

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092221\
 Data File : BF125549.D
 Acq On : 22 Sep 2021 9:34
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Sep 22 10:48:15 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 21 19:21:36 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	252605	20.000	ng	0.00
21) Naphthalene-d8	8.181	136	995283	20.000	ng	# 0.00
39) Acenaphthene-d10	9.939	164	525512	20.000	ng	0.00
64) Phenanthrene-d10	11.422	188	848428	20.000	ng	# 0.00
76) Chrysene-d12	14.063	240	473798	20.000	ng	0.00
86) Perylene-d12	15.539	264	435701	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	1161750	72.728	ng	0.00
7) Phenol-d6	6.522	99	1494546	72.566	ng	0.00
23) Nitrobenzene-d5	7.463	82	1253887	80.354	ng	0.00
42) 2,4,6-Tribromophenol	10.728	330	404070	78.503	ng	0.00
45) 2-Fluorobiphenyl	9.263	172	2352952	80.454	ng	0.00
79) Terphenyl-d14	13.010	244	2154290	73.423	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.693	88	235403	36.551	ng	# 100
3) Pyridine	3.452	79	632857	36.671	ng	# 72
4) n-Nitrosodimethylamine	3.404	42	239035	35.118	ng	# 1
6) Aniline	6.563	93	869629	36.676	ng	94
8) 2-Chlorophenol	6.687	128	673949	36.947	ng	97
9) Benzaldehyde	6.451	77	401326	40.091	ng	92
10) Phenol	6.534	94	760865	36.295	ng	90
11) bis(2-Chloroethyl)ether	6.634	93	610820	36.446	ng	98
12) 1,3-Dichlorobenzene	6.840	146	722137	36.834	ng	98
13) 1,4-Dichlorobenzene	6.916	146	727291	36.701	ng	97
14) 1,2-Dichlorobenzene	7.075	146	679878	36.734	ng	98
15) Benzyl Alcohol	7.034	79	538610	36.544	ng	94
16) 2,2'-oxybis(1-Chloropr...	7.175	45	817417	35.447	ng	89
17) 2-Methylphenol	7.145	107	534755	36.727	ng	95
18) Hexachloroethane	7.416	117	259335	36.754	ng	90
19) n-Nitroso-di-n-propyla...	7.316	70	442847	35.926	ng	# 69
20) 3+4-Methylphenols	7.298	107	654535	36.154	ng	92
22) Acetophenone	7.310	105	850356	36.927	ng	# 89
24) Nitrobenzene	7.481	77	655900	39.666	ng	96
25) Isophorone	7.722	82	1238331	36.876	ng	93
26) 2-Nitrophenol	7.792	139	336709	45.221	ng	88
27) 2,4-Dimethylphenol	7.828	122	565129	37.870	ng	94
28) bis(2-Chloroethoxy)met...	7.928	93	714223	37.019	ng	97
29) 2,4-Dichlorophenol	8.034	162	578860	38.372	ng	97
30) 1,2,4-Trichlorobenzene	8.122	180	594971	37.422	ng	97
31) Naphthalene	8.204	128	1888943	36.955	ng	99
32) Benzoic acid	7.945	122	418228	46.158	ng	98
33) 4-Chloroaniline	8.245	127	781313	37.292	ng	99
34) Hexachlorobutadiene	8.322	225	337722	37.671	ng	98
35) Caprolactam	8.622	113	159308	38.506	ng	# 85
36) 4-Chloro-3-methylphenol	8.722	107	563628	38.034	ng	97
37) 2-Methylnaphthalene	8.892	142	1292777	37.340	ng	99
38) 1-Methylnaphthalene	8.992	142	1223239	37.656	ng	99
40) 1,2,4,5-Tetrachloroben...	9.063	216	527077	37.250	ng	99
41) Hexachlorocyclopentadiene	9.051	237	291594	36.850	ng	99
43) 2,4,6-Trichlorophenol	9.169	196	406344	38.115	ng	98

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44) 2,4,5-Trichlorophenol	9.204	196	433970	38.789	ng	94
46) 1,1'-Biphenyl	9.357	154	1464513	36.987	ng	95
47) 2-Chloronaphthalene	9.386	162	1228422	37.353	ng	96
48) 2-Nitroaniline	9.475	65	308318	43.936	ng	94
49) Acenaphthylene	9.798	152	1813820	37.528	ng	97
50) Dimethylphthalate	9.663	163	1353746	37.430	ng	99
51) 2,6-Dinitrotoluene	9.716	165	309428	43.003	ng	96
52) Acenaphthene	9.975	154	1134681	37.191	ng	95
53) 3-Nitroaniline	9.886	138	322497	42.200	ng	90
54) 2,4-Dinitrophenol	9.986	184	121849	47.004	ng	# 79
55) Dibenzofuran	10.145	168	1669117	36.866	ng	97
56) 4-Nitrophenol	10.033	139	262790	42.390	ng	# 86
57) 2,4-Dinitrotoluene	10.122	165	389453	45.163	ng	# 76
58) Fluorene	10.486	166	1220605	35.053	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.257	232	330103	39.283	ng	# 90
60) Diethylphthalate	10.357	149	1357086	38.072	ng	99
61) 4-Chlorophenyl-phenyle...	10.475	204	571262	36.888	ng	93
62) 4-Nitroaniline	10.498	138	291941	42.273	ng	# 77
63) Azobenzene	10.639	77	1262205	36.444	ng	86
65) 4,6-Dinitro-2-methylph...	10.528	198	187533	45.534	ng	# 73
66) n-Nitrosodiphenylamine	10.598	169	1148887	37.298	ng	95
67) 4-Bromophenyl-phenylether	10.969	248	363767	37.518	ng	97
68) Hexachlorobenzene	11.033	284	407876	37.328	ng	# 89
69) Atrazine	11.122	200	327404	39.525	ng	93
70) Pentachlorophenol	11.222	266	250852	40.266	ng	95
71) Phenanthrene	11.451	178	1767250	37.082	ng	97
72) Anthracene	11.498	178	1768328	37.223	ng	97
73) Carbazole	11.651	167	1647045	36.658	ng	98
74) Di-n-butylphthalate	11.986	149	1966871	36.886	ng	99
75) Fluoranthene	12.633	202	1685814	36.056	ng	96
77) Benzidine	12.751	184	528547	39.343	ng	99
78) Pyrene	12.863	202	1711552	38.548	ng	96
80) Butylbenzylphthalate	13.480	149	700399	39.039	ng	96
81) Benzo(a)anthracene	14.051	228	1204045	37.332	ng	99
82) 3,3'-Dichlorobenzidine	14.010	252	379506	38.404	ng	98
83) Chrysene	14.086	228	1239841	38.055	ng	96
84) Bis(2-ethylhexyl)phtha...	14.045	149	869215	39.631	ng	99
85) Di-n-octyl phthalate	14.657	149	1464639	40.162	ng	# 89
87) Indeno(1,2,3-cd)pyrene	17.033	276	1240676	39.667	ng	98
88) Benzo(b)fluoranthene	15.104	252	1189851	37.340	ng	98
89) Benzo(k)fluoranthene	15.139	252	1055152	36.207	ng	97
90) Benzo(a)pyrene	15.480	252	1033875	37.279	ng	96
91) Dibenzo(a,h)anthracene	17.051	278	1052599	39.708	ng	98
92) Benzo(g,h,i)perylene	17.486	276	1037326	39.649	ng	93

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

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