

Data Path : Z:\HPCHEM1\BNA M\DATA\BM072017\
 Data File : BM010908.D
 Acq On : 20 Jul 2017 14:47
 Operator : SJ/JU
 Sample : IDOC-W-2
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 IDOC-W-2

Quant Time: Jul 21 03:16:13 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	310715	20.00	ng	-0.01
21) Naphthalene-d8	10.20	136	1208799	20.00	ng	-0.01
38) Acenaphthene-d10	14.09	164	746264	20.00	ng	0.00
63) Phenanthrene-d10	16.84	188	1652178	20.00	ng	0.00
75) Chrysene-d12	21.06	240	1445038	20.00	ng	0.00
86) Perylene-d12	23.19	264	1141087	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.08	112	2495093	133.02	ng	0.00
7) Phenol-d6	6.65	99	3275974	126.69	ng	0.00
23) Nitrobenzene-d5	8.59	82	2467432	78.46	ng	0.00
41) 2,4,6-Tribromophenol	15.59	330	1379685	120.98	ng	0.00
44) 2-Fluorobiphenyl	12.70	172	5052529	84.23	ng	0.00
78) Terphenyl-d14	19.50	244	6088069	81.49	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.10	88	251576	30.396	ng	# 100
3) Pyridine	3.46	79	751761	33.229	ng	# 87
4) n-Nitrosodimethylamine	3.38	42	441152	38.154	ng	# 70
6) Aniline	6.79	93	1090046	33.650	ng	95
8) 2-Chlorophenol	7.02	128	756324	40.725	ng	91
9) Benzaldehyde	6.60	77	443137	24.680	ng	85
10) Phenol	6.67	94	1112317	40.684	ng	95
11) bis(2-Chloroethyl)ether	6.89	93	827417	39.294	ng	95
12) 1,3-Dichlorobenzene	7.33	146	868146	37.826	ng	# 86
13) 1,4-Dichlorobenzene	7.47	146	878540	37.093	ng	# 91
14) 1,2-Dichlorobenzene	7.79	146	838762	37.164	ng	# 90
15) Benzyl Alcohol	7.69	79	981749	40.095	ng	# 84
16) 2,2'-oxybis(1-Chloropropan	7.97	45	731055	40.045	ng	75
17) 2-Methylphenol	7.89	107	724007	40.946	ng	# 80
18) Hexachloroethane	8.50	117	377129	37.962	ng	# 79
19) n-Nitroso-di-n-propylamine	8.25	70	766262	37.408	ng	# 93
20) 3+4-Methylphenols	8.22	107	978865	39.846	ng	85
22) Acetophenone	8.26	105	1366578	37.548	ng	# 95
24) Nitrobenzene	8.63	77	1228706	38.123	ng	# 89
25) Isophorone	9.15	82	2006594	38.948	ng	# 87
26) 2-Nitrophenol	9.33	139	421045	39.260	ng	# 32
27) 2,4-Dimethylphenol	9.40	122	725250	37.889	ng	# 66
28) bis(2-Chloroethoxy)methane	9.64	93	1144083	39.614	ng	97
29) 2,4-Dichlorophenol	9.86	162	833577	38.552	ng	95
30) 1,2,4-Trichlorobenzene	10.07	180	994305	37.615	ng	98
31) Naphthalene	10.25	128	2393926	38.949	ng	99
32) Benzoic acid	9.55	122	533032	35.492	ng	# 66
33) 4-Chloroaniline	10.37	127	316278	12.772	ng	# 79
34) Hexachlorobutadiene	10.54	225	760995	37.102	ng	95
35) Caprolactam	11.16	113	214506	37.263	ng	# 83
36) 4-Chloro-3-methylphenol	11.50	107	966921	36.969	ng	# 76
37) 2-Methylnaphthalene	11.87	142	1784385	40.419	ng	# 87
39) 1,2,4,5-Tetrachlorobenzene	12.26	216	1244648	38.263	ng	# 100
40) Hexachlorocyclopentadiene	12.23	237	1980417	106.203	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.50	196	759791	37.933	nq	96
43) 2,4,5-Trichlorophenol	12.57	196	797749	40.341	nq #	88
45) 1,1'-Biphenyl	12.92	154	2394004	36.823	nq	98
46) 2-Chloronaphthalene	12.95	162	1942553	40.749	nq	93
47) 2-Nitroaniline	13.17	65	664638	42.412	nq #	73
48) Acenaphthylene	13.81	152	2824749	39.840	nq	99
49) Dimethylphthalate	13.57	163	2369837	37.328	nq #	99
50) 2,6-Dinitrotoluene	13.68	165	502713	40.369	nq #	63
51) Acenaphthene	14.16	154	1688479	38.973	nq	97
52) 3-Nitroaniline	14.01	138	338668	29.916	nq #	39
53) 2,4-Dinitrophenol	14.22	184	659252	76.583	nq	93
54) Dibenzofuran	14.50	168	2944274	42.012	nq #	91
55) 4-Nitrophenol	14.33	139	687632	69.938	nq #	1
56) 2,4-Dinitrotoluene	14.47	165	678453	39.263	nq	94
57) Fluorene	15.15	166	2325939	38.608	nq	98
58) 2,3,4,6-Tetrachlorophenol	14.73	232	790221	43.006	nq #	100
59) Diethylphthalate	14.94	149	2349789	37.477	nq	100
60) 4-Chlorophenyl-phenylether	15.15	204	1408723	38.066	nq	98
61) 4-Nitroaniline	15.18	138	419077	35.417	nq #	1
62) Azobenzene	15.44	77	2783285	43.715	nq	88
64) 4,6-Dinitro-2-methylphenol	15.24	198	425951	39.097	nq	93
65) n-Nitrosodiphenylamine	15.37	169	1956171	41.378	nq	98
66) 4-Bromophenyl-phenylether	16.05	248	921680	44.018	nq #	93
67) Hexachlorobenzene	16.16	284	935526	39.074	nq #	92
68) Atrazine	16.34	200	773460	39.506	nq	98
69) Pentachlorophenol	16.50	266	1144644	74.576	nq	99
70) Phenanthrene	16.89	178	3575108	41.595	nq	99
71) Anthracene	16.98	178	3568850	41.830	nq	98
72) Carbazole	17.26	167	2838945	37.781	nq	99
73) Di-n-butylphthalate	17.84	149	3479803	39.143	nq #	96
74) Fluoranthene	18.92	202	4017264	37.739	nq	99
76) Benzidine	19.12	184	1862179	48.156	nq	99
77) Pyrene	19.28	202	4020059	47.989	nq	99
79) Butylbenzylphthalate	20.22	149	1306261	44.539	nq #	71
80) Benzo(a)anthracene	21.04	228	3464174	41.372	nq	99
81) 3,3'-Dichlorobenzidine	20.99	252	986596	29.967	nq #	96
82) Chrysene	21.10	228	3204679	40.187	nq	99
83) Bis(2-ethylhexyl)phthalate	21.00	149	1794183	42.009	nq	99
84) Di-n-octyl phthalate	21.86	149	2715696	38.809	nq	100
85) Indeno(1,2,3-cd)pyrene	25.30	276	3048053	33.334	nq #	100
87) Benzo(b)fluoranthene	22.56	252	3034946	41.871	nq #	92
88) Benzo(k)fluoranthene	22.60	252	2925330	42.254	nq #	95
89) Benzo(a)pyrene	23.10	252	2790766	42.300	nq #	96
90) Dibenzo(a,h)anthracene	25.30	278	2539022	39.651	nq #	93
91) Benzo(g,h,i)perylene	25.93	276	2510076	39.876	nq #	83

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

