

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061218\
 Data File : BF106364.D
 Acq On : 13 Jun 2018 00:31
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 6/13/2018 4:56:27 PM

Quant Time: Jun 13 04:17:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 11 15:00:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	102482	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	359033	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	144715	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	309161	20.00	ng	0.00
75) Chrysene-d12	14.16	240	314236	20.00	ng	0.00
86) Perylene-d12	15.69	264	230256	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	575344	95.02	ng	0.04
7) Phenol-d6	6.62	99	646913	84.56	ng	0.04
23) Nitrobenzene-d5	7.54	82	553693	90.72	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	135192	97.09	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	859646	119.21	ng	0.00
78) Terphenyl-d14	13.09	244	1129939	89.31	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.80	88	144999	46.271	ng	96
3) Pyridine	3.59	79	377212	43.195	ng	99
4) n-Nitrosodimethylamine	3.55	42	166646	41.656	ng	94
6) Aniline	6.64	93	410589	36.756	ng	# 59
8) 2-Chlorophenol	6.78	128	277087	42.554	ng	98
9) Benzaldehyde	6.53	77	231309	39.863	ng	99
10) Phenol	6.64	94	367495	44.004	ng	# 33
11) bis(2-Chloroethyl)ether	6.71	93	300183	41.987	ng	97
12) 1,3-Dichlorobenzene	6.92	146	326063	42.546	ng	99
13) 1,4-Dichlorobenzene	7.00	146	332536	42.759	ng	98
14) 1,2-Dichlorobenzene	7.15	146	301941	41.243	ng	98
15) Benzyl Alcohol	7.12	79	243828	37.218	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.24	45	315579	30.793	ng	# 17
17) 2-Methylphenol	7.25	107	235534	40.543	ng	# 91
18) Hexachloroethane	7.49	117	89373	32.252	ng	92
19) n-Nitroso-di-n-propylamine	7.38	70	210843	36.737	ng	98
20) 3+4-Methylphenols	7.40	107	307297	41.362	ng	93
22) Acetophenone	7.38	105	399165	44.117	ng	# 98
24) Nitrobenzene	7.55	77	286263	43.155	ng	93
25) Isophorone	7.80	82	467449	39.213	ng	98
26) 2-Nitrophenol	7.87	139	123142	47.827	ng	97
27) 2,4-Dimethylphenol	7.92	122	203146	40.254	ng	98
28) bis(2-Chloroethoxy)methane	8.00	93	326825	42.283	ng	98
29) 2,4-Dichlorophenol	8.14	162	205117	42.131	ng	99
30) 1,2,4-Trichlorobenzene	8.20	180	227847	41.738	ng	96
31) Naphthalene	8.28	128	690373	42.536	ng	100
32) Benzoic acid	8.01	122	89136m	27.600	ng	
33) 4-Chloroaniline	8.33	127	281592	39.150	ng	99
34) Hexachlorobutadiene	8.39	225	156613	44.656	ng	98
35) Caprolactam	8.68	113	59352	36.041	ng	88
36) 4-Chloro-3-methylphenol	8.83	107	195735	36.345	ng	98
37) 2-Methylnaphthalene	8.97	142	429964	38.895	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.14	216	218528	46.517	ng	# 96
42) 2,4,6-Trichlorophenol	9.25	196	137671	45.985	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	9.32	196	122489	37.952	nq	95
45) 1,1'-Biphenyl	9.44	154	513702	45.830	nq	98
46) 2-Chloronaphthalene	9.46	162	392998	43.812	nq	98
47) 2-Nitroaniline	9.56	65	136653	48.427	nq	89
48) Acenaphthylene	9.88	152	617173	44.951	nq	99
49) Dimethylphthalate	9.73	163	465098	44.985	nq	# 98
50) 2,6-Dinitrotoluene	9.80	165	89567	42.410	nq	99
51) Acenaphthene	10.05	154	371486	41.896	nq	99
52) 3-Nitroaniline	9.97	138	119023	47.484	nq	97
53) 2,4-Dinitrophenol	10.07	184	5985	14.122	nq	# 20
54) Dibenzofuran	10.22	168	525135	43.981	nq	99
55) 4-Nitrophenol	10.19	139	65687	34.189	nq	91
56) 2,4-Dinitrotoluene	10.20	165	111674	42.108	nq	# 90
57) Fluorene	10.57	166	425195	44.947	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.35	232	110916	43.360	nq	# 84
59) Diethylphthalate	10.42	149	456021	44.437	nq	97
60) 4-Chlorophenyl-phenylether	10.55	204	218822	43.979	nq	91
61) 4-Nitroaniline	10.58	138	129553	53.122	nq	100
62) Azobenzene	10.71	77	432585	40.426	nq	98
64) 4,6-Dinitro-2-methylphenol	10.61	198	19985	16.706	nq	87
65) n-Nitrosodiphenylamine	10.67	169	409202	39.383	nq	94
66) 4-Bromophenyl-phenylether	11.05	248	143868	40.914	nq	# 83
67) Hexachlorobenzene	11.11	284	150915	41.630	nq	# 88
68) Atrazine	11.20	200	123752	38.648	nq	99
69) Pentachlorophenol	11.32	266	65555	29.372	nq	96
70) Phenanthrene	11.53	178	666210	44.001	nq	99
71) Anthracene	11.58	178	680373	44.855	nq	99
72) Carbazole	11.74	167	633979	45.273	nq	100
73) Di-n-butylphthalate	12.06	149	737572	47.325	nq	100
74) Fluoranthene	12.72	202	810696	50.208	nq	98
76) Benzidine	12.85	184	490381	46.889	nq	99
77) Pyrene	12.95	202	846576	43.032	nq	98
79) Butylbenzylphthalate	13.57	149	380874	51.720	nq	94
80) Benzo(a)anthracene	14.15	228	789690	42.229	nq	99
81) 3,3'-Dichlorobenzidine	14.11	252	309541	49.061	nq	# 94
82) Chrysene	14.18	228	714251	41.835	nq	99
83) Bis(2-ethylhexyl)phthalate	14.12	149	535621	50.727	nq	99
84) Di-n-octyl phthalate	14.74	149	892589	58.637	nq	98
85) Indeno(1,2,3-cd)pyrene	17.26	276	522607	38.354	nq	94
87) Benzo(b)fluoranthene	15.24	252	597410m	42.921	nq	
88) Benzo(k)fluoranthene	15.27	252	590351	42.827	nq	# 97
89) Benzo(a)pyrene	15.62	252	528136	42.186	nq	98
90) Dibenzo(a,h)anthracene	17.28	278	450806	43.192	nq	96
91) Benzo(g,h,i)perylene	17.74	276	415593	40.972	nq	93

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

