

Data Path : Z:\HPCHEM1\BNA M\DATA\BM072617\
 Data File : BM010985.D
 Acq On : 26 Jul 2017 23:48
 Operator : SJ/JU
 Sample : PB100891BSD
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB100891BSD

Manual Integrations
 APPROVED

Sohil
 7/27/2017 6:21:39 PM

Quant Time: Jul 27 09:57:53 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	7.39	152	281601	20.00	ng	0.00
21) Naphthalene-d8	10.15	136	1170471	20.00	ng	0.00
38) Acenaphthene-d10	14.05	164	785613	20.00	ng	0.00
63) Phenanthrene-d10	16.81	188	1909166	20.00	ng	0.00
75) Chrysene-d12	21.03	240	1658693	20.00	ng	0.00
86) Perylene-d12	23.14	264	1096714	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.03	112	2971874	178.40	ng	0.00
7) Phenol-d6	6.60	99	3989748	168.57	ng	0.01
23) Nitrobenzene-d5	8.55	82	3118264	99.99	ng	0.00
41) 2,4,6-Tribromophenol	15.56	330	2026595	160.95	ng	0.00
44) 2-Fluorobiphenyl	12.66	172	6696793	106.91	ng	0.00
78) Terphenyl-d14	19.47	244	9277723	110.80	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.04	88	276792	37.007	ng	# 100
3) Pyridine	3.41	79	861916	42.189	ng	# 89
4) n-Nitrosodimethylamine	3.33	42	505829	48.585	ng	# 69
6) Aniline	6.75	93	1350030	45.257	ng	91
8) 2-Chlorophenol	6.97	128	895126	52.926	ng	85
9) Benzaldehyde	6.55	77	454549	29.682	ng	84
10) Phenol	6.63	94	1372392	53.384	ng	97
11) bis(2-Chloroethyl)ether	6.85	93	992265	49.537	ng	94
12) 1,3-Dichlorobenzene	7.28	146	1004554	48.737	ng	# 84
13) 1,4-Dichlorobenzene	7.43	146	1027226	48.620	ng	# 90
14) 1,2-Dichlorobenzene	7.74	146	983585	48.691	ng	# 90
15) Benzyl Alcohol	7.65	79	1256548	55.341	ng	# 83
16) 2,2'-oxybis(1-Chloropropan	7.93	45	905178	50.218	ng	78
17) 2-Methylphenol	7.85	107	901231	54.852	ng	# 80
18) Hexachloroethane	8.45	117	441648	47.934	ng	# 82
19) n-Nitroso-di-n-propylamine	8.21	70	997659	49.602	ng	92
20) 3+4-Methylphenols	8.17	107	1202521	52.651	ng	# 85
22) Acetophenone	8.22	105	1672196	47.660	ng	# 95
24) Nitrobenzene	8.59	77	1543988	47.777	ng	# 89
25) Isophorone	9.11	82	2518775	48.440	ng	# 87
26) 2-Nitrophenol	9.28	139	533572	51.891	ng	# 27
27) 2,4-Dimethylphenol	9.36	122	931968	51.486	ng	# 77
28) bis(2-Chloroethoxy)methane	9.60	93	1451405	50.046	ng	# 98
29) 2,4-Dichlorophenol	9.81	162	1051699	51.704	ng	95
30) 1,2,4-Trichlorobenzene	10.02	180	1212285	48.202	ng	100
31) Naphthalene	10.20	128	2927522	49.718	ng	100
32) Benzoic acid	9.54	122	643985	44.261	ng	# 65
33) 4-Chloroaniline	10.33	127	537026	22.862	ng	# 72
34) Hexachlorobutadiene	10.49	225	962106	48.570	ng	98
35) Caprolactam	11.13	113	286593m	49.684	ng	
36) 4-Chloro-3-methylphenol	11.46	107	1279087	48.700	ng	# 72
37) 2-Methylnaphthalene	11.83	142	2293651	53.025	ng	# 88
39) 1,2,4,5-Tetrachlorobenzene	12.21	216	1638544	48.546	ng	# 100
40) Hexachlorocyclopentadiene	12.19	237	2546145	129.461	ng	99

