

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070518\
 Data File : BM015790.D
 Acq On : 06 Jul 2018 07:43
 Operator : SJ/JU
 Sample : PB110755BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB110755BS

Manual Integrations
 APPROVED

Sohil
 7/6/2018 5:21:59 PM

Quant Time: Jul 06 08:52:40 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 05 17:39:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.31	152	126514	20.00	ng	0.00
21) Naphthalene-d8	11.14	136	539019	20.00	ng	0.00
38) Acenaphthene-d10	14.93	164	313100	20.00	ng	0.00
63) Phenanthrene-d10	17.66	188	675761	20.00	ng	0.00
75) Chrysene-d12	21.76	240	634543	20.00	ng	0.00
86) Perylene-d12	24.17	264	480407	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.81	112	759539	102.96	ng	0.00
7) Phenol-d6	7.46	99	1001649	99.18	ng	0.00
23) Nitrobenzene-d5	9.50	82	1084029	98.12	ng	0.00
41) 2,4,6-Tribromophenol	16.41	330	430785	105.30	ng	0.00
44) 2-Fluorobiphenyl	13.56	172	2550546	101.13	ng	0.00
78) Terphenyl-d14	20.22	244	3828786	111.92	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.56	88	102192	33.495	ng	# 100
3) Pyridine	4.00	79	277153	34.425	ng	# 79
4) n-Nitrosodimethylamine	3.92	42	261740	41.752	ng	# 42
6) Aniline	7.63	93	388473	31.727	ng	94
8) 2-Chlorophenol	7.87	128	438556	49.533	ng	95
9) Benzaldehyde	7.45	77	211612	33.244	ng	90
10) Phenol	7.49	94	507355	49.997	ng	93
11) bis(2-Chloroethyl)ether	7.73	93	352733	42.730	ng	88
12) 1,3-Dichlorobenzene	8.19	146	426315	43.429	ng	# 92
13) 1,4-Dichlorobenzene	8.35	146	437361	43.085	ng	# 94
14) 1,2-Dichlorobenzene	8.67	146	434282	44.236	ng	94
15) Benzyl Alcohol	8.56	79	406523	46.365	ng	90
16) 2,2'-oxybis(1-Chloropropan	8.84	45	565509	62.727	ng	99
17) 2-Methylphenol	8.76	107	359606	51.150	ng	# 86
18) Hexachloroethane	9.39	117	184421	45.946	ng	96
19) n-Nitroso-di-n-propylamine	9.13	70	339513	41.830	ng	# 78
20) 3+4-Methylphenols	9.09	107	481852	49.342	ng	89
22) Acetophenone	9.15	105	634534	45.190	ng	# 87
24) Nitrobenzene	9.55	77	520294	48.382	ng	97
25) Isophorone	10.06	82	884939	52.481	ng	95
26) 2-Nitrophenol	10.26	139	233302	50.382	ng	# 63
27) 2,4-Dimethylphenol	10.30	122	406765	57.823	ng	# 86
28) bis(2-Chloroethoxy)methane	10.54	93	529022	48.156	ng	99
29) 2,4-Dichlorophenol	10.79	162	432440	55.504	ng	98
30) 1,2,4-Trichlorobenzene	11.00	180	434239	47.986	ng	97
31) Naphthalene	11.19	128	1329743	47.583	ng	99
32) Benzoic acid	10.50	122	277989m	44.633	ng	
33) 4-Chloroaniline	11.30	127	213391	18.729	ng	# 89
34) Hexachlorobutadiene	11.45	225	284673	47.136	ng	97
35) Caprolactam	12.13	113	124650m	48.241	ng	
36) 4-Chloro-3-methylphenol	12.42	107	476945	53.351	ng	89
37) 2-Methylnaphthalene	12.78	142	1029153	51.769	ng	# 94
39) 1,2,4,5-Tetrachlorobenzene	13.14	216	511017	50.852	ng	# 100
40) Hexachlorocyclopentadiene	13.10	237	608753	123.849	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.38	196	342691	53.948	nq	98
43) 2,4,5-Trichlorophenol	13.46	196	361230	52.643	nq #	86
45) 1,1'-Biphenyl	13.77	154	1321330	48.811	nq	99
46) 2-Chloronaphthalene	13.82	162	1014311	50.114	nq #	91
47) 2-Nitroaniline	14.03	65	348736	49.408	nq	83
48) Acenaphthylene	14.66	152	1669320	50.700	nq	99
49) Dimethylphthalate	14.39	163	1328494	48.111	nq	99
50) 2,6-Dinitrotoluene	14.52	165	279628	51.937	nq #	76
51) Acenaphthene	15.00	154	967271	47.211	nq	97
52) 3-Nitroaniline	14.85	138	160763	27.280	nq #	81
53) 2,4-Dinitrophenol	15.07	184	281330	102.282	nq	94
54) Dibenzofuran	15.33	168	1572434	49.807	nq #	88
55) 4-Nitrophenol	15.16	139	456034	104.855	nq #	69
56) 2,4-Dinitrotoluene	15.30	165	399319	51.373	nq	95
57) Fluorene	15.97	166	1286924	49.432	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.55	232	320697	55.384	nq #	100
59) Diethylphthalate	15.73	149	1413604	47.771	nq	97
60) 4-Chlorophenyl-phenylether	15.95	204	651936	48.665	nq	90
61) 4-Nitroaniline	16.01	138	265925	43.499	nq #	39
62) Azobenzene	16.25	77	1465804	52.845	nq	81
64) 4,6-Dinitro-2-methylphenol	16.07	198	196856	47.919	nq	89
65) n-Nitrosodiphenylamine	16.17	169	1100834	47.542	nq	98
66) 4-Bromophenyl-phenylether	16.84	248	391388	51.874	nq #	83
67) Hexachlorobenzene	16.96	284	423770	50.263	nq #	77
68) Atrazine	17.10	200	400133	52.536	nq	99
69) Pentachlorophenol	17.31	266	476342	91.305	nq	98
70) Phenanthrene	17.70	178	1916234	49.565	nq	99
71) Anthracene	17.79	178	1959675	51.615	nq	99
72) Carbazole	18.06	167	1729650	48.888	nq	99
73) Di-n-butylphthalate	18.57	149	2343521	49.144	nq #	98
74) Fluoranthene	19.68	202	2127660	47.872	nq	99
76) Benzidine	19.86	184	754568	40.587	nq	99
77) Pyrene	20.04	202	2137692	54.082	nq	100
79) Butylbenzylphthalate	20.89	149	1013677	53.053	nq #	86
80) Benzo(a)anthracene	21.75	228	2006596	49.876	nq	99
81) 3,3'-Dichlorobenzidine	21.67	252	362958	25.109	nq	98
82) Chrysene	21.80	228	1845955	49.255	nq	100
83) Bis(2-ethylhexyl)phthalate	21.63	149	1432688	50.748	nq	96
84) Di-n-octyl phthalate	22.55	149	2224000	48.313	nq #	91
85) Indeno(1,2,3-cd)pyrene	26.65	276	1476194	42.395	nq #	93
87) Benzo(b)fluoranthene	23.44	252	1666473	51.957	nq #	96
88) Benzo(k)fluoranthene	23.49	252	1577627	50.635	nq	99
89) Benzo(a)pyrene	24.07	252	1472802	52.164	nq	100
90) Dibenzo(a,h)anthracene	26.65	278	1259203	49.569	nq	98
91) Benzo(g,h,i)perylene	27.41	276	1140138	51.081	nq #	94

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

