

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF021221\
 Data File : BF123197.D
 Acq On : 12 Feb 2021 14:49
 Operator : JU/CG
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

mohammad
 2/15/2021 11:25:53 AM

Quant Time: Feb 12 15:18:46 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF021221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 12 14:20:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	162062	20.00	ng	0.00
21) Naphthalene-d8	8.157	136	606698	20.00	ng	0.00
39) Acenaphthene-d10	9.916	164	301488	20.00	ng	0.00
64) Phenanthrene-d10	11.404	188	547035	20.00	ng	0.00
76) Chrysene-d12	14.057	240	444043	20.00	ng	# 0.00
86) Perylene-d12	15.533	264	417089	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	1127530	107.32	ng	0.00
7) Phenol-d6	6.504	99	1502721	105.52	ng	0.00
23) Nitrobenzene-d5	7.439	82	1298177	107.69	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	358407	109.51	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	2054672	101.31	ng	0.00
79) Terphenyl-d14	12.998	244	2377290	105.19	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.622	88	269835	51.62	ng	98
3) Pyridine	3.375	79	718091	51.37	ng	96
4) n-Nitrosodimethylamine	3.340	42	310316	52.75	ng	94
6) Aniline	6.539	93	923302	52.68	ng	99
8) 2-Chlorophenol	6.657	128	636383	55.88	ng	97
9) Benzaldehyde	6.422	77	388915	59.74	ng	99
10) Phenol	6.516	94	819158	58.00	ng	95
11) bis(2-Chloroethyl)ether	6.610	93	581600	49.48	ng	99
12) 1,3-Dichlorobenzene	6.816	146	660683	49.75	ng	97
13) 1,4-Dichlorobenzene	6.892	146	672439	48.62	ng	98
14) 1,2-Dichlorobenzene	7.045	146	624262	48.24	ng	97
15) Benzyl Alcohol	7.016	79	562959	56.39	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.151	45	738965	40.77	ng	98
17) 2-Methylphenol	7.122	107	520679	53.73	ng	99
18) Hexachloroethane	7.386	117	249526	51.51	ng	91
19) n-Nitroso-di-n-propyla...	7.292	70	430852	50.31	ng	99
20) 3+4-Methylphenols	7.281	107	633355	55.49	ng	94
22) Acetophenone	7.286	105	783287	49.55	ng	96
24) Nitrobenzene	7.457	77	685316	55.18	ng	98
25) Isophorone	7.698	82	1199364	54.31	ng	99
26) 2-Nitrophenol	7.775	139	343568	66.89	ng	93
27) 2,4-Dimethylphenol	7.810	122	497663	53.70	ng	99
28) bis(2-Chloroethoxy)met...	7.904	93	644421	42.77	ng	99
29) 2,4-Dichlorophenol	8.016	162	496147	51.19	ng	96
30) 1,2,4-Trichlorobenzene	8.098	180	527836	48.34	ng	98
31) Naphthalene	8.180	128	1755125	52.74	ng	99
32) Benzoic acid	7.945	122	442291	65.06	ng	99
33) 4-Chloroaniline	8.228	127	750067	50.79	ng	98
34) Hexachlorobutadiene	8.298	225	285778	45.31	ng	100
35) Caprolactam	8.610	113	177720	54.26	ng	99
36) 4-Chloro-3-methylphenol	8.710	107	576636	57.22	ng	97
37) 2-Methylnaphthalene	8.875	142	1173023	53.10	ng	99
38) 1-Methylnaphthalene	8.975	142	1097351	52.46	ng	100
40) 1,2,4,5-Tetrachloroben...	9.039	216	445820	47.53	ng	97
41) Hexachlorocyclopentadiene	9.027	237	232723	52.27	ng	97
43) 2,4,6-Trichlorophenol	9.151	196	346631	53.91	ng	98

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44) 2,4,5-Trichlorophenol	9.186	196	379311	56.41	ng	#	94
46) 1,1'-Biphenyl	9.339	154	1272162	51.44	ng		99
47) 2-Chloronaphthalene	9.363	162	1034109	49.95	ng		97
48) 2-Nitroaniline	9.457	65	361074	57.55	ng		88
49) Acenaphthylene	9.780	152	1650152	54.46	ng		99
50) Dimethylphthalate	9.639	163	1163432	49.19	ng		99
51) 2,6-Dinitrotoluene	9.698	165	290280	57.61	ng	#	80
52) Acenaphthene	9.951	154	1104464	56.75	ng		99
53) 3-Nitroaniline	9.869	138	350934	58.90	ng		92
54) 2,4-Dinitrophenol	9.974	184	171664	90.36	ng		95
55) Dibenzofuran	10.127	168	1439568	50.83	ng		97
56) 4-Nitrophenol	10.027	139	296300	68.80	ng		92
57) 2,4-Dinitrotoluene	10.104	165	395559	60.37	ng		92
58) Fluorene	10.469	166	1114317	49.25	ng		100
59) 2,3,4,6-Tetrachlorophenol	10.239	232	300064	52.62	ng	#	85
60) Diethylphthalate	10.339	149	1175352	50.53	ng		98
61) 4-Chlorophenyl-phenyle...	10.457	204	505053	47.51	ng		91
62) 4-Nitroaniline	10.486	138	348752	58.26	ng		99
63) Azobenzene	10.621	77	1201965	52.81	ng		98
65) 4,6-Dinitro-2-methylph...	10.516	198	225365	81.98	ng		87
66) n-Nitrosodiphenylamine	10.580	169	1015288	51.49	ng		99
67) 4-Bromophenyl-phenylether	10.951	248	325425	46.71	ng		97
68) Hexachlorobenzene	11.021	284	353979	47.37	ng		97
69) Atrazine	11.104	200	298073	54.74	ng		98
70) Pentachlorophenol	11.210	266	235919	58.25	ng		98
71) Phenanthrene	11.433	178	1710940	50.36	ng		100
72) Anthracene	11.486	178	1715724	52.64	ng		100
73) Carbazole	11.639	167	1639289	54.20	ng		99
74) Di-n-butylphthalate	11.968	149	1774009	49.46	ng		100
75) Fluoranthene	12.627	202	1755955	51.71	ng		99
78) Pyrene	12.857	202	1800846	53.72	ng		100
80) Butylbenzylphthalate	13.474	149	810403	56.18	ng		97
81) Benzo(a)anthracene	14.045	228	1643819	54.70	ng		98
82) 3,3'-Dichlorobenzidine	14.009	252	572894	60.57	ng	#	97
83) Chrysene	14.086	228	1629184	54.17	ng		99
84) Bis(2-ethylhexyl)phtha...	14.033	149	1033896	53.94	ng		100
85) Di-n-octyl phthalate	14.651	149	1796015	55.80	ng		100
87) Indeno(1,2,3-cd)pyrene	17.027	276	1478646	53.24	ng	#	94
88) Benzo(b)fluoranthene	15.104	252	1721094m	57.17	ng		
89) Benzo(k)fluoranthene	15.139	252	1435693	51.32	ng		100
90) Benzo(a)pyrene	15.474	252	1456470	55.09	ng	#	98
91) Dibenzo(a,h)anthracene	17.050	278	1253277	54.82	ng		97
92) Benzo(g,h,i)perylene	17.480	276	1209620	54.60	ng		95

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

