

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072018\
 Data File : BF107524.D
 Acq On : 20 Jul 2018 15:03
 Operator : JU/SJ
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

Sohil
 7/23/2018 12:44:26 PM

Quant Time: Jul 20 15:58:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 20 14:36:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	339120	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	1368438	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	606198	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	1172491	20.00	ng	0.00
75) Chrysene-d12	14.17	240	858469	20.00	ng	0.00
86) Perylene-d12	15.69	264	724401	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	2367928	106.05	ng	0.00
7) Phenol-d6	6.62	99	2763395	95.72	ng	0.00
23) Nitrobenzene-d5	7.55	82	2429620	94.68	ng	0.00
41) 2,4,6-Tribromophenol	10.82	330	803195	150.80	ng	0.00
44) 2-Fluorobiphenyl	9.34	172	3706062	87.85	ng	0.00
78) Terphenyl-d14	13.10	244	3493244	101.10	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.86	88	578620	52.489	ng	# 82
3) Pyridine	3.64	79	1596003	53.678	ng	# 85
4) n-Nitrosodimethylamine	3.62	42	648412	53.109	ng	95
6) Aniline	6.65	93	1832533	48.019	ng	# 88
8) 2-Chlorophenol	6.77	128	1250463	55.376	ng	94
9) Benzaldehyde	6.54	77	789859	34.011	ng	98
10) Phenol	6.64	94	1619545	52.328	ng	80
11) bis(2-Chloroethyl)ether	6.72	93	1365616	54.068	ng	92
12) 1,3-Dichlorobenzene	6.92	146	1419088	51.762	ng	98
13) 1,4-Dichlorobenzene	7.00	146	1449244	53.327	ng	97
14) 1,2-Dichlorobenzene	7.15	146	1254409	49.331	ng	98
15) Benzyl Alcohol	7.12	79	951586	32.822	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.25	45	2197430	81.531	ng	87
17) 2-Methylphenol	7.23	107	1075210	50.079	ng	96
18) Hexachloroethane	7.49	117	530750	52.487	ng	97
19) n-Nitroso-di-n-propylamine	7.40	70	876595	42.409	ng	99
20) 3+4-Methylphenols	7.39	107	1175756	43.110	ng	# 72
22) Acetophenone	7.40	105	1643538	43.713	ng	# 90
24) Nitrobenzene	7.57	77	1347296	46.717	ng	97
25) Isophorone	7.81	82	2410867	47.396	ng	98
26) 2-Nitrophenol	7.88	139	668422	68.182	ng	95
27) 2,4-Dimethylphenol	7.91	122	1046606	51.196	ng	100
28) bis(2-Chloroethoxy)methane	8.01	93	1560970	49.000	ng	98
29) 2,4-Dichlorophenol	8.12	162	1086122	51.054	ng	98
30) 1,2,4-Trichlorobenzene	8.20	180	1148111	44.567	ng	99
31) Naphthalene	8.28	128	3038984	45.720	ng	# 93
32) Benzoic acid	8.07	122	814485	64.918	ng	96
33) 4-Chloroaniline	8.34	127	1324852	46.023	ng	96
34) Hexachlorobutadiene	8.39	225	697727	40.477	ng	99
35) Caprolactam	8.73	113	350747m	50.685	ng	
36) 4-Chloro-3-methylphenol	8.82	107	1179897	42.447	ng	98
37) 2-Methylnaphthalene	8.98	142	2135347	48.737	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.14	216	1093247	46.344	ng	98
40) Hexachlorocyclopentadiene	9.12	237	654401	53.113	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.25	196	753822	51.192	nq	99
43) 2,4,5-Trichlorophenol	9.30	196	777710	55.208	nq	98
45) 1,1'-Biphenyl	9.44	154	2689929	51.323	nq	93
46) 2-Chloronaphthalene	9.47	162	2115266	51.099	nq	95
47) 2-Nitroaniline	9.57	65	764713	60.365	nq	89
48) Acenaphthylene	9.89	152	3036775	52.567	nq	# 90
49) Dimethylphthalate	9.74	163	2461907	51.115	nq	97
50) 2,6-Dinitrotoluene	9.81	165	557905	60.702	nq	96
51) Acenaphthene	10.06	154	1929506	50.383	nq	96
52) 3-Nitroaniline	9.99	138	690693	71.547	nq	98
53) 2,4-Dinitrophenol	10.09	184	228541	70.580	nq	# 20
54) Dibenzofuran	10.23	168	2795033	47.229	nq	89
55) 4-Nitrophenol	10.15	139	526395	67.386	nq	96
56) 2,4-Dinitrotoluene	10.22	165	686105	57.272	nq	95
57) Fluorene	10.58	166	2083301	53.152	nq	95
58) 2,3,4,6-Tetrachlorophenol	10.35	232	661005	48.441	nq	91
59) Diethylphthalate	10.44	149	2474724	52.887	nq	# 92
60) 4-Chlorophenyl-phenylether	10.56	204	1145787	49.077	nq	98
61) 4-Nitroaniline	10.61	138	558913	60.836	nq	98
62) Azobenzene	10.72	77	2333233	42.461	nq	87
64) 4,6-Dinitro-2-methylphenol	10.63	198	360700	64.579	nq	83
65) n-Nitrosodiphenylamine	10.69	169	2074756	56.027	nq	96
66) 4-Bromophenyl-phenylether	11.05	248	756475	56.393	nq	91
67) Hexachlorobenzene	11.12	284	831486	66.962	nq	# 85
68) Atrazine	11.21	200	669226	47.823	nq	95
69) Pentachlorophenol	11.32	266	580097	63.217	nq	99
70) Phenanthrene	11.54	178	2858123m	48.914	nq	
71) Anthracene	11.59	178	2862467	49.184	nq	# 91
72) Carbazole	11.75	167	2971776	53.936	nq	# 92
73) Di-n-butylphthalate	12.07	149	3312667	52.315	nq	# 93
74) Fluoranthene	12.74	202	3022499	45.510	nq	94
76) Benzidine	12.85	184	1260334	37.829	nq	96
77) Pyrene	12.97	202	3089282	52.919	nq	92
79) Butylbenzylphthalate	13.57	149	1575687	74.100	nq	# 93
80) Benzo(a)anthracene	14.15	228	2707313	51.542	nq	94
81) 3,3'-Dichlorobenzidine	14.12	252	789579	44.821	nq	# 96
82) Chrysene	14.19	228	2539452	50.611	nq	92
83) Bis(2-ethylhexyl)phthalate	14.12	149	1947858	63.923	nq	# 95
84) Di-n-octyl phthalate	14.74	149	3007417	63.489	nq	# 97
85) Indeno(1,2,3-cd)pyrene	17.28	276	2233463	62.535	nq	96
87) Benzo(b)fluoranthene	15.24	252	2323300	54.794	nq	97
88) Benzo(k)fluoranthene	15.28	252	2095381	50.639	nq	97
89) Benzo(a)pyrene	15.64	252	2074684	53.311	nq	98
90) Dibenzo(a,h)anthracene	17.30	278	1852838	63.084	nq	96
91) Benzo(g,h,i)perylene	17.77	276	1866829	62.552	nq	95

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

