

Data Path : Z:\HPCHEM1\BNA M\DATA\BM080717\
 Data File : BM011167.D
 Acq On : 08 Aug 2017 00:08
 Operator : SJ/JU
 Sample : I4592-02MSD
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 15-50-RT3460-RT4215-RT3451MSD

Manual Integrations
 APPROVED

Sohil
 8/10/2017 6:13:07 PM

Quant Time: Aug 08 03:17:01 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	7.31	152	161291	20.00	ng	-0.08
21) Naphthalene-d8	10.07	136	698134	20.00	ng	-0.08
38) Acenaphthene-d10	13.98	164	582330	20.00	ng	-0.07
63) Phenanthrene-d10	16.74	188	1786715	20.00	ng	-0.07
75) Chrysene-d12	20.97	240	2227368	20.00	ng	-0.06
86) Perylene-d12	23.05	264	1994956	20.00	ng	-0.09

System Monitoring Compounds

5) 2-Fluorophenol	4.96	112	1537626	161.15	ng	-0.07
7) Phenol-d6	6.52	99	2194268	161.86	ng	-0.07
23) Nitrobenzene-d5	8.46	82	2040331	109.69	ng	-0.08
41) 2,4,6-Tribromophenol	15.49	330	1736809	186.08	ng	-0.06
44) 2-Fluorobiphenyl	12.59	172	4562458	98.26	ng	-0.08
78) Terphenyl-d14	19.41	244	10244020	91.10	ng	-0.06

