

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM052324\
 Data File : BM045751.D
 Acq On : 23 May 2024 13:27
 Operator : MA/JU
 Sample : PB161074BSD
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB161074BSD

Quant Time: May 24 02:51:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM052224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 23 04:12:56 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.563	152	321762	20.000	ng	0.00
21) Naphthalene-d8	10.334	136	1397357	20.000	ng	0.00
39) Acenaphthene-d10	14.198	164	858350	20.000	ng	0.00
64) Phenanthrene-d10	16.945	188	1852558	20.000	ng	0.00
76) Chrysene-d12	21.162	240	1518618	20.000	ng	0.00
86) Perylene-d12	23.956	264	1742007	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.252	112	1789831	99.175	ng	0.00
7) Phenol-d6	6.840	99	2302908	95.047	ng	0.00
23) Nitrobenzene-d5	8.728	82	2164033	84.243	ng	0.00
42) 2,4,6-Tribromophenol	15.698	330	864769	95.119	ng	0.00
45) 2-Fluorobiphenyl	12.822	172	4523657	84.704	ng	0.00
79) Terphenyl-d14	19.586	244	7048377	85.425	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.193	88	253592	33.854	ng	99
3) Pyridine	3.587	79	664530	32.388	ng	98
4) n-Nitrosodimethylamine	3.499	42	498138	43.035	ng	98
6) Aniline	6.940	93	947385	40.882	ng	99
8) 2-Chlorophenol	7.169	128	992640	50.221	ng	98
9) Benzaldehyde	6.740	77	186472	12.634	ng	98
10) Phenol	6.863	94	1194651	48.788	ng	99
11) bis(2-Chloroethyl)ether	7.022	93	921504	43.422	ng	99
12) 1,3-Dichlorobenzene	7.451	146	987333	41.915	ng	99
13) 1,4-Dichlorobenzene	7.598	146	998367	42.027	ng	99
14) 1,2-Dichlorobenzene	7.916	146	960595	43.278	ng	99
15) Benzyl Alcohol	7.851	79	894076	47.252	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.104	45	1486314	45.125	ng	99
17) 2-Methylphenol	8.075	107	823547	48.518	ng	99
18) Hexachloroethane	8.634	117	381249	41.344	ng	94
19) n-Nitroso-di-n-propyla...	8.398	70	717073	42.095	ng	99
20) 3+4-Methylphenols	8.410	107	1051061	46.501	ng	96
22) Acetophenone	8.398	105	1348296	41.720	ng	# 99
24) Nitrobenzene	8.769	77	1112662	42.588	ng	98
25) Isophorone	9.304	82	2045319	42.773	ng	99
26) 2-Nitrophenol	9.475	139	513089	44.745	ng	99
27) 2,4-Dimethylphenol	9.581	122	939566	64.422	ng	99
28) bis(2-Chloroethoxy)met...	9.781	93	1284216	43.617	ng	100
29) 2,4-Dichlorophenol	10.022	162	957684	48.018	ng	99
30) 1,2,4-Trichlorobenzene	10.192	180	951590	42.135	ng	99
31) Naphthalene	10.381	128	2871875	41.484	ng	100
32) Benzoic acid	9.816	122	557546	41.492	ng	99
33) 4-Chloroaniline	10.551	127	703451	26.619	ng	99
34) Hexachlorobutadiene	10.669	225	557664	40.276	ng	99
35) Caprolactam	11.422	113	309208m	47.936	ng	
36) 4-Chloro-3-methylphenol	11.704	107	1015371	44.771	ng	99
37) 2-Methylnaphthalene	11.998	142	2003617	41.549	ng	99
38) 1-Methylnaphthalene	12.222	142	1891379	39.752	ng	100
40) 1,2,4,5-Tetrachloroben...	12.369	216	980947	44.314	ng	98
41) Hexachlorocyclopentadiene	12.357	237	1401608	151.285	ng	99
43) 2,4,6-Trichlorophenol	12.645	196	744424	49.830	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.733	196	800845	48.166	ng	99
46) 1,1'-Biphenyl	13.033	154	2551281	44.095	ng	100
47) 2-Chloronaphthalene	13.063	162	1973387	43.582	ng	99
48) 2-Nitroaniline	13.316	65	729829	51.081	ng	99
49) Acenaphthylene	13.922	152	3177502	46.279	ng	100
50) Dimethylphthalate	13.698	163	2666128	45.919	ng	100
51) 2,6-Dinitrotoluene	13.804	165	580472	46.581	ng	96
52) Acenaphthene	14.263	154	1928719	44.812	ng	99
53) 3-Nitroaniline	14.157	138	440665	34.671	ng	99
54) 2,4-Dinitrophenol	14.339	184	697717	95.975	ng	99
55) Dibenzofuran	14.604	168	3014297	44.426	ng	99
56) 4-Nitrophenol	14.539	139	1124319	104.047	ng	99
57) 2,4-Dinitrotoluene	14.592	165	781410	47.220	ng	99
58) Fluorene	15.251	166	2462937	44.413	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.851	232	702352	50.836	ng	100
60) Diethylphthalate	15.063	149	2702603	44.547	ng	99
61) 4-Chlorophenyl-phenyle...	15.257	204	1205812	44.432	ng	98
62) 4-Nitroaniline	15.339	138	600001	47.525	ng	97
63) Azobenzene	15.545	77	2631114	44.757	ng	99
65) 4,6-Dinitro-2-methylph...	15.351	198	440860	47.563	ng	95
66) n-Nitrosodiphenylamine	15.480	169	2176147	45.884	ng	99
67) 4-Bromophenyl-phenylether	16.145	248	786081	44.660	ng	99
68) Hexachlorobenzene	16.245	284	898452	42.851	ng	96
69) Atrazine	16.451	200	820817	56.677	ng	99
70) Pentachlorophenol	16.615	266	1174260	91.805	ng	99
71) Phenanthrene	16.986	178	3929465	44.939	ng	99
72) Anthracene	17.074	178	3927245	45.683	ng	100
73) Carbazole	17.368	167	3853219	44.793	ng	100
74) Di-n-butylphthalate	17.933	149	4766157	44.620	ng	100
75) Fluoranthene	19.004	202	4604930	43.191	ng	99
77) Benzidine	19.221	184	1323031	51.019	ng	100
78) Pyrene	19.368	202	4762791	44.459	ng	100
80) Butylbenzylphthalate	20.292	149	2161469	46.580	ng	99
81) Benzo(a)anthracene	21.145	228	4368086	45.160	ng	100
82) 3,3'-Dichlorobenzidine	21.092	252	1342032	40.487	ng	100
83) Chrysene	21.203	228	4010885	44.209	ng	100
84) Bis(2-ethylhexyl)phtha...	21.092	149	2965220	45.509	ng	100
85) Di-n-octyl phthalate	22.150	149	5289499	46.550	ng	100
87) Indeno(1,2,3-cd)pyrene	27.056	276	4874226	48.628	ng	100
88) Benzo(b)fluoranthene	23.080	252	4596252	46.527	ng	100
89) Benzo(k)fluoranthene	23.139	252	4405200	46.981	ng	100
90) Benzo(a)pyrene	23.821	252	4024721	46.886	ng	100
91) Dibenzo(a,h)anthracene	27.115	278	4083509	49.180	ng	99
92) Benzo(g,h,i)perylene	28.026	276	4065411	47.200	ng	100

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

