

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040221\
 Data File : BG048565.D
 Acq On : 2 Apr 2021 12:32
 Operator : CG/JU
 Sample : PB135464BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB135464BS

Manual Integrations
 APPROVED

mohammad
 4/5/2021 11:54:34 AM

Quant Time: Apr 02 13:14:21 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 02 10:20:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.093	152	22077	20.00	ng	0.00
21) Naphthalene-d8	10.925	136	92834	20.00	ng	# 0.00
39) Acenaphthene-d10	14.750	164	65996	20.00	ng	0.00
64) Phenanthrene-d10	17.506	188	171856	20.00	ng	# 0.00
76) Chrysene-d12	21.801	240	158606	20.00	ng	# 0.00
86) Perylene-d12	25.144	264	125295	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.643	112	122541	98.00	ng	0.00
7) Phenol-d6	7.247	99	160048	88.25	ng	0.00
23) Nitrobenzene-d5	9.268	82	134788	70.30	ng	0.00
42) 2,4,6-Tribromophenol	16.243	330	76001	87.61	ng	0.00
45) 2-Fluorobiphenyl	13.369	172	320958	72.28	ng	0.00
79) Terphenyl-d14	20.114	244	574140	66.20	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.505	88	17914	35.61	ng	# 92
3) Pyridine	3.910	79	49407	39.17	ng	96
4) n-Nitrosodimethylamine	3.816	42	37988	46.10	ng	93
6) Aniline	7.412	93	70396	33.80	ng	# 94
8) 2-Chlorophenol	7.658	128	69278	49.02	ng	94
9) Benzaldehyde	7.224	77	9982	8.61	ng	97
10) Phenol	7.277	94	87404	48.14	ng	98
11) bis(2-Chloroethyl)ether	7.512	93	57283	45.47	ng	94
12) 1,3-Dichlorobenzene	7.987	146	73085	45.91	ng	97
13) 1,4-Dichlorobenzene	8.134	146	75249	46.88	ng	98
14) 1,2-Dichlorobenzene	8.458	146	70985	46.18	ng	95
15) Benzyl Alcohol	8.334	79	69296	46.65	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.628	45	52837	43.48	ng	96
17) 2-Methylphenol	8.534	107	59059	50.27	ng	86
18) Hexachloroethane	9.192	117	27540	46.18	ng	# 82
19) n-Nitroso-di-n-propyla...	8.904	70	52428	41.94	ng	99
20) 3+4-Methylphenols	8.869	107	80670	49.47	ng	96
22) Acetophenone	8.922	105	103807	43.37	ng	# 97
24) Nitrobenzene	9.309	77	83435	44.34	ng	99
25) Isophorone	9.832	82	147065	41.53	ng	# 97
26) 2-Nitrophenol	10.026	139	39894	46.14	ng	98
27) 2,4-Dimethylphenol	10.079	122	70566	51.87	ng	97
28) bis(2-Chloroethoxy)met...	10.320	93	81494	50.21	ng	98
29) 2,4-Dichlorophenol	10.567	162	71551	46.45	ng	94
30) 1,2,4-Trichlorobenzene	10.784	180	80657	45.56	ng	95
31) Naphthalene	10.978	128	212112	44.44	ng	99
32) Benzoic acid	10.214	122	46918	44.83	ng	94
33) 4-Chloroaniline	11.078	127	53081	26.35	ng	95
34) Hexachlorobutadiene	11.260	225	54738	42.34	ng	94
35) Caprolactam	11.854	113	25854m	46.00	ng	
36) 4-Chloro-3-methylphenol	12.200	107	81284	46.99	ng	99
37) 2-Methylnaphthalene	12.582	142	164759	43.36	ng	94
38) 1-Methylnaphthalene	12.800	142	155712	42.97	ng	98
40) 1,2,4,5-Tetrachloroben...	12.946	216	92829	45.37	ng	97
41) Hexachlorocyclopentadiene	12.923	237	107062	90.34	ng	93
43) 2,4,6-Trichlorophenol	13.181	196	65755	46.76	ng	94

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.258	196	70723	47.01	ng	94
46) 1,1'-Biphenyl	13.581	154	215926	44.78	ng	98
47) 2-Chloronaphthalene	13.628	162	167125	45.49	ng	96
48) 2-Nitroaniline	13.828	65	57794	49.49	ng	90
49) Acenaphthylene	14.474	152	257413	44.70	ng	98
50) Dimethylphthalate	14.198	163	229972	46.95	ng	# 98
51) 2,6-Dinitrotoluene	14.315	165	53023	47.32	ng	96
52) Acenaphthene	14.815	154	216855m	52.05	ng	
53) 3-Nitroaniline	14.650	138	37741	34.34	ng	81
54) 2,4-Dinitrophenol	14.856	184	69802	110.77	ng	98
55) Dibenzofuran	15.150	168	270807	44.51	ng	# 94
56) 4-Nitrophenol	14.956	139	91713	103.39	ng	87
57) 2,4-Dinitrotoluene	15.103	165	72626	45.81	ng	91
58) Fluorene	15.796	166	226887	44.55	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.373	232	70883	48.30	ng	99
60) Diethylphthalate	15.561	149	237136	47.11	ng	97
61) 4-Chlorophenyl-phenyle...	15.790	204	122341	44.65	ng	96
62) 4-Nitroaniline	15.814	138	53158	46.25	ng	97
63) Azobenzene	16.084	77	210322	42.48	ng	97
65) 4,6-Dinitro-2-methylph...	15.867	198	50384	48.79	ng	90
66) n-Nitrosodiphenylamine	16.002	169	204982	41.92	ng	97
67) 4-Bromophenyl-phenylether	16.683	248	82761	43.44	ng	95
68) Hexachlorobenzene	16.807	284	84767	42.61	ng	96
69) Atrazine	16.948	200	75159	37.69	ng	90
70) Pentachlorophenol	17.147	266	119321	96.51	ng	91
71) Phenanthrene	17.547	178	385645	43.23	ng	98
72) Anthracene	17.641	178	390480	43.92	ng	99
73) Carbazole	17.905	167	364802	43.27	ng	98
74) Di-n-butylphthalate	18.458	149	439472	46.66	ng	99
75) Fluoranthene	19.556	202	490372	44.34	ng	99
77) Benzidine	19.733	184	132391	24.65	ng	97
78) Pyrene	19.921	202	470580	43.15	ng	99
80) Butylbenzylphthalate	20.796	149	180695	44.61	ng	90
81) Benzo(a)anthracene	21.783	228	437020	41.24	ng	99
82) 3,3'-Dichlorobenzidine	21.695	252	136512	37.50	ng	92
83) Chrysene	21.848	228	420902	41.62	ng	94
84) Bis(2-ethylhexyl)phtha...	21.677	149	255076	45.21	ng	100
85) Di-n-octyl phthalate	22.935	149	426719	43.39	ng	# 97
87) Indeno(1,2,3-cd)pyrene	28.986	276	491590	55.96	ng	# 92
88) Benzo(b)fluoranthene	24.080	252	436722	53.67	ng	# 98
89) Benzo(k)fluoranthene	24.157	252	435077	55.61	ng	98
90) Benzo(a)pyrene	24.991	252	396393	52.85	ng	100
91) Dibenzo(a,h)anthracene	29.063	278	420287	56.06	ng	99
92) Benzo(g,h,i)perylene	30.191	276	408785	55.67	ng	# 92

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

