

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120518\
 Data File : BM017928.D
 Acq On : 05 Dec 2018 07:12
 Operator : JU/SJ
 Sample : PB115307BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB115307BS

Manual Integrations
 APPROVED

mohammad
 12/5/2018 4:50:23 PM

Quant Time: Dec 05 10:09:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM111218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 12 18:37:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	126676	20.00	ng	-0.04
21) Naphthalene-d8	10.35	136	503571	20.00	ng	-0.04
39) Acenaphthene-d10	14.23	164	266448	20.00	ng	-0.04
64) Phenanthrene-d10	16.99	188	558003	20.00	ng	-0.03
76) Chrysene-d12	21.19	240	612978	20.00	ng	-0.02
87) Perylene-d12	23.37	264	693987	20.00	ng	-0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.21	112	573261	76.37	ng	-0.03
7) Phenol-d6	6.77	99	744243	70.99	ng	-0.04
23) Nitrobenzene-d5	8.73	82	668827	68.57	ng	-0.04
42) 2,4,6-Tribromophenol	15.73	330	276059	72.11	ng	-0.03
45) 2-Fluorobiphenyl	12.84	172	1565274	76.21	ng	-0.04
79) Terphenyl-d14	19.63	244	2056283	76.03	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.21	88	95059	30.408	ng	# 86
3) Pyridine	3.59	79	259234	28.270	ng	# 84
4) n-Nitrosodimethylamine	3.50	42	113599	28.123	ng	# 75
6) Aniline	6.92	93	193149	14.857	ng	96
8) 2-Chlorophenol	7.15	128	315443	35.115	ng	94
9) Benzaldehyde	6.74	77	140500	18.398	ng	94
10) Phenