

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031022\
 Data File : BF127288.D
 Acq On : 10 Mar 2022 15:04
 Operator : CG\JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Quant Time: Mar 10 15:42:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 10 14:28:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	82297	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	269733	20.000	ng	# 0.00
39) Acenaphthene-d10	9.916	164	162734	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	299257	20.000	ng	# 0.00
76) Chrysene-d12	14.057	240	235011	20.000	ng	# 0.00
86) Perylene-d12	15.539	264	195165	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	472181	84.977	ng	0.00
7) Phenol-d6	6.510	99	638038	85.506	ng	0.00
23) Nitrobenzene-d5	7.440	82	623438	92.452	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	166781	92.136	ng	0.00
45) 2-Fluorobiphenyl	9.234	172	1078225	88.711	ng	0.00
79) Terphenyl-d14	12.992	244	1250454	87.794	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.634	88	132590	47.510	ng	# 82
3) Pyridine	3.405	79	370211	50.611	ng	# 76
4) n-Nitrosodimethylamine	3.381	42	203178	50.270	ng	# 63
6) Aniline	6.540	93	389295	43.998	ng	93
8) 2-Chlorophenol	6.657	128	239904	42.844	ng	89
9) Benzaldehyde	6.428	77	161780	37.632	ng	95
10) Phenol	6.522	94	346783	42.193	ng	74
11) bis(2-Chloroethyl)ether	6.610	93	262977	43.742	ng	91
12) 1,3-Dichlorobenzene	6.810	146	273207	43.533	ng	97
13) 1,4-Dichlorobenzene	6.887	146	273955	43.567	ng	98
14) 1,2-Dichlorobenzene	7.040	146	252374	42.112	ng	99
15) Benzyl Alcohol	7.016	79	250087	45.070	ng	90
16) 2,2'-oxybis(1-Chloropr...	7.140	45	490394	42.887	ng	99
17) 2-Methylphenol	7.128	107	206113	43.495	ng	# 88
18) Hexachloroethane	7.381	117	103897	43.273	ng	# 79
19) n-Nitroso-di-n-propyla...	7.287	70	197482	42.510	ng	# 86
20) 3+4-Methylphenols	7.281	107	248891	41.006	ng	91
22) Acetophenone	7.287	105	336351	44.471	ng	# 88
24) Nitrobenzene	7.457	77	296834	46.642	ng	92
25) Isophorone	7.698	82	514310	47.916	ng	# 98
26) 2-Nitrophenol	7.775	139	124789	47.892	ng	# 79
27) 2,4-Dimethylphenol	7.804	122	186803	47.425	ng	91
28) bis(2-Chloroethoxy)met...	7.904	93	322436	47.134	ng	99
29) 2,4-Dichlorophenol	8.016	162	214948	47.368	ng	98
30) 1,2,4-Trichlorobenzene	8.092	180	249663	48.297	ng	98
31) Naphthalene	8.175	128	682382	45.529	ng	100
32) Benzoic acid	7.951	122	137534	66.987	ng	88
33) 4-Chloroaniline	8.228	127	269558	46.215	ng	# 91
34) Hexachlorobutadiene	8.287	225	167297	48.405	ng	98
35) Caprolactam	8.616	113	64403	48.508	ng	# 68
36) 4-Chloro-3-methylphenol	8.716	107	241734	49.380	ng	98
37) 2-Methylnaphthalene	8.869	142	459912	47.272	ng	96
38) 1-Methylnaphthalene	8.969	142	445495	47.714	ng	99
40) 1,2,4,5-Tetrachloroben...	9.034	216	242256	45.139	ng	# 95
41) Hexachlorocyclopentadiene	9.016	237	99907	68.727	ng	94
43) 2,4,6-Trichlorophenol	9.145	196	160319	48.323	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.192	196	169340	47.892	ng	# 90
46) 1,1'-Biphenyl	9.334	154	569223	44.278	ng	97
47) 2-Chloronaphthalene	9.363	162	450501	45.351	ng	98
48) 2-Nitroaniline	9.463	65	179566	47.662	ng	91
49) Acenaphthylene	9.781	152	686333	44.662	ng	98
50) Dimethylphthalate	9.639	163	565654	46.979	ng	# 99
51) 2,6-Dinitrotoluene	9.704	165	115132	47.545	ng	# 90
52) Acenaphthene	9.951	154	409181	44.676	ng	96
53) 3-Nitroaniline	9.881	138	120764	45.638	ng	98
54) 2,4-Dinitrophenol	9.986	184	62652	64.375	ng	93
55) Dibenzofuran	10.122	168	624109	43.680	ng	97
56) 4-Nitrophenol	10.045	139	83210	53.881	ng	# 81
57) 2,4-Dinitrotoluene	10.110	165	151782	47.445	ng	# 92
58) Fluorene	10.469	166	485466	43.578	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.239	232	148038	51.266	ng	# 78
60) Diethylphthalate	10.334	149	514104	46.431	ng	98
61) 4-Chlorophenyl-phenyle...	10.457	204	269338	45.177	ng	# 87
62) 4-Nitroaniline	10.498	138	121359	46.012	ng	# 77
63) Azobenzene	10.616	77	571671	44.617	ng	88
65) 4,6-Dinitro-2-methylph...	10.522	198	93979	51.671	ng	84
66) n-Nitrosodiphenylamine	10.581	169	422692	43.194	ng	99
67) 4-Bromophenyl-phenylether	10.945	248	166890	44.174	ng	96
68) Hexachlorobenzene	11.010	284	179696	43.700	ng	# 93
69) Atrazine	11.104	200	150347	44.232	ng	95
70) Pentachlorophenol	11.210	266	83171	59.283	ng	98
71) Phenanthrene	11.433	178	722634	43.653	ng	100
72) Anthracene	11.486	178	720824	43.843	ng	99
73) Carbazole	11.639	167	687309	44.728	ng	99
74) Di-n-butylphthalate	11.963	149	813430	45.332	ng	100
75) Fluoranthene	12.622	202	822807	44.507	ng	92
77) Benzidine	12.745	184	251680	42.179	ng	96
78) Pyrene	12.857	202	842430	44.063	ng	99
80) Butylbenzylphthalate	13.463	149	340916	47.258	ng	# 97
81) Benzo(a)anthracene	14.045	228	749499	45.856	ng	99
82) 3,3'-Dichlorobenzidine	14.010	252	221627	42.477	ng	# 95
83) Chrysene	14.086	228	695814	45.730	ng	99
84) Bis(2-ethylhexyl)phtha...	14.016	149	452777	46.441	ng	# 99
85) Di-n-octyl phthalate	14.633	149	742648	50.817	ng	96
87) Indeno(1,2,3-cd)pyrene	17.057	276	541203	45.310	ng	# 83
88) Benzo(b)fluoranthene	15.104	252	628222	47.972	ng	# 93
89) Benzo(k)fluoranthene	15.139	252	558099	44.340	ng	# 92
90) Benzo(a)pyrene	15.480	252	484221	47.068	ng	# 92
91) Dibenzo(a,h)anthracene	17.068	278	448622	44.806	ng	# 90
92) Benzo(g,h,i)perylene	17.515	276	418733	43.020	ng	# 83

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

SSTDICC050

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