

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121718\
 Data File : BF111541.D
 Acq On : 17 Dec 2018 16:11
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 12/18/2018 5:35:42 PM

Quant Time: Dec 18 13:03:03 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	162819	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	802676	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	391912	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	842974	20.00	ng	0.00
76) Chrysene-d12	13.96	240	622783	20.00	ng	0.00
87) Perylene-d12	15.39	264	508224	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	790260	80.77	ng	0.00
7) Phenol-d6	6.45	99	1004625	77.19	ng	0.00
23) Nitrobenzene-d5	7.36	82	1132604	84.03	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	272932	77.74	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1964447	77.42	ng	0.00
79) Terphenyl-d14	12.90	244	2011752	75.85	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.52	88	206167	43.627	ng	98
3) Pyridine	3.25	79	533202	44.925	ng	94
4) n-Nitrosodimethylamine	3.19	42	247431	45.897	ng	91
6) Aniline	6.46	93	619800	38.391	ng	# 27
8) 2-Chlorophenol	6.59	128	473684	40.517	ng	95
9) Benzaldehyde	6.34	77	276791	35.063	ng	96
10) Phenol	6.46	94	559219	38.571	ng	# 68
11) bis(2-Chloroethyl)ether	6.53	93	442737	38.954	ng	99
12) 1,3-Dichlorobenzene	6.73	146	524419	40.750	ng	98
13) 1,4-Dichlorobenzene	6.81	146	527745	40.402	ng	98
14) 1,2-Dichlorobenzene	6.96	146	498521	40.567	ng	98
15) Benzyl Alcohol	6.94	79	401380	40.531	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.07	45	887188	39.979	ng	88
17) 2-Methylphenol	7.06	107	377965	40.172	ng	# 92
18) Hexachloroethane	7.30	117	225348	46.356	ng	96
19) n-Nitroso-di-n-propylamine	7.21	70	375533	43.735	ng	94
20) 3+4-Methylphenols	7.21	107	537114	45.510	ng	# 81
22) Acetophenone	7.20	105	741969	41.408	ng	# 90
24) Nitrobenzene	7.37	77	569780	41.437	ng	98
25) Isophorone	7.62	82	903805	39.628	ng	99
26) 2-Nitrophenol	7.69	139	306255	42.947	ng	94
27) 2,4-Dimethylphenol	7.74	122	421866	40.808	ng	99
28) bis(2-Chloroethoxy)methane	7.82	93	612905	38.927	ng	97
29) 2,4-Dichlorophenol	7.94	162	452357	40.990	ng	98
30) 1,2,4-Trichlorobenzene	8.02	180	464269	40.187	ng	100
31) Naphthalene	8.10	128	1473102	39.238	ng	96
32) Benzoic acid	7.87	122	348579m	39.019	ng	
33) 4-Chloroaniline	8.15	127	649191	38.887	ng	97
34) Hexachlorobutadiene	8.22	225	260276	40.893	ng	98
35) Caprolactam	8.53	113	158626m	38.715	ng	
36) 4-Chloro-3-methylphenol	8.64	107	463568	38.162	ng	96
37) 2-Methylnaphthalene	8.79	142	955467	37.610	ng	99
38) 1-Methylnaphthalene	8.89	142	914032	37.242	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.96	216	418752	38.962	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.95	237	206190	37.366	nq	99
43) 2,4,6-Trichlorophenol	9.07	196	300298	39.477	nq	99
44) 2,4,5-Trichlorophenol	9.12	196	305252	37.710	nq	98
46) 1,1'-Biphenyl	9.25	154	1266897	38.151	nq	95
47) 2-Chloronaphthalene	9.28	162	983278	38.955	nq	95
48) 2-Nitroaniline	9.37	65	308461	37.637	nq	89
49) Acenaphthylene	9.69	152	1421842	37.949	nq	95
50) Dimethylphthalate	9.56	163	1069027	37.876	nq	98
51) 2,6-Dinitrotoluene	9.62	165	246453	38.792	nq	96
52) Acenaphthene	9.86	154	941965	39.777	nq	98
53) 3-Nitroaniline	9.79	138	295161	38.250	nq	91
54) 2,4-Dinitrophenol	9.89	184	100185	41.449	nq	# 79
55) Dibenzofuran	10.04	168	1317201	38.320	nq	95
56) 4-Nitrophenol	9.96	139	201912	37.809	nq	94
57) 2,4-Dinitrotoluene	10.02	165	321325	38.535	nq	98
58) Fluorene	10.38	166	958836	38.457	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.16	232	242700	38.375	nq	99
60) Diethylphthalate	10.25	149	1072918	37.528	nq	97
61) 4-Chlorophenyl-phenylether	10.37	204	469854	38.518	nq	97
62) 4-Nitroaniline	10.40	138	299662	39.228	nq	100
63) Azobenzene	10.53	77	1025045	36.367	nq	97
65) 4,6-Dinitro-2-methylphenol	10.43	198	152856	32.119	nq	# 75
66) n-Nitrosodiphenylamine	10.49	169	965987m	33.230	nq	
67) 4-Bromophenyl-phenylether	10.86	248	311200m	34.252	nq	
68) Hexachlorobenzene	10.93	284	303485	32.473	nq	96
69) Atrazine	11.02	200	304812m	36.282	nq	
70) Pentachlorophenol	11.13	266	209511	36.549	nq	99
71) Phenanthrene	11.34	178	1675637	38.576	nq	94
72) Anthracene	11.39	178	1713415	38.571	nq	96
73) Carbazole	11.55	167	1786256	38.885	nq	96
74) Di-n-butylphthalate	11.87	149	2043278	39.939	nq	96
75) Fluoranthene	12.53	202	1673755	39.388	nq	98
77) Benzidine	12.64	184	871806	38.081	nq	98
78) Pyrene	12.76	202	1722751	39.235	nq	97
80) Butylbenzylphthalate	13.37	149	855553	39.975	nq	89
81) Benzo(a)anthracene	13.94	228	1307151	38.601	nq	98
82) 3,3'-Dichlorobenzidine	13.90	252	551768	40.081	nq	97
83) Chrysene	13.98	228	1361958	38.179	nq	98
84) Bis(2-ethylhexyl)phthalate	13.93	149	1095197	40.064	nq	99
85) Di-n-octyl phthalate	14.55	149	1863142	40.407	nq	97
86) Indeno(1,2,3-cd)pyrene	16.82	276	939359	36.485	nq	97
88) Benzo(b)fluoranthene	14.98	252	1148885	38.774	nq	99
89) Benzo(k)fluoranthene	15.01	252	1141733	40.327	nq	99
90) Benzo(a)pyrene	15.33	252	1051728	39.362	nq	98
91) Dibenzo(a,h)anthracene	16.83	278	788643	37.417	nq	96
92) Benzo(g,h,i)perylene	17.24	276	784739	37.930	nq	95

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

