

Data Path : Z:\HPCHEM1\BNA M\DATA\BM072017\
 Data File : BM010912.D
 Acq On : 20 Jul 2017 17:14
 Operator : SJ/JU
 Sample : PB100784BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB100784BS

Manual Integrations
 APPROVED

Sohil
 7/21/2017 1:28:37 PM

Quant Time: Jul 21 04:01:24 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM071217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 12 14:21:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	226728	20.00	ng	-0.01
21) Naphthalene-d8	10.20	136	954444	20.00	ng	-0.01
38) Acenaphthene-d10	14.09	164	689540	20.00	ng	0.00
63) Phenanthrene-d10	16.84	188	1761756	20.00	ng	0.00
75) Chrysene-d12	21.06	240	1552960	20.00	ng	0.00
86) Perylene-d12	23.19	264	1259345	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.08	112	1853959	135.45	ng	0.00
7) Phenol-d6	6.64	99	2483714	131.63	ng	0.00
23) Nitrobenzene-d5	8.59	82	1986583	80.00	ng	0.00
41) 2,4,6-Tribromophenol	15.60	330	1497059	142.07	ng	0.00
44) 2-Fluorobiphenyl	12.70	172	4680830	84.45	ng	0.00
78) Terphenyl-d14	19.50	244	7172990	89.33	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.09	88	178214	29.508	ng	# 100
3) Pyridine	3.46	79	539406	32.675	ng	# 87
4) n-Nitrosodimethylamine	3.38	42	343981	40.770	ng	# 72
6) Aniline	6.79	93	844831	35.742	ng	96
8) 2-Chlorophenol	7.02	128	570978	42.134	ng	83
9) Benzaldehyde	6.60	77	332205	25.355	ng	80
10) Phenol	6.67	94	861909	43.203	ng	98
11) bis(2-Chloroethyl)ether	6.89	93	600610	39.089	ng	95
12) 1,3-Dichlorobenzene	7.33	146	625669	37.360	ng	# 81
13) 1,4-Dichlorobenzene	7.47	146	640562	37.064	ng	# 91
14) 1,2-Dichlorobenzene	7.79	146	630863	38.307	ng	# 92
15) Benzyl Alcohol	7.69	79	798462	44.689	ng	# 83
16) 2,2'-oxybis(1-Chloropropan	7.97	45	552099	41.445	ng	80
17) 2-Methylphenol	7.89	107	581440	45.064	ng	# 78
18) Hexachloroethane	8.50	117	288607	39.813	ng	# 80
19) n-Nitroso-di-n-propylamine	8.25	70	627659	41.992	ng	93
20) 3+4-Methylphenols	8.22	107	776280	43.305	ng	# 83
22) Acetophenone	8.26	105	1071900	37.300	ng	# 94
24) Nitrobenzene	8.63	77	993560	39.043	ng	# 89
25) Isophorone	9.15	82	1654775	40.679	ng	# 85
26) 2-Nitrophenol	9.33	139	351541	41.514	ng	# 29
27) 2,4-Dimethylphenol	9.40	122	610195	40.374	ng	# 70
28) bis(2-Chloroethoxy)methane	9.64	93	932695	40.901	ng	100
29) 2,4-Dichlorophenol	9.86	162	730817	42.806	ng	95
30) 1,2,4-Trichlorobenzene	10.07	180	820039	39.290	ng	97
31) Naphthalene	10.25	128	1953615	40.255	ng	99
32) Benzoic acid	9.56	122	440534	37.150	ng	# 63
33) 4-Chloroaniline	10.37	127	298139	15.248	ng	# 76
34) Hexachlorobutadiene	10.54	225	632340	39.045	ng	98
35) Caprolactam	11.16	113	210189m	46.243	ng	
36) 4-Chloro-3-methylphenol	11.51	107	879523	42.589	ng	# 76
