

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG030918\
 Data File : BG033071.D
 Acq On : 9 Mar 2018 14:21
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/12/2018 3:51:23 PM

Quant Time: Mar 09 16:08:25 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 06 17:38:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.91	152	53091	20.00	ng	0.00
21) Naphthalene-d8	11.80	136	252845	20.00	ng	0.00
38) Acenaphthene-d10	15.53	164	146754	20.00	ng	0.00
63) Phenanthrene-d10	18.26	188	341519	20.00	ng	0.00
75) Chrysene-d12	22.61	240	313557	20.00	ng	0.00
86) Perylene-d12	26.41	264	344186	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	6.32	112	258236	77.82	ng	0.00
7) Phenol-d6	8.00	99	361879	78.24	ng	0.00
23) Nitrobenzene-d5	10.11	82	333865	90.86	ng	0.00
41) 2,4,6-Tribromophenol	17.00	330	136430	88.41	ng	0.00
44) 2-Fluorobiphenyl	14.16	172	726264	71.37	ng	0.00
78) Terphenyl-d14	20.77	244	996070	71.37	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.96	88	52245	36.276	ng	94
3) Pyridine	4.43	79	153374	38.638	ng	90
4) n-Nitrosodimethylamine	4.33	42	72452	45.525	ng	# 83
6) Aniline	8.20	93	232714	37.168	ng	99
8) 2-Chlorophenol	8.46	128	140557	38.697	ng	94
9) Benzaldehyde	8.01	77	106933	40.112	ng	96
10) Phenol	8.03	94	185336	39.748	ng	96
11) bis(2-Chloroethyl)ether	8.29	93	139954	36.707	ng	95
12) 1,3-Dichlorobenzene	8.80	146	155609	36.577	ng	98
13) 1,4-Dichlorobenzene	8.95	146	156868	36.631	ng	99
14) 1,2-Dichlorobenzene	9.28	146	150839	36.757	ng	96
15) Benzyl Alcohol	9.14	79	134345	39.508	ng	94
16) 2,2'-oxybis(1-Chloropropan	9.44	45	275972	42.104	ng	99
17) 2-Methylphenol	9.34	107	130472	37.946	ng	94
18) Hexachloroethane	10.04	117	55942	38.367	ng	87
19) n-Nitroso-di-n-propylamine	9.72	70	122793	37.603	ng	# 90
20) 3+4-Methylphenols	9.67	107	184802	38.655	ng	93
22) Acetophenone	9.76	105	231169	36.202	ng	# 97
24) Nitrobenzene	10.15	77	174774	43.963	ng	93
25) Isophorone	10.68	82	312740	35.239	ng	95
26) 2-Nitrophenol	10.88	139	79225	53.460	ng	96
27) 2,4-Dimethylphenol	10.91	122	134470	34.879	ng	92
28) bis(2-Chloroethoxy)methane	11.15	93	178778	34.580	ng	96
29) 2,4-Dichlorophenol	11.42	162	134693	38.494	ng	95
30) 1,2,4-Trichlorobenzene	11.65	180	145953	35.584	ng	96
31) Naphthalene	11.85	128	474543	35.917	ng	98
32) Benzoic acid	11.02	122	107807	45.841	ng	# 87
33) 4-Chloroaniline	11.93	127	213704	36.175	ng	97
34) Hexachlorobutadiene	12.11	225	85260	37.059	ng	98
35) Caprolactam	12.68	113	55503	39.615	ng	98
36) 4-Chloro-3-methylphenol	12.99	107	167554	41.149	ng	95
37) 2-Methylnaphthalene	13.40	142	343327	35.918	ng	93
39) 1,2,4,5-Tetrachlorobenzene	13.75	216	160257	36.786	ng	# 96
40) Hexachlorocyclopentadiene	13.72	237	85353	38.444	ng	90

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.97	196	108015	39.316	nq	97
43) 2,4,5-Trichlorophenol	14.04	196	128068	40.277	nq	97
45) 1,1'-Biphenyl	14.37	154	453528	35.365	nq	94
46) 2-Chloronaphthalene	14.43	162	338455	35.331	nq	94
47) 2-Nitroaniline	14.60	65	128474	52.118	nq	98
48) Acenaphthylene	15.26	152	571986	36.470	nq	99
49) Dimethylphthalate	14.95	163	451952	36.270	nq	100
50) 2,6-Dinitrotoluene	15.08	165	99488	48.194	nq	89
51) Acenaphthene	15.59	154	368658	37.743	nq	98
52) 3-Nitroaniline	15.41	138	113871	45.159	nq	# 88
53) 2,4-Dinitrophenol	15.61	184	51583	79.428	nq	# 1
54) Dibenzofuran	15.92	168	504232	36.255	nq	93
55) 4-Nitrophenol	15.68	139	72055	39.267	nq	88
56) 2,4-Dinitrotoluene	15.86	165	140226	55.733	nq	87
57) Fluorene	16.56	166	418555	36.462	nq	100
58) 2,3,4,6-Tetrachlorophenol	16.13	232	103879m	42.078	nq	
59) Diethylphthalate	16.29	149	487178	37.068	nq	97
60) 4-Chlorophenyl-phenylether	16.53	204	208839	37.190	nq	92
61) 4-Nitroaniline	16.56	138	108828	41.188	nq	91
62) Azobenzene	16.83	77	438516	37.296	nq	96
64) 4,6-Dinitro-2-methylphenol	16.61	198	68182	61.532	nq	93
65) n-Nitrosodiphenylamine	16.75	169	387801	34.905	nq	98
66) 4-Bromophenyl-phenylether	17.43	248	130034	36.009	nq	# 88
67) Hexachlorobenzene	17.56	284	140252	35.939	nq	# 93
68) Atrazine	17.66	200	128938	35.258	nq	98
69) Pentachlorophenol	17.89	266	93514	38.043	nq	98
70) Phenanthrene	18.30	178	678105	35.590	nq	99
71) Anthracene	18.39	178	681269	35.615	nq	99
72) Carbazole	18.64	167	650188	34.485	nq	97
73) Di-n-butylphthalate	19.14	149	835352	36.115	nq	99
74) Fluoranthene	20.25	202	758703	36.048	nq	95
76) Benzidine	20.41	184	354397	33.158	nq	97
77) Pyrene	20.61	202	775060	35.051	nq	97
79) Butylbenzylphthalate	21.46	149	383060	39.073	nq	94
80) Benzo(a)anthracene	22.59	228	726869	36.150	nq	99
81) 3,3'-Dichlorobenzidine	22.47	252	263532	35.388	nq	# 97
82) Chrysene	22.67	228	698292	36.356	nq	98
83) Bis(2-ethylhexyl)phthalate	22.41	149	554706	38.224	nq	95
84) Di-n-octyl phthalate	23.82	149	886276	40.067	nq	98
85) Indeno(1,2,3-cd)pyrene	30.84	276	777400	34.705	nq	# 95
87) Benzo(b)fluoranthene	25.19	252	730517	35.896	nq	# 97
88) Benzo(k)fluoranthene	25.27	252	726208	35.906	nq	# 97
89) Benzo(a)pyrene	26.23	252	692322	35.767	nq	# 95
90) Dibenzo(a,h)anthracene	30.91	278	611073	31.364	nq	# 95
91) Benzo(g,h,i)perylene	32.22	276	637220	33.178	nq	# 92

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

