

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020222\
 Data File : BF126782.D
 Acq On : 02 Feb 2022 13:02
 Operator : CG/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Quant Time: Feb 02 14:34:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF020222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 02 14:30:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.951	152	133529	20.000	ng	0.00
21) Naphthalene-d8	8.234	136	507159	20.000	ng	0.00
39) Acenaphthene-d10	9.992	164	284658	20.000	ng	0.00
64) Phenanthrene-d10	11.486	188	485202	20.000	ng	0.00
76) Chrysene-d12	14.133	240	329458	20.000	ng	# 0.00
86) Perylene-d12	15.645	264	232584	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.563	112	907912	89.473	ng	0.00
7) Phenol-d6	6.575	99	1287727	91.338	ng	0.00
23) Nitrobenzene-d5	7.522	82	1303310	100.586	ng	0.00
42) 2,4,6-Tribromophenol	10.786	330	287849	108.053	ng	0.00
45) 2-Fluorobiphenyl	9.310	172	1623841	87.365	ng	0.00
79) Terphenyl-d14	13.074	244	1870098	95.206	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.793	88	236776	47.344	ng	# 87
3) Pyridine	3.557	79	660071	47.116	ng	# 83
4) n-Nitrosodimethylamine	3.534	42	207056	41.771	ng	# 69
6) Aniline	6.616	93	856796	48.497	ng	99
8) 2-Chlorophenol	6.740	128	420380	45.167	ng	99
9) Benzaldehyde	6.504	77	372051	33.900	ng	87
10) Phenol	6.587	94	688631	45.924	ng	96
11) bis(2-Chloroethyl)ether	6.687	93	558170	45.871	ng	95
12) 1,3-Dichlorobenzene	6.893	146	477655	46.229	ng	# 95
13) 1,4-Dichlorobenzene	6.969	146	478831	45.892	ng	94
14) 1,2-Dichlorobenzene	7.122	146	446037	45.345	ng	95
15) Benzyl Alcohol	7.093	79	531882	42.316	ng	92
16) 2,2'-oxybis(1-Chloropr...	7.222	45	607046	43.727	ng	81
17) 2-Methylphenol	7.192	107	427654	46.736	ng	# 93
18) Hexachloroethane	7.463	117	189443	44.967	ng	# 92
19) n-Nitroso-di-n-propyla...	7.369	70	426143	41.414	ng	# 83
20) 3+4-Methylphenols	7.351	107	525575	44.297	ng	# 84
22) Acetophenone	7.363	105	677627	43.356	ng	# 88
24) Nitrobenzene	7.540	77	646703	48.347	ng	# 88
25) Isophorone	7.775	82	1133297	45.100	ng	95
26) 2-Nitrophenol	7.851	139	225869	64.732	ng	# 79
27) 2,4-Dimethylphenol	7.881	122	379722	45.857	ng	84
28) bis(2-Chloroethoxy)met...	7.981	93	709402	45.265	ng	# 97
29) 2,4-Dichlorophenol	8.087	162	373307	47.526	ng	96
30) 1,2,4-Trichlorobenzene	8.175	180	395073	46.902	ng	99
31) Naphthalene	8.257	128	1249342	45.600	ng	100
32) Benzoic acid	8.004	122	321690	68.485	ng	# 68
33) 4-Chloroaniline	8.304	127	515797	47.121	ng	# 91
34) Hexachlorobutadiene	8.369	225	246162	45.942	ng	98
35) Caprolactam	8.687	113	143371	51.113	ng	# 76
36) 4-Chloro-3-methylphenol	8.781	107	472065	45.958	ng	89
37) 2-Methylnaphthalene	8.951	142	777336	45.621	ng	97
38) 1-Methylnaphthalene	9.045	142	749532	45.385	ng	99
40) 1,2,4,5-Tetrachloroben...	9.110	216	382586	47.918	ng	96
41) Hexachlorocyclopentadiene	9.098	237	235527	48.451	ng	96
43) 2,4,6-Trichlorophenol	9.222	196	276120	48.512	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.263	196	284118	48.198	ng	# 93
46) 1,1'-Biphenyl	9.416	154	968541	45.163	ng	98
47) 2-Chloronaphthalene	9.439	162	774585	45.973	ng	99
48) 2-Nitroaniline	9.534	65	337948	58.092	ng	92
49) Acenaphthylene	9.857	152	1208302	45.717	ng	99
50) Dimethylphthalate	9.716	163	924575	45.795	ng	# 99
51) 2,6-Dinitrotoluene	9.775	165	196512	56.328	ng	# 76
52) Acenaphthene	10.033	154	770600	47.691	ng	98
53) 3-Nitroaniline	9.951	138	227133	56.943	ng	# 80
54) 2,4-Dinitrophenol	10.051	184	116069	91.314	ng	88
55) Dibenzofuran	10.204	168	1109571	45.695	ng	99
56) 4-Nitrophenol	10.098	139	175388	55.219	ng	# 74
57) 2,4-Dinitrotoluene	10.181	165	266881	63.129	ng	94
58) Fluorene	10.545	166	823772	44.933	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.316	232	247371	51.328	ng	# 83
60) Diethylphthalate	10.416	149	894567	44.574	ng	97
61) 4-Chlorophenyl-phenyle...	10.533	204	417514	45.599	ng	93
62) 4-Nitroaniline	10.569	138	218686	56.491	ng	# 73
63) Azobenzene	10.698	77	1263692	42.548	ng	92
65) 4,6-Dinitro-2-methylph...	10.592	198	159390	77.478	ng	89
66) n-Nitrosodiphenylamine	10.657	169	726587	44.363	ng	99
67) 4-Bromophenyl-phenylether	11.028	248	264119	47.232	ng	96
68) Hexachlorobenzene	11.092	284	288161	48.299	ng	# 87
69) Atrazine	11.180	200	241269	46.125	ng	95
70) Pentachlorophenol	11.280	266	182004	52.461	ng	98
71) Phenanthrene	11.516	178	1238032	44.662	ng	100
72) Anthracene	11.563	178	1261845	45.370	ng	100
73) Carbazole	11.716	167	1180708	45.217	ng	99
74) Di-n-butylphthalate	12.039	149	1379021	43.796	ng	# 97
75) Fluoranthene	12.704	202	1261901	46.088	ng	98
77) Benzidine	12.822	184	328681	28.919	ng	99
78) Pyrene	12.933	202	1281141	48.086	ng	99
80) Butylbenzylphthalate	13.545	149	552110	47.699	ng	# 91
81) Benzo(a)anthracene	14.121	228	1050237	46.900	ng	98
82) 3,3'-Dichlorobenzidine	14.086	252	300967	40.913	ng	# 96
83) Chrysene	14.163	228	991140	45.999	ng	99
84) Bis(2-ethylhexyl)phtha...	14.098	149	699429	44.618	ng	# 99
85) Di-n-octyl phthalate	14.721	149	978590	40.774	ng	94
87) Indeno(1,2,3-cd)pyrene	17.204	276	733149	51.018	ng	# 88
88) Benzo(b)fluoranthene	15.198	252	725399	47.031	ng	# 93
89) Benzo(k)fluoranthene	15.233	252	755542	49.891	ng	# 94
90) Benzo(a)pyrene	15.586	252	599865	47.801	ng	# 94
91) Dibenzo(a,h)anthracene	17.221	278	601263	50.506	ng	# 92
92) Benzo(g,h,i)perylene	17.680	276	608222	52.104	ng	# 88

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

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