

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092121\
 Data File : BF125531.D
 Acq On : 21 Sep 2021 11:43
 Operator : JU/CG
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICCC040

Quant Time: Sep 21 13:29:20 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 13:26:09 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	230822	20.000	ng	0.00
21) Naphthalene-d8	8.180	136	912858	20.000	ng	# 0.00
39) Acenaphthene-d10	9.939	164	467669	20.000	ng	0.00
64) Phenanthrene-d10	11.421	188	754868	20.000	ng	# 0.00
76) Chrysene-d12	14.062	240	422384	20.000	ng	0.00
86) Perylene-d12	15.539	264	376383	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	1076330	74.091	ng	0.00
7) Phenol-d6	6.522	99	1384477	71.706	ng	0.00
23) Nitrobenzene-d5	7.463	82	1109054	59.986	ng	0.00
42) 2,4,6-Tribromophenol	10.727	330	348545	61.192	ng	0.00
45) 2-Fluorobiphenyl	9.263	172	2129940	66.830	ng	0.00
79) Terphenyl-d14	13.010	244	1967439	75.530	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.693	88	220513	30.740	ng	# 100
3) Pyridine	3.451	79	595753	32.029	ng	# 75
4) n-Nitrosodimethylamine	3.404	42	233055	25.173	ng	# 14
6) Aniline	6.563	93	814031	36.715	ng	94
8) 2-Chlorophenol	6.686	128	626119	41.103	ng	98
9) Benzaldehyde	6.451	77	370660	31.645	ng	93
10) Phenol	6.534	94	707245	32.326	ng	89
11) bis(2-Chloroethyl)ether	6.634	93	565971	35.899	ng	100
12) 1,3-Dichlorobenzene	6.839	146	672002	38.219	ng	98
13) 1,4-Dichlorobenzene	6.916	146	675602	38.111	ng	98
14) 1,2-Dichlorobenzene	7.075	146	628311	36.976	ng	99
15) Benzyl Alcohol	7.033	79	503784	36.429	ng	95
16) 2,2'-oxybis(1-Chloropr...	7.175	45	784221	27.081	ng	91
17) 2-Methylphenol	7.139	107	496204	38.537	ng	97
18) Hexachloroethane	7.416	117	241855	38.757	ng	93
19) n-Nitroso-di-n-propyla...	7.316	70	407902	35.152	ng	# 71
20) 3+4-Methylphenols	7.298	107	612596	39.657	ng	90
22) Acetophenone	7.310	105	778784	34.618	ng	# 90
24) Nitrobenzene	7.481	77	585994	29.974	ng	94
25) Isophorone	7.722	82	1136903	32.230	ng	92
26) 2-Nitrophenol	7.798	139	277068	31.458	ng	# 93
27) 2,4-Dimethylphenol	7.828	122	505699	39.729	ng	96
28) bis(2-Chloroethoxy)met...	7.928	93	659160	34.510	ng	97
29) 2,4-Dichlorophenol	8.033	162	521631	35.835	ng	98
30) 1,2,4-Trichlorobenzene	8.122	180	546078	35.149	ng	97
31) Naphthalene	8.204	128	1732051	38.094	ng	99
32) Benzoic acid	7.939	122	319680	37.008	ng	98
33) 4-Chloroaniline	8.245	127	717170	38.336	ng	98
34) Hexachlorobutadiene	8.322	225	307200	30.358	ng	99
35) Caprolactam	8.622	113	142839	32.394	ng	# 82
36) 4-Chloro-3-methylphenol	8.722	107	507341	32.837	ng	98
37) 2-Methylnaphthalene	8.898	142	1164320	36.885	ng	97
38) 1-Methylnaphthalene	8.992	142	1097807	37.320	ng	99
40) 1,2,4,5-Tetrachloroben...	9.063	216	473005	31.642	ng	99
41) Hexachlorocyclopentadiene	9.051	237	273455	57.240	ng	98
43) 2,4,6-Trichlorophenol	9.169	196	360510	35.139	ng	97

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44) 2,4,5-Trichlorophenol	9.204	196	390681	35.339	ng	91
46) 1,1'-Biphenyl	9.363	154	1322248	37.706	ng	97
47) 2-Chloronaphthalene	9.386	162	1105484	38.178	ng	97
48) 2-Nitroaniline	9.475	65	255406	25.188	ng	96
49) Acenaphthylene	9.804	152	1621730	37.864	ng	98
50) Dimethylphthalate	9.663	163	1193678	35.797	ng	99
51) 2,6-Dinitrotoluene	9.716	165	260912	34.821	ng	96
52) Acenaphthene	9.974	154	1018410	39.237	ng	96
53) 3-Nitroaniline	9.886	138	273149	33.229	ng	92
54) 2,4-Dinitrophenol	9.986	184	89739	22.771	ng #	86
55) Dibenzofuran	10.145	168	1493037	38.506	ng	97
56) 4-Nitrophenol	10.027	139	223836	39.957	ng	85
57) 2,4-Dinitrotoluene	10.122	165	321761	33.369	ng #	78
58) Fluorene	10.486	166	1085891	35.268	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.257	232	287542	32.555	ng #	90
60) Diethylphthalate	10.357	149	1196469	37.267	ng	99
61) 4-Chlorophenyl-phenyle...	10.480	204	498490	32.185	ng	90
62) 4-Nitroaniline	10.498	138	243692	29.213	ng #	79
63) Azobenzene	10.639	77	1161482	33.942	ng	86
65) 4,6-Dinitro-2-methylph...	10.527	198	148030	29.881	ng #	76
66) n-Nitrosodiphenylamine	10.598	169	1031125	40.384	ng	95
67) 4-Bromophenyl-phenylether	10.969	248	319803	34.921	ng	97
68) Hexachlorobenzene	11.033	284	367237	37.437	ng	90
69) Atrazine	11.121	200	287116	37.589	ng	94
70) Pentachlorophenol	11.221	266	219886	49.486	ng	96
71) Phenanthrene	11.451	178	1562311	38.680	ng	96
72) Anthracene	11.498	178	1572717	38.814	ng	97
73) Carbazole	11.651	167	1484934	39.305	ng	98
74) Di-n-butylphthalate	11.986	149	1767641	40.953	ng	99
75) Fluoranthene	12.633	202	1544811	35.059	ng	96
77) Benzidine	12.751	184	493392	38.059	ng	98
78) Pyrene	12.863	202	1542417	44.210	ng	96
80) Butylbenzylphthalate	13.480	149	626251	44.938	ng	94
81) Benzo(a)anthracene	14.051	228	1073414	35.237	ng	99
82) 3,3'-Dichlorobenzidine	14.009	252	336322	36.277	ng	98
83) Chrysene	14.086	228	1096557	37.192	ng	96
84) Bis(2-ethylhexyl)phtha...	14.045	149	759350	41.149	ng	100
85) Di-n-octyl phthalate	14.657	149	1281934	43.145	ng #	90
87) Indeno(1,2,3-cd)pyrene	17.033	276	1089294	44.620	ng	98
88) Benzo(b)fluoranthene	15.104	252	1041824	36.496	ng	99
89) Benzo(k)fluoranthene	15.139	252	934986	35.593	ng	97
90) Benzo(a)pyrene	15.480	252	906480	37.433	ng	96
91) Dibenzo(a,h)anthracene	17.050	278	921755	44.278	ng	98
92) Benzo(g,h,i)perylene	17.486	276	908649	44.385	ng	93

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

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