

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012117\
 Data File : BG025568.D
 Acq On : 21 Jan 2017 11:23
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 1/23/2017 5:00:09 PM

Quant Time: Jan 23 02:26:22 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 10 11:26:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.18	152	64412	20.00	ng	-0.02
21) Naphthalene-d8	11.00	136	276928	20.00	ng	-0.02
38) Acenaphthene-d10	14.81	164	220516	20.00	ng	-0.02
63) Phenanthrene-d10	17.55	188	502750	20.00	ng	-0.02
75) Chrysene-d12	21.85	240	736192	20.00	ng	-0.02
86) Perylene-d12	25.22	264	804611	20.00	ng	-0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.72	112	302078	85.24	ng	-0.02
7) Phenol-d6	7.34	99	434150	86.99	ng	-0.02
23) Nitrobenzene-d5	9.36	82	524454	83.66	ng	-0.02
41) 2,4,6-Tribromophenol	16.30	330	270596	76.22	ng	-0.02
44) 2-Fluorobiphenyl	13.43	172	1107576	81.32	ng	-0.02
78) Terphenyl-d14	20.14	244	2068110	74.48	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.55	88	69945	46.26	ng	91
3) Pyridine	3.97	79	193247	44.09	ng	97
4) n-Nitrosodimethylamine	3.88	42	115653	44.45	ng	# 90
6) Aniline	7.50	93	292069	43.19	ng	98
8) 2-Chlorophenol	7.74	128	165167	41.32	ng	96
9) Benzaldehyde	7.32	77	140823	38.56	ng	95
10) Phenol	7.37	94	236087	42.67	ng	98
11) bis(2-Chloroethyl)ether	7.59	93	176695	41.45	ng	93
12) 1,3-Dichlorobenzene	8.06	146	198071	40.95	ng	99
13) 1,4-Dichlorobenzene	8.21	146	208560	41.60	ng	98
14) 1,2-Dichlorobenzene	8.54	146	194128	40.58	ng	93
15) Benzyl Alcohol	8.43	79	213684	44.85	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.70	45	364584	46.18	ng	97
17) 2-Methylphenol	8.63	107	158888	43.73	ng	92
18) Hexachloroethane	9.26	117	80583	41.85	ng	94
19) n-Nitroso-di-n-propylamine	8.98	70	186017	44.13	ng	92
20) 3+4-Methylphenols	8.96	107	220998	43.95	ng	97
22) Acetophenone	9.01	105	313610	40.76	ng	# 95
24) Nitrobenzene	9.41	77	285890	41.58	ng	97
25) Isophorone	9.92	82	486884	42.90	ng	# 96
26) 2-Nitrophenol	10.12	139	103659	39.42	ng	89
27) 2,4-Dimethylphenol	10.16	122	173967	40.24	ng	100
28) bis(2-Chloroethoxy)methane	10.39	93	241401	40.96	ng	98
29) 2,4-Dichlorophenol	10.66	162	192438	41.60	ng	94
30) 1,2,4-Trichlorobenzene	10.86	180	222195	39.76	ng	95
31) Naphthalene	11.06	128	568970	41.14	ng	98
32) Benzoic acid	10.32	122	120298	34.60	ng	92
33) 4-Chloroaniline	11.18	127	245907	42.28	ng	# 95
34) Hexachlorobutadiene	11.32	225	173767	38.09	ng	95
35) Caprolactam	11.94	113	67764	38.98	ng	# 88
36) 4-Chloro-3-methylphenol	12.29	107	226055	39.59	ng	97
37) 2-Methylnaphthalene	12.65	142	430544	42.01	ng	98
39) 1,2,4,5-Tetrachlorobenzene	13.01	216	299144	41.09	ng	97
40) Hexachlorocyclopentadiene	12.97	237	98046	41.46	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.26	196	182413	39.79	nq	95
43) 2,4,5-Trichlorophenol	13.34	196	184649	38.48	nq	94
45) 1,1'-Biphenyl	13.64	154	607188	40.94	nq	97
46) 2-Chloronaphthalene	13.70	162	477325	41.19	nq	96
47) 2-Nitroaniline	13.91	65	215188	44.00	nq	97
48) Acenaphthylene	14.54	152	742916	41.37	nq	95
49) Dimethylphthalate	14.25	163	655171	40.75	nq	96
50) 2,6-Dinitrotoluene	14.40	165	146633	41.75	nq	96
51) Acenaphthene	14.87	154	479166	40.79	nq	98
52) 3-Nitroaniline	14.73	138	144102	42.68	nq	92
53) 2,4-Dinitrophenol	14.96	184	86395	40.21	nq	93
54) Dibenzofuran	15.21	168	753033	40.55	nq	96
55) 4-Nitrophenol	15.06	139	98092	38.90	nq	91
56) 2,4-Dinitrotoluene	15.19	165	208377	40.75	nq	# 89
57) Fluorene	15.85	166	634631	40.82	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.43	232	188074	36.59	nq	97
59) Diethylphthalate	15.60	149	692368	41.07	nq	98
60) 4-Chlorophenyl-phenylether	15.83	204	352193	39.97	nq	98
61) 4-Nitroaniline	15.90	138	146086	43.00	nq	# 80
62) Azobenzene	16.13	77	707315	43.51	nq	98
64) 4,6-Dinitro-2-methylphenol	15.95	198	145592	41.82	nq	90
65) n-Nitrosodiphenylamine	16.05	169	569308	41.89	nq	97
66) 4-Bromophenyl-phenylether	16.73	248	259187	41.79	nq	97
67) Hexachlorobenzene	16.86	284	289277	40.64	nq	# 90
68) Atrazine	16.99	200	193704	32.53	nq	100
69) Pentachlorophenol	17.21	266	127432	34.54	nq	93
70) Phenanthrene	17.60	178	1065568	41.13	nq	95
71) Anthracene	17.69	178	1083925	41.18	nq	98
72) Carbazole	17.97	167	989146	42.01	nq	95
73) Di-n-butylphthalate	18.48	149	1211983	41.84	nq	96
74) Fluoranthene	19.59	202	1472733	43.04	nq	95
76) Benzidine	19.78	184	578162	31.48	nq	98
77) Pyrene	19.96	202	1503934	39.09	nq	96
79) Butylbenzylphthalate	20.81	149	625127	40.46	nq	98
80) Benzo(a)anthracene	21.82	228	1662015	40.47	nq	98
81) 3,3'-Dichlorobenzidine	21.73	252	666040	41.76	nq	100
82) Chrysene	21.89	228	1594133	40.48	nq	97
83) Bis(2-ethylhexyl)phthalate	21.67	149	910656	42.36	nq	# 97
84) Di-n-octyl phthalate	22.91	149	1561619	43.75	nq	97
85) Indeno(1,2,3-cd)pyrene	29.13	276	2186358	43.68	nq	# 96
87) Benzo(b)fluoranthene	24.14	252	1792568m	40.83	nq	
88) Benzo(k)fluoranthene	24.21	252	1721656	40.61	nq	98
89) Benzo(a)pyrene	25.07	252	1742384	41.09	nq	97
90) Dibenzo(a,h)anthracene	29.17	278	1865521	41.51	nq	98
91) Benzo(g,h,i)perylene	30.36	276	1841550	41.24	nq	99

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

