

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021617\
 Data File : BF093000.D
 Acq On : 17 Feb 2017 00:05
 Operator : SJ/MA
 Sample : I1865-01MSD
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TILCON-SCREENINGMSD

Quant Time: Feb 17 02:50:31 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 16 23:55:37 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	135588	20.00	ng	0.01
21) Naphthalene-d8	8.14	136	567262	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	306126	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	492264	20.00	ng	0.00
75) Chrysene-d12	14.02	240	361305	20.00	ng	0.00
86) Perylene-d12	15.44	264	292135	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	1048731	132.70	ng	0.01
7) Phenol-d6	6.50	99	1509770	148.47	ng	0.00
23) Nitrobenzene-d5	7.44	82	944803	92.75	ng	0.01
41) 2,4,6-Tribromophenol	10.69	330	318347	143.78	ng	0.01
44) 2-Fluorobiphenyl	9.22	172	1404022	83.09	ng	0.00
78) Terphenyl-d14	12.97	244	1321304	85.93	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.48	88	149759	35.07	ng	# 86
3) Pyridine	3.21	79	429689	39.57	ng	87
4) n-Nitrosodimethylamine	3.16	42	220496	43.25	ng	90
6) Aniline	6.52	93	421072	30.36	ng	84
8) 2-Chlorophenol	6.65	128	401825	46.09	ng	98
9) Benzaldehyde	6.41	77	80218	11.01	ng	99
10) Phenol	6.52	94	504974	42.44	ng	92
11) bis(2-Chloroethyl)ether	6.60	93	408435	43.01	ng	97
12) 1,3-Dichlorobenzene	6.80	146	420565m	41.18	ng	
13) 1,4-Dichlorobenzene	6.88	146	428499	41.46	ng	98
14) 1,2-Dichlorobenzene	7.04	146	412322	41.72	ng	95
15) Benzyl Alcohol	7.01	79	344204	45.72	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.14	45	718408	43.55	ng	94
17) 2-Methylphenol	7.13	107	355237	47.49	ng	96
18) Hexachloroethane	7.37	117	145655	41.28	ng	# 84
19) n-Nitroso-di-n-propylamine	7.28	70	280257	43.30	ng	94
20) 3+4-Methylphenols	7.28	107	390403	43.65	ng	# 78
22) Acetophenone	7.28	105	513123	39.62	ng	# 92
24) Nitrobenzene	7.45	77	424004	40.54	ng	97
25) Isophorone	7.69	82	806128	42.47	ng	97
26) 2-Nitrophenol	7.77	139	234064	47.08	ng	87
27) 2,4-Dimethylphenol	7.80	122	350370	40.12	ng	98
28) bis(2-Chloroethoxy)methane	7.90	93	517379	41.15	ng	99
29) 2,4-Dichlorophenol	8.01	162	358368	44.00	ng	93
30) 1,2,4-Trichlorobenzene	8.09	180	347156	39.56	ng	96
31) Naphthalene	8.17	128	1129669	39.28	ng	99
32) Benzoic acid	7.89	122	31045	12.50	ng	94
33) 4-Chloroaniline	8.21	127	249928	22.16	ng	# 92
34) Hexachlorobutadiene	8.28	225	176836	36.62	ng	100
35) Caprolactam	8.59	113	111235m	43.42	ng	
36) 4-Chloro-3-methylphenol	8.70	107	351901	41.45	ng	100
37) 2-Methylnaphthalene	8.85	142	737313	39.93	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.02	216	341421	39.73	ng	99
40) Hexachlorocyclopentadiene	9.01	237	266717	68.79	ng	98

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42) 2,4,6-Trichlorophenol	9.14	196	246278	43.62	ng	93
43) 2,4,5-Trichlorophenol	9.18	196	253543	40.98	ng #	87
45) 1,1'-Biphenyl	9.32	154	924189	41.69	ng	98
46) 2-Chloronaphthalene	9.34	162	688746	39.72	ng	96
47) 2-Nitroaniline	9.45	65	266834	45.77	ng	88
48) Acenaphthylene	9.76	152	1107084	36.96	ng	99
49) Dimethylphthalate	9.63	163	1397068	67.75	ng	100
50) 2,6-Dinitrotoluene	9.69	165	176099	39.55	ng #	78
51) Acenaphthene	9.93	154	713514	39.84	ng	96
52) 3-Nitroaniline	9.86	138	170060	33.37	ng	83
53) 2,4-Dinitrophenol	9.96	184	112800	51.58	ng #	59
54) Dibenzofuran	10.10	168	979475	40.89	ng	98
55) 4-Nitrophenol	10.03	139	321206	87.51	ng	87
56) 2,4-Dinitrotoluene	10.09	165	231132	38.99	ng	91
57) Fluorene	10.44	166	665756	36.12	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.22	232	196933	47.72	ng	97
59) Diethylphthalate	10.33	149	832828	42.86	ng	95
60) 4-Chlorophenyl-phenylether	10.44	204	348035	39.31	ng #	74
61) 4-Nitroaniline	10.48	138	222439	42.62	ng	82
62) Azobenzene	10.60	77	927736	46.21	ng	89
64) 4,6-Dinitro-2-methylphenol	10.50	198	127338	42.80	ng #	64
65) n-Nitrosodiphenylamine	10.56	169	628151	39.94	ng	100
66) 4-Bromophenyl-phenylether	10.92	248	199545	38.80	ng	92
67) Hexachlorobenzene	10.99	284	195711	38.51	ng #	87
68) Atrazine	11.08	200	188600	37.37	ng	99
69) Pentachlorophenol	11.18	266	212967	81.07	ng	96
70) Phenanthrene	11.40	178	1046316	37.10	ng	98
71) Anthracene	11.46	178	1218706	43.03	ng	98
72) Carbazole	11.61	167	1019659	37.66	ng	99
73) Di-n-butylphthalate	11.94	149	1200835	41.01	ng #	98
74) Fluoranthene	12.59	202	1087727	37.39	ng	97
76) Benzidine	12.72	184	295512	19.19	ng	96
77) Pyrene	12.82	202	1097601	40.59	ng	99
79) Butylbenzylphthalate	13.44	149	535846	48.80	ng	89
80) Benzo(a)anthracene	14.01	228	833515	37.72	ng	99
81) 3,3'-Dichlorobenzidine	13.96	252	244290	31.01	ng	97
82) Chrysene	14.04	228	821653	39.71	ng	98
83) Bis(2-ethylhexyl)phthalate	14.00	149	706092	47.02	ng #	96
84) Di-n-octyl phthalate	14.60	149	1101549	45.05	ng	98
85) Indeno(1,2,3-cd)pyrene	16.85	276	744860	46.48	ng #	94
87) Benzo(b)fluoranthene	15.04	252	798273m	41.52	ng	
88) Benzo(k)fluoranthene	15.06	252	614232m	37.00	ng	
89) Benzo(a)pyrene	15.38	252	687166	41.31	ng #	96
90) Dibenzo(a,h)anthracene	16.87	278	619655	46.10	ng	99
91) Benzo(g,h,i)perylene	17.28	276	633505	46.65	ng #	92

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

