

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101117\
 Data File : BF099358.D
 Acq On : 11 Oct 2017 21:36
 Operator : SJ/JU
 Sample : I5710-02MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-29BMS

Manual Integrations
 APPROVED

Sohil
 10/13/2017 12:07:17 PM

Quant Time: Oct 12 01:23:57 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 16:25:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	108411	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	443943	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	199840	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	340142	20.00	ng	0.00
75) Chrysene-d12	14.02	240	185414	20.00	ng	0.00
86) Perylene-d12	15.48	264	175586	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.49	112	867997	127.46	ng	0.02
7) Phenol-d6	6.49	99	1052492	125.28	ng	0.00
23) Nitrobenzene-d5	7.42	82	631724	91.26	ng	0.00
41) 2,4,6-Tribromophenol	10.69	330	210381	128.66	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	1079159	90.15	ng	0.00
78) Terphenyl-d14	12.96	244	830369	101.85	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.69	88	114939	35.332	ng	96
3) Pyridine	3.50	79	80841	9.186	ng	98
4) n-Nitrosodimethylamine	3.41	42	136864	36.675	ng	82
6) Aniline	6.52	93	352679	31.104	ng	98
8) 2-Chlorophenol	6.65	128	328737	42.625	ng	92
9) Benzaldehyde	6.41	77	123900	25.425	ng	93
10) Phenol	6.50	94	428122	45.566	ng	99
11) bis(2-Chloroethyl)ether	6.59	93	309991	42.440	ng	98
12) 1,3-Dichlorobenzene	6.79	146	348346	43.129	ng	98
13) 1,4-Dichlorobenzene	6.87	146	354931	43.191	ng	99
14) 1,2-Dichlorobenzene	7.02	146	328100	43.210	ng	99
15) Benzyl Alcohol	7.00	79	256375	41.598	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.13	45	439354	38.607	ng	97
17) 2-Methylphenol	7.11	107	276288	43.783	ng	96
18) Hexachloroethane	7.36	117	120850	40.914	ng	99
19) n-Nitroso-di-n-propylamine	7.27	70	203687	37.975	ng	95
20) 3+4-Methylphenols	7.26	107	321521	40.276	ng	# 78
22) Acetophenone	7.27	105	431198	40.089	ng	# 95
24) Nitrobenzene	7.44	77	325709	41.477	ng	96
25) Isophorone	7.67	82	605951	42.338	ng	99
26) 2-Nitrophenol	7.75	139	185384	49.093	ng	92
27) 2,4-Dimethylphenol	7.79	122	292539	41.021	ng	98
28) bis(2-Chloroethoxy)methane	7.88	93	392613	44.019	ng	99
29) 2,4-Dichlorophenol	7.99	162	279891	45.713	ng	97
30) 1,2,4-Trichlorobenzene	8.07	180	276031	45.075	ng	99
31) Naphthalene	8.16	128	926074	44.415	ng	99
32) Benzoic acid	7.92	122	139083	23.743	ng	94
33) 4-Chloroaniline	8.21	127	291978	33.533	ng	# 94
34) Hexachlorobutadiene	8.27	225	136942	43.416	ng	98
35) Caprolactam	8.59	113	80472m	40.973	ng	
36) 4-Chloro-3-methylphenol	8.69	107	287536	41.781	ng	94
37) 2-Methylnaphthalene	8.85	142	630690	46.530	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.02	216	253253	42.494	ng	99
40) Hexachlorocyclopentadiene	9.00	237	164809	60.511	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.13	196	179971	42.626	nq	100
43) 2,4,5-Trichlorophenol	9.17	196	190068	45.096	nq	93
45) 1,1'-Biphenyl	9.31	154	733747	42.722	nq	99
46) 2-Chloronaphthalene	9.34	162	579193	44.915	nq	97
47) 2-Nitroaniline	9.44	65	168089	42.570	nq	90
48) Acenaphthylene	9.76	152	869206	44.031	nq	99
49) Dimethylphthalate	9.62	163	826442	54.794	nq	99
50) 2,6-Dinitrotoluene	9.68	165	153607	46.780	nq	89
51) Acenaphthene	9.93	154	509975	40.782	nq	99
52) 3-Nitroaniline	9.85	138	141710	37.012	nq	92
53) 2,4-Dinitrophenol	9.96	184	112776	66.990	nq	88
54) Dibenzofuran	10.10	168	759310	45.605	nq	99
55) 4-Nitrophenol	10.02	139	224091	69.218	nq	# 79
56) 2,4-Dinitrotoluene	10.09	165	198176	46.503	nq	# 93
57) Fluorene	10.44	166	603282	45.081	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.22	232	138363	47.523	nq	96
59) Diethylphthalate	10.31	149	627760	42.594	nq	98
60) 4-Chlorophenyl-phenylether	10.43	204	270828	45.053	nq	97
61) 4-Nitroaniline	10.47	138	166651	45.116	nq	88
62) Azobenzene	10.59	77	581437	42.906	nq	91
64) 4,6-Dinitro-2-methylphenol	10.50	198	86305	41.703	nq	# 78
65) n-Nitrosodiphenylamine	10.55	169	541911	45.989	nq	99
66) 4-Bromophenyl-phenylether	10.92	248	158604	47.637	nq	98
67) Hexachlorobenzene	10.99	284	159137	44.752	nq	98
68) Atrazine	11.07	200	158391	44.676	nq	98
69) Pentachlorophenol	11.19	266	143945	65.042	nq	99
70) Phenanthrene	11.40	178	838138	46.034	nq	99
71) Anthracene	11.46	178	851465	45.706	nq	99
72) Carbazole	11.61	167	786582	43.117	nq	100
73) Di-n-butylphthalate	11.93	149	946940	43.124	nq	99
74) Fluoranthene	12.59	202	783319	42.487	nq	97
76) Benzidine	12.71	184	92655	13.511	nq	97
77) Pyrene	12.82	202	789230	55.821	nq	99
79) Butylbenzylphthalate	13.43	149	332584	50.052	nq	91
80) Benzo(a)anthracene	14.00	228	525156	46.775	nq	99
81) 3,3'-Dichlorobenzidine	13.96	252	149057	36.216	nq	98
82) Chrysene	14.04	228	493105	46.133	nq	99
83) Bis(2-ethylhexyl)phthalate	13.98	149	423629	46.301	nq	99
84) Di-n-octyl phthalate	14.59	149	600767	40.778	nq	# 95
85) Indeno(1,2,3-cd)pyrene	16.97	276	559934	65.486	nq	96
87) Benzo(b)fluoranthene	15.05	252	476189	44.109	nq	97
88) Benzo(k)fluoranthene	15.08	252	426016	39.953	nq	99
89) Benzo(a)pyrene	15.42	252	456832	46.099	nq	# 96
90) Dibenzo(a,h)anthracene	16.98	278	461067	58.137	nq	97
91) Benzo(g,h,i)perylene	17.42	276	489273	62.412	nq	96

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

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