

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
 Data File : BM015896.D
 Acq On : 11 Jul 2018 19:37
 Operator : SJ/JU
 Sample : PB111002BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB111002BS

Manual Integrations
 APPROVED

Sohil
 7/12/2018 6:46:55 PM

Quant Time: Jul 12 04:02:46 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	8.29	152	90513	20.00	ng	-0.02
21) Naphthalene-d8	11.12	136	361109	20.00	ng	-0.02
38) Acenaphthene-d10	14.92	164	184398	20.00	ng	-0.02
63) Phenanthrene-d10	17.64	188	349831	20.00	ng	-0.02
75) Chrysene-d12	21.75	240	299160	20.00	ng	-0.02
86) Perylene-d12	24.14	264	287879	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.80	112	658421	124.75	ng	-0.02
7) Phenol-d6	7.45	99	832062	115.16	ng	-0.02
23) Nitrobenzene-d5	9.48	82	581812	78.61	ng	-0.02
41) 2,4,6-Tribromophenol	16.40	330	261513	108.54	ng	-0.02
44) 2-Fluorobiphenyl	13.54	172	1165943	78.50	ng	-0.02
78) Terphenyl-d14	20.20	244	1309790	81.21	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.56	88	67348	30.854	ng	# 100
3) Pyridine	4.00	79	174509	30.297	ng	# 75
4) n-Nitrosodimethylamine	3.91	42	162178	36.160	ng	# 41
6) Aniline	7.62	93	203920	23.279	ng	91
8) 2-Chlorophenol	7.85	128	237300	37.463	ng	96
9) Benzaldehyde	7.43	77	143309	31.469	ng	90
10) Phenol	7.47	94	273443	37.664	ng	94
11) bis(2-Chloroethyl)ether	7.71	93	199558	33.790	ng	88
12) 1,3-Dichlorobenzene	8.17	146	250867	35.721	ng	# 92
13) 1,4-Dichlorobenzene	8.33	146	253725	34.936	ng	# 95
14) 1,2-Dichlorobenzene	8.65	146	246306	35.068	ng	# 94
15) Benzyl Alcohol	8.55	79	206239	32.878	ng	89
16) 2,2'-oxybis(1-Chloropropan	8.81	45	311585	48.308	ng	98
17) 2-Methylphenol	8.74	107	186291	37.037	ng	# 86
18) Hexachloroethane	9.37	117	108695	37.850	ng	96
19) n-Nitroso-di-n-propylamine	9.10	70	178081	30.667	ng	# 76
20) 3+4-Methylphenols	9.07	107	243899	34.909	ng	88
22) Acetophenone	9.13	105	337531	35.881	ng	# 87
24) Nitrobenzene	9.52	77	282087	39.155	ng	97
25) Isophorone	10.04	82	450193	39.853	ng	93
26) 2-Nitrophenol	10.23	139	119546	38.535	ng	# 57
27) 2,4-Dimethylphenol	10.28	122	200677	42.581	ng	# 85
28) bis(2-Chloroethoxy)methane	10.52	93	270871	36.805	ng	100
29) 2,4-Dichlorophenol	10.77	162	209841	40.203	ng	97
30) 1,2,4-Trichlorobenzene	10.98	180	228816	37.743	ng	97
31) Naphthalene	11.17	128	697694	37.266	ng	99
32) Benzoic acid	10.45	122	120070m	28.776	ng	
33) 4-Chloroaniline	11.29	127	99984	13.099	ng	# 91
34) Hexachlorobutadiene	11.43	225	149695	36.998	ng	97
35) Caprolactam	12.09	113	51838m	29.946	ng	
36) 4-Chloro-3-methylphenol	12.39	107	215451	35.974	ng	88
37) 2-Methylnaphthalene	12.76	142	503791	37.827	ng	# 93
39) 1,2,4,5-Tetrachlorobenzene	13.12	216	236965	40.039	ng	# 100
40) Hexachlorocyclopentadiene	13.08	237	263982	93.688	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.36	196	150121	40.128	ng	98
43) 2,4,5-Trichlorophenol	13.44	196	152518	37.740	ng	# 84
45) 1,1'-Biphenyl	13.75	154	616068	38.642	ng	98
46) 2-Chloronaphthalene	13.80	162	473993	39.764	ng	# 91
47) 2-Nitroaniline	14.02	65	151042	36.335	ng	# 79
48) Acenaphthylene	14.64	152	749389	38.646	ng	99
49) Dimethylphthalate	14.37	163	574622	35.334	ng	99
50) 2,6-Dinitrotoluene	14.50	165	117842	37.164	ng	# 62
51) Acenaphthene	14.98	154	434869	36.040	ng	98
52) 3-Nitroaniline	14.83	138	74642	21.507	ng	# 75
53) 2,4-Dinitrophenol	15.05	184	98552	63.674	ng	# 81
54) Dibenzofuran	15.31	168	686032	36.897	ng	# 86
55) 4-Nitrophenol	15.13	139	166458	64.987	ng	# 77
56) 2,4-Dinitrotoluene	15.29	165	158188	34.555	ng	100
57) Fluorene	15.95	166	539927	35.214	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.53	232	125023	36.661	ng	# 100
59) Diethylphthalate	15.71	149	586497	33.654	ng	96
60) 4-Chlorophenyl-phenylether	15.93	204	266551	33.785	ng	# 86
61) 4-Nitroaniline	15.99	138	101834	28.284	ng	# 32
62) Azobenzene	16.23	77	638442	39.082	ng	79
64) 4,6-Dinitro-2-methylphenol	16.05	198	69948	32.890	ng	87
65) n-Nitrosodiphenylamine	16.15	169	446560	37.253	ng	97
66) 4-Bromophenyl-phenylether	16.83	248	153894	39.400	ng	# 83
67) Hexachlorobenzene	16.94	284	168773	38.668	ng	# 72
68) Atrazine	17.09	200	151173	38.341	ng	98
69) Pentachlorophenol	17.29	266	168616	62.432	ng	99
70) Phenanthrene	17.68	178	750528	37.500	ng	98
71) Anthracene	17.77	178	763006	38.820	ng	99
72) Carbazole	18.04	167	653275	35.668	ng	98
73) Di-n-butylphthalate	18.56	149	880179	35.654	ng	# 97
74) Fluoranthene	19.66	202	768917	33.419	ng	99
76) Benzidine	19.84	184	355195	40.524	ng	98
77) Pyrene	20.02	202	773303	41.497	ng	98
79) Butylbenzylphthalate	20.87	149	356936	39.624	ng	# 82
80) Benzo(a)anthracene	21.73	228	717116	37.808	ng	99
81) 3,3'-Dichlorobenzidine	21.66	252	173800	25.503	ng	99
82) Chrysene	21.79	228	669981	37.918	ng	100
83) Bis(2-ethylhexyl)phthalate	21.62	149	510490	38.354	ng	94
84) Di-n-octyl phthalate	22.53	149	820354	37.800	ng	# 89
85) Indeno(1,2,3-cd)pyrene	26.61	276	810590	49.378	ng	# 94
87) Benzo(b)fluoranthene	23.41	252	688411	35.817	ng	# 96
88) Benzo(k)fluoranthene	23.46	252	663086	35.515	ng	98
89) Benzo(a)pyrene	24.04	252	655917	38.768	ng	99
90) Dibenzo(a,h)anthracene	26.61	278	692748	45.508	ng	# 97
91) Benzo(g,h,i)perylene	27.37	276	665975	49.792	ng	# 94

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

