

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG073018\
 Data File : BG035984.D
 Acq On : 30 Jul 2018 20:13
 Operator : JU/SJ
 Sample : PB111523BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111523BS

Manual Integrations
 APPROVED

Sohil
 7/31/2018 12:20:51 PM

Quant Time: Jul 31 01:48:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 25 11:04:14 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	31414	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	124847	20.00	ng	0.00
38) Acenaphthene-d10	14.65	164	96757	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	276220	20.00	ng	0.00
75) Chrysene-d12	21.66	240	308316	20.00	ng	0.00
86) Perylene-d12	24.85	264	298887	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	237949	125.76	ng	0.00
7) Phenol-d6	7.20	99	356579	126.31	ng	0.00
23) Nitrobenzene-d5	9.18	82	231439	77.44	ng	0.00
41) 2,4,6-Tribromophenol	16.14	330	173468	136.02	ng	0.00
44) 2-Fluorobiphenyl	13.27	172	542619	75.13	ng	0.00
78) Terphenyl-d14	20.00	244	1159994	82.93	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.51	88	25247	26.216	ng	# 77
3) Pyridine	3.91	79	71584	28.473	ng	# 76
4) n-Nitrosodimethylamine	3.82	42	32925	30.500	ng	# 80
6) Aniline	7.35	93	96628	26.408	ng	97
8) 2-Chlorophenol	7.59	128	78651	40.870	ng	98
9) Benzaldehyde	7.17	77	58505	33.776	ng	97
10) Phenol	7.22	94	114960	40.307	ng	96
11) bis(2-Chloroethyl)ether	7.44	93	75657	33.872	ng	87
12) 1,3-Dichlorobenzene	7.91	146	84026	36.157	ng	94
13) 1,4-Dichlorobenzene	8.06	146	86716	36.725	ng	89
14) 1,2-Dichlorobenzene	8.38	146	82605	36.417	ng	99
15) Benzyl Alcohol	8.26	79	88501	34.522	ng	89
16) 2,2'-oxybis(1-Chloropropan	8.55	45	87408	46.677	ng	81
17) 2-Methylphenol	8.46	107	79308	40.898	ng	# 88
18) Hexachloroethane	9.10	117	33176	36.748	ng	96
19) n-Nitroso-di-n-propylamine	8.82	70	74788	31.460	ng	# 80
20) 3+4-Methylphenols	8.79	107	106347	39.532	ng	93
22) Acetophenone	8.85	105	136754	35.224	ng	# 91
24) Nitrobenzene	9.23	77	111883	37.225	ng	93
25) Isophorone	9.74	82	210816	38.000	ng	100
26) 2-Nitrophenol	9.94	139	44917	42.079	ng	# 89
27) 2,4-Dimethylphenol	10.00	122	81547	45.131	ng	91
28) bis(2-Chloroethoxy)methane	10.23	93	113551	37.145	ng	97
29) 2,4-Dichlorophenol	10.48	162	92136	44.670	ng	96
30) 1,2,4-Trichlorobenzene	10.69	180	99048	39.162	ng	95
31) Naphthalene	10.87	128	246103	37.842	ng	98
32) Benzoic acid	10.17	122	22159m	18.217	ng	
33) 4-Chloroaniline	10.99	127	49809	17.573	ng	# 92
34) Hexachlorobutadiene	11.15	225	75379	38.221	ng	99
35) Caprolactam	11.76	113	33724m	38.101	ng	
36) 4-Chloro-3-methylphenol	12.11	107	115490	41.773	ng	93
37) 2-Methylnaphthalene	12.48	142	197054	40.671	ng	94
39) 1,2,4,5-Tetrachlorobenzene	12.85	216	134644	36.869	ng	98
40) Hexachlorocyclopentadiene	12.82	237	150865	78.771	ng	91

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.09	196	86761	40.625	nq	97
43) 2,4,5-Trichlorophenol	13.16	196	96478	40.381	nq	94
45) 1,1'-Biphenyl	13.48	154	279986	37.324	nq	99
46) 2-Chloronaphthalene	13.53	162	223071	38.230	nq	98
47) 2-Nitroaniline	13.73	65	79540	38.558	nq	100
48) Acenaphthylene	14.37	152	378452	38.719	nq	97
49) Dimethylphthalate	14.10	163	312234	36.832	nq	98
50) 2,6-Dinitrotoluene	14.22	165	72899	42.228	nq	94
51) Acenaphthene	14.71	154	212643	34.924	nq	98
52) 3-Nitroaniline	14.56	138	46430	26.315	nq	# 96
53) 2,4-Dinitrophenol	14.76	184	39630	47.551	nq	# 65
54) Dibenzofuran	15.04	168	372259	38.443	nq	93
55) 4-Nitrophenol	14.87	139	99910	79.512	nq	# 86
56) 2,4-Dinitrotoluene	15.01	165	109335	44.147	nq	90
57) Fluorene	15.70	166	315630	38.889	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.27	232	101833	44.455	nq	93
59) Diethylphthalate	15.46	149	328082	37.637	nq	98
60) 4-Chlorophenyl-phenylether	15.68	204	181092	37.963	nq	96
61) 4-Nitroaniline	15.72	138	75366	40.835	nq	# 76
62) Azobenzene	15.98	77	328690	38.228	nq	94
64) 4,6-Dinitro-2-methylphenol	15.78	198	53631	34.879	nq	81
65) n-Nitrosodiphenylamine	15.90	169	285625	36.329	nq	98
66) 4-Bromophenyl-phenylether	16.58	248	123325	38.484	nq	91
67) Hexachlorobenzene	16.70	284	130479	39.537	nq	# 91
68) Atrazine	16.85	200	138311	42.727	nq	95
69) Pentachlorophenol	17.05	266	111377	55.409	nq	97
70) Phenanthrene	17.43	178	548271	39.779	nq	99
71) Anthracene	17.52	178	562278	40.970	nq	97
72) Carbazole	17.80	167	512584	39.328	nq	98
73) Di-n-butylphthalate	18.34	149	589963	38.305	nq	# 98
74) Fluoranthene	19.44	202	733411	39.504	nq	97
76) Benzidine	19.62	184	258654	31.910	nq	98
77) Pyrene	19.80	202	733111	37.605	nq	98
79) Butylbenzylphthalate	20.68	149	276834	39.709	nq	96
80) Benzo(a)anthracene	21.63	228	745828	38.818	nq	98
81) 3,3'-Dichlorobenzidine	21.55	252	181850	27.145	nq	# 99
82) Chrysene	21.70	228	700167	39.325	nq	98
83) Bis(2-ethylhexyl)phthalate	21.53	149	375054	39.572	nq	99
84) Di-n-octyl phthalate	22.73	149	641531	40.426	nq	# 92
85) Indeno(1,2,3-cd)pyrene	28.49	276	773637	39.247	nq	# 85
87) Benzo(b)fluoranthene	23.84	252	694097	38.208	nq	# 96
88) Benzo(k)fluoranthene	23.90	252	704659	39.677	nq	# 98
89) Benzo(a)pyrene	24.70	252	682852	40.404	nq	# 97
90) Dibenzo(a,h)anthracene	28.55	278	644151	38.823	nq	95
91) Benzo(g,h,i)perylene	29.63	276	640240	39.433	nq	# 93

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

