

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042017\
 Data File : BF094537.D
 Acq On : 20 Apr 2017 11:47
 Operator : SJ/MA
 Sample : PB98356BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB98356BS

Manual Integrations
 APPROVED

mohammad
 4/21/2017 11:35:56 AM

Quant Time: Apr 20 17:40:13 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 17 14:59:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.77	152	114691	20.00	ng	0.00
21) Naphthalene-d8	7.26	136	451402	20.00	ng	0.00
38) Acenaphthene-d10	9.40	164	205085	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	388100	20.00	ng	0.00
75) Chrysene-d12	14.47	240	290844	20.00	ng	0.00
86) Perylene-d12	16.08	264	220336	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	4.32	112	864104	125.00	ng	0.02
7) Phenol-d6	5.40	99	1053029	129.21	ng	0.00
23) Nitrobenzene-d5	6.42	82	651773	89.16	ng	0.00
41) 2,4,6-Tribromophenol	10.38	330	238692	148.80	ng	0.00
44) 2-Fluorobiphenyl	8.58	172	1219648	87.50	ng	0.00
78) Terphenyl-d14	13.19	244	1170434	107.57	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.11	88	107405	26.72	ng	97
3) Pyridine	2.64	79	311262	35.42	ng	99
4) n-Nitrosodimethylamine	2.58	42	147515	40.57	ng	90
6) Aniline	5.40	93	322905	29.81	ng	# 80
8) 2-Chlorophenol	5.54	128	335570	42.66	ng	93
9) Benzaldehyde	5.27	77	52046	8.40	ng	97
10) Phenol	5.41	94	402390	40.19	ng	97
11) bis(2-Chloroethyl)ether	5.48	93	316566	43.28	ng	93
12) 1,3-Dichlorobenzene	5.70	146	347844	39.91	ng	99
13) 1,4-Dichlorobenzene	5.79	146	355617	40.56	ng	98
14) 1,2-Dichlorobenzene	5.96	146	327621	40.56	ng	98
15) Benzyl Alcohol	5.95	79	252211	41.72	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.10	45	398705	41.58	ng	78
17) 2-Methylphenol	6.09	107	263399	42.51	ng	94
18) Hexachloroethane	6.35	117	122507	40.76	ng	# 85
19) n-Nitroso-di-n-propylamine	6.25	70	234581	43.44	ng	97
20) 3+4-Methylphenols	6.27	107	368408	43.76	ng	# 62
22) Acetophenone	6.24	105	466451	42.50	ng	# 85
24) Nitrobenzene	6.44	77	330385	42.92	ng	92
25) Isophorone	6.73	82	595260	43.93	ng	# 94
26) 2-Nitrophenol	6.81	139	187448	45.32	ng	98
27) 2,4-Dimethylphenol	6.89	122	295473	43.98	ng	98
28) bis(2-Chloroethoxy)methane	6.99	93	382500	43.64	ng	99
29) 2,4-Dichlorophenol	7.11	162	281918	45.11	ng	99
30) 1,2,4-Trichlorobenzene	7.20	180	275360	42.62	ng	96
31) Naphthalene	7.29	128	920117	42.40	ng	100
32) Benzoic acid	7.07	122	180341	50.76	ng	95
33) 4-Chloroaniline	7.37	127	91590	10.15	ng	# 92
34) Hexachlorobutadiene	7.44	225	138713	43.06	ng	99
35) Caprolactam	7.83	113	96615m	49.21	ng	
36) 4-Chloro-3-methylphenol	8.00	107	303251	46.30	ng	89
37) 2-Methylnaphthalene	8.13	142	645192	43.46	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.34	216	247415	42.37	ng	99
40) Hexachlorocyclopentadiene	8.32	237	239833	101.52	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.49	196	188246	45.90	nq	97
43) 2,4,5-Trichlorophenol	8.55	196	198150	47.05	nq #	92
45) 1,1'-Biphenyl	8.70	154	724621	43.25	nq	97
46) 2-Chloronaphthalene	8.72	162	583072	43.76	nq	95
47) 2-Nitroaniline	8.85	65	175323	45.74	nq #	85
48) Acenaphthylene	9.22	152	906758	43.66	nq	98
49) Dimethylphthalate	9.10	163	726668	47.55	nq	99
50) 2,6-Dinitrotoluene	9.17	165	163892	47.77	nq	98
51) Acenaphthene	9.44	154	549652	44.18	nq	100
52) 3-Nitroaniline	9.37	138	93155	23.47	nq	86
53) 2,4-Dinitrophenol	9.51	184	146807	78.95	nq #	68
54) Dibenzofuran	9.65	168	819543	45.33	nq	98
55) 4-Nitrophenol	9.63	139	246645m	87.96	nq	
56) 2,4-Dinitrotoluene	9.66	165	211323	50.20	nq #	84
57) Fluorene	10.07	166	649264	46.11	nq	100
58) 2,3,4,6-Tetrachlorophenol	9.82	232	149310	49.46	nq	96
59) Diethylphthalate	9.96	149	709477	49.20	nq	99
60) 4-Chlorophenyl-phenylether	10.08	204	280534	47.88	nq	93
61) 4-Nitroaniline	10.12	138	164848	40.86	nq	98
62) Azobenzene	10.27	77	662353	47.31	nq	96
64) 4,6-Dinitro-2-methylphenol	10.17	198	110179	43.33	nq #	63
65) n-Nitrosodiphenylamine	10.23	169	604541	44.79	nq	98
66) 4-Bromophenyl-phenylether	10.68	248	175740	45.60	nq	96
67) Hexachlorobenzene	10.75	284	174172	44.85	nq #	88
68) Atrazine	10.91	200	185410	45.92	nq	95
69) Pentachlorophenol	11.00	266	165903	107.58	nq	98
70) Phenanthrene	11.25	178	973573	44.55	nq	99
71) Anthracene	11.31	178	1017069	44.99	nq	99
72) Carbazole	11.52	167	937751	44.20	nq	98
73) Di-n-butylphthalate	11.97	149	1174330	50.27	nq	99
74) Fluoranthene	12.71	202	1021670	45.11	nq	99
76) Benzidine	12.88	184	259151	22.01	nq #	91
77) Pyrene	12.98	202	1015928	50.00	nq	100
79) Butylbenzylphthalate	13.80	149	489892	55.01	nq #	85
80) Benzo(a)anthracene	14.45	228	779145	46.09	nq	100
81) 3,3'-Dichlorobenzidine	14.43	252	171163	28.60	nq #	96
82) Chrysene	14.50	228	712115	46.09	nq	99
83) Bis(2-ethylhexyl)phthalate	14.52	149	677940	56.70	nq #	99
84) Di-n-octyl phthalate	15.27	149	1018503	50.78	nq	97
85) Indeno(1,2,3-cd)pyrene	17.37	276	659220	44.85	nq	99
87) Benzo(b)fluoranthene	15.69	252	632356	46.92	nq	98
88) Benzo(k)fluoranthene	15.72	252	576896	45.23	nq	95
89) Benzo(a)pyrene	16.03	252	575215	45.87	nq	98
90) Dibenzo(a,h)anthracene	17.38	278	536193	51.50	nq	97
91) Benzo(g,h,i)perylene	17.74	276	566056	53.40	nq	99

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

