-Methylphenol	8.688	107	115620	47.39	ng	91
18) Hexachloroethane	9.346	117	50514	45.22	ng	95
19) n-Nitroso-di-n-propyla...	9.058	70	121208	43.60	ng	# 88
20) 3+4-Methylphenols	9.017	107	159906	46.56	ng	97
22) Acetophenone	9.082	105	201102	39.98	ng	# 97
24) Nitrobenzene	9.470	77	173598	40.85	ng	95
25) Isophorone	9.987	82	308763	40.73	ng	# 95
26) 2-Nitrophenol	10.186	139	71518	45.83	ng	# 93
27) 2,4-Dimethylphenol	10.228	122	127078	47.13	ng	94
28) bis(2-Chloroethoxy)met...	10.468	93	173364	39.99	ng	92
29) 2,4-Dichlorophenol	10.715	162	121961	44.38	ng	98
30) 1,2,4-Trichlorobenzene	10.938	180	127065	40.71	ng	93
31) Naphthalene	11.132	128	391354	41.32	ng	99
32) Benzoic acid	10.374	122	66570	33.27	ng	# 77
33) 4-Chloroaniline	11.232	127	161113	40.55	ng	98
34) Hexachlorobutadiene	11.403	225	78720	40.17	ng	95
35) Caprolactam	12.008	113	46098m	43.83	ng	
36) 4-Chloro-3-methylphenol	12.331	107	149434	43.51	ng	# 89
37) 2-Methylnaphthalene	12.719	142	279672	42.80	ng	92
38) 1-Methylnaphthalene	12.936	142	257313	40.61	ng	93
40) 1,2,4,5-Tetrachloroben...	13.077	216	139422	41.23	ng	97
41) Hexachlorocyclopentadiene	13.054	237	176285	90.19	ng	93
43) 2,4,6-Trichlorophenol	13.312	196	102378	42.96	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.383	196	109411	43.91	ng	93
46) 1,1'-Biphenyl	13.712	154	338996	40.58	ng	93
47) 2-Chloronaphthalene	13.759	162	269332	41.99	ng	98
48) 2-Nitroaniline	13.958	65	118006	46.25	ng	95
49) Acenaphthylene	14.599	152	445911	42.43	ng	97
50) Dimethylphthalate	14.317	163	367116	42.41	ng	99
51) 2,6-Dinitrotoluene	14.446	165	80739	46.57	ng	90
52) Acenaphthene	14.940	154	316342m	46.84	ng	
53) 3-Nitroaniline	14.775	138	81750	39.86	ng	87
54) 2,4-Dinitrophenol	14.981	184	86024	79.98	ng	93
55) Dibenzofuran	15.269	168	424062	40.60	ng	97
56) 4-Nitrophenol	15.075	139	139075	84.30	ng	# 79
57) 2,4-Dinitrotoluene	15.228	165	114688	46.93	ng	95
58) Fluorene	15.921	166	340621	42.46	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.492	232	89696	41.33	ng	97
60) Diethylphthalate	15.668	149	383002	42.11	ng	96
61) 4-Chlorophenyl-phenyle...	15.903	204	182899	41.66	ng	98
62) 4-Nitroaniline	15.938	138	90259	42.30	ng	# 65
63) Azobenzene	16.197	77	426317	40.49	ng	82
65) 4,6-Dinitro-2-methylph...	15.991	198	61256	40.51	ng	86
66) n-Nitrosodiphenylamine	16.121	169	305117	42.65	ng	97
67) 4-Bromophenyl-phenylether	16.796	248	109027	42.97	ng	100
68) Hexachlorobenzene	16.920	284	109039	43.24	ng	# 94
69) Atrazine	17.055	200	124081	44.03	ng	98
70) Pentachlorophenol	17.260	266	129972	75.72	ng	99
71) Phenanthrene	17.666	178	555908	41.90	ng	98
72) Anthracene	17.754	178	566841	43.00	ng	98
73) Carbazole	18.018	167	525285	39.98	ng	98
74) Di-n-butylphthalate	18.553	149	651228	42.23	ng	# 97
75) Fluoranthene	19.658	202	634677	38.92	ng	96
77) Benzidine	19.834	184	542766	85.82	ng	97
78) Pyrene	20.022	202	636274	45.96	ng	97
80) Butylbenzylphthalate	20.886	149	260180	44.43	ng	94
81) Benzo(a)anthracene	21.902	228	553393	42.86	ng	99
82) 3,3'-Dichlorobenzidine	21.808	252	191492	44.75	ng	# 96
83) Chrysene	21.973	228	532322	43.83	ng	96
84) Bis(2-ethylhexyl)phtha...	21.773	149	365820	43.97	ng	# 96
85) Di-n-octyl phthalate	23.048	149	621615	44.79	ng	96
87) Indeno(1,2,3-cd)pyrene	29.264	276	598698	45.26	ng	# 92
88) Benzo(b)fluoranthene	24.246	252	549802	47.27	ng	# 95
89) Benzo(k)fluoranthene	24.317	252	524158	45.81	ng	# 97
90) Benzo(a)pyrene	25.181	252	529077	53.50	ng	# 95
91) Dibenzo(a,h)anthracene	29.329	278	508619	46.48	ng	# 93
92) Benzo(g,h,i)perylene	30.504	276	493519	46.05	ng	# 92

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

