

Data Path : Z:\HPCHEM1\BNA M\DATA\BM081117\
 Data File : BM011207.D
 Acq On : 11 Aug 2017 18:41
 Operator : SJ/JU
 Sample : PB101388BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB101388BS

Manual Integrations
 APPROVED

sohil
 8/14/2017 6:57:10 PM

Quant Time: Aug 11 23:24:46 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM072617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 26 17:29:06 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.29	152	161483	20.00	ng	-0.10
21) Naphthalene-d8	10.05	136	642488	20.00	ng	-0.10
38) Acenaphthene-d10	13.96	164	453723	20.00	ng	-0.09
63) Phenanthrene-d10	16.72	188	1099194	20.00	ng	-0.09
75) Chrysene-d12	20.94	240	1137350	20.00	ng	-0.08
86) Perylene-d12	23.02	264	1152911	20.00	ng	-0.12

System Monitoring Compounds

5) 2-Fluorophenol	4.93	112	843292	88.28	ng	-0.09
7) Phenol-d6	6.50	99	1162130	85.62	ng	-0.09
23) Nitrobenzene-d5	8.45	82	1560208	91.14	ng	-0.09
41) 2,4,6-Tribromophenol	15.47	330	549039	75.50	ng	-0.09
44) 2-Fluorobiphenyl	12.57	172	2940048	81.27	ng	-0.09
78) Terphenyl-d14	19.39	244	4289635	74.71	ng	-0.08

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.93	88	166197	38.749	ng	# 100
3) Pyridine	3.30	79	482834	41.213	ng	# 88
4) n-Nitrosodimethylamine	3.22	42	286372	47.966	ng	# 68
6) Aniline	6.64	93	554662	32.425	ng	# 78
8) 2-Chlorophenol	6.87	128	379053	39.084	ng	# 75
9) Benzaldehyde	6.46	77	237645	27.061	ng	# 75
10) Phenol	6.53	94	598777	40.617	ng	# 95
11) bis(2-Chloroethyl)ether	6.75	93	471118	41.015	ng	# 87
12) 1,3-Dichlorobenzene	7.18	146	466407	39.460	ng	# 77
13) 1,4-Dichlorobenzene	7.33	146	479080	39.542	ng	# 87
14) 1,2-Dichlorobenzene	7.64	146	452912	39.099	ng	# 88
15) Benzyl Alcohol	7.55	79	605967	46.540	ng	# 79
16) 2,2'-oxybis(1-Chloropropan	7.82	45	501035	48.473	ng	# 75
17) 2-Methylphenol	7.75	107	393724	41.788	ng	# 75
18) Hexachloroethane	8.35	117	221868	41.992	ng	# 84
19) n-Nitroso-di-n-propylamine	8.10	70	537436	46.597	ng	# 90
20) 3+4-Methylphenols	8.08	107	514579	39.290	ng	# 79
22) Acetophenone	8.12	105	770116	39.987	ng	# 91
24) Nitrobenzene	8.49	77	784454	44.222	ng	# 82
25) Isophorone	9.00	82	1295489	45.388	ng	# 83
26) 2-Nitrophenol	9.18	139	216272	38.317	ng	# 1
27) 2,4-Dimethylphenol	9.26	122	373299	37.570	ng	# 64
28) bis(2-Chloroethoxy)methane	9.50	93	676386	42.489	ng	# 97
29) 2,4-Dichlorophenol	9.72	162	430087	38.520	ng	# 89
30) 1,2,4-Trichlorobenzene	9.92	180	567981	41.142	ng	# 99
31) Naphthalene	10.10	128	1275481	39.462	ng	# 99
32) Benzoic acid	9.40	122	239362	29.970	ng	# 51
33) 4-Chloroaniline	10.23	127	206576	16.021	ng	# 71
34) Hexachlorobutadiene	10.39	225	465606	42.821	ng	# 96
35) Caprolactam	11.02	113	112900	35.656	ng	# 90
36) 4-Chloro-3-methylphenol	11.37	107	541876	37.586	ng	# 71
37) 2-Methylnaphthalene	11.73	142	978794	41.223	ng	# 84
39) 1,2,4,5-Tetrachlorobenzene	12.11	216	751100	38.531	ng	# 100
40) Hexachlorocyclopentadiene	12.09	237	1170234	103.026	ng	# 99

