

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011619\
 Data File : BF112070.D
 Acq On : 16 Jan 2019 16:41
 Operator : JU/SJ
 Sample : SSTDICC050
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

Sohil
 1/17/2019 2:07:56 PM

Quant Time: Jan 16 17:29:06 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 16 16:57:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	295210	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	1098707	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	534038	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	1021688	20.00	ng	0.00
76) Chrysene-d12	14.03	240	823343	20.00	ng	0.00
87) Perylene-d12	15.50	264	716676	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.49	112	1557397	90.26	ng	0.00
7) Phenol-d6	6.51	99	1920721	86.36	ng	0.00
23) Nitrobenzene-d5	7.43	82	1693537	82.38	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	604951	87.09	ng	0.00
45) 2-Fluorobiphenyl	9.23	172	3253672	88.84	ng	0.00
79) Terphenyl-d14	12.97	244	3568099	82.26	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.63	88	369882	49.023	ng	99
3) Pyridine	3.39	79	978635	50.042	ng	98
4) n-Nitrosodimethylamine	3.35	42	386389	39.711	ng	100
6) Aniline	6.53	93	1169644	43.495	ng	# 65
8) 2-Chlorophenol	6.66	128	922690	48.156	ng	98
9) Benzaldehyde	6.41	77	486512	34.973	ng	99
10) Phenol	6.53	94	1063908	44.013	ng	99
11) bis(2-Chloroethyl)ether	6.60	93	874073	46.741	ng	98
12) 1,3-Dichlorobenzene	6.80	146	1039653	46.805	ng	99
13) 1,4-Dichlorobenzene	6.88	146	1059203	46.991	ng	99
14) 1,2-Dichlorobenzene	7.04	146	970076	45.023	ng	98
15) Benzyl Alcohol	7.01	79	743043	42.226	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.14	45	1222842	38.353	ng	95
17) 2-Methylphenol	7.13	107	690276	44.825	ng	97
18) Hexachloroethane	7.37	117	365916	46.495	ng	99
19) n-Nitroso-di-n-propylamine	7.28	70	604172	41.810	ng	99
20) 3+4-Methylphenols	7.28	107	831108	43.365	ng	# 82
22) Acetophenone	7.27	105	1203293	43.259	ng	98
24) Nitrobenzene	7.45	77	906491	43.590	ng	94
25) Isophorone	7.69	82	1499586	45.047	ng	98
26) 2-Nitrophenol	7.76	139	432491	42.238	ng	98
27) 2,4-Dimethylphenol	7.80	122	705755	47.670	ng	100
28) bis(2-Chloroethoxy)methane	7.89	93	1030242	46.365	ng	100
29) 2,4-Dichlorophenol	8.01	162	785696	46.152	ng	99
30) 1,2,4-Trichlorobenzene	8.09	180	887594	47.261	ng	99
31) Naphthalene	8.17	128	2386687	45.760	ng	99
32) Benzoic acid	7.94	122	348208m	36.243	ng	
33) 4-Chloroaniline	8.22	127	1072374	47.563	ng	98
34) Hexachlorobutadiene	8.29	225	493595	42.217	ng	98
35) Caprolactam	8.60	113	268858	48.337	ng	90
36) 4-Chloro-3-methylphenol	8.71	107	748294	44.208	ng	98
37) 2-Methylnaphthalene	8.86	142	1629557	50.422	ng	99
38) 1-Methylnaphthalene	8.96	142	1568675	50.606	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.03	216	852368	50.065	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.02	237	324628	32.977	ng	99
43) 2,4,6-Trichlorophenol	9.15	196	547812	47.439	ng	99
44) 2,4,5-Trichlorophenol	9.19	196	571622	47.246	ng	97
46) 1,1'-Biphenyl	9.33	154	2124435	47.557	ng	98
47) 2-Chloronaphthalene	9.35	162	1711516	48.817	ng	98
48) 2-Nitroaniline	9.45	65	427232	39.858	ng	92
49) Acenaphthylene	9.77	152	2463702	47.531	ng	100
50) Dimethylphthalate	9.63	163	1915155	47.739	ng	100
51) 2,6-Dinitrotoluene	9.69	165	419721	46.786	ng	88
52) Acenaphthene	9.94	154	1493281	44.688	ng	100
53) 3-Nitroaniline	9.86	138	462243	48.115	ng	90
54) 2,4-Dinitrophenol	9.96	184	97478	25.661	ng	91
55) Dibenzofuran	10.11	168	2287795	47.077	ng	97
56) 4-Nitrophenol	10.03	139	299061	44.056	ng	95
57) 2,4-Dinitrotoluene	10.09	165	510362	44.027	ng	90
58) Fluorene	10.45	166	1671946	47.282	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.23	232	439781	43.142	ng	98
60) Diethylphthalate	10.32	149	1844018	47.498	ng	98
61) 4-Chlorophenyl-phenylether	10.44	204	920505	46.421	ng	96
62) 4-Nitroaniline	10.47	138	454490	49.876	ng	97
63) Azobenzene	10.60	77	1780471	48.007	ng	97
65) 4,6-Dinitro-2-methylphenol	10.51	198	213938	33.936	ng	94
66) n-Nitrosodiphenylamine	10.56	169	1636756	47.739	ng	99
67) 4-Bromophenyl-phenylether	10.93	248	590217	44.601	ng	93
68) Hexachlorobenzene	11.01	284	630245	42.974	ng	94
69) Atrazine	11.09	200	531169	44.095	ng	98
70) Pentachlorophenol	11.20	266	295914	35.458	ng	99
71) Phenanthrene	11.42	178	2442483	44.714	ng	98
72) Anthracene	11.47	178	2458717	44.340	ng	98
73) Carbazole	11.62	167	2468214	45.760	ng	98
74) Di-n-butylphthalate	11.94	149	2798772	44.671	ng	99
75) Fluoranthene	12.60	202	2555764	45.576	ng	98
77) Benzidine	12.72	184	1121724	45.595	ng	98
78) Pyrene	12.83	202	2638505	46.575	ng	99
80) Butylbenzylphthalate	13.44	149	1213044	50.648	ng	97
81) Benzo(a)anthracene	14.02	228	2281371	48.146	ng	99
82) 3,3'-Dichlorobenzidine	13.98	252	845355	48.324	ng	98
83) Chrysene	14.06	228	2166029	47.942	ng	99
84) Bis(2-ethylhexyl)phthalate	14.00	149	1697413	53.966	ng	99
85) Di-n-octyl phthalate	14.62	149	2613093	58.052	ng	99
86) Indeno(1,2,3-cd)pyrene	17.00	276	1934072	58.719	ng	100
88) Benzo(b)fluoranthene	15.07	252	2085510	48.646	ng	98
89) Benzo(k)fluoranthene	15.11	252	1890702	46.918	ng	99
90) Benzo(a)pyrene	15.44	252	1855538	49.225	ng	98
91) Dibenzo(a,h)anthracene	17.01	278	1603368	52.885	ng	99
92) Benzo(g,h,i)perylene	17.44	276	1595286	53.797	ng	99

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

