

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091015\
 Data File : BF081569.D
 Acq On : 10 Sep 2015 15:58
 Operator : UM/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Sep 11 12:14:04 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 10 16:31:29 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.26	152	45830	20.00	ng	0.00
21) Naphthalene-d8	11.17	136	182830	20.00	ng	0.00
38) Acenaphthene-d10	15.30	164	88530	20.00	ng	0.00
63) Phenanthrene-d10	18.81	188	178753	20.00	ng	0.00
75) Chrysene-d12	23.42	240	138477	20.00	ng	0.00
86) Perylene-d12	24.85	264	81498	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	233913	79.19	ng	0.00
7) Phenol-d6	7.69	99	317146	81.11	ng	0.00
23) Nitrobenzene-d5	9.57	82	312571	82.48	ng	0.00
41) 2,4,6-Tribromophenol	17.21	330	65963	83.68	ng	0.00
44) 2-Fluorobiphenyl	13.82	172	560834	81.18	ng	0.00
78) Terphenyl-d14	22.15	244	517219	86.95	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.60	88	58051	43.76	ng	# 91
3) Pyridine	2.17	79	141565	38.70	ng	# 77
4) n-Nitrosodimethylamine	2.09	42	88825	43.71	ng	# 59
6) Aniline	7.58	93	211236	38.26	ng	# 84
8) 2-Chlorophenol	7.82	128	130004	39.94	ng	# 80
9) Benzaldehyde	7.29	77	90791	37.06	ng	# 75
10) Phenol	7.71	94	170458	37.59	ng	# 51
11) bis(2-Chloroethyl)ether	7.81	93	132600	40.53	ng	# 69
12) 1,3-Dichlorobenzene	8.11	146	136211	39.17	ng	# 88
13) 1,4-Dichlorobenzene	8.30	146	136885	38.66	ng	# 88
14) 1,2-Dichlorobenzene	8.62	146	129630	40.66	ng	# 90
15) Benzyl Alcohol	8.72	79	129941	40.51	ng	# 72
16) 2,2'-oxybis(1-Chloropropan	9.03	45	281778	44.94	ng	92
17) 2-Methylphenol	9.05	107	105838	40.75	ng	# 79
18) Hexachloroethane	9.36	117	55467	40.26	ng	# 88
19) n-Nitroso-di-n-propylamine	9.35	70	112325	42.21	ng	# 88
20) 3+4-Methylphenols	9.43	107	139136	39.73	ng	85
22) Acetophenone	9.27	105	195415	39.13	ng	# 74
24) Nitrobenzene	9.61	77	163597	41.57	ng	# 63
25) Isophorone	10.23	82	289157	41.02	ng	# 84
26) 2-Nitrophenol	10.34	139	67483	39.34	ng	# 33
27) 2,4-Dimethylphenol	10.62	122	116650	39.38	ng	# 76
28) bis(2-Chloroethoxy)methane	10.80	93	167599	41.17	ng	# 91
29) 2,4-Dichlorophenol	10.95	162	102166	39.77	ng	91
30) 1,2,4-Trichlorobenzene	11.08	180	114374	40.98	ng	98
31) Naphthalene	11.21	128	390874	40.09	ng	98
32) Benzoic acid	11.09	122	42320	21.49	ng	# 59
33) 4-Chloroaniline	11.46	127	157781	39.67	ng	# 85
34) Hexachlorobutadiene	11.57	225	73227	43.12	ng	97
35) Caprolactam	12.46	113	28689	36.41	ng	# 6
36) 4-Chloro-3-methylphenol	12.77	107	126928	44.14	ng	# 74
37) 2-Methylnaphthalene	12.87	142	256702	40.46	ng	# 87
39) 1,2,4,5-Tetrachlorobenzene	13.27	216	111045	39.66	ng	99
40) Hexachlorocyclopentadiene	13.26	237	60257	43.85	ng	98

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42) 2,4,6-Trichlorophenol	13.62	196	79212	40.99	ng	98
43) 2,4,5-Trichlorophenol	13.73	196	80287	40.73	ng #	89
45) 1,1'-Biphenyl	14.01	154	309937	38.05	ng	94
46) 2-Chloronaphthalene	14.00	162	231659	39.38	ng #	88
47) 2-Nitroaniline	14.36	65	99654	41.57	ng #	62
48) Acenaphthylene	14.96	152	405602	38.87	ng	97
49) Dimethylphthalate	14.90	163	279415	40.62	ng #	97
50) 2,6-Dinitrotoluene	14.98	165	57714	36.97	ng #	16
51) Acenaphthene	15.38	154	226740	38.20	ng	90
52) 3-Nitroaniline	15.36	138	65250	36.71	ng #	50
53) 2,4-Dinitrophenol	15.64	184	24284	34.64	ng #	36
54) Dibenzofuran	15.81	168	341102	39.35	ng #	85
55) 4-Nitrophenol	15.98	139	45484	40.74	ng #	11
56) 2,4-Dinitrotoluene	15.93	165	79928	40.21	ng #	51
57) Fluorene	16.62	166	273762	41.62	ng	96
58) 2,3,4,6-Tetrachlorophenol	16.17	232	63775	42.38	ng	91
59) Diethylphthalate	16.60	149	303914	42.00	ng	95
60) 4-Chlorophenyl-phenylether	16.71	204	133302	43.48	ng #	81
61) 4-Nitroaniline	16.82	138	64272	35.80	ng #	12
62) Azobenzene	17.09	77	335695	41.73	ng	84
64) 4,6-Dinitro-2-methylphenol	16.89	198	38335	38.34	ng #	69
65) n-Nitrosodiphenylamine	17.03	169	245270	37.71	ng	94
66) 4-Bromophenyl-phenylether	17.85	248	73780	40.82	ng #	84
67) Hexachlorobenzene	17.90	284	73733	39.02	ng #	83
68) Atrazine	18.45	200	74211	39.43	ng #	78
69) Pentachlorophenol	18.45	266	36596	44.17	ng	96
70) Phenanthrene	18.87	178	387490	38.99	ng	97
71) Anthracene	19.00	178	388273	38.03	ng	98
72) Carbazole	19.46	167	363298	37.92	ng	95
73) Di-n-butylphthalate	20.52	149	483314	42.72	ng #	97
74) Fluoranthene	21.48	202	421991	39.23	ng	86
76) Benzidine	21.78	184	147869	24.87	ng	98
77) Pyrene	21.82	202	432296	42.14	ng	98
79) Butylbenzylphthalate	22.84	149	183190	41.06	ng #	73
80) Benzo(a)anthracene	23.41	228	343792	38.98	ng	96
81) 3,3'-Dichlorobenzidine	23.41	252	96602	32.47	ng	98
82) Chrysene	23.44	228	284525	35.99	ng	95
83) Bis(2-ethylhexyl)phthalate	23.52	149	237238	40.30	ng #	89
84) Di-n-octyl phthalate	24.18	149	336938	34.76	ng #	93
85) Indeno(1,2,3-cd)pyrene	25.91	276	211301	36.43	ng #	84
87) Benzo(b)fluoranthene	24.50	252	200115m	34.39	ng	
88) Benzo(k)fluoranthene	24.54	252	219301m	45.42	ng	
89) Benzo(a)pyrene	24.80	252	195046	39.56	ng #	83
90) Dibenzo(a,h)anthracene	25.92	278	177309	46.54	ng #	81
91) Benzo(g,h,i)perylene	26.22	276	177376	48.78	ng #	75

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

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