

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112421\
 Data File : BM033255.D
 Acq On : 24 Nov 2021 19:59
 Operator : CG/JU
 Sample : M4812-02MS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SETTLING-TANK-01MS

Quant Time: Nov 26 04:10:12 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 24 15:16:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.937	152	125295	20.000	ng	0.00
21) Naphthalene-d8	10.736	136	473275	20.000	ng	# 0.00
39) Acenaphthene-d10	14.560	164	279146	20.000	ng	0.00
64) Phenanthrene-d10	17.295	188	560902	20.000	ng	# 0.00
76) Chrysene-d12	21.459	240	591174	20.000	ng	0.00
86) Perylene-d12	23.788	264	620366	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.519	112	986733	130.051	ng	0.01
7) Phenol-d6	7.107	99	1225368	120.589	ng	0.01
23) Nitrobenzene-d5	9.101	82	942736	92.473	ng	0.00
42) 2,4,6-Tribromophenol	16.048	330	438097	143.971	ng	0.00
45) 2-Fluorobiphenyl	13.183	172	1850801	93.506	ng	0.00
79) Terphenyl-d14	19.906	244	2657997	88.906	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.437	88	130267	40.271	ng	# 29
3) Pyridine	3.837	79	277116	33.725	ng	# 81
4) n-Nitrosodimethylamine	3.743	42	202631	48.555	ng	# 67
6) Aniline	7.272	93	245127	21.423	ng	97
8) 2-Chlorophenol	7.507	128	406042	51.399	ng	98
9) Benzaldehyde	7.084	77	255207	44.011	ng	94
10) Phenol	7.137	94	485517	47.245	ng	88
11) bis(2-Chloroethyl)ether	7.366	93	410565	52.218	ng	100
12) 1,3-Dichlorobenzene	7.825	146	441036	47.515	ng	95
13) 1,4-Dichlorobenzene	7.972	146	452776	47.683	ng	98
14) 1,2-Dichlorobenzene	8.289	146	432648	47.706	ng	97
15) Benzyl Alcohol	8.184	79	411035	51.977	ng	92
16) 2,2'-oxybis(1-Chloropr...	8.460	45	683818	49.070	ng	94
17) 2-Methylphenol	8.384	107	341629	50.359	ng	94
18) Hexachloroethane	9.019	117	167022	48.445	ng	93
19) n-Nitroso-di-n-propyla...	8.742	70	340921	50.459	ng	93
20) 3+4-Methylphenols	8.707	107	456311	49.659	ng	95
22) Acetophenone	8.766	105	614849	50.479	ng	# 92
24) Nitrobenzene	9.142	77	519440	52.889	ng	97
25) Isophorone	9.666	82	853047	53.505	ng	99
26) 2-Nitrophenol	9.848	139	211164	57.643	ng	# 85
27) 2,4-Dimethylphenol	9.907	122	369191	59.013	ng	99
28) bis(2-Chloroethoxy)met...	10.142	93	521503	50.208	ng	99
29) 2,4-Dichlorophenol	10.383	162	373825	52.750	ng	96
30) 1,2,4-Trichlorobenzene	10.595	180	410993	48.861	ng	98
31) Naphthalene	10.789	128	1258116	50.395	ng	99
32) Benzoic acid	10.007	122	72951	21.237	ng	93
33) 4-Chloroaniline	10.901	127	54150	5.640	ng	97
34) Hexachlorobutadiene	11.060	225	243610	46.851	ng	97
35) Caprolactam	11.695	113	100615	74.173	ng	97
36) 4-Chloro-3-methylphenol	12.019	107	411824	51.116	ng	99
37) 2-Methylnaphthalene	12.389	142	884423	52.036	ng	94
38) 1-Methylnaphthalene	12.607	142	815516	49.181	ng	# 99
40) 1,2,4,5-Tetrachloroben...	12.754	216	430218	51.350	ng	97
41) Hexachlorocyclopentadiene	12.730	237	443493	106.622	ng	97
43) 2,4,6-Trichlorophenol	12.995	196	289093	54.014	ng	94

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.072	196	313854	53.526	ng	# 93
46) 1,1'-Biphenyl	13.395	154	1071455	51.050	ng	99
47) 2-Chloronaphthalene	13.442	162	837837	52.243	ng	98
48) 2-Nitroaniline	13.648	65	281666	54.337	ng	92
49) Acenaphthylene	14.283	152	1343414	53.637	ng	99
50) Dimethylphthalate	14.019	163	1026206	51.726	ng	99
51) 2,6-Dinitrotoluene	14.136	165	220431	55.500	ng	# 75
52) Acenaphthene	14.624	154	791689	51.441	ng	98
53) 3-Nitroaniline	14.466	138	115882	27.380	ng	99
54) 2,4-Dinitrophenol	14.671	184	62864	30.123	ng	# 69
55) Dibenzofuran	14.960	168	1270065	49.628	ng	93
56) 4-Nitrophenol	14.777	139	374544	110.417	ng	# 83
57) 2,4-Dinitrotoluene	14.918	165	300342	57.727	ng	# 71
58) Fluorene	15.607	166	1012513	50.527	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.183	232	294105	57.332	ng	# 82
60) Diethylphthalate	15.377	149	1008512	51.235	ng	96
61) 4-Chlorophenyl-phenyle...	15.595	204	503054	49.229	ng	90
62) 4-Nitroaniline	15.630	138	229561	52.304	ng	# 79
63) Azobenzene	15.889	77	1115111	51.174	ng	90
65) 4,6-Dinitro-2-methylph...	15.677	198	66430	22.866	ng	85
66) n-Nitrosodiphenylamine	15.813	169	879917	53.472	ng	97
67) 4-Bromophenyl-phenylether	16.489	248	302534	51.597	ng	92
68) Hexachlorobenzene	16.601	284	326506	50.733	ng	# 83
69) Atrazine	16.760	200	318248	92.570	ng	97
70) Pentachlorophenol	16.948	266	374141	98.659	ng	98
71) Phenanthrene	17.342	178	1585293	51.418	ng	99
72) Anthracene	17.430	178	1625032	54.635	ng	100
73) Carbazole	17.701	167	1534540	51.860	ng	97
74) Di-n-butylphthalate	18.254	149	1790769	58.260	ng	# 97
75) Fluoranthene	19.342	202	1866895	52.692	ng	97
77) Benzidine	19.530	184	854800	128.604	ng	99
78) Pyrene	19.706	202	1917411	49.271	ng	100
80) Butylbenzylphthalate	20.595	149	861233	59.166	ng	96
81) Benzo(a)anthracene	21.442	228	1897900	50.018	ng	99
82) 3,3'-Dichlorobenzidine	21.371	252	390190	22.767	ng	100
83) Chrysene	21.494	228	1847105	49.974	ng	99
84) Bis(2-ethylhexyl)phtha...	21.359	149	1275269	62.213	ng	95
85) Di-n-octyl phthalate	22.265	149	2254424	64.770	ng	98
87) Indeno(1,2,3-cd)pyrene	26.182	276	2214598	51.994	ng	98
88) Benzo(b)fluoranthene	23.083	252	1985394	52.693	ng	99
89) Benzo(k)fluoranthene	23.130	252	1919698	50.777	ng	98
90) Benzo(a)pyrene	23.688	252	1918827	60.806	ng	99
91) Dibenzo(a,h)anthracene	26.200	278	1920637	53.853	ng	97
92) Benzo(g,h,i)perylene	26.924	276	1896153	53.302	ng	97

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

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