

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032522\
 Data File : BP009555.D
 Acq On : 25 Mar 2022 16:44
 Operator : CG/JU
 Sample : N2091-01MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MHBMSD

Quant Time: Mar 26 03:00:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 25 06:34:38 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	235893	20.000	ng	0.00
21) Naphthalene-d8	10.557	136	910665	20.000	ng	0.00
39) Acenaphthene-d10	14.416	164	554670	20.000	ng	0.00
64) Phenanthrene-d10	17.175	188	1119708	20.000	ng	-0.01
76) Chrysene-d12	21.292	240	1157982	20.000	ng	0.00
86) Perylene-d12	23.686	264	1322011	20.000	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	1681339	124.442	ng	0.00
7) Phenol-d6	6.957	99	2148031	117.561	ng	0.00
23) Nitrobenzene-d5	8.922	82	1306945	76.265	ng	-0.01
42) 2,4,6-Tribromophenol	15.916	330	878225	95.949	ng	0.00
45) 2-Fluorobiphenyl	13.034	172	2893009	72.305	ng	0.00
79) Terphenyl-d14	19.757	244	4063199	62.980	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	3.305	88	252473	47.178	ng 98
3) Pyridine	3.699	79	714912	49.602	ng 100
4) n-Nitrosodimethylamine	3.605	42	292320	55.070	ng 99
6) Aniline	7.099	93	980038	47.061	ng 98
8) 2-Chlorophenol	7.340	128	798313	51.816	ng 98
9) Benzaldehyde	6.916	77	397180	46.839	ng 97
10) Phenol	6.987	94	999943	54.550	ng 99
11) bis(2-Chloroethyl)ether	7.199	93	749358	51.654	ng 97
12) 1,3-Dichlorobenzene	7.657	146	866170	49.812	ng 98
13) 1,4-Dichlorobenzene	7.804	146	888066	49.951	ng 100
14) 1,2-Dichlorobenzene	8.116	146	844071	49.120	ng 99
15) Benzyl Alcohol	8.016	79	667336	45.809	ng 99
16) 2,2'-oxybis(1-Chloropr...	8.299	45	844230	53.684	ng 97
17) 2-Methylphenol	8.228	107	642091	50.830	ng 99
18) Hexachloroethane	8.840	117	308804	49.562	ng 96
19) n-Nitroso-di-n-propyla...	8.581	70	564020	48.804	ng 99
20) 3+4-Methylphenols	8.557	107	872235	50.217	ng 99
22) Acetophenone	8.593	105	1119109	49.651	ng # 98
24) Nitrobenzene	8.969	77	840537	51.922	ng 98
25) Isophorone	9.498	82	1547147	52.928	ng 98
26) 2-Nitrophenol	9.681	139	415967	49.296	ng 95
27) 2,4-Dimethylphenol	9.751	122	695882	58.444	ng 100
28) bis(2-Chloroethoxy)met...	9.981	93	949852	49.906	ng 98
29) 2,4-Dichlorophenol	10.222	162	742325	51.111	ng 99
30) 1,2,4-Trichlorobenzene	10.428	180	803893	47.774	ng 99
31) Naphthalene	10.610	128	2344894	49.989	ng 100
32) Benzoic acid	9.934	122	445368	47.842	ng 94
33) 4-Chloroaniline	10.728	127	668249	34.911	ng 98
34) Hexachlorobutadiene	10.898	225	515065	45.175	ng 98
35) Caprolactam	11.522	113	222928	45.480	ng 98
36) 4-Chloro-3-methylphenol	11.869	107	740947	47.476	ng 100
37) 2-Methylnaphthalene	12.228	142	1611620	48.967	ng 99
38) 1-Methylnaphthalene	12.445	142	1549891	48.340	ng 99
40) 1,2,4,5-Tetrachloroben...	12.604	216	904236	49.134	ng 99
41) Hexachlorocyclopentadiene	12.581	237	676138	84.001	ng 100
43) 2,4,6-Trichlorophenol	12.851	196	592741	48.593	ng 99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.928	196	648187	48.933	ng	99
46) 1,1'-Biphenyl	13.245	154	2084361	50.373	ng	99
47) 2-Chloronaphthalene	13.286	162	1632084	50.098	ng	99
48) 2-Nitroaniline	13.492	65	446140	50.281	ng	100
49) Acenaphthylene	14.134	152	2654797	52.789	ng	100
50) Dimethylphthalate	13.881	163	1978208	47.068	ng	100
51) 2,6-Dinitrotoluene	13.992	165	440939	50.073	ng	99
52) Acenaphthene	14.481	154	1490247	49.606	ng	100
53) 3-Nitroaniline	14.328	138	359748	38.678	ng	# 95
54) 2,4-Dinitrophenol	14.545	184	506441	91.842	ng	98
55) Dibenzofuran	14.816	168	2412194	47.795	ng	99
56) 4-Nitrophenol	14.663	139	713108	102.854	ng	98
57) 2,4-Dinitrotoluene	14.792	165	593276	48.998	ng	98
58) Fluorene	15.469	166	1902168	47.922	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.051	232	548913	45.710	ng	94
60) Diethylphthalate	15.251	149	1942920	46.189	ng	99
61) 4-Chlorophenyl-phenyle...	15.463	204	1000686	45.766	ng	97
62) 4-Nitroaniline	15.498	138	446777	45.738	ng	98
63) Azobenzene	15.757	77	1800110	49.194	ng	99
65) 4,6-Dinitro-2-methylph...	15.563	198	334502	46.459	ng	94
66) n-Nitrosodiphenylamine	15.680	169	1676276	50.589	ng	99
67) 4-Bromophenyl-phenylether	16.363	248	635828	46.406	ng	96
68) Hexachlorobenzene	16.486	284	736964	46.735	ng	95
69) Atrazine	16.645	200	611699	52.506	ng	99
70) Pentachlorophenol	16.833	266	837891	99.401	ng	99
71) Phenanthrene	17.222	178	2998394	50.025	ng	100
72) Anthracene	17.310	178	3047327	51.494	ng	99
73) Carbazole	17.586	167	2792672	48.253	ng	99
74) Di-n-butylphthalate	18.151	149	3275796	48.301	ng	100
75) Fluoranthene	19.204	202	3579163	49.324	ng	98
77) Benzidine	19.386	184	1708065	60.178	ng	99
78) Pyrene	19.557	202	3699311	48.480	ng	100
80) Butylbenzylphthalate	20.439	149	1495961	46.125	ng	98
81) Benzo(a)anthracene	21.280	228	3807288	50.044	ng	100
82) 3,3'-Dichlorobenzidine	21.215	252	1354349	46.082	ng	# 98
83) Chrysene	21.327	228	3533788	49.053	ng	99
84) Bis(2-ethylhexyl)phtha...	21.215	149	2173952	48.089	ng	100
85) Di-n-octyl phthalate	22.133	149	3864619	49.596	ng	98
87) Indeno(1,2,3-cd)pyrene	26.180	276	5016597	50.017	ng	97
88) Benzo(b)fluoranthene	22.957	252	4146843	52.545	ng	99
89) Benzo(k)fluoranthene	23.004	252	4068142	52.648	ng	99
90) Benzo(a)pyrene	23.580	252	4093600	60.626	ng	98
91) Dibenzo(a,h)anthracene	26.192	278	4398243	52.616	ng	97
92) Benzo(g,h,i)perylene	26.945	276	4392462	53.415	ng	96

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

