

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110217\
 Data File : BF100097.D
 Acq On : 2 Nov 2017 20:50
 Operator : SJ/JU
 Sample : I6249-03MS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-2MS

Manual Integrations
 APPROVED

Sohil
 11/3/2017 1:23:24 PM

Quant Time: Nov 03 11:19:52 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 18:13:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	86750	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	350753	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	154746	20.00	ng	-0.01
63) Phenanthrene-d10	11.30	188	239652	20.00	ng	-0.01
75) Chrysene-d12	13.95	240	151432	20.00	ng	-0.02
86) Perylene-d12	15.41	264	155013	20.00	ng	-0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.39	112	435798	78.87	ng	0.00
7) Phenol-d6	6.42	99	500477	76.39	ng	0.00
23) Nitrobenzene-d5	7.34	82	281139	51.60	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	101765	72.16	ng	-0.01
44) 2-Fluorobiphenyl	9.12	172	545205	54.36	ng	-0.01
78) Terphenyl-d14	12.87	244	362919	52.37	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.50	88	52613	21.562	ng	# 89
3) Pyridine	3.29	79	116161	19.796	ng	# 87
4) n-Nitrosodimethylamine	3.23	42	44236	21.102	ng	# 65
6) Aniline	6.44	93	122931	14.471	ng	# 70
8) 2-Chlorophenol	6.56	128	157183	26.690	ng	# 92
9) Benzaldehyde	6.33	77	58913	15.048	ng	# 94
10) Phenol	6.43	94	200337	27.850	ng	# 97
11) bis(2-Chloroethyl)ether	6.51	93	144315	25.980	ng	# 93
12) 1,3-Dichlorobenzene	6.71	146	167259m	25.295	ng	# 97
13) 1,4-Dichlorobenzene	6.79	146	170665	25.431	ng	# 97
14) 1,2-Dichlorobenzene	6.94	146	160590	26.256	ng	# 99
15) Benzyl Alcohol	6.92	79	109010	26.162	ng	# 15
16) 2,2'-oxybis(1-Chloropropan	7.04	45	176949	28.676	ng	# 94
17) 2-Methylphenol	7.04	107	125519	25.481	ng	# 98
18) Hexachloroethane	7.27	117	51004	22.780	ng	# 87
19) n-Nitroso-di-n-propylamine	7.19	70	98006	25.526	ng	# 95
20) 3+4-Methylphenols	7.20	107	167529	29.207	ng	# 92
22) Acetophenone	7.19	105	208861	26.152	ng	# 87
24) Nitrobenzene	7.36	77	142134	25.892	ng	# 97
25) Isophorone	7.59	82	267891	27.098	ng	# 82
26) 2-Nitrophenol	7.68	139	82949	26.382	ng	# 93
27) 2,4-Dimethylphenol	7.72	122	140042	30.230	ng	# 99
28) bis(2-Chloroethoxy)methane	7.80	93	179854	25.977	ng	# 96
29) 2,4-Dichlorophenol	7.93	162	132724	27.579	ng	# 99
30) 1,2,4-Trichlorobenzene	7.99	180	131222	25.520	ng	# 99
31) Naphthalene	8.07	128	438891	25.922	ng	# 87
32) Benzoic acid	7.84	122	16940	4.154	ng	# 94
33) 4-Chloroaniline	8.14	127	70490	10.082	ng	# 99
34) Hexachlorobutadiene	8.18	225	65274	24.680	ng	# 75
35) Caprolactam	8.50	113	39559	26.139	ng	# 91
36) 4-Chloro-3-methylphenol	8.63	107	127419	25.941	ng	# 98
37) 2-Methylnaphthalene	8.76	142	298777	26.333	ng	# 97
39) 1,2,4,5-Tetrachlorobenzene	8.93	216	121299	24.270	ng	# 82
40) Hexachlorocyclopentadiene	8.91	237	686	10.831	ng	

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.06	196	81984	27.119	nq	99
43) 2,4,5-Trichlorophenol	9.10	196	86420	26.938	nq	95
45) 1,1'-Biphenyl	9.23	154	346782	24.772	nq	97
46) 2-Chloronaphthalene	9.26	162	268918	26.451	nq	95
47) 2-Nitroaniline	9.37	65	69265	25.958	nq #	75
48) Acenaphthylene	9.67	152	393759	25.146	nq	99
49) Dimethylphthalate	9.53	163	408588	33.342	nq	99
50) 2,6-Dinitrotoluene	9.60	165	66205	25.699	nq	90
51) Acenaphthene	9.84	154	234784	24.539	nq	99
52) 3-Nitroaniline	9.78	138	55059	18.138	nq #	90
53) 2,4-Dinitrophenol	9.93	184	4748	10.093	nq #	81
54) Dibenzofuran	10.02	168	353484	26.173	nq	96
55) 4-Nitrophenol	9.97	139	68978m	43.490	nq	
56) 2,4-Dinitrotoluene	10.02	165	85617	27.596	nq #	83
57) Fluorene	10.36	166	283130	25.320	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.15	232	64074	28.486	nq #	88
59) Diethylphthalate	10.23	149	298606	24.952	nq	96
60) 4-Chlorophenyl-phenylether	10.35	204	129144	24.540	nq	91
61) 4-Nitroaniline	10.40	138	62394	21.544	nq	85
62) Azobenzene	10.50	77	245163	24.439	nq	91
64) 4,6-Dinitro-2-methylphenol	10.45	198	11451	7.601	nq #	73
65) n-Nitrosodiphenylamine	10.47	169	248473	28.087	nq	98
66) 4-Bromophenyl-phenylether	10.83	248	71662	28.404	nq	90
67) Hexachlorobenzene	10.91	284	69354	26.870	nq #	88
68) Atrazine	10.99	200	74162	27.908	nq	98
69) Pentachlorophenol	11.12	266	50404	45.701	nq	98
70) Phenanthrene	11.32	178	358152	26.481	nq	98
71) Anthracene	11.37	178	365609	26.632	nq	98
72) Carbazole	11.54	167	333385	25.824	nq	98
73) Di-n-butylphthalate	11.84	149	438604	26.636	nq #	98
74) Fluoranthene	12.52	202	321587	22.880	nq	98
76) Benzidine	12.64	184	155604	23.483	nq	98
77) Pyrene	12.74	202	319193	27.720	nq	98
79) Butylbenzylphthalate	13.34	149	145078	25.586	nq	91
80) Benzo(a)anthracene	13.94	228	246763	25.705	nq	98
81) 3,3'-Dichlorobenzidine	13.89	252	80328	23.052	nq	99
82) Chrysene	13.97	228	237622	26.329	nq	98
83) Bis(2-ethylhexyl)phthalate	13.90	149	193639	24.318	nq	100
84) Di-n-octyl phthalate	14.51	149	323104	23.641	nq #	92
85) Indeno(1,2,3-cd)pyrene	16.87	276	220033	26.557	nq	97
87) Benzo(b)fluoranthene	14.99	252	246013	25.133	nq	98
88) Benzo(k)fluoranthene	15.01	252	235339	25.497	nq	99
89) Benzo(a)pyrene	15.35	252	232150	26.403	nq	99
90) Dibenzo(a,h)anthracene	16.87	278	188603	24.140	nq	97
91) Benzo(g,h,i)perylene	17.32	276	173468	22.694	nq	96

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

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