

Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\
 Data File : BG032565.D
 Acq On : 6 Feb 2018 13:48
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 2/7/2018 4:14:23 PM

Quant Time: Feb 07 03:15:12 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 18:57:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.71	152	17938	20.00	ng	0.00
21) Naphthalene-d8	11.58	136	75892	20.00	ng	0.00
38) Acenaphthene-d10	15.34	164	63017	20.00	ng	0.00
63) Phenanthrene-d10	18.08	188	163630	20.00	ng	0.00
75) Chrysene-d12	22.41	240	194636	20.00	ng	0.00
86) Perylene-d12	26.07	264	233489	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.14	112	68990	77.25	ng	0.00
7) Phenol-d6	7.82	99	92856	77.92	ng	0.00
23) Nitrobenzene-d5	9.91	82	89198	73.41	ng	0.00
41) 2,4,6-Tribromophenol	16.82	330	106279	88.60	ng	0.00
44) 2-Fluorobiphenyl	13.97	172	436027	82.24	ng	0.00
78) Terphenyl-d14	20.61	244	748625	82.39	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.81	88	13361	44.466	ng	# 73
3) Pyridine	4.26	79	38011	46.849	ng	# 74
4) n-Nitrosodimethylamine	4.17	42	19656m	46.539	ng	
6) Aniline	8.01	93	57252	38.531	ng	93
8) 2-Chlorophenol	8.26	128	43700	41.395	ng	# 81
9) Benzaldehyde	7.81	77	22344m	36.301	ng	
10) Phenol	7.85	94	45731	40.408	ng	95
11) bis(2-Chloroethyl)ether	8.10	93	34991	40.578	ng	# 69
12) 1,3-Dichlorobenzene	8.59	146	59623	41.763	ng	95
13) 1,4-Dichlorobenzene	8.75	146	55460	38.758	ng	98
14) 1,2-Dichlorobenzene	9.07	146	56789	40.513	ng	97
15) Benzyl Alcohol	8.95	79	34724	35.354	ng	92
16) 2,2'-oxybis(1-Chloropropan	9.23	45	31519	38.573	ng	61
17) 2-Methylphenol	9.15	107	35069	38.153	ng	94
18) Hexachloroethane	9.83	117	19809	41.232	ng	85
19) n-Nitroso-di-n-propylamine	9.52	70	28703	36.325	ng	# 87
20) 3+4-Methylphenols	9.49	107	51345	39.823	ng	88
22) Acetophenone	9.55	105	66947	37.632	ng	# 88
24) Nitrobenzene	9.95	77	42890	36.505	ng	# 86
25) Isophorone	10.47	82	79308	38.334	ng	# 83
26) 2-Nitrophenol	10.68	139	29199	41.112	ng	# 78
27) 2,4-Dimethylphenol	10.71	122	37926	37.204	ng	95
28) bis(2-Chloroethoxy)methane	10.95	93	45007	39.671	ng	98
29) 2,4-Dichlorophenol	11.22	162	49799	36.876	ng	88
30) 1,2,4-Trichlorobenzene	11.44	180	67884	37.716	ng	92
31) Naphthalene	11.64	128	153672	41.459	ng	99
32) Benzoic acid	10.85	122	30534	41.342	ng	# 81
33) 4-Chloroaniline	11.74	127	65360	39.535	ng	92
34) Hexachlorobutadiene	11.91	225	53704	38.350	ng	93
35) Caprolactam	12.49	113	15170	39.131	ng	# 32
36) 4-Chloro-3-methylphenol	12.82	107	57050	44.537	ng	# 90
37) 2-Methylnaphthalene	13.20	142	146934	47.185	ng	97
39) 1,2,4,5-Tetrachlorobenzene	13.56	216	116173	39.406	ng	# 97
40) Hexachlorocyclopentadiene	13.53	237	70555	38.303	ng	90

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.79	196	53867	33.591	nq	96
43) 2,4,5-Trichlorophenol	13.87	196	74545	39.877	nq #	91
45) 1,1'-Biphenyl	14.18	154	220778	41.468	nq	96
46) 2-Chloronaphthalene	14.23	162	171539	41.108	nq	98
47) 2-Nitroaniline	14.43	65	30514m	33.067	nq	
48) Acenaphthylene	15.07	152	263163	41.717	nq	98
49) Dimethylphthalate	14.77	163	215039	37.009	nq	98
50) 2,6-Dinitrotoluene	14.90	165	50146	41.119	nq #	75
51) Acenaphthene	15.41	154	181038	41.380	nq	96
52) 3-Nitroaniline	15.24	138	43402	40.744	nq #	73
53) 2,4-Dinitrophenol	15.44	184	34500	47.099	nq #	58
54) Dibenzofuran	15.74	168	265604	41.017	nq	98
55) 4-Nitrophenol	15.53	139	29098	36.487	nq #	76
56) 2,4-Dinitrotoluene	15.68	165	70010	40.319	nq #	65
57) Fluorene	16.38	166	216769	40.285	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.95	232	67955	36.858	nq #	89
59) Diethylphthalate	16.12	149	230981	40.816	nq	96
60) 4-Chlorophenyl-phenylether	16.36	204	130521	38.466	nq	96
61) 4-Nitroaniline	16.40	138	47522	41.623	nq #	79
62) Azobenzene	16.65	77	138177	40.858	nq	85
64) 4,6-Dinitro-2-methylphenol	16.44	198	46645	38.158	nq #	66
65) n-Nitrosodiphenylamine	16.57	169	198364	40.621	nq	94
66) 4-Bromophenyl-phenylether	17.25	248	99385	40.998	nq	95
67) Hexachlorobenzene	17.38	284	113132	42.171	nq #	89
68) Atrazine	17.49	200	83382	39.776	nq	97
69) Pentachlorophenol	17.72	266	64896	38.616	nq	95
70) Phenanthrene	18.12	178	365721	40.078	nq	98
71) Anthracene	18.21	178	380088	41.832	nq	97
72) Carbazole	18.47	167	311304	38.777	nq	97
73) Di-n-butylphthalate	18.97	149	384124	43.217	nq	99
74) Fluoranthene	20.09	202	461759	38.586	nq	95
76) Benzidine	20.24	184	170188	44.530	nq	99
77) Pyrene	20.44	202	483744	40.522	nq	99
79) Butylbenzylphthalate	21.30	149	166470	44.210	nq #	78
80) Benzo(a)anthracene	22.38	228	477474	39.571	nq	100
81) 3,3'-Dichlorobenzidine	22.28	252	194197	41.539	nq #	96
82) Chrysene	22.45	228	472538	40.551	nq	97
83) Bis(2-ethylhexyl)phthalate	22.22	149	237499	44.483	nq #	97
84) Di-n-octyl phthalate	23.58	149	383881	45.331	nq #	89
85) Indeno(1,2,3-cd)pyrene	30.32	276	704912	47.819	nq #	85
87) Benzo(b)fluoranthene	24.90	252	514142	36.443	nq #	95
88) Benzo(k)fluoranthene	24.97	252	517168	37.005	nq #	97
89) Benzo(a)pyrene	25.90	252	523160	38.897	nq #	96
90) Dibenzo(a,h)anthracene	30.40	278	592254	43.045	nq #	94
91) Benzo(g,h,i)perylene	31.66	276	576451	42.210	nq #	91

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

