

Data Path : Z:\HPCHEM1\BNA M\DATA\BM081617\
 Data File : BM011259.D
 Acq On : 16 Aug 2017 18:07
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Quant Time: Aug 17 03:53:11 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	121774	20.00	ng	-0.01
21) Naphthalene-d8	10.03	136	469594	20.00	ng	-0.01
38) Acenaphthene-d10	13.94	164	346127	20.00	ng	-0.01
63) Phenanthrene-d10	16.70	188	979059	20.00	ng	-0.01
75) Chrysene-d12	20.93	240	1290212	20.00	ng	-0.01
86) Perylene-d12	23.00	264	1124270	20.00	ng	-0.02

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System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	4.92	112	638637	88.65	ng	-0.01
7) Phenol-d6	6.48	99	938353	91.68	ng	-0.01
23) Nitrobenzene-d5	8.42	82	1319086	105.43	ng	-0.02
41) 2,4,6-Tribromophenol	15.45	330	488933	88.13	ng	-0.02
44) 2-Fluorobiphenyl	12.55	172	2347376	85.05	ng	-0.01
78) Terphenyl-d14	19.37	244	5295416	81.30	ng	-0.01

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Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.91	88	155531	48.087	ng	# 100
3) Pyridine	3.29	79	426639	48.292	ng	# 87
4) n-Nitrosodimethylamine	3.21	42	241535	53.648	ng	# 69
6) Aniline	6.62	93	559987	43.411	ng	# 84
8) 2-Chlorophenol	6.85	128	305329	41.748	ng	# 67
9) Benzaldehyde	6.44	77	282489	42.657	ng	# 72
10) Phenol	6.50	94	514669m	46.296	ng	# 87
11) bis(2-Chloroethyl)ether	6.73	93	386943	44.671	ng	# 78
12) 1,3-Dichlorobenzene	7.16	146	375922	42.176	ng	# 86
13) 1,4-Dichlorobenzene	7.31	146	384611	42.097	ng	# 91
14) 1,2-Dichlorobenzene	7.62	146	349241	39.980	ng	# 80
15) Benzyl Alcohol	7.53	79	459332	46.781	ng	# 77
16) 2,2'-oxybis(1-Chloropropan	7.82	45	386429	49.576	ng	# 74
17) 2-Methylphenol	7.73	107	323630	45.550	ng	# 82
18) Hexachloroethane	8.33	117	176051	44.186	ng	# 87
19) n-Nitroso-di-n-propylamine	8.09	70	459707	52.854	ng	# 79
20) 3+4-Methylphenols	8.06	107	438643	44.413	ng	# 89
22) Acetophenone	8.09	105	650022	46.178	ng	# 80
24) Nitrobenzene	8.46	77	679458	52.406	ng	# 84
25) Isophorone	8.99	82	1080987	51.817	ng	# 2
26) 2-Nitrophenol	9.16	139	180284	43.701	ng	# 65
27) 2,4-Dimethylphenol	9.25	122	326291	44.929	ng	# 97
28) bis(2-Chloroethoxy)methane	9.48	93	565040	48.563	ng	# 90
29) 2,4-Dichlorophenol	9.69	162	364582	44.676	ng	# 98
30) 1,2,4-Trichlorobenzene	9.90	180	457724	45.363	ng	# 97
31) Naphthalene	10.08	128	996747	42.193	ng	# 46
32) Benzoic acid	9.40	122	251876	43.149	ng	# 65
33) 4-Chloroaniline	10.21	127	372019	39.476	ng	# 99
34) Hexachlorobutadiene	10.37	225	373019	46.937	ng	# 87
35) Caprolactam	11.00	113	117468	50.758	ng	# 68
36) 4-Chloro-3-methylphenol	11.35	107	514184	48.796	ng	# 86
37) 2-Methylnaphthalene	11.70	142	766474	44.166	ng	# 100
39) 1,2,4,5-Tetrachlorobenzene	12.09	216	643477	43.271	ng	# 96
40) Hexachlorocyclopentadiene	12.07	237	364763	42.096	ng	

