

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012517\
 Data File : BF092485.D
 Acq On : 25 Jan 2017 19:13
 Operator : SJ/MA
 Sample : PB96282BS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB96282BS

Quant Time: Jan 26 14:54:40 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 24 13:48:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	159274	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	644665	20.00	ng	-0.01
38) Acenaphthene-d10	10.03	164	361339	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	599743	20.00	ng	-0.01
75) Chrysene-d12	14.16	240	520154	20.00	ng	-0.01
86) Perylene-d12	15.63	264	431805	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.57	112	1108139	108.25	ng	0.02
7) Phenol-d6	6.59	99	1287303	98.72	ng	0.00
23) Nitrobenzene-d5	7.53	82	874787	74.12	ng	-0.01
41) 2,4,6-Tribromophenol	10.82	330	288462	117.15	ng	0.00
44) 2-Fluorobiphenyl	9.34	172	1363449	69.78	ng	-0.01
78) Terphenyl-d14	13.10	244	1404042	71.88	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.70	88	162629	29.87	ng	# 78
3) Pyridine	3.45	79	445775	28.76	ng	# 83
4) n-Nitrosodimethylamine	3.40	42	243146	31.98	ng	82
6) Aniline	6.62	93	246864	12.69	ng	# 45
8) 2-Chlorophenol	6.75	128	388337	36.11	ng	99
9) Benzaldehyde	6.51	77	207562	22.43	ng	97
10) Phenol	6.60	94	551548	35.04	ng	# 71
11) bis(2-Chloroethyl)ether	6.70	93	395960	33.62	ng	97
12) 1,3-Dichlorobenzene	6.91	146	434120	34.14	ng	98
13) 1,4-Dichlorobenzene	6.98	146	415552	32.47	ng	99
14) 1,2-Dichlorobenzene	7.14	146	412732	33.99	ng	96
15) Benzyl Alcohol	7.10	79	343530	34.96	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.25	45	753231	32.34	ng	93
17) 2-Methylphenol	7.22	107	340246	37.03	ng	96
18) Hexachloroethane	7.48	117	143273	32.50	ng	90
19) n-Nitroso-di-n-propylamine	7.38	70	275153	30.99	ng	91
20) 3+4-Methylphenols	7.37	107	392507	35.21	ng	93
22) Acetophenone	7.38	105	531264	34.60	ng	# 98
24) Nitrobenzene	7.55	77	419975	33.97	ng	99
25) Isophorone	7.79	82	780230	33.55	ng	95
26) 2-Nitrophenol	7.87	139	199157	38.44	ng	97
27) 2,4-Dimethylphenol	7.90	122	374091	34.72	ng	97
28) bis(2-Chloroethoxy)methane	8.01	93	497229	32.57	ng	100
29) 2,4-Dichlorophenol	8.11	162	333832	35.74	ng	92
30) 1,2,4-Trichlorobenzene	8.20	180	356855	35.34	ng	96
31) Naphthalene	8.28	128	1081623	32.78	ng	99
32) Benzoic acid	8.02	122	257080	33.49	ng	99
33) 4-Chloroaniline	8.33	127	115161	8.25	ng	99
34) Hexachlorobutadiene	8.41	225	179073	32.43	ng	97
35) Caprolactam	8.69	113	116538m	37.36	ng	
36) 4-Chloro-3-methylphenol	8.81	107	351416	34.89	ng	97
37) 2-Methylnaphthalene	8.98	142	745640	35.70	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.15	216	344041	37.75	ng	99
40) Hexachlorocyclopentadiene	9.14	237	341807	74.88	ng	98

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42) 2,4,6-Trichlorophenol	9.25	196	252863	36.21	nq	90
43) 2,4,5-Trichlorophenol	9.30	196	224695	35.16	nq #	82
45) 1,1'-Biphenyl	9.45	154	960946	36.79	nq	98
46) 2-Chloronaphthalene	9.47	162	704567	33.02	nq	98
47) 2-Nitroaniline	9.56	65	245403	34.16	nq	94
48) Acenaphthylene	9.89	152	1151498	36.80	nq	97
49) Dimethylphthalate	9.74	163	791065	34.43	nq	98
50) 2,6-Dinitrotoluene	9.80	165	176568	37.59	nq #	59
51) Acenaphthene	10.06	154	749902	38.57	nq	99
52) 3-Nitroaniline	9.97	138	81789	13.90	nq	87
53) 2,4-Dinitrophenol	10.08	184	145854	57.23	nq #	59
54) Dibenzofuran	10.24	168	1040174	39.40	nq	92
55) 4-Nitrophenol	10.13	139	357928	76.62	nq	89
56) 2,4-Dinitrotoluene	10.21	165	275984	44.27	nq #	76
57) Fluorene	10.58	166	753251	38.48	nq	97
58) 2,3,4,6-Tetrachlorophenol	10.35	232	204568	42.73	nq	97
59) Diethylphthalate	10.45	149	848944	38.53	nq	96
60) 4-Chlorophenyl-phenylether	10.57	204	336295	35.42	nq	87
61) 4-Nitroaniline	10.59	138	183387	31.17	nq	96
62) Azobenzene	10.73	77	934194	37.26	nq	97
64) 4,6-Dinitro-2-methylphenol	10.61	198	108077	34.94	nq #	34
65) n-Nitrosodiphenylamine	10.68	169	648499	34.62	nq	100
66) 4-Bromophenyl-phenylether	11.06	248	208780	34.48	nq	94
67) Hexachlorobenzene	11.13	284	213233	34.84	nq #	89
68) Atrazine	11.22	200	234299	36.55	nq	95
69) Pentachlorophenol	11.31	266	265615	67.87	nq	97
70) Phenanthrene	11.54	178	1146695	34.72	nq	99
71) Anthracene	11.58	178	1164534	36.08	nq	99
72) Carbazole	11.74	167	1138190	36.93	nq	98
73) Di-n-butylphthalate	12.08	149	1265617	35.91	nq #	98
74) Fluoranthene	12.73	202	1229678	38.06	nq	96
76) Benzidine	12.84	184	258928	13.41	nq	99
77) Pyrene	12.96	202	1240973	32.93	nq	99
79) Butylbenzylphthalate	13.57	149	543542	35.60	nq	99
80) Benzo(a)anthracene	14.15	228	994658	35.58	nq	99
81) 3,3'-Dichlorobenzidine	14.11	252	247928	23.36	nq #	93
82) Chrysene	14.18	228	932792	31.24	nq	99
83) Bis(2-ethylhexyl)phthalate	14.13	149	682864	35.51	nq	99
84) Di-n-octyl phthalate	14.75	149	1354913	38.81	nq	97
85) Indeno(1,2,3-cd)pyrene	17.11	276	822284	37.23	nq #	94
87) Benzo(b)fluoranthene	15.20	252	926462	33.42	nq	98
88) Benzo(k)fluoranthene	15.23	252	946564m	41.06	nq	
89) Benzo(a)pyrene	15.56	252	888884	37.90	nq #	96
90) Dibenzo(a,h)anthracene	17.13	278	686997	36.23	nq	97
91) Benzo(g,h,i)perylene	17.55	276	678280	35.37	nq #	90

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

