

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG030518\
 Data File : BG032995.D
 Acq On : 5 Mar 2018 18:29
 Operator : SJ/JU
 Sample : PB107049BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB107049BS

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:26:30 PM

Quant Time: Mar 06 02:13:55 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 01 17:38:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.91	152	58293	20.00	ng	-0.01
21) Naphthalene-d8	11.80	136	259056	20.00	ng	0.00
38) Acenaphthene-d10	15.53	164	155042	20.00	ng	0.00
63) Phenanthrene-d10	18.26	188	340916	20.00	ng	0.00
75) Chrysene-d12	22.61	240	304821	20.00	ng	-0.01
86) Perylene-d12	26.40	264	327726	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	6.33	112	444774	122.07	ng	0.00
7) Phenol-d6	8.01	99	598594	117.86	ng	0.00
23) Nitrobenzene-d5	10.11	82	354688	93.85	ng	0.00
41) 2,4,6-Tribromophenol	17.00	330	198925	122.02	ng	0.00
44) 2-Fluorobiphenyl	14.16	172	775572	72.15	ng	-0.01
78) Terphenyl-d14	20.77	244	1091122	80.42	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.96	88	50912	32.196	ng	95
3) Pyridine	4.42	79	139453	31.996	ng	93
4) n-Nitrosodimethylamine	4.32	42	82811	47.390	ng	# 84
6) Aniline	8.20	93	88579	12.885	ng	97
8) 2-Chlorophenol	8.46	128	167190	41.921	ng	96
9) Benzaldehyde	8.01	77	57624	19.686	ng	99
10) Phenol	8.04	94	218529	42.684	ng	95
11) bis(2-Chloroethyl)ether	8.29	93	150201	35.879	ng	98
12) 1,3-Dichlorobenzene	8.80	146	164462	35.208	ng	95
13) 1,4-Dichlorobenzene	8.95	146	165802	35.262	ng	95
14) 1,2-Dichlorobenzene	9.29	146	156542	34.743	ng	95
15) Benzyl Alcohol	9.15	79	147909	39.615	ng	89
16) 2,2'-oxybis(1-Chloropropan	9.44	45	289761	40.263	ng	99
17) 2-Methylphenol	9.34	107	146931	38.920	ng	96
18) Hexachloroethane	10.04	117	60301	37.666	ng	# 84
19) n-Nitroso-di-n-propylamine	9.72	70	128374	35.804	ng	# 90
20) 3+4-Methylphenols	9.67	107	204498	38.958	ng	93
22) Acetophenone	9.76	105	235547	36.003	ng	# 98
24) Nitrobenzene	10.16	77	182000	44.578	ng	94
25) Isophorone	10.68	82	348657	38.345	ng	95
26) 2-Nitrophenol	10.88	139	85113	55.305	ng	98
27) 2,4-Dimethylphenol	10.91	122	180259	45.635	ng	91
28) bis(2-Chloroethoxy)methane	11.15	93	212612	40.138	ng	97
29) 2,4-Dichlorophenol	11.42	162	157305	43.879	ng	96
30) 1,2,4-Trichlorobenzene	11.65	180	147278	35.047	ng	97
31) Naphthalene	11.85	128	505110	37.314	ng	98
32) Benzoic acid	11.01	122	107631	45.034	ng	89
33) 4-Chloroaniline	11.94	127	80827	13.354	ng	96
34) Hexachlorobutadiene	12.11	225	82233	34.886	ng	99
35) Caprolactam	12.68	113	62377	43.454	ng	92
36) 4-Chloro-3-methylphenol	12.99	107	190035	45.551	ng	93
37) 2-Methylnaphthalene	13.40	142	374975	38.288	ng	92
39) 1,2,4,5-Tetrachlorobenzene	13.75	216	157700	34.264	ng	# 96
40) Hexachlorocyclopentadiene	13.73	237	195676	83.423	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.97	196	121712	41.934	nq	96
43) 2,4,5-Trichlorophenol	14.04	196	133626	39.778	nq	# 95
45) 1,1'-Biphenyl	14.37	154	475073	35.065	nq	96
46) 2-Chloronaphthalene	14.43	162	357663	35.340	nq	98
47) 2-Nitroaniline	14.61	65	131418	50.775	nq	94
48) Acenaphthylene	15.26	152	592691	35.770	nq	99
49) Dimethylphthalate	14.95	163	470060	35.707	nq	99
50) 2,6-Dinitrotoluene	15.08	165	105090	48.187	nq	93
51) Acenaphthene	15.60	154	400081	38.771	nq	97
52) 3-Nitroaniline	15.41	138	61973	26.921	nq	# 85
53) 2,4-Dinitrophenol	15.61	184	79936	104.127	nq	# 40
54) Dibenzofuran	15.92	168	554612	37.746	nq	95
55) 4-Nitrophenol	15.68	139	173659	80.506	nq	85
56) 2,4-Dinitrotoluene	15.85	165	147226	55.476	nq	# 88
57) Fluorene	16.56	166	445385	36.726	nq	98
58) 2,3,4,6-Tetrachlorophenol	16.13	232	113912m	43.676	nq	
59) Diethylphthalate	16.29	149	507420	36.544	nq	98
60) 4-Chlorophenyl-phenylether	16.54	204	219646	37.024	nq	91
61) 4-Nitroaniline	16.56	138	107930	38.989	nq	84
62) Azobenzene	16.83	77	471650	37.970	nq	96
64) 4,6-Dinitro-2-methylphenol	16.61	198	63211	58.480	nq	95
65) n-Nitrosodiphenylamine	16.75	169	411945	37.144	nq	99
66) 4-Bromophenyl-phenylether	17.43	248	131494	36.477	nq	# 86
67) Hexachlorobenzene	17.56	284	135758	34.849	nq	# 90
68) Atrazine	17.66	200	141730	38.824	nq	96
69) Pentachlorophenol	17.89	266	181133	73.819	nq	100
70) Phenanthrene	18.30	178	707455	37.196	nq	99
71) Anthracene	18.39	178	728424	38.148	nq	99
72) Carbazole	18.64	167	676730	35.956	nq	# 97
73) Di-n-butylphthalate	19.14	149	859668	37.232	nq	99
74) Fluoranthene	20.25	202	785920	37.407	nq	94
76) Benzidine	20.40	184	113916	10.964	nq	97
77) Pyrene	20.61	202	785261	36.530	nq	97
79) Butylbenzylphthalate	21.46	149	398446	41.807	nq	92
80) Benzo(a)anthracene	22.59	228	748628	38.299	nq	99
81) 3,3'-Dichlorobenzidine	22.47	252	120691	16.671	nq	# 96
82) Chrysene	22.66	228	696814	37.319	nq	99
83) Bis(2-ethylhexyl)phthalate	22.42	149	569552	40.372	nq	96
84) Di-n-octyl phthalate	23.83	149	981098	45.625	nq	99
85) Indeno(1,2,3-cd)pyrene	30.82	276	802414	36.849	nq	97
87) Benzo(b)fluoranthene	25.18	252	735697	37.967	nq	# 96
88) Benzo(k)fluoranthene	25.26	252	733255	38.075	nq	# 96
89) Benzo(a)pyrene	26.23	252	719864	39.058	nq	# 97
90) Dibenzo(a,h)anthracene	30.90	278	636881	34.330	nq	# 94
91) Benzo(g,h,i)perylene	32.22	276	673660	36.837	nq	# 92

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

