

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101317\
 Data File : BF099449.D
 Acq On : 13 Oct 2017 19:44
 Operator : SJ/JU
 Sample : PB103199BSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB103199BSD

Manual Integrations
 APPROVED

Sohil
 10/15/2017 12:39:59 PM

Quant Time: Oct 13 23:19:21 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	100965	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	405660	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	163806	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	240702	20.00	ng	0.00
75) Chrysene-d12	14.02	240	167191	20.00	ng	0.00
86) Perylene-d12	15.49	264	189079	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.49	112	806738	127.20	ng	0.02
7) Phenol-d6	6.50	99	959486	122.63	ng	0.01
23) Nitrobenzene-d5	7.42	82	574037	90.75	ng	0.00
41) 2,4,6-Tribromophenol	10.69	330	162925	121.55	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	1014436	103.39	ng	0.00
78) Terphenyl-d14	12.96	244	642385	87.38	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.69	88	105967	34.98	ng	94
3) Pyridine	3.48	79	299653	36.56	ng	94
4) n-Nitrosodimethylamine	3.42	42	123177	35.44	ng	82
6) Aniline	6.52	93	250933	23.76	ng	99
8) 2-Chlorophenol	6.65	128	298267	41.53	ng	92
9) Benzaldehyde	6.41	77	102355	22.55	ng	93
10) Phenol	6.51	94	360935	41.25	ng	97
11) bis(2-Chloroethyl)ether	6.59	93	275294	40.47	ng	98
12) 1,3-Dichlorobenzene	6.79	146	317284	42.18	ng	97
13) 1,4-Dichlorobenzene	6.87	146	320316	41.85	ng	99
14) 1,2-Dichlorobenzene	7.03	146	296636	41.95	ng	96
15) Benzyl Alcohol	7.00	79	223573	38.95	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.13	45	387631	36.57	ng	93
17) 2-Methylphenol	7.11	107	244169	41.55	ng	98
18) Hexachloroethane	7.36	117	109065	39.65	ng	97
19) n-Nitroso-di-n-propylamine	7.27	70	179610	35.96	ng	93
20) 3+4-Methylphenols	7.26	107	288953	38.87	ng	# 73
22) Acetophenone	7.27	105	376468	38.30	ng	# 97
24) Nitrobenzene	7.44	77	288467	40.20	ng	98
25) Isophorone	7.67	82	530190	40.54	ng	99
26) 2-Nitrophenol	7.76	139	163729	47.45	ng	# 86
27) 2,4-Dimethylphenol	7.79	122	258433	39.66	ng	95
28) bis(2-Chloroethoxy)methane	7.89	93	344646	42.29	ng	100
29) 2,4-Dichlorophenol	8.00	162	240687	43.02	ng	97
30) 1,2,4-Trichlorobenzene	8.07	180	242335	43.31	ng	99
31) Naphthalene	8.16	128	812556	42.65	ng	99
32) Benzoic acid	7.92	122	137769	25.74	ng	97
33) 4-Chloroaniline	8.21	127	78586	9.88	ng	95
34) Hexachlorobutadiene	8.27	225	118614	41.15	ng	99
35) Caprolactam	8.59	113	76053m	42.38	ng	
36) 4-Chloro-3-methylphenol	8.70	107	245902	39.10	ng	95
37) 2-Methylnaphthalene	8.85	142	551501	44.53	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.02	216	216167	44.25	ng	98
40) Hexachlorocyclopentadiene	9.00	237	85462	38.28	ng	96

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42) 2,4,6-Trichlorophenol	9.13	196	148085	42.79	ng	99
43) 2,4,5-Trichlorophenol	9.17	196	154525	44.73	ng	95
45) 1,1'-Biphenyl	9.32	154	625554	44.43	ng	98
46) 2-Chloronaphthalene	9.34	162	486700	46.04	ng	98
47) 2-Nitroaniline	9.44	65	140406	43.38	ng	98
48) Acenaphthylene	9.76	152	722358	44.64	ng	99
49) Dimethylphthalate	9.61	163	564238	45.64	ng	99
50) 2,6-Dinitrotoluene	9.68	165	125590	46.66	ng	89
51) Acenaphthene	9.93	154	420569	41.03	ng	99
52) 3-Nitroaniline	9.85	138	69809	22.24	ng	# 87
53) 2,4-Dinitrophenol	9.96	184	63391	47.77	ng	97
54) Dibenzofuran	10.10	168	627567	45.98	ng	100
55) 4-Nitrophenol	10.02	139	152030	57.29	ng	84
56) 2,4-Dinitrotoluene	10.09	165	153457	43.93	ng	92
57) Fluorene	10.44	166	478797	43.65	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.22	232	103887	43.53	ng	96
59) Diethylphthalate	10.31	149	518581	42.93	ng	98
60) 4-Chlorophenyl-phenylether	10.43	204	218955	44.44	ng	98
61) 4-Nitroaniline	10.47	138	115068	38.00	ng	86
62) Azobenzene	10.59	77	470113	42.32	ng	91
64) 4,6-Dinitro-2-methylphenol	10.50	198	49573	33.85	ng	82
65) n-Nitrosodiphenylamine	10.55	169	424910	50.96	ng	98
66) 4-Bromophenyl-phenylether	10.92	248	120444	51.12	ng	# 89
67) Hexachlorobenzene	10.99	284	114169	45.37	ng	99
68) Atrazine	11.07	200	124957	49.81	ng	99
69) Pentachlorophenol	11.19	266	91625	58.51	ng	99
70) Phenanthrene	11.40	178	608616	47.24	ng	98
71) Anthracene	11.46	178	624867	47.40	ng	99
72) Carbazole	11.61	167	556711	43.12	ng	99
73) Di-n-butylphthalate	11.93	149	694371	44.69	ng	98
74) Fluoranthene	12.59	202	541017	41.47	ng	99
76) Benzidine	12.72	184	261928	42.36	ng	97
77) Pyrene	12.82	202	545181	42.76	ng	100
79) Butylbenzylphthalate	13.43	149	247958	41.38	ng	98
80) Benzo(a)anthracene	14.01	228	469913	46.42	ng	99
81) 3,3'-Dichlorobenzidine	13.97	252	140493	37.86	ng	99
82) Chrysene	14.05	228	451932	46.89	ng	99
83) Bis(2-ethylhexyl)phthalate	13.98	149	333416	40.41	ng	# 100
84) Di-n-octyl phthalate	14.59	149	606923	45.69	ng	95
85) Indeno(1,2,3-cd)pyrene	16.99	276	523423	67.89	ng	96
87) Benzo(b)fluoranthene	15.06	252	492255	42.34	ng	98
88) Benzo(k)fluoranthene	15.09	252	510928	44.50	ng	98
89) Benzo(a)pyrene	15.43	252	502352	47.08	ng	# 96
90) Dibenzo(a,h)anthracene	17.00	278	437831	51.27	ng	97
91) Benzo(g,h,i)perylene	17.44	276	435344	51.57	ng	95

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

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