

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121321\
 Data File : BP008379.D
 Acq On : 13 Dec 2021 14:51
 Operator : CG/JU
 Sample : M5017-29
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP3

Quant Time: Dec 14 02:19:13 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 11 03:53:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	13	20.000	ng	# 0.00
21) Naphthalene-d8	10.563	136	11	20.000	ng	# 0.00
39) Acenaphthene-d10	14.439	164	37	20.000	ng	# 0.02
64) Phenanthrene-d10	17.216	188	13	20.000	ng	# 0.04
76) Chrysene-d12	21.333	240	108	20.000	ng	# 0.05
86) Perylene-d12	23.721	264	344	20.000	ng	# 0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.346	112	411478	509629.568	ng	-0.01
7) Phenol-d6	6.946	99	528328	484730.651	ng	0.00
23) Nitrobenzene-d5	8.910	82	350355	1448341.827	ng	-0.02
42) 2,4,6-Tribromophenol	15.910	330	244013	515968.395	ng	0.00
45) 2-Fluorobiphenyl	13.022	172	711221	279209.181	ng	-0.01
79) Terphenyl-d14	19.745	244	1417917	274383.152	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.263	88	133	353.080	ng	# 1
3) Pyridine	3.652	79	235	233.744	ng	# 1
4) n-Nitrosodimethylamine	3.581	42	237	565.170	ng	# 1
6) Aniline	7.104	93	91	71.225	ng	# 1
8) 2-Chlorophenol	7.446	128	34	41.136	ng	# 86
9) Benzaldehyde	6.946	77	1045	1942.236	ng	# 10
10) Phenol	6.969	94	7802	6965.778	ng	# 23
11) bis(2-Chloroethyl)ether	7.187	93	64	71.489	ng	# 30
12) 1,3-Dichlorobenzene	7.693	146	29	30.397	ng	# 96
13) 1,4-Dichlorobenzene	7.969	146	22	22.745	ng	# 93
14) 1,2-Dichlorobenzene	7.969	146	22	22.870	ng	# 95
15) Benzyl Alcohol	8.034	79	169	195.700	ng	# 46
16) 2,2'-oxybis(1-Chloropr...	8.393	45	201	148.799	ng	# 80
17) 2-Methylphenol	8.228	107	31	42.543	ng	# 1
18) Hexachloroethane	8.840	117	42	116.778	ng	# 9
19) n-Nitroso-di-n-propyla...	8.593	70	104	137.897	ng	# 1
20) 3+4-Methylphenols	8.557	107	33	32.814	ng	# 1
22) Acetophenone	8.598	105	80	265.886	ng	# 1
24) Nitrobenzene	8.998	77	113	481.688	ng	# 52
25) Isophorone	9.516	82	43	101.605	ng	# 85
26) 2-Nitrophenol	9.681	139	28	276.637	ng	# 92
27) 2,4-Dimethylphenol	9.751	122	5	33.031	ng	# 1
28) bis(2-Chloroethoxy)met...	9.992	93	27	100.103	ng	# 1
29) 2,4-Dichlorophenol	10.234	162	16	88.501	ng	# 1
30) 1,2,4-Trichlorobenzene	10.416	180	24	114.375	ng	# 36
31) Naphthalene	10.610	128	99	175.689	ng	# 74
32) Benzoic acid	9.951	122	12	96.913	ng	# 1
33) 4-Chloroaniline	10.734	127	14	59.887	ng	# 1
34) Hexachlorobutadiene	10.910	225	14	97.493	ng	# 35
35) Caprolactam	11.545	113	24	597.627	ng	# 1
36) 4-Chloro-3-methylphenol	11.869	107	30	145.160	ng	# 34
37) 2-Methylnaphthalene	12.234	142	6	15.055	ng	# 35
38) 1-Methylnaphthalene	12.445	142	58	149.510	ng	# 76
40) 1,2,4,5-Tetrachloroben...	12.610	216	21	17.472	ng	# 27
41) Hexachlorocyclopentadiene	12.451	237	5	10.821	ng	# 78
43) 2,4,6-Trichlorophenol	12.875	196	8	9.824	ng	# 64

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.969	196	26	29.797	ng	# 10
46) 1,1'-Biphenyl	13.016	154	1439	536.039	ng	# 1
47) 2-Chloronaphthalene	13.069	162	14	6.604	ng	# 88
48) 2-Nitroaniline	13.528	65	79	108.137	ng	# 34
49) Acenaphthylene	14.151	152	225	68.795	ng	# 80
50) Dimethylphthalate	13.886	163	367	132.090	ng	# 81
51) 2,6-Dinitrotoluene	14.010	165	36	62.432	ng	# 81
52) Acenaphthene	14.510	154	14	7.183	ng	# 49
53) 3-Nitroaniline	14.363	138	7	11.979	ng	# 30
54) 2,4-Dinitrophenol	14.586	184	29	85.060	ng	# 68
55) Dibenzofuran	14.851	168	56	16.813	ng	# 82
56) 4-Nitrophenol	14.869	139	73	162.979	ng	# 93
57) 2,4-Dinitrotoluene	14.828	165	28	35.383	ng	# 1
58) Fluorene	15.504	166	15	5.543	ng	# 1
59) 2,3,4,6-Tetrachlorophenol	15.069	232	159	204.918	ng	# 31
60) Diethylphthalate	15.304	149	45	16.321	ng	# 51
61) 4-Chlorophenyl-phenylether	15.486	204	31	20.885	ng	# 52
62) 4-Nitroaniline	15.316	138	16	26.386	ng	# 91
63) Azobenzene	15.710	77	75	26.593	ng	# 82
65) 4,6-Dinitro-2-methylph...	15.598	198	23	270.495	ng	# 1
66) n-Nitrosodiphenylamine	15.904	169	8128	21584.343	ng	# 50
67) 4-Bromophenyl-phenylether	16.469	248	47	325.128	ng	# 89
68) Hexachlorobenzene	16.527	284	20	129.430	ng	# 38
69) Atrazine	16.686	200	33	337.657	ng	# 91
70) Pentachlorophenol	16.786	266	14	152.364	ng	# 92
71) Phenanthrene	17.222	178	2607	3798.220	ng	# 98
72) Anthracene	17.222	178	2607	3827.273	ng	# 98
73) Carbazole	17.604	167	1006	1512.917	ng	# 91
74) Di-n-butylphthalate	18.180	149	241	289.530	ng	# 92
75) Fluoranthene	19.198	202	6528	7060.074	ng	# 100
78) Pyrene	19.551	202	6826	1061.943	ng	# 93
80) Butylbenzylphthalate	20.421	149	1243	404.175	ng	# 89
81) Benzo(a)anthracene	21.268	228	10136	1416.083	ng	# 94
82) 3,3'-Dichlorobenzidine	21.239	252	380	112.680	ng	# 46
83) Chrysene	21.315	228	10163	1492.097	ng	# 91
84) Bis(2-ethylhexyl)phtha...	21.280	149	1216	260.420	ng	# 73
85) Di-n-octyl phthalate	22.151	149	845	106.529	ng	# 56
87) Indeno(1,2,3-cd)pyrene	26.139	276	9334	372.966	ng	# 95
88) Benzo(b)fluoranthene	22.980	252	18453	889.814	ng	# 83
89) Benzo(k)fluoranthene	22.980	252	18453	908.891	ng	# 82
90) Benzo(a)pyrene	23.556	252	12664	714.501	ng	# 88
91) Dibenzo(a,h)anthracene	26.156	278	5784	278.284	ng	# 80
92) Benzo(g,h,i)perylene	26.897	276	7381	352.011	ng	# 87

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

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