

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091923\
 Data File : BM041912.D
 Acq On : 19 Sep 2023 20:49
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Sep 20 02:10:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM091823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 19 04:49:34 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	316905	20.000	ng	0.00
21) Naphthalene-d8	10.780	136	1386394	20.000	ng	0.00
39) Acenaphthene-d10	14.610	164	878176	20.000	ng	0.00
64) Phenanthrene-d10	17.362	188	1877265	20.000	ng	0.00
76) Chrysene-d12	21.539	240	1674034	20.000	ng	0.00
86) Perylene-d12	23.956	264	1769492	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	1524646	80.133	ng	0.00
7) Phenol-d6	7.110	99	2150105	81.389	ng	0.00
23) Nitrobenzene-d5	9.157	82	2225881	81.745	ng	0.00
42) 2,4,6-Tribromophenol	16.104	330	856357	78.156	ng	0.00
45) 2-Fluorobiphenyl	13.233	172	4809325	80.382	ng	0.00
79) Terphenyl-d14	19.974	244	7463730	81.203	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.410	88	307263	39.593	ng	100
3) Pyridine	3.816	79	963630	39.971	ng	98
4) n-Nitrosodimethylamine	3.746	42	471080	41.109	ng	99
6) Aniline	7.304	93	1397161	39.919	ng	99
8) 2-Chlorophenol	7.522	128	844428	39.864	ng	98
9) Benzaldehyde	7.128	77	622588	41.686	ng	99
10) Phenol	7.139	94	1097764	40.483	ng	99
11) bis(2-Chloroethyl)ether	7.398	93	905333	40.025	ng	99
12) 1,3-Dichlorobenzene	7.845	146	897781	39.693	ng	99
13) 1,4-Dichlorobenzene	7.998	146	908906	39.548	ng	99
14) 1,2-Dichlorobenzene	8.316	146	872641	39.502	ng	100
15) Benzyl Alcohol	8.216	79	842604	40.401	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.492	45	1390028	40.182	ng	99
17) 2-Methylphenol	8.392	107	793191	40.227	ng	99
18) Hexachloroethane	9.033	117	340941	39.467	ng	98
19) n-Nitroso-di-n-propyla...	8.781	70	803095	40.895	ng	99
20) 3+4-Methylphenols	8.728	107	1092639	40.371	ng	98
22) Acetophenone	8.810	105	1406247	40.575	ng	# 99
24) Nitrobenzene	9.204	77	1097295	40.600	ng	100
25) Isophorone	9.710	82	2151156	40.027	ng	100
26) 2-Nitrophenol	9.898	139	493826	40.567	ng	99
27) 2,4-Dimethylphenol	9.933	122	865145	39.518	ng	100
28) bis(2-Chloroethoxy)met...	10.192	93	1251806	39.853	ng	100
29) 2,4-Dichlorophenol	10.416	162	861192	40.648	ng	98
30) 1,2,4-Trichlorobenzene	10.633	180	867383	39.876	ng	98
31) Naphthalene	10.833	128	2777112	39.789	ng	100
32) Benzoic acid	10.086	122	701725	40.682	ng	100
33) 4-Chloroaniline	10.963	127	1249638	40.002	ng	99
34) Hexachlorobutadiene	11.063	225	521565	39.256	ng	99
35) Caprolactam	11.798	113	302017	39.885	ng	96
36) 4-Chloro-3-methylphenol	12.057	107	943821	39.973	ng	99
37) 2-Methylnaphthalene	12.439	142	2057956	40.102	ng	99
38) 1-Methylnaphthalene	12.657	142	1927537	39.993	ng	100
40) 1,2,4,5-Tetrachloroben...	12.786	216	1010322	39.670	ng	99
41) Hexachlorocyclopentadiene	12.739	237	566232	39.541	ng	99
43) 2,4,6-Trichlorophenol	13.039	196	696631	40.193	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.104	196	818047	40.329	ng	99
46) 1,1'-Biphenyl	13.445	154	2605537	40.213	ng	99
47) 2-Chloronaphthalene	13.498	162	1886822	39.987	ng	99
48) 2-Nitroaniline	13.721	65	696861	41.330	ng	98
49) Acenaphthylene	14.345	152	3150960	40.238	ng	100
50) Dimethylphthalate	14.074	163	2576653	39.688	ng	100
51) 2,6-Dinitrotoluene	14.215	165	553858	40.256	ng	98
52) Acenaphthene	14.680	154	1882767	39.975	ng	100
53) 3-Nitroaniline	14.551	138	627034	40.719	ng	98
54) 2,4-Dinitrophenol	14.751	184	343278	39.950	ng	98
55) Dibenzofuran	15.010	168	3025174	39.790	ng	100
56) 4-Nitrophenol	14.833	139	500788	41.004	ng	98
57) 2,4-Dinitrotoluene	14.998	165	751063	41.261	ng	100
58) Fluorene	15.662	166	2402421	40.081	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.227	232	665350	39.963	ng	99
60) Diethylphthalate	15.421	149	2577026	39.904	ng	99
61) 4-Chlorophenyl-phenyle...	15.651	204	1254890	39.907	ng	98
62) 4-Nitroaniline	15.721	138	609617	41.167	ng	99
63) Azobenzene	15.945	77	2667311	40.740	ng	99
65) 4,6-Dinitro-2-methylph...	15.745	198	451458	41.880	ng	98
66) n-Nitrosodiphenylamine	15.874	169	2134935	40.267	ng	100
67) 4-Bromophenyl-phenylether	16.545	248	744145	38.467	ng	98
68) Hexachlorobenzene	16.633	284	813466	39.715	ng	98
69) Atrazine	16.815	200	632564	38.259	ng	100
70) Pentachlorophenol	16.992	266	624675	40.365	ng	99
71) Phenanthrene	17.409	178	3762195	40.184	ng	100
72) Anthracene	17.498	178	3861555	40.432	ng	100
73) Carbazole	17.780	167	3534724	40.363	ng	99
74) Di-n-butylphthalate	18.303	149	4575362	40.696	ng	100
75) Fluoranthene	19.415	202	4448519	40.472	ng	99
77) Benzidine	19.615	184	1340041	39.217	ng	99
78) Pyrene	19.780	202	4646231	39.888	ng	100
80) Butylbenzylphthalate	20.656	149	2067239	39.903	ng	97
81) Benzo(a)anthracene	21.521	228	4398335	40.285	ng	100
82) 3,3'-Dichlorobenzidine	21.456	252	1564081	40.678	ng	100
83) Chrysene	21.574	228	4115196	40.212	ng	100
84) Bis(2-ethylhexyl)phtha...	21.397	149	3069179	39.934	ng	99
85) Di-n-octyl phthalate	22.327	149	5261564	40.670	ng	100
87) Indeno(1,2,3-cd)pyrene	26.474	276	4243042	40.161	ng	100
88) Benzo(b)fluoranthene	23.209	252	4030945	39.019	ng	100
89) Benzo(k)fluoranthene	23.262	252	4137064	40.525	ng	100
90) Benzo(a)pyrene	23.850	252	3869763	39.920	ng	100
91) Dibenzo(a,h)anthracene	26.485	278	3565388	40.480	ng	99
92) Benzo(g,h,i)perylene	27.250	276	3369444	40.038	ng	99

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

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