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.96	88	172793	40.335	ng	# 100
3) Pyridine	3.34	79	472657	40.393	ng	# 88
4) n-Nitrosodimethylamine	3.25	42	319742	53.619	ng	# 75
6) Aniline	6.67	93	510084	29.854	ng	# 83
8) 2-Chlorophenol	6.89	128	480007	49.552	ng	# 75
9) Benzaldehyde	6.48	77	298761	34.061	ng	# 76
10) Phenol	6.55	94	727959	49.438	ng	# 89
11) bis(2-Chloroethyl)ether	6.77	93	592537	51.646	ng	# 90
12) 1,3-Dichlorobenzene	7.20	146	497557	42.145	ng	# 75
13) 1,4-Dichlorobenzene	7.35	146	512085	42.317	ng	# 91
14) 1,2-Dichlorobenzene	7.66	146	510062	44.085	ng	# 87
15) Benzyl Alcohol	7.56	79	806472	62.013	ng	# 79
16) 2,2'-oxybis(1-Chloropropan	7.85	45	569904	55.202	ng	# 76
17) 2-Methylphenol	7.77	107	528529	56.163	ng	# 72
18) Hexachloroethane	8.37	117	230994	43.771	ng	# 80
19) n-Nitroso-di-n-propylamine	8.13	70	714325	62.007	ng	# 90
20) 3+4-Methylphenols	8.10	107	713454	54.539	ng	# 79
22) Acetophenone	8.13	105	1014794	48.492	ng	# 89
24) Nitrobenzene	8.50	77	1003069	52.039	ng	# 83
25) Isophorone	9.03	82	1784704	57.545	ng	# 85
26) 2-Nitrophenol	9.20	139	296305	48.312	ng	# 18
27) 2,4-Dimethylphenol	9.28	122	537471	49.781	ng	# 65
28) bis(2-Chloroethoxy)methane	9.52	93	893979	51.681	ng	# 98
29) 2,4-Dichlorophenol	9.73	162	615790	50.756	ng	# 89
30) 1,2,4-Trichlorobenzene	9.94	180	702890	46.856	ng	# 98
31) Naphthalene	10.12	128	1697582	48.336	ng	# 98
32) Benzoic acid	9.45	122	361271	41.629	ng	# 46
33) 4-Chloroaniline	10.27	127	230746	16.470	ng	# 69
34) Hexachlorobutadiene	10.41	225	556834	47.129	ng	# 96
35) Caprolactam	11.05	113	199834m	58.082	ng	#
36) 4-Chloro-3-methylphenol	11.39	107	907866	57.952	ng	# 69
37) 2-Methylnaphthalene	11.75	142	1414017	54.806	ng	# 87
39) 1,2,4,5-Tetrachlorobenzene	12.13	216	1048617	41.913	ng	# 100
40) Hexachlorocyclopentadiene	12.11	237	322255	22.105	ng	# 99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.39	196	724273	47.829	nq	98
43) 2,4,5-Trichlorophenol	12.46	196	783395	50.229	nq #	85
45) 1,1'-Biphenyl	12.80	154	2130035	42.302	nq	97
46) 2-Chloronaphthalene	12.83	162	1691028	45.581	nq #	94
47) 2-Nitroaniline	13.06	65	790992	60.267	nq #	62
48) Acenaphthylene	13.69	152	2699260	48.753	nq	98
49) Dimethylphthalate	13.46	163	2557790	51.345	nq #	99
50) 2,6-Dinitrotoluene	13.57	165	550874	54.684	nq #	43
51) Acenaphthene	14.05	154	1693439	49.397	nq	100
52) 3-Nitroaniline	13.91	138	392068	44.850	nq #	23
53) 2,4-Dinitrophenol	14.12	184	768893	107.398	nq #	80
54) Dibenzofuran	14.39	168	3043428	54.789	nq #	90
55) 4-Nitrophenol	14.24	139	656795	85.517	nq #	85
56) 2,4-Dinitrotoluene	14.37	165	830813	58.748	nq #	79
57) Fluorene	15.04	166	2605338	53.870	nq	99
58) 2,3,4,6-Tetrachlorophenol	14.62	232	878272	60.078	nq #	100
59) Diethylphthalate	14.85	149	2773787	54.521	nq	99
60) 4-Chlorophenyl-phenylether	15.05	204	1543287	52.853	nq	96
61) 4-Nitroaniline	15.08	138	351790	37.887	nq #	1
62) Azobenzene	15.34	77	3642948	66.905	nq	86
64) 4,6-Dinitro-2-methylphenol	15.14	198	563408	46.579	nq	94
65) n-Nitrosodiphenylamine	15.27	169	2332114	45.609	nq	98
66) 4-Bromophenyl-phenylether	15.94	248	1071784	46.662	nq #	91
67) Hexachlorobenzene	16.04	284	1136529	43.788	nq #	86
68) Atrazine	16.24	200	1045112	51.003	nq	97
69) Pentachlorophenol	16.40	266	1568091	95.162	nq	98
70) Phenanthrene	16.78	178	4794534	51.248	nq	100
71) Anthracene	16.87	178	4782966	51.262	nq	99
72) Carbazole	17.16	167	3892980	47.709	nq	98
73) Di-n-butylphthalate	17.74	149	5130129	51.627	nq #	96
74) Fluoranthene	18.82	202	6481135	55.901	nq	99
76) Benzidine	19.03	184	1225944m	23.007	nq	
77) Pyrene	19.18	202	6716061	50.746	nq	100
79) Butylbenzylphthalate	20.13	149	2480721	52.711	nq #	67
80) Benzo(a)anthracene	20.95	228	6558931	51.572	nq	99
81) 3,3'-Dichlorobenzidine	20.90	252	1659861	33.269	nq #	98
82) Chrysene	21.00	228	6163684	50.488	nq	99
83) Bis(2-ethylhexyl)phthalate	20.92	149	3564664	51.487	nq #	99
84) Di-n-octyl phthalate	21.76	149	5918196	52.668	nq	97
85) Indeno(1,2,3-cd)pyrene	25.09	276	5992531	47.396	nq #	96
87) Benzo(b)fluoranthene	22.44	252	6645628	51.903	nq #	92
88) Benzo(k)fluoranthene	22.49	252	6188946	50.432	nq #	95
89) Benzo(a)pyrene	22.96	252	5960440	51.446	nq #	96
90) Dibenzo(a,h)anthracene	25.10	278	4756265	44.284	nq #	91
91) Benzo(g,h,i)perylene	25.71	276	4704031	44.602	nq #	85

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

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