

Data Path : U:\HPCHEM1\BNA M\DATA\BM082217\
 Data File : BM011333.D
 Acq On : 22 Aug 2017 18:45
 Operator : SJ/JU
 Sample : I4892-03MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 HR-06-081817MSD

Manual Integrations
 APPROVED

Sohil
 8/23/2017 7:21:32 PM

Quant Time: Aug 23 03:34:53 2017
 Quant Method : Z:\HPCHEM1\BNA M\Methods\8270-BM072617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 16 02:25:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.27	152	135770	20.00	ng	-0.02
21) Naphthalene-d8	10.02	136	501050	20.00	ng	-0.02
38) Acenaphthene-d10	13.93	164	365222	20.00	ng	-0.02
63) Phenanthrene-d10	16.70	188	993367	20.00	ng	-0.02
75) Chrysene-d12	20.93	240	1333867	20.00	ng	-0.02
86) Perylene-d12	22.99	264	1176675	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	4.91	112	1050708	130.82	ng	-0.02
7) Phenol-d6	6.48	99	1417254	124.20	ng	-0.02
23) Nitrobenzene-d5	8.42	82	1491654	111.73	ng	-0.02
41) 2,4,6-Tribromophenol	15.45	330	806522	137.78	ng	-0.02
44) 2-Fluorobiphenyl	12.54	172	2555475	87.75	ng	-0.02
78) Terphenyl-d14	19.37	244	5025838	74.64	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.92	88	155428	43.101	ng	# 100
3) Pyridine	3.29	79	339297	34.446	ng	# 88
4) n-Nitrosodimethylamine	3.20	42	268122	53.415	ng	# 67
6) Aniline	6.62	93	330023	22.946	ng	# 85
8) 2-Chlorophenol	6.84	128	328731	40.314	ng	# 61
9) Benzaldehyde	6.43	77	318321	43.113	ng	# 67
10) Phenol	6.50	94	486627	39.261	ng	# 89
11) bis(2-Chloroethyl)ether	6.72	93	435206	45.064	ng	# 82
12) 1,3-Dichlorobenzene	7.16	146	378036	38.041	ng	# 74
13) 1,4-Dichlorobenzene	7.30	146	382883	37.588	ng	# 89
14) 1,2-Dichlorobenzene	7.61	146	363058	37.277	ng	# 87
15) Benzyl Alcohol	7.52	79	554819	50.681	ng	# 75
16) 2,2'-oxybis(1-Chloropropan	7.80	45	466591	53.690	ng	# 79
17) 2-Methylphenol	7.73	107	346965	43.800	ng	# 69
18) Hexachloroethane	8.32	117	175117	39.421	ng	# 87
19) n-Nitroso-di-n-propylamine	8.08	70	510922	52.687	ng	# 89
20) 3+4-Methylphenols	8.05	107	448309	40.712	ng	# 76
22) Acetophenone	8.09	105	693274	46.159	ng	# 86
24) Nitrobenzene	8.46	77	738119	53.356	ng	# 80
25) Isophorone	8.98	82	1188223	53.382	ng	# 81
26) 2-Nitrophenol	9.16	139	182434	41.446	ng	# 1
27) 2,4-Dimethylphenol	9.23	122	331396	42.767	ng	# 61
28) bis(2-Chloroethoxy)methane	9.47	93	605636	48.784	ng	# 97
29) 2,4-Dichlorophenol	9.69	162	375973	43.179	ng	# 85
30) 1,2,4-Trichlorobenzene	9.89	180	466302	43.312	ng	# 96
31) Naphthalene	10.07	128	1098651	43.587	ng	# 100
32) Benzoic acid	9.39	122	208466	33.470	ng	# 34
33) 4-Chloroaniline	10.21	127	148354	14.754	ng	# 60
34) Hexachlorobutadiene	10.36	225	372000	43.870	ng	# 93
35) Caprolactam	11.00	113	87681m	35.509	ng	#
36) 4-Chloro-3-methylphenol	11.34	107	490235	43.602	ng	# 73