37) 2-Methylnaphthalene	11.87	142	1552021	44.525	ng	# 88
39) 1,2,4,5-Tetrachlorobenzene	12.25	216	1120699	37.287	ng	# 100
40) Hexachlorocyclopentadiene	12.23	237	1728852	100.340	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.50	196	737698	39.860	nq	98
43) 2,4,5-Trichlorophenol	12.57	196	784312	42.924	nq	# 87
45) 1,1'-Biphenyl	12.92	154	2224223	37.025	nq	98
46) 2-Chloronaphthalene	12.95	162	1760135	39.960	nq	94
47) 2-Nitroaniline	13.17	65	658431	45.472	nq	# 70
48) Acenaphthylene	13.81	152	2711654	41.391	nq	98
49) Dimethylphthalate	13.57	163	2398955	40.895	nq	# 98
50) 2,6-Dinitrotoluene	13.68	165	502983	43.714	nq	# 56
51) Acenaphthene	14.16	154	1663160	41.547	nq	98
52) 3-Nitroaniline	14.01	138	338202	32.332	nq	# 43
53) 2,4-Dinitrophenol	14.22	184	698302	87.792	nq	# 88
54) Dibenzofuran	14.49	168	2915011	45.017	nq	# 89
55) 4-Nitrophenol	14.33	139	725058	79.812	nq	# 1
56) 2,4-Dinitrotoluene	14.47	165	725274	45.426	nq	# 94
57) Fluorene	15.15	166	2415096	43.386	nq	98
58) 2,3,4,6-Tetrachlorophenol	14.73	232	813796	47.933	nq	# 100
59) Diethylphthalate	14.94	149	2511466	43.350	nq	98
60) 4-Chlorophenyl-phenylether	15.15	204	1440374	42.124	nq	96
61) 4-Nitroaniline	15.19	138	461624	42.223	nq	# 1
62) Azobenzene	15.44	77	2959258	50.303	nq	88
64) 4,6-Dinitro-2-methylphenol	15.24	198	475978	40.972	nq	93
65) n-Nitrosodiphenylamine	15.37	169	2110523	41.866	nq	97
66) 4-Bromophenyl-phenylether	16.04	248	968697	43.386	nq	# 86
67) Hexachlorobenzene	16.16	284	1018068	39.876	nq	# 91
68) Atrazine	16.34	200	866704	41.515	nq	96
69) Pentachlorophenol	16.50	266	1292207	78.954	nq	97
70) Phenanthrene	16.89	178	4010640	43.760	nq	100
71) Anthracene	16.98	178	4029002	44.286	nq	99
72) Carbazole	17.26	167	3253953	40.611	nq	100
73) Di-n-butylphthalate	17.84	149	4007711	42.277	nq	# 97
74) Fluoranthene	18.92	202	4629054	40.781	nq	97
76) Benzidine	19.12	184	2369052	57.007	nq	99
77) Pyrene	19.29	202	4751431	52.777	nq	100
79) Butylbenzylphthalate	20.22	149	1550176	49.182	nq	# 74
80) Benzo(a)anthracene	21.04	228	3982165	44.254	nq	98
81) 3,3'-Dichlorobenzidine	20.99	252	1046712	29.583	nq	# 97
82) Chrysene	21.10	228	3658501	42.690	nq	99
83) Bis(2-ethylhexyl)phthalate	21.00	149	2056312	44.801	nq	# 99
84) Di-n-octyl phthalate	21.86	149	3070975	40.836	nq	99
85) Indeno(1,2,3-cd)pyrene	25.30	276	3887062	39.555	nq	# 100
87) Benzo(b)fluoranthene	22.56	252	3479606	43.498	nq	# 92
88) Benzo(k)fluoranthene	22.60	252	3391483	44.388	nq	# 94
89) Benzo(a)pyrene	23.10	252	3321708	45.620	nq	# 96
90) Dibenzo(a,h)anthracene	25.31	278	3201464	45.301	nq	# 91
91) Benzo(g,h,i)perylene	25.94	276	3192324	45.953	nq	# 87

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

