

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121117\
 Data File : BG031468.D
 Acq On : 11 Dec 2017 23:37
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 12/12/2017 3:55:19 PM

Quant Time: Dec 12 04:41:52 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 11 17:25:55 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	52747	20.00	ng	0.00
21) Naphthalene-d8	10.92	136	245797	20.00	ng	0.00
38) Acenaphthene-d10	14.73	164	174091	20.00	ng	0.00
63) Phenanthrene-d10	17.47	188	439206	20.00	ng	0.00
75) Chrysene-d12	21.74	240	477909	20.00	ng	0.00
86) Perylene-d12	25.02	264	474856	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.67	112	239928	80.63	ng	0.00
7) Phenol-d6	7.27	99	322578	78.08	ng	0.00
23) Nitrobenzene-d5	9.27	82	300192	75.28	ng	0.00
41) 2,4,6-Tribromophenol	16.22	330	164759	80.78	ng	0.00
44) 2-Fluorobiphenyl	13.35	172	880166	74.21	ng	0.00
78) Terphenyl-d14	20.06	244	1579475	75.19	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	52055	36.662	ng	85
3) Pyridine	3.92	79	145689	38.971	ng	89
4) n-Nitrosodimethylamine	3.84	42	66751	37.908	ng	# 91
6) Aniline	7.43	93	208485	38.321	ng	95
8) 2-Chlorophenol	7.67	128	130457	39.302	ng	90
9) Benzaldehyde	7.25	77	89087	37.204	ng	86
10) Phenol	7.30	94	168259	39.648	ng	90
11) bis(2-Chloroethyl)ether	7.52	93	133850	38.643	ng	89
12) 1,3-Dichlorobenzene	8.00	146	151184	38.300	ng	96
13) 1,4-Dichlorobenzene	8.15	146	154474	37.654	ng	96
14) 1,2-Dichlorobenzene	8.46	146	148462	37.989	ng	96
15) Benzyl Alcohol	8.35	79	116484	36.045	ng	91
16) 2,2'-oxybis(1-Chloropropan	8.63	45	169858m	43.651	ng	
17) 2-Methylphenol	8.56	107	113730	39.314	ng	97
18) Hexachloroethane	9.19	117	53257	38.512	ng	96
19) n-Nitroso-di-n-propylamine	8.91	70	117189	36.041	ng	# 95
20) 3+4-Methylphenols	8.89	107	158605	36.804	ng	94
22) Acetophenone	8.93	105	227070	38.508	ng	# 93
24) Nitrobenzene	9.32	77	154299	37.759	ng	92
25) Isophorone	9.84	82	286846	38.061	ng	# 92
26) 2-Nitrophenol	10.03	139	78631	38.914	ng	89
27) 2,4-Dimethylphenol	10.09	122	119862	39.049	ng	92
28) bis(2-Chloroethoxy)methane	10.31	93	190066	39.064	ng	96
29) 2,4-Dichlorophenol	10.58	162	145189	40.147	ng	89
30) 1,2,4-Trichlorobenzene	10.78	180	175072	37.898	ng	98
31) Naphthalene	10.97	128	450185	37.281	ng	96
32) Benzoic acid	10.25	122	53916	34.628	ng	87
33) 4-Chloroaniline	11.08	127	200134	38.171	ng	93
34) Hexachlorobutadiene	11.25	225	115635	37.737	ng	97
35) Caprolactam	11.85	113	55474	39.064	ng	# 74
36) 4-Chloro-3-methylphenol	12.21	107	155928	37.773	ng	# 86
37) 2-Methylnaphthalene	12.57	142	351009	37.543	ng	93
39) 1,2,4,5-Tetrachlorobenzene	12.93	216	254106	37.347	ng	# 96
40) Hexachlorocyclopentadiene	12.91	237	53995	34.751	ng	90

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.18	196	136902	39.383	nq	99
43) 2,4,5-Trichlorophenol	13.26	196	149935	40.041	nq #	94
45) 1,1'-Biphenyl	13.56	154	532662	37.252	nq	99
46) 2-Chloronaphthalene	13.62	162	387999	37.074	nq	97
47) 2-Nitroaniline	13.82	65	108339	36.840	nq #	84
48) Acenaphthylene	14.46	152	630884	37.464	nq	98
49) Dimethylphthalate	14.18	163	540651	37.512	nq	99
50) 2,6-Dinitrotoluene	14.30	165	116621	37.855	nq #	82
51) Acenaphthene	14.80	154	398373	37.413	nq	98
52) 3-Nitroaniline	14.64	138	121347	38.641	nq #	80
53) 2,4-Dinitrophenol	14.86	184	34576	32.903	nq #	50
54) Dibenzofuran	15.13	168	627114	36.908	nq	98
55) 4-Nitrophenol	14.97	139	75850	37.705	nq #	83
56) 2,4-Dinitrotoluene	15.10	165	166288	38.003	nq #	76
57) Fluorene	15.78	166	507641	37.287	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.36	232	139189	39.545	nq #	92
59) Diethylphthalate	15.53	149	557442	38.585	nq	98
60) 4-Chlorophenyl-phenylether	15.76	204	308541	37.042	nq	93
61) 4-Nitroaniline	15.80	138	127987	38.836	nq	90
62) Azobenzene	16.05	77	445081	37.191	nq	95
64) 4,6-Dinitro-2-methylphenol	15.87	198	94322	36.761	nq	79
65) n-Nitrosodiphenylamine	15.98	169	498520	38.778	nq	97
66) 4-Bromophenyl-phenylether	16.65	248	198486	38.864	nq	97
67) Hexachlorobenzene	16.78	284	200624	38.042	nq	96
68) Atrazine	16.92	200	190521	40.531	nq	98
69) Pentachlorophenol	17.13	266	84268	36.550	nq	98
70) Phenanthrene	17.52	178	865344	37.725	nq	99
71) Anthracene	17.61	178	859291	37.878	nq	100
72) Carbazole	17.88	167	799280	38.933	nq	99
73) Di-n-butylphthalate	18.41	149	950298	40.083	nq	100
74) Fluoranthene	19.52	202	1108833	38.627	nq	98
76) Benzidine	19.69	184	441595	39.705	nq	98
77) Pyrene	19.88	202	1134695	38.211	nq	98
79) Butylbenzylphthalate	20.74	149	431782	41.448	nq #	83
80) Benzo(a)anthracene	21.72	228	1097098	38.033	nq	99
81) 3,3'-Dichlorobenzidine	21.63	252	415172	41.354	nq	97
82) Chrysene	21.79	228	1053792	38.107	nq	97
83) Bis(2-ethylhexyl)phthalate	21.60	149	609985	40.699	nq	98
84) Di-n-octyl phthalate	22.82	149	1018909	41.214	nq	95
85) Indeno(1,2,3-cd)pyrene	28.79	276	1188164	40.204	nq	99
87) Benzo(b)fluoranthene	23.98	252	1078221	37.980	nq	99
88) Benzo(k)fluoranthene	24.05	252	1097713	37.889	nq	98
89) Benzo(a)pyrene	24.87	252	1032348	38.333	nq	99
90) Dibenzo(a,h)anthracene	28.83	278	1003471	38.965	nq	99
91) Benzo(g,h,i)perylene	29.97	276	979884	38.702	nq	98

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

