

Data Path : Z:\HPCHEM1\BNA M\DATA\BM083117\
 Data File : BM011396.D
 Acq On : 01 Sep 2017 05:46
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Sep 01 08:20:55 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM083117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 31 18:39:59 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	478958	20.00	ng	0.00
21) Naphthalene-d8	10.80	136	2266032	20.00	ng	0.00
38) Acenaphthene-d10	14.62	164	1331131	20.00	ng	0.00
63) Phenanthrene-d10	17.36	188	3011876	20.00	ng	0.00
75) Chrysene-d12	21.52	240	3590477	20.00	ng	-0.01
86) Perylene-d12	23.90	264	3431550	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.54	112	2188375	76.91	ng	0.00
7) Phenol-d6	7.14	99	3177818	80.54	ng	0.00
23) Nitrobenzene-d5	9.15	82	2805247	81.58	ng	0.00
41) 2,4,6-Tribromophenol	16.10	330	1200539	77.95	ng	0.00
44) 2-Fluorobiphenyl	13.25	172	7302188	79.08	ng	0.00
78) Terphenyl-d14	19.97	244	10929394	75.53	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.43	88	451220	38.048	ng	# 100
3) Pyridine	3.84	79	1419583	39.088	ng	# 84
4) n-Nitrosodimethylamine	3.75	42	749852	40.437	ng	# 68
6) Aniline	7.31	93	2052199	38.298	ng	95
8) 2-Chlorophenol	7.55	128	1311807	39.106	ng	91
9) Benzaldehyde	7.13	77	877986	38.282	ng	98
10) Phenol	7.16	94	1706976	39.354	ng	# 75
11) bis(2-Chloroethyl)ether	7.40	93	1340221	38.659	ng	# 83
12) 1,3-Dichlorobenzene	7.87	146	1461615	38.264	ng	97
13) 1,4-Dichlorobenzene	8.02	146	1497734	38.556	ng	97
14) 1,2-Dichlorobenzene	8.35	146	1460416	38.842	ng	97
15) Benzyl Alcohol	8.22	79	1115253	39.096	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.52	45	2291289	39.523	ng	92
17) 2-Methylphenol	8.42	107	1100321	39.618	ng	98
18) Hexachloroethane	9.07	117	515552	38.768	ng	93
19) n-Nitroso-di-n-propylamine	8.80	70	1142987	39.454	ng	# 85
20) 3+4-Methylphenols	8.75	107	1527302	39.634	ng	99
22) Acetophenone	8.81	105	2110136	37.475	ng	# 90
24) Nitrobenzene	9.19	77	1502550	39.655	ng	93
25) Isophorone	9.72	82	2839422	36.852	ng	94
26) 2-Nitrophenol	9.90	139	600366	37.964	ng	# 84
27) 2,4-Dimethylphenol	9.96	122	1346107	37.469	ng	96
28) bis(2-Chloroethoxy)methane	10.20	93	1999302	38.025	ng	100
29) 2,4-Dichlorophenol	10.44	162	1302673	38.756	ng	98
30) 1,2,4-Trichlorobenzene	10.66	180	1402365	37.689	ng	99
31) Naphthalene	10.85	128	4715526	38.543	ng	100
32) Benzoic acid	10.09	122	879922	37.708	ng	100
33) 4-Chloroaniline	10.95	127	2012911	37.609	ng	# 91
34) Hexachlorobutadiene	11.13	225	787481	37.144	ng	97
35) Caprolactam	11.74	113	444999	36.761	ng	# 57
36) 4-Chloro-3-methylphenol	12.06	107	1526721	38.672	ng	91
37) 2-Methylnaphthalene	12.45	142	3326014	39.058	ng	98
39) 1,2,4,5-Tetrachlorobenzene	12.82	216	1526684	38.112	ng	# 100
40) Hexachlorocyclopentadiene	12.80	237	686463	37.382	ng	98

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42) 2,4,6-Trichlorophenol	13.05	196	1015863	38.310	ng	98
43) 2,4,5-Trichlorophenol	13.12	196	1042784	38.722	ng	# 89
45) 1,1'-Biphenyl	13.46	154	4543639	38.449	ng	97
46) 2-Chloronaphthalene	13.50	162	3341821	38.068	ng	95
47) 2-Nitroaniline	13.70	65	1080294	39.829	ng	83
48) Acenaphthylene	14.35	152	5246679	38.358	ng	100
49) Dimethylphthalate	14.07	163	3936701	37.219	ng	99
50) 2,6-Dinitrotoluene	14.19	165	792157	39.554	ng	93
51) Acenaphthene	14.69	154	3489779	39.291	ng	100
52) 3-Nitroaniline	14.52	138	1041963	40.055	ng	85
53) 2,4-Dinitrophenol	14.72	184	285137	42.911	ng	91
54) Dibenzofuran	15.02	168	4741101	38.711	ng	96
55) 4-Nitrophenol	14.81	139	798751	39.455	ng	# 48
56) 2,4-Dinitrotoluene	14.97	165	1055139	38.447	ng	# 81
57) Fluorene	15.66	166	4263729	39.932	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.24	232	862393	39.575	ng	# 100
59) Diethylphthalate	15.43	149	4091236	37.542	ng	99
60) 4-Chlorophenyl-phenylether	15.66	204	1982102	39.798	ng	96
61) 4-Nitroaniline	15.68	138	1121732	41.481	ng	84
62) Azobenzene	15.95	77	3860237	39.304	ng	93
64) 4,6-Dinitro-2-methylphenol	15.73	198	463095	41.601	ng	90
65) n-Nitrosodiphenylamine	15.87	169	3655486	38.517	ng	99
66) 4-Bromophenyl-phenylether	16.55	248	1195538	37.737	ng	# 90
67) Hexachlorobenzene	16.66	284	1381482	37.158	ng	# 88
68) Atrazine	16.81	200	1030787	37.059	ng	97
69) Pentachlorophenol	17.00	266	810950	39.732	ng	98
70) Phenanthrene	17.40	178	6630482	39.530	ng	98
71) Anthracene	17.49	178	6644904	39.397	ng	98
72) Carbazole	17.76	167	6386418	39.291	ng	99
73) Di-n-butylphthalate	18.32	149	7531220	39.882	ng	100
74) Fluoranthene	19.41	202	7774329	40.621	ng	96
76) Benzidine	19.59	184	2690771	37.971	ng	99
77) Pyrene	19.77	202	8217587	37.350	ng	98
79) Butylbenzylphthalate	20.66	149	3553386	36.558	ng	94
80) Benzo(a)anthracene	21.51	228	7715278	38.420	ng	98
81) 3,3'-Dichlorobenzidine	21.44	252	3113062	40.832	ng	99
82) Chrysene	21.56	228	7302530	38.581	ng	97
83) Bis(2-ethylhexyl)phthalate	21.43	149	5457128	38.296	ng	99
84) Di-n-octyl phthalate	22.34	149	9575338	39.862	ng	# 95
85) Indeno(1,2,3-cd)pyrene	26.36	276	7681970	43.510	ng	# 88
87) Benzo(b)fluoranthene	23.18	252	7883578	37.420	ng	99
88) Benzo(k)fluoranthene	23.23	252	7824104	38.734	ng	97
89) Benzo(a)pyrene	23.80	252	7373503	38.804	ng	# 97
90) Dibenzo(a,h)anthracene	26.37	278	6781116	41.698	ng	99
91) Benzo(g,h,i)perylene	27.11	276	6322921	40.232	ng	99

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

