

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070918\
 Data File : BM015844.D
 Acq On : 09 Jul 2018 15:45
 Operator : SJ/JU
 Sample : PB110833BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB110833BS

Manual Integrations
 APPROVED

Sohil
 7/10/2018 2:50:02 PM

Quant Time: Jul 10 02:07:19 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	8.30	152	78391	20.00	ng	-0.02
21) Naphthalene-d8	11.13	136	353778	20.00	ng	-0.02
38) Acenaphthene-d10	14.92	164	210126	20.00	ng	-0.01
63) Phenanthrene-d10	17.64	188	481080	20.00	ng	-0.01
75) Chrysene-d12	21.76	240	529048	20.00	ng	-0.01
86) Perylene-d12	24.16	264	521187	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.80	112	597795	130.78	ng	-0.01
7) Phenol-d6	7.45	99	798597	127.62	ng	-0.01
23) Nitrobenzene-d5	9.49	82	547695	75.54	ng	-0.02
41) 2,4,6-Tribromophenol	16.40	330	345060	125.68	ng	-0.01
44) 2-Fluorobiphenyl	13.54	172	1231287	72.75	ng	-0.02
78) Terphenyl-d14	20.21	244	2120075	74.33	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.55	88	51545	27.266	ng	# 100
3) Pyridine	4.00	79	128771	25.813	ng	# 74
4) n-Nitrosodimethylamine	3.91	42	128172	32.997	ng	# 41
6) Aniline	7.62	93	142032	18.721	ng	# 91
8) 2-Chlorophenol	7.86	128	219400	39.993	ng	95
9) Benzaldehyde	7.44	77	91106	23.099	ng	92
10) Phenol	7.48	94	258943	41.182	ng	93
11) bis(2-Chloroethyl)ether	7.72	93	177262	34.656	ng	87
12) 1,3-Dichlorobenzene	8.19	146	208571	34.291	ng	# 90
13) 1,4-Dichlorobenzene	8.33	146	214317	34.073	ng	94
14) 1,2-Dichlorobenzene	8.66	146	211216	34.722	ng	# 95
15) Benzyl Alcohol	8.55	79	199023	36.634	ng	# 87
16) 2,2'-oxybis(1-Chloropropan	8.83	45	285274	51.068	ng	98
17) 2-Methylphenol	8.75	107	178172	40.900	ng	# 87
18) Hexachloroethane	9.39	117	88955	35.766	ng	97
19) n-Nitroso-di-n-propylamine	9.12	70	169411	33.686	ng	# 81
20) 3+4-Methylphenols	9.08	107	239239	39.537	ng	90
22) Acetophenone	9.15	105	315987	34.287	ng	# 87
24) Nitrobenzene	9.53	77	261860	37.100	ng	99
25) Isophorone	10.05	82	429178	38.780	ng	# 95
26) 2-Nitrophenol	10.25	139	110542	36.371	ng	# 60
27) 2,4-Dimethylphenol	10.29	122	212768	46.082	ng	# 85
28) bis(2-Chloroethoxy)methane	10.53	93	263072	36.486	ng	98
29) 2,4-Dichlorophenol	10.77	162	211404	41.341	ng	97
30) 1,2,4-Trichlorobenzene	10.99	180	212043	35.701	ng	96
31) Naphthalene	11.18	128	661610	36.071	ng	100
32) Benzoic acid	10.45	122	131533m	32.177	ng	
33) 4-Chloroaniline	11.30	127	123760	16.549	ng	# 88
34) Hexachlorobutadiene	11.43	225	137982	34.810	ng	98
35) Caprolactam	12.10	113	62571m	36.896	ng	
36) 4-Chloro-3-methylphenol	12.40	107	237485	40.475	ng	86
37) 2-Methylnaphthalene	12.77	142	504172	38.640	ng	# 93
39) 1,2,4,5-Tetrachlorobenzene	13.13	216	242441	35.949	ng	# 100
40) Hexachlorocyclopentadiene	13.09	237	233779	74.920	ng	98

