

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050517\
 Data File : BM009901.D
 Acq On : 05 May 2017 14:08
 Operator : SJ/MA
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

mohammad
 5/8/2017 2:49:13 PM

Quant Time: May 05 14:41:45 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM050517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 05 14:20:43 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	107990	20.00	ng	0.00
21) Naphthalene-d8	10.58	136	468269	20.00	ng	0.00
38) Acenaphthene-d10	14.44	164	292680	20.00	ng	0.00
63) Phenanthrene-d10	17.20	188	692519	20.00	ng	0.00
75) Chrysene-d12	21.39	240	881159	20.00	ng	0.00
86) Perylene-d12	23.69	264	761299	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.37	112	1141795	84.90	ng	0.00
7) Phenol-d6	6.97	99	1796202	87.54	ng	0.00
23) Nitrobenzene-d5	8.96	82	2191046	85.49	ng	0.00
41) 2,4,6-Tribromophenol	15.94	330	722366	92.35	ng	0.00
44) 2-Fluorobiphenyl	13.06	172	3922678	88.66	ng	0.00
78) Terphenyl-d14	19.82	244	7863647	93.79	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.29	88	190179	71.97	ng	# 100
3) Pyridine	3.69	79	694640	77.14	ng	# 73
4) n-Nitrosodimethylamine	3.61	42	458102	78.13	ng	# 58
6) Aniline	7.13	93	1152792	84.54	ng	# 88
8) 2-Chlorophenol	7.36	128	637650	85.23	ng	# 81
9) Benzaldehyde	6.94	77	588496	69.97	ng	# 81
10) Phenol	7.00	94	962068	85.33	ng	# 93
11) bis(2-Chloroethyl)ether	7.22	93	755075	83.93	ng	# 79
12) 1,3-Dichlorobenzene	7.67	146	688024	82.59	ng	# 85
13) 1,4-Dichlorobenzene	7.82	146	710876	82.82	ng	# 93
14) 1,2-Dichlorobenzene	8.13	146	692021	84.09	ng	# 93
15) Benzyl Alcohol	8.04	79	817610	85.78	ng	# 87
16) 2,2'-oxybis(1-Chloropropan	8.30	45	1371285	82.21	ng	# 94
17) 2-Methylphenol	8.24	107	652198	85.99	ng	# 86
18) Hexachloroethane	8.85	117	317405	83.57	ng	# 94
19) n-Nitroso-di-n-propylamine	8.60	70	831314	85.10	ng	# 85
20) 3+4-Methylphenols	8.57	107	919873	88.69	ng	# 87
22) Acetophenone	8.62	105	1251193	86.38	ng	# 86
24) Nitrobenzene	9.00	77	1114136	83.46	ng	# 89
25) Isophorone	9.53	82	1925026	85.10	ng	# 88
26) 2-Nitrophenol	9.70	139	386988	90.79	ng	# 44
27) 2,4-Dimethylphenol	9.77	122	652622	84.09	ng	# 78
28) bis(2-Chloroethoxy)methane	10.00	93	1132295	86.18	ng	# 98
29) 2,4-Dichlorophenol	10.25	162	645336	87.22	ng	# 95
30) 1,2,4-Trichlorobenzene	10.45	180	739085	85.16	ng	# 95
31) Naphthalene	10.63	128	2152941	84.25	ng	# 99
32) Benzoic acid	10.00	122	468299	93.86	ng	# 73
33) 4-Chloroaniline	10.76	127	938026	87.53	ng	# 76
34) Hexachlorobutadiene	10.89	225	511168	84.55	ng	# 99
35) Caprolactam	11.61	113	247595m	87.84	ng	# 81
36) 4-Chloro-3-methylphenol	11.90	107	859597	85.26	ng	# 92
37) 2-Methylnaphthalene	12.25	142	1543123	85.71	ng	# 100
39) 1,2,4,5-Tetrachlorobenzene	12.62	216	906221	87.17	ng	# 97
40) Hexachlorocyclopentadiene	12.58	237	275348	145.51	ng	# 97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.87	196	598009	89.37	ng	97
43) 2,4,5-Trichlorophenol	12.96	196	636031	88.15	ng #	88
45) 1,1'-Biphenyl	13.27	154	2140504	86.96	ng	95
46) 2-Chloronaphthalene	13.32	162	1703751	87.69	ng	96
47) 2-Nitroaniline	13.54	65	815016	86.71	ng #	65
48) Acenaphthylene	14.16	152	2660592	87.16	ng	100
49) Dimethylphthalate	13.91	163	2300176	82.92	ng #	98
50) 2,6-Dinitrotoluene	14.04	165	464693	86.44	ng #	43
51) Acenaphthene	14.51	154	1615038	85.50	ng	99
52) 3-Nitroaniline	14.38	138	499527	85.79	ng #	53
53) 2,4-Dinitrophenol	14.60	184	257179	108.91	ng #	79
54) Dibenzofuran	14.84	168	2499253	86.68	ng	94
55) 4-Nitrophenol	14.70	139	348050	94.11	ng #	6
56) 2,4-Dinitrotoluene	14.84	165	697994	89.88	ng #	66
57) Fluorene	15.50	166	2239648	88.29	ng	100
58) 2,3,4,6-Tetrachlorophenol	15.08	232	488823	88.52	ng #	100
59) Diethylphthalate	15.27	149	2489861	84.23	ng	98
60) 4-Chlorophenyl-phenylether	15.49	204	1196126	88.96	ng	98
61) 4-Nitroaniline	15.55	138	504063	86.54	ng #	13
62) Azobenzene	15.78	77	2807894	85.45	ng	81
64) 4,6-Dinitro-2-methylphenol	15.60	198	402980	97.00	ng #	79
65) n-Nitrosodiphenylamine	15.71	169	1919129	88.83	ng	100
66) 4-Bromophenyl-phenylether	16.38	248	720167	90.58	ng #	89
67) Hexachlorobenzene	16.50	284	792610	86.97	ng #	88
68) Atrazine	16.67	200	721174	86.84	ng	98
69) Pentachlorophenol	16.86	266	348112	104.73	ng	98
70) Phenanthrene	17.24	178	3484798	88.19	ng	99
71) Anthracene	17.33	178	3550324	89.26	ng	99
72) Carbazole	17.61	167	3139116	86.67	ng	100
73) Di-n-butylphthalate	18.15	149	4612410	89.03	ng #	97
74) Fluoranthene	19.26	202	4366564	87.45	ng	99
76) Benzidine	19.45	184	2148396	80.96	ng	100
77) Pyrene	19.63	202	4583556	88.39	ng	99
79) Butylbenzylphthalate	20.51	149	2209475	88.35	ng #	72
80) Benzo(a)anthracene	21.37	228	4518019	86.05	ng	99
81) 3,3'-Dichlorobenzidine	21.31	252	1730134	82.37	ng #	98
82) Chrysene	21.43	228	4300385	87.31	ng	99
83) Bis(2-ethylhexyl)phthalate	21.27	149	3378478	87.74	ng	98
84) Di-n-octyl phthalate	22.17	149	5791791	86.41	ng #	85
85) Indeno(1,2,3-cd)pyrene	26.07	276	4063619	71.08	ng #	100
87) Benzo(b)fluoranthene	23.00	252	4222706	88.34	ng #	95
88) Benzo(k)fluoranthene	23.05	252	4252192m	92.10	ng	
89) Benzo(a)pyrene	23.60	252	3925715	87.39	ng	99
90) Dibenzo(a,h)anthracene	26.07	278	3691548	80.98	ng #	96
91) Benzo(g,h,i)perylene	26.79	276	3226206	74.59	ng #	92

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

