

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061115\
 Data File : BF079910.D
 Acq On : 12 Jun 2015 00:00
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

apatel
 6/13/2015 8:44:16 AM

Quant Time: Jun 12 09:14:45 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF061115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 12 08:55:39 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	46096	20.00	ng	0.00
21) Naphthalene-d8	8.68	136	189094	20.00	ng	0.00
38) Acenaphthene-d10	10.46	164	97037	20.00	ng	0.00
63) Phenanthrene-d10	11.96	188	217058	20.00	ng	-0.01
75) Chrysene-d12	14.85	240	240248	20.00	ng	-0.15
86) Perylene-d12	16.71	264	228069	20.00	ng	-0.22

System Monitoring Compounds

5) 2-Fluorophenol	6.00	112	197705	78.65	ng	0.00
7) Phenol-d6	6.99	99	279904	81.73	ng	0.00
23) Nitrobenzene-d5	7.95	82	240426	83.89	ng	0.00
41) 2,4,6-Tribromophenol	11.26	330	73471	76.43	ng	0.00
44) 2-Fluorobiphenyl	9.76	172	540978	78.07	ng	0.00
78) Terphenyl-d14	13.61	244	799862	76.76	ng	-0.08

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.32	88	41986	36.71	ng	97
3) Pyridine	4.11	79	128073	43.70	ng	97
4) n-Nitrosodimethylamine	4.03	42	55296	38.29	ng	# 95
6) Aniline	7.05	93	194111	39.16	ng	99
8) 2-Chlorophenol	7.18	128	122690	40.16	ng	97
9) Benzaldehyde	6.94	77	72375	36.76	ng	98
10) Phenol	7.00	94	158157	41.99	ng	99
11) bis(2-Chloroethyl)ether	7.11	93	122538	40.43	ng	98
12) 1,3-Dichlorobenzene	7.34	146	143165	40.17	ng	99
13) 1,4-Dichlorobenzene	7.42	146	142501	38.49	ng	98
14) 1,2-Dichlorobenzene	7.56	146	127287	38.71	ng	99
15) Benzyl Alcohol	7.52	79	107818	39.26	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.66	45	177318	39.50	ng	99
17) 2-Methylphenol	7.62	107	106755	39.99	ng	98
18) Hexachloroethane	7.92	117	45347	39.56	ng	99
19) n-Nitroso-di-n-propylamine	7.78	70	96860	39.58	ng	97
20) 3+4-Methylphenols	7.77	107	137659	40.35	ng	99
22) Acetophenone	7.79	105	183586	41.11	ng	99
24) Nitrobenzene	7.96	77	124495	40.84	ng	98
25) Isophorone	8.20	82	267065	40.04	ng	99
26) 2-Nitrophenol	8.30	139	39912	37.88	ng	94
27) 2,4-Dimethylphenol	8.31	122	116564	39.00	ng	99
28) bis(2-Chloroethoxy)methane	8.41	93	154919	39.12	ng	99
29) 2,4-Dichlorophenol	8.52	162	101944	40.72	ng	97
30) 1,2,4-Trichlorobenzene	8.63	180	123643	38.86	ng	98
31) Naphthalene	8.71	128	414043	40.06	ng	99
32) Benzoic acid	8.36	122	35224m	33.35	ng	
33) 4-Chloroaniline	8.74	127	169666m	42.29	ng	
34) Hexachlorobutadiene	8.82	225	67386	38.15	ng	95
35) Caprolactam	9.08	113	34716	43.84	ng	93
36) 4-Chloro-3-methylphenol	9.21	107	118637	41.48	ng	97
37) 2-Methylnaphthalene	9.40	142	283289	40.57	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.58	216	125606	38.65	ng	99
40) Hexachlorocyclopentadiene	9.56	237	50921	44.59	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.68	196	79699	39.80	nq	99
43) 2,4,5-Trichlorophenol	9.71	196	94503	42.75	nq	97
45) 1,1'-Biphenyl	9.86	154	313509	39.92	nq	98
46) 2-Chloronaphthalene	9.90	162	270258	40.42	nq	99
47) 2-Nitroaniline	9.98	65	59999	42.33	nq	95
48) Acenaphthylene	10.32	152	462197	40.95	nq	100
49) Dimethylphthalate	10.15	163	302126	38.86	nq	100
50) 2,6-Dinitrotoluene	10.22	165	59929	42.55	nq	92
51) Acenaphthene	10.49	154	253581	39.53	nq	99
52) 3-Nitroaniline	10.39	138	69597	40.76	nq	94
53) 2,4-Dinitrophenol	10.49	184	8646	45.56	nq	80
54) Dibenzofuran	10.66	168	375358	40.05	nq	98
55) 4-Nitrophenol	10.52	139	51580	42.80	nq	87
56) 2,4-Dinitrotoluene	10.63	165	69091	37.83	nq	95
57) Fluorene	11.02	166	321318	39.33	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.78	232	74823	44.66	nq	97
59) Diethylphthalate	10.86	149	322051	41.43	nq	98
60) 4-Chlorophenyl-phenylether	10.99	204	146913	40.27	nq	94
61) 4-Nitroaniline	11.00	138	77480	43.11	nq	93
62) Azobenzene	11.15	77	310258	41.48	nq	97
64) 4,6-Dinitro-2-methylphenol	11.04	198	19967	42.33	nq	95
65) n-Nitrosodiphenylamine	11.11	169	286870	40.94	nq	99
66) 4-Bromophenyl-phenylether	11.50	248	92393	37.98	nq	# 92
67) Hexachlorobenzene	11.58	284	105604	39.80	nq	# 92
68) Atrazine	11.62	200	81649	37.86	nq	91
69) Pentachlorophenol	11.76	266	42508	38.64	nq	96
70) Phenanthrene	11.99	178	513813	39.72	nq	100
71) Anthracene	12.04	178	503299	40.21	nq	99
72) Carbazole	12.19	167	472347	38.91	nq	99
73) Di-n-butylphthalate	12.51	149	495293	37.45	nq	99
74) Fluoranthene	13.22	202	557795	39.00	nq	97
76) Benzidine	13.34	184	281108	42.25	nq	99
77) Pyrene	13.47	202	585271	38.79	nq	100
79) Butylbenzylphthalate	14.14	149	214416	42.08	nq	# 86
80) Benzo(a)anthracene	14.83	228	583911	38.27	nq	98
81) 3,3'-Dichlorobenzidine	14.78	252	193535	39.68	nq	99
82) Chrysene	14.88	228	556961	41.18	nq	99
83) Bis(2-ethylhexyl)phthalate	14.80	149	336497	41.26	nq	99
84) Di-n-octyl phthalate	15.57	149	543651	40.88	nq	99
85) Indeno(1,2,3-cd)pyrene	18.61	276	641321	40.14	nq	99
87) Benzo(b)fluoranthene	16.16	252	546017	39.07	nq	98
88) Benzo(k)fluoranthene	16.19	252	596828	41.88	nq	99
89) Benzo(a)pyrene	16.63	252	536222	40.35	nq	# 96
90) Dibenzo(a,h)anthracene	18.63	278	509799	39.02	nq	98
91) Benzo(g,h,i)perylene	19.18	276	519489	38.43	nq	99

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

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