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42) 2,4,6-Trichlorophenol	12.37	196	448906	38.047	nq	96
43) 2,4,5-Trichlorophenol	12.45	196	457526	37.650	nq #	86
45) 1,1'-Biphenyl	12.78	154	1377160	35.102	nq	95
46) 2-Chloronaphthalene	12.82	162	1082161	37.437	nq #	94
47) 2-Nitroaniline	13.04	65	447991	43.808	nq #	62
48) Acenaphthylene	13.67	152	1631832	37.827	nq	100
49) Dimethylphthalate	13.44	163	1396484	35.979	nq #	99
50) 2,6-Dinitrotoluene	13.56	165	295253	37.617	nq #	52
51) Acenaphthene	14.03	154	997734	37.353	nq	99
52) 3-Nitroaniline	13.89	138	180947	26.566	nq #	27
53) 2,4-Dinitrophenol	14.10	184	372195	66.723	nq #	75
54) Dibenzofuran	14.37	168	1747671	40.380	nq #	91
55) 4-Nitrophenol	14.22	139	312809	52.273	nq #	75
56) 2,4-Dinitrotoluene	14.35	165	413790	37.553	nq #	75
57) Fluorene	15.02	166	1408860	37.388	nq	100
58) 2,3,4,6-Tetrachlorophenol	14.60	232	464897	40.815	nq #	100
59) Diethylphthalate	14.83	149	1468601	37.049	nq	99
60) 4-Chlorophenyl-phenylether	15.03	204	871583	38.310	nq	95
61) 4-Nitroaniline	15.06	138	176985	24.464	nq #	1
62) Azobenzene	15.32	77	2040911	48.107	nq	86
64) 4,6-Dinitro-2-methylphenol	15.12	198	270614	36.366	nq	92
65) n-Nitrosodiphenylamine	15.25	169	1222184	38.852	nq	99
66) 4-Bromophenyl-phenylether	15.93	248	577629	40.878	nq #	90
67) Hexachlorobenzene	16.03	284	584720	36.619	nq #	87
68) Atrazine	16.22	200	514874	40.843	nq	96
69) Pentachlorophenol	16.38	266	742146	73.209	nq	99
70) Phenanthrene	16.76	178	2328603	40.458	nq	99
71) Anthracene	16.85	178	2302930	40.120	nq	98
72) Carbazole	17.13	167	1791755	35.693	nq	99
73) Di-n-butylphthalate	17.72	149	2356512	38.548	nq #	95
74) Fluoranthene	18.80	202	2780678	38.986	nq	98
76) Benzidine	19.01	184	1191105	43.777	nq	99
77) Pyrene	19.16	202	2817395	41.690	nq	98
79) Butylbenzylphthalate	20.11	149	1000147	41.619	nq #	64
80) Benzo(a)anthracene	20.93	228	2686027	41.361	nq	97
81) 3,3'-Dichlorobenzidine	20.88	252	758162	29.759	nq #	97
82) Chrysene	20.98	228	2507243	40.220	nq	99
83) Bis(2-ethylhexyl)phthalate	20.90	149	1409734	39.876	nq #	96
84) Di-n-octyl phthalate	21.74	149	2223741	38.756	nq #	96
85) Indeno(1,2,3-cd)pyrene	25.05	276	3091144	47.879	nq #	96
87) Benzo(b)fluoranthene	22.42	252	2723205	36.802	nq #	91
88) Benzo(k)fluoranthene	22.46	252	2672166	37.678	nq #	94
89) Benzo(a)pyrene	22.93	252	2642784	39.470	nq #	96
90) Dibenzo(a,h)anthracene	25.06	278	2456607	39.578	nq #	92
91) Benzo(g,h,i)perylene	25.66	276	2581907m	42.361	nq	

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