37) 2-Methylnaphthalene	11.70	142	860441	46.468	ng	# 86
39) 1,2,4,5-Tetrachlorobenzene	12.09	216	647026	41.235	ng	# 100
40) Hexachlorocyclopentadiene	12.06	237	825476	90.284	ng	# 95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.34	196	391324	41.204	nq	96
43) 2,4,5-Trichlorophenol	12.42	196	422171	43.159	nq #	83
45) 1,1'-Biphenyl	12.75	154	1212729	38.402	nq	92
46) 2-Chloronaphthalene	12.79	162	972042	41.776	nq #	92
47) 2-Nitroaniline	13.02	65	474262	57.615	nq #	61
48) Acenaphthylene	13.65	152	1439157	41.445	nq	99
49) Dimethylphthalate	13.42	163	1262515	40.409	nq #	99
50) 2,6-Dinitrotoluene	13.53	165	271286	42.939	nq #	36
51) Acenaphthene	14.00	154	912470	42.439	nq	99
52) 3-Nitroaniline	13.86	138	213205	38.888	nq #	8
53) 2,4-Dinitrophenol	14.07	184	371592	82.758	nq #	73
54) Dibenzofuran	14.34	168	1609432	46.197	nq #	91
55) 4-Nitrophenol	14.20	139	328236	68.143	nq #	75
56) 2,4-Dinitrotoluene	14.33	165	399947	45.092	nq #	64
57) Fluorene	15.00	166	1310076	43.191	nq	99
58) 2,3,4,6-Tetrachlorophenol	14.58	232	449549	49.032	nq #	100
59) Diethylphthalate	14.80	149	1354141	42.439	nq	99
60) 4-Chlorophenyl-phenylether	15.00	204	802193	43.804	nq	97
61) 4-Nitroaniline	15.04	138	210507	36.148	nq #	1
62) Azobenzene	15.30	77	2042542	59.812	nq	85
64) 4,6-Dinitro-2-methylphenol	15.10	198	270218	40.181	nq	91
65) n-Nitrosodiphenylamine	15.23	169	1173517	41.279	nq	98
66) 4-Bromophenyl-phenylether	15.90	248	540450	42.321	nq #	92
67) Hexachlorobenzene	16.00	284	549811	38.101	nq #	87
68) Atrazine	16.20	200	503555	44.200	nq	94
69) Pentachlorophenol	16.36	266	740570	80.836	nq	97
70) Phenanthrene	16.74	178	2357052	45.315	nq	99
71) Anthracene	16.83	178	2412989	46.515	nq	99
72) Carbazole	17.11	167	2000416	44.095	nq	99
73) Di-n-butylphthalate	17.70	149	2391094	43.281	nq #	95
74) Fluoranthene	18.77	202	3185411	49.418	nq	98
76) Benzidine	18.99	184	779469	24.427	nq	98
77) Pyrene	19.14	202	3294454	41.567	nq	99
79) Butylbenzylphthalate	20.09	149	1184292	42.021	nq #	55
80) Benzo(a)anthracene	20.91	228	3302977	43.367	nq	99
81) 3,3'-Dichlorobenzidine	20.86	252	897738	30.046	nq #	96
82) Chrysene	20.96	228	3113957	42.593	nq	99
83) Bis(2-ethylhexyl)phthalate	20.88	149	1674882	40.396	nq #	98
84) Di-n-octyl phthalate	21.72	149	2830832	42.068	nq #	94
85) Indeno(1,2,3-cd)pyrene	25.00	276	3375130	44.576	nq #	96
87) Benzo(b)fluoranthene	22.39	252	3327243	44.057	nq #	92
88) Benzo(k)fluoranthene	22.43	252	3086608	42.643	nq #	94
89) Benzo(a)pyrene	22.90	252	3066564	44.875	nq #	96
90) Dibenzo(a,h)anthracene	25.02	278	2776037	43.821	nq #	90
91) Benzo(g,h,i)perylene	25.61	276	2813505	45.228	nq #	85

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