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42) 2,4,6-Trichlorophenol	12.46	196	1048051	51.302	nq	99
43) 2,4,5-Trichlorophenol	12.53	196	1099404	52.250	nq	# 87
45) 1,1'-Biphenyl	12.87	154	3230822	47.561	nq	97
46) 2-Chloronaphthalene	12.91	162	2517694	50.303	nq	94
47) 2-Nitroaniline	13.13	65	938956	53.029	nq	# 70
48) Acenaphthylene	13.77	152	3821110	51.157	nq	99
49) Dimethylphthalate	13.53	163	3302152	49.135	nq	# 98
50) 2,6-Dinitrotoluene	13.65	165	682996	50.256	nq	# 61
51) Acenaphthene	14.12	154	2318013	50.120	nq	99
52) 3-Nitroaniline	13.97	138	475369	40.308	nq	# 40
53) 2,4-Dinitrophenol	14.19	184	960116	99.406	nq	# 88
54) Dibenzofuran	14.46	168	4043842	53.962	nq	# 90
55) 4-Nitrophenol	14.30	139	1024686	98.895	nq	97
56) 2,4-Dinitrotoluene	14.44	165	997033	52.258	nq	93
57) Fluorene	15.11	166	3324782	50.957	nq	99
58) 2,3,4,6-Tetrachlorophenol	14.69	232	1103520	55.954	nq	# 100
59) Diethylphthalate	14.91	149	3446428	50.214	nq	99
60) 4-Chlorophenyl-phenylether	15.12	204	2015628	51.167	nq	98
61) 4-Nitroaniline	15.15	138	619081	49.422	nq	# 1
62) Azobenzene	15.41	77	4043656	55.048	nq	89
64) 4,6-Dinitro-2-methylphenol	15.20	198	638639	49.412	nq	97
65) n-Nitrosodiphenylamine	15.33	169	2901319	53.101	nq	98
66) 4-Bromophenyl-phenylether	16.01	248	1343408	54.736	nq	# 89
67) Hexachlorobenzene	16.12	284	1378884	49.719	nq	# 91
68) Atrazine	16.30	200	1186909	54.208	nq	98
69) Pentachlorophenol	16.46	266	1735144	98.546	nq	96
70) Phenanthrene	16.85	178	5397907	53.996	nq	100
71) Anthracene	16.94	178	5343060	53.592	nq	99
72) Carbazole	17.22	167	4299323	49.310	nq	99
73) Di-n-butylphthalate	17.80	149	5488986	51.696	nq	# 96
74) Fluoranthene	18.88	202	6212852	50.150	nq	98
76) Benzidine	19.09	184	2938632	74.057	nq	99
77) Pyrene	19.24	202	6084280	61.733	nq	99
79) Butylbenzylphthalate	20.19	149	2102492	59.991	nq	# 72
80) Benzo(a)anthracene	21.01	228	4996550	52.756	nq	99
81) 3,3'-Dichlorobenzidine	20.96	252	1281845	34.500	nq	# 98
82) Chrysene	21.06	228	4588383	50.470	nq	100
83) Bis(2-ethylhexyl)phthalate	20.97	149	2871406	55.693	nq	# 98
84) Di-n-octyl phthalate	21.82	149	4083986	48.805	nq	100
85) Indeno(1,2,3-cd)pyrene	25.21	276	3300714	35.056	nq	# 96
87) Benzo(b)fluoranthene	22.52	252	3946192	56.063	nq	# 93
88) Benzo(k)fluoranthene	22.56	252	3585314	53.144	nq	# 95
89) Benzo(a)pyrene	23.05	252	3379322	53.057	nq	# 96
90) Dibenzo(a,h)anthracene	25.23	278	2731756	46.266	nq	# 90
91) Benzo(g,h,i)perylene	25.85	276	2706052	46.673	nq	# 85

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

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