

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012821\
 Data File : BG048055.D
 Acq On : 28 Jan 2021 12:45
 Operator : CG/JU
 Sample : PB134332BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB134332BSD

Manual Integrations
 APPROVED

mohammad
 1/29/2021 10:33:04 AM

Quant Time: Jan 28 13:42:05 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 25 19:39:42 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.123	152	39798	20.00	ng	# 0.00
21) Naphthalene-d8	10.937	136	155004	20.00	ng	0.00
39) Acenaphthene-d10	14.745	164	113890	20.00	ng	0.00
64) Phenanthrene-d10	17.488	188	285557	20.00	ng	# 0.00
76) Chrysene-d12	21.766	240	333070	20.00	ng	# 0.00
86) Perylene-d12	25.044	264	328324	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.679	112	324779	125.42	ng	0.00
7) Phenol-d6	7.271	99	451376	114.64	ng	0.00
23) Nitrobenzene-d5	9.286	82	265959	67.89	ng	0.00
42) 2,4,6-Tribromophenol	16.231	330	215698	113.71	ng	0.00
45) 2-Fluorobiphenyl	13.370	172	593219	67.34	ng	-0.02
79) Terphenyl-d14	20.085	244	1210636	63.02	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.564	88	45424	39.35	ng	# 96
3) Pyridine	3.963	79	134327	38.77	ng	88
4) n-Nitrosodimethylamine	3.869	42	70586	44.09	ng	# 77
6) Aniline	7.441	93	152579	32.08	ng	92
8) 2-Chlorophenol	7.688	128	127290	46.96	ng	91
9) Benzaldehyde	7.253	77	20913	10.35	ng	87
10) Phenol	7.300	94	174566	46.25	ng	81
11) bis(2-Chloroethyl)ether	7.535	93	128492	43.47	ng	96
12) 1,3-Dichlorobenzene	8.011	146	140123	42.83	ng	# 91
13) 1,4-Dichlorobenzene	8.158	146	141700	43.20	ng	92
14) 1,2-Dichlorobenzene	8.481	146	135283	43.21	ng	96
15) Benzyl Alcohol	8.358	79	142588	42.80	ng	# 86
16) 2,2'-oxybis(1-Chloropr...	8.652	45	158017	38.09	ng	96
17) 2-Methylphenol	8.558	107	113701	44.58	ng	# 87
18) Hexachloroethane	9.216	117	53907	42.13	ng	96
19) n-Nitroso-di-n-propyla...	8.928	70	114038	39.68	ng	98
20) 3+4-Methylphenols	8.881	107	153855	43.60	ng	89
22) Acetophenone	8.945	105	201857	42.26	ng	# 93
24) Nitrobenzene	9.327	77	175072	44.62	ng	# 82
25) Isophorone	9.856	82	294782	41.93	ng	# 93
26) 2-Nitrophenol	10.044	139	73113	45.29	ng	# 72
27) 2,4-Dimethylphenol	10.091	122	122272	52.13	ng	94
28) bis(2-Chloroethoxy)met...	10.332	93	167626	39.63	ng	98
29) 2,4-Dichlorophenol	10.579	162	138890	46.02	ng	96
30) 1,2,4-Trichlorobenzene	10.796	180	159423	43.69	ng	95
31) Naphthalene	10.990	128	384687	42.92	ng	99
32) Benzoic acid	10.215	122	92807	44.18	ng	# 76
33) 4-Chloroaniline	11.084	127	82965	20.25	ng	# 92
34) Hexachlorobutadiene	11.272	225	113088	43.64	ng	98
35) Caprolactam	11.854	113	44000m	38.30	ng	
36) 4-Chloro-3-methylphenol	12.195	107	147071	43.43	ng	90
37) 2-Methylnaphthalene	12.582	142	288570	42.63	ng	# 94
38) 1-Methylnaphthalene	12.800	142	270535	41.89	ng	# 99
40) 1,2,4,5-Tetrachloroben...	12.947	216	192559	44.35	ng	# 97
41) Hexachlorocyclopentadiene	12.929	237	267218	108.65	ng	100
43) 2,4,6-Trichlorophenol	13.176	196	132738	43.83	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.252	196	141780	45.18	ng	# 95
46) 1,1'-Biphenyl	13.581	154	373086	42.37	ng	99
47) 2-Chloronaphthalene	13.628	162	312116	40.73	ng	97
48) 2-Nitroaniline	13.822	65	108143	38.89	ng	# 83
49) Acenaphthylene	14.468	152	472185	41.75	ng	99
50) Dimethylphthalate	14.198	163	407273	39.51	ng	98
51) 2,6-Dinitrotoluene	14.310	165	94017	43.87	ng	# 71
52) Acenaphthene	14.809	154	365037	47.68	ng	96
53) 3-Nitroaniline	14.639	138	64549	27.46	ng	# 77
54) 2,4-Dinitrophenol	14.844	184	128907	96.20	ng	# 67
55) Dibenzofuran	15.144	168	481819	41.49	ng	95
56) 4-Nitrophenol	14.938	139	159401	86.91	ng	# 54
57) 2,4-Dinitrotoluene	15.097	165	132448	42.80	ng	# 81
58) Fluorene	15.790	166	390253	41.80	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.361	232	137090	43.69	ng	# 95
60) Diethylphthalate	15.555	149	417658	39.41	ng	97
61) 4-Chlorophenyl-phenyle...	15.779	204	234773	41.07	ng	98
62) 4-Nitroaniline	15.802	138	90837	35.47	ng	# 67
63) Azobenzene	16.072	77	395362	39.63	ng	90
65) 4,6-Dinitro-2-methylph...	15.861	198	91042	49.40	ng	91
66) n-Nitrosodiphenylamine	15.996	169	360854	42.05	ng	99
67) 4-Bromophenyl-phenylether	16.678	248	159195	41.72	ng	97
68) Hexachlorobenzene	16.795	284	170378	41.06	ng	98
69) Atrazine	16.936	200	134521	41.27	ng	97
70) Pentachlorophenol	17.136	266	242683	96.31	ng	97
71) Phenanthrene	17.530	178	689856	41.87	ng	97
72) Anthracene	17.624	178	703854	43.44	ng	97
73) Carbazole	17.888	167	654820	43.31	ng	99
74) Di-n-butylphthalate	18.440	149	760787	41.20	ng	# 98
75) Fluoranthene	19.533	202	977790	45.26	ng	96
77) Benzidine	19.703	184	221699	23.07	ng	100
78) Pyrene	19.891	202	990980	40.63	ng	97
80) Butylbenzylphthalate	20.773	149	365318	39.21	ng	90
81) Benzo(a)anthracene	21.742	228	1002138	41.40	ng	99
82) 3,3'-Dichlorobenzidine	21.654	252	289050	35.44	ng	# 99
83) Chrysene	21.813	228	965994	42.00	ng	99
84) Bis(2-ethylhexyl)phtha...	21.648	149	522025	40.91	ng	98
85) Di-n-octyl phthalate	22.882	149	882624	41.35	ng	97
87) Indeno(1,2,3-cd)pyrene	28.799	276	1184290	44.84	ng	# 89
88) Benzo(b)fluoranthene	24.004	252	1055592	45.24	ng	# 97
89) Benzo(k)fluoranthene	24.075	252	1002292	44.20	ng	# 97
90) Benzo(a)pyrene	24.897	252	939077	42.85	ng	# 96
91) Dibenzo(a,h)anthracene	28.863	278	973941	44.69	ng	# 95
92) Benzo(g,h,i)perylene	29.968	276	979021	46.63	ng	# 93

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