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.35	196	397880	44.205	nq	96
43) 2,4,5-Trichlorophenol	12.42	196	415029	44.770	nq	# 84
45) 1,1'-Biphenyl	12.76	154	1258920	42.064	nq	95
46) 2-Chloronaphthalene	12.80	162	933097	42.315	nq	94
47) 2-Nitroaniline	13.02	65	456160	58.473	nq	# 55
48) Acenaphthylene	13.66	152	1455269	44.221	nq	99
49) Dimethylphthalate	13.43	163	1290383	43.580	nq	# 99
50) 2,6-Dinitrotoluene	13.54	165	270797	45.226	nq	# 35
51) Acenaphthene	14.01	154	897135	44.027	nq	99
52) 3-Nitroaniline	13.87	138	209990	40.414	nq	# 7
53) 2,4-Dinitrophenol	14.08	184	205439	48.278	nq	# 81
54) Dibenzofuran	14.35	168	1467888	44.459	nq	# 90
55) 4-Nitrophenol	14.20	139	193319	42.348	nq	# 83
56) 2,4-Dinitrotoluene	14.34	165	401414	47.754	nq	# 71
57) Fluorene	15.00	166	1302933	45.325	nq	99
58) 2,3,4,6-Tetrachlorophenol	14.59	232	399365	45.961	nq	# 100
59) Diethylphthalate	14.81	149	1421656	47.013	nq	99
60) 4-Chlorophenyl-phenylether	15.02	204	766045	44.138	nq	99
61) 4-Nitroaniline	15.04	138	203628	36.896	nq	# 1
62) Azobenzene	15.30	77	1757693	54.311	nq	84
64) 4,6-Dinitro-2-methylphenol	15.10	198	305807	46.138	nq	94
65) n-Nitrosodiphenylamine	15.23	169	1155596	41.243	nq	98
66) 4-Bromophenyl-phenylether	15.91	248	515757	40.978	nq	# 88
67) Hexachlorobenzene	16.01	284	578553	40.679	nq	# 85
68) Atrazine	16.21	200	459982	40.966	nq	94
69) Pentachlorophenol	16.36	266	388650	43.043	nq	94
70) Phenanthrene	16.74	178	2204873	43.009	nq	100
71) Anthracene	16.84	178	2209155	43.208	nq	99
72) Carbazole	17.12	167	1993167	44.577	nq	99
73) Di-n-butylphthalate	17.71	149	2504204	45.991	nq	# 96
74) Fluoranthene	18.78	202	2982509	46.946	nq	98
76) Benzidine	19.00	184	326396m	10.575	nq	
77) Pyrene	19.14	202	3126248	40.779	nq	99
79) Butylbenzylphthalate	20.09	149	1162125	42.629	nq	# 56
80) Benzo(a)anthracene	20.92	228	3148664	42.740	nq	99
81) 3,3'-Dichlorobenzidine	20.87	252	1144198	39.591	nq	# 97
82) Chrysene	20.97	228	3000632	42.432	nq	100
83) Bis(2-ethylhexyl)phthalate	20.89	149	1678540	41.854	nq	97
84) Di-n-octyl phthalate	21.73	149	2805454	43.101	nq	# 94
85) Indeno(1,2,3-cd)pyrene	25.01	276	3039870	41.507	nq	# 96
87) Benzo(b)fluoranthene	22.40	252	3008394	41.692	nq	# 90
88) Benzo(k)fluoranthene	22.44	252	3008329	43.498	nq	# 95
89) Benzo(a)pyrene	22.91	252	2792624	42.771	nq	# 97
90) Dibenzo(a,h)anthracene	25.03	278	2526209	41.736	nq	# 90
91) Benzo(g,h,i)perylene	25.63	276	2497621	42.022	nq	# 84

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

