

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080521\
 Data File : BP006513.D
 Acq On : 05 Aug 2021 19:13
 Operator : CG/JU
 Sample : PB138213BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB138213BS

Quant Time: Aug 06 01:10:47 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 03 15:41:30 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.758	152	237611	20.000	ng	0.00
21) Naphthalene-d8	10.540	136	998146	20.000	ng	# 0.00
39) Acenaphthene-d10	14.387	164	554349	20.000	ng	0.00
64) Phenanthrene-d10	17.139	188	1059752	20.000	ng	# 0.00
76) Chrysene-d12	21.239	240	1018989	20.000	ng	# 0.00
86) Perylene-d12	23.563	264	1056962	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	2326915	150.414	ng	0.00
7) Phenol-d6	6.940	99	2950334	134.256	ng	0.00
23) Nitrobenzene-d5	8.916	82	1944310	89.911	ng	0.00
42) 2,4,6-Tribromophenol	15.881	330	851261	146.932	ng	0.00
45) 2-Fluorobiphenyl	13.016	172	3366769	90.865	ng	0.00
79) Terphenyl-d14	19.716	244	4822368	94.829	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.305	88	240416	35.696	ng	90
3) Pyridine	3.693	79	614204	31.028	ng	95
4) n-Nitrosodimethylamine	3.611	42	330497	44.235	ng	# 79
6) Aniline	7.093	93	1091916	45.743	ng	95
8) 2-Chlorophenol	7.322	128	830807	49.635	ng	90
9) Benzaldehyde	6.905	77	422392	31.935	ng	91
10) Phenol	6.969	94	1079487	48.989	ng	99
11) bis(2-Chloroethyl)ether	7.193	93	809049	48.354	ng	91
12) 1,3-Dichlorobenzene	7.646	146	822987	47.176	ng	96
13) 1,4-Dichlorobenzene	7.793	146	842268	47.572	ng	98
14) 1,2-Dichlorobenzene	8.105	146	807859	47.552	ng	97
15) Benzyl Alcohol	8.005	79	804328	48.807	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.287	45	943590	47.391	ng	92
17) 2-Methylphenol	8.205	107	695981	50.020	ng	99
18) Hexachloroethane	8.828	117	336636	48.037	ng	89
19) n-Nitroso-di-n-propyla...	8.569	70	671344	45.535	ng	96
20) 3+4-Methylphenols	8.534	107	942024	49.740	ng	98
22) Acetophenone	8.581	105	1244647	47.021	ng	# 98
24) Nitrobenzene	8.957	77	994672	44.249	ng	92
25) Isophorone	9.487	82	1701454	43.806	ng	95
26) 2-Nitrophenol	9.663	139	416770	50.945	ng	93
27) 2,4-Dimethylphenol	9.728	122	751576	54.863	ng	97
28) bis(2-Chloroethoxy)met...	9.969	93	1026288	49.384	ng	98
29) 2,4-Dichlorophenol	10.193	162	673356	48.550	ng	94
30) 1,2,4-Trichlorobenzene	10.404	180	695994	46.700	ng	96
31) Naphthalene	10.593	128	2445536	45.596	ng	99
32) Benzoic acid	9.881	122	542862	51.224	ng	89
33) 4-Chloroaniline	10.710	127	639719	29.474	ng	95
34) Hexachlorobutadiene	10.875	225	401500	46.231	ng	98
35) Caprolactam	11.510	113	248194	49.628	ng	# 73
36) 4-Chloro-3-methylphenol	11.840	107	835018	47.386	ng	93
37) 2-Methylnaphthalene	12.204	142	1660434	45.686	ng	99
38) 1-Methylnaphthalene	12.422	142	1529892	44.295	ng	98
40) 1,2,4,5-Tetrachloroben...	12.575	216	699608	48.219	ng	# 95
41) Hexachlorocyclopentadiene	12.551	237	999623	129.903	ng	98
43) 2,4,6-Trichlorophenol	12.816	196	489626	49.423	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.887	196	543946	49.603	ng	# 90
46) 1,1'-Biphenyl	13.222	154	1918820	45.731	ng	96
47) 2-Chloronaphthalene	13.263	162	1500728	46.445	ng	95
48) 2-Nitroaniline	13.475	65	575076	50.432	ng	88
49) Acenaphthylene	14.110	152	2485804	47.711	ng	100
50) Dimethylphthalate	13.863	163	1891757	46.881	ng	99
51) 2,6-Dinitrotoluene	13.975	165	419369	46.575	ng	# 83
52) Acenaphthene	14.451	154	1437681	45.340	ng	98
53) 3-Nitroaniline	14.304	138	380792	38.147	ng	84
54) 2,4-Dinitrophenol	14.510	184	472843	100.419	ng	# 85
55) Dibenzofuran	14.787	168	2228434	44.679	ng	95
56) 4-Nitrophenol	14.622	139	883030	101.877	ng	87
57) 2,4-Dinitrotoluene	14.763	165	586872	48.600	ng	# 84
58) Fluorene	15.440	166	1838200	46.110	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.016	232	446686	48.606	ng	# 73
60) Diethylphthalate	15.228	149	2031050	47.922	ng	97
61) 4-Chlorophenyl-phenylether	15.440	204	839676	46.510	ng	90
62) 4-Nitroaniline	15.469	138	513421	48.471	ng	90
63) Azobenzene	15.734	77	2274202	46.024	ng	93
65) 4,6-Dinitro-2-methylph...	15.522	198	286659	44.270	ng	73
66) n-Nitrosodiphenylamine	15.657	169	1587994	47.674	ng	100
67) 4-Bromophenyl-phenylether	16.334	248	494011	48.584	ng	# 87
68) Hexachlorobenzene	16.439	284	556130	48.061	ng	# 88
69) Atrazine	16.616	200	521468	50.169	ng	95
70) Pentachlorophenol	16.787	266	789334	107.514	ng	99
71) Phenanthrene	17.181	178	2810783	47.022	ng	99
72) Anthracene	17.275	178	2870479	49.696	ng	100
73) Carbazole	17.551	167	2729412	46.327	ng	98
74) Di-n-butylphthalate	18.116	149	3561585	53.060	ng	99
75) Fluoranthene	19.157	202	3166587	47.197	ng	94
77) Benzidine	19.345	184	1760195	69.031	ng	99
78) Pyrene	19.510	202	3256867	47.851	ng	99
80) Butylbenzylphthalate	20.398	149	1651022	55.164	ng	# 85
81) Benzo(a)anthracene	21.222	228	3070517	46.108	ng	100
82) 3,3'-Dichlorobenzidine	21.163	252	1085816	62.107	ng	# 98
83) Chrysene	21.275	228	3017700	46.742	ng	99
84) Bis(2-ethylhexyl)phtha...	21.163	149	2490923	55.383	ng	# 97
85) Di-n-octyl phthalate	22.069	149	4240841	57.207	ng	# 94
87) Indeno(1,2,3-cd)pyrene	25.974	276	3735536	48.741	ng	# 89
88) Benzo(b)fluoranthene	22.857	252	3237995	47.136	ng	# 95
89) Benzo(k)fluoranthene	22.904	252	3170000	48.063	ng	# 95
90) Benzo(a)pyrene	23.463	252	3156970	51.237	ng	# 94
91) Dibenzo(a,h)anthracene	25.992	278	3188846	48.123	ng	# 93
92) Benzo(g,h,i)perylene	26.715	276	3140089	49.450	ng	# 89

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

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