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070918\
 Data File : BM015844.D
 Acq On : 09 Jul 2018 15:45
 Operator : SJ/JU
 Sample : PB110833BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB110833BS

Manual Integrations
 APPROVED

Sohil
 7/10/2018 2:50:02 PM

Quant Time: Jul 10 02:07:19 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)
42) 2,4,6-Trichlorophenol	13.37	196	162560	38.132	ng	99
43) 2,4,5-Trichlorophenol	13.44	196	171863	37.320	ng	# 82
45) 1,1'-Biphenyl	13.76	154	648311	35.685	ng	99
46) 2-Chloronaphthalene	13.81	162	498183	36.676	ng	# 91
47) 2-Nitroaniline	14.02	65	177778	37.531	ng	# 79
48) Acenaphthylene	14.65	152	811419	36.721	ng	99
49) Dimethylphthalate	14.37	163	656372	35.419	ng	100
50) 2,6-Dinitrotoluene	14.51	165	136127	37.674	ng	# 70
51) Acenaphthene	14.99	154	477553	34.731	ng	98
52) 3-Nitroaniline	14.84	138	81076	20.500	ng	# 77
53) 2,4-Dinitrophenol	15.06	184	124793	69.977	ng	# 84
54) Dibenzofuran	15.32	168	778037	36.722	ng	# 85
55) 4-Nitrophenol	15.14	139	232591	79.687	ng	# 73
56) 2,4-Dinitrotoluene	15.29	165	200342	38.405	ng	98
57) Fluorene	15.96	166	639621	36.608	ng	98
58) 2,3,4,6-Tetrachlorophenol	15.54	232	154304	39.707	ng	# 100
59) Diethylphthalate	15.72	149	717360	36.123	ng	95
60) 4-Chlorophenyl-phenylether	15.94	204	312341	34.741	ng	# 88
61) 4-Nitroaniline	16.00	138	150321	36.639	ng	# 35
62) Azobenzene	16.24	77	756344	40.630	ng	81
64) 4,6-Dinitro-2-methylphenol	16.06	198	94571	32.336	ng	87
65) n-Nitrosodiphenylamine	16.16	169	553959	33.605	ng	98
66) 4-Bromophenyl-phenylether	16.83	248	190213	35.412	ng	# 82
67) Hexachlorobenzene	16.96	284	208464	34.731	ng	# 78
68) Atrazine	17.09	200	207614	38.290	ng	97
69) Pentachlorophenol	17.30	266	235610	63.438	ng	98
70) Phenanthrene	17.69	178	999981	36.333	ng	99
71) Anthracene	17.78	178	1012922	37.476	ng	99
72) Carbazole	18.05	167	934454	37.100	ng	98
73) Di-n-butylphthalate	18.57	149	1229432	36.215	ng	# 98
74) Fluoranthene	19.67	202	1157833	36.593	ng	98
76) Benzidine	19.84	184	356934	23.027	ng	97
77) Pyrene	20.03	202	1184598	35.945	ng	99
79) Butylbenzylphthalate	20.87	149	588769	36.959	ng	# 82
80) Benzo(a)anthracene	21.74	228	1243213	37.063	ng	99
81) 3,3'-Dichlorobenzidine	21.66	252	219383	18.203	ng	99
82) Chrysene	21.79	228	1147015	36.708	ng	98
83) Bis(2-ethylhexyl)phthalate	21.62	149	862482	36.642	ng	95
84) Di-n-octyl phthalate	22.54	149	1497975	39.030	ng	# 90
85) Indeno(1,2,3-cd)pyrene	26.63	276	1329879	45.809	ng	# 93
87) Benzo(b)fluoranthene	23.43	252	1227738	35.283	ng	# 95
88) Benzo(k)fluoranthene	23.47	252	1170320	34.623	ng	98
89) Benzo(a)pyrene	24.05	252	1152938	37.640	ng	99
90) Dibenzo(a,h)anthracene	26.63	278	1138118	41.297	ng	97
91) Benzo(g,h,i)perylene	27.39	276	1037371	42.841	ng	# 94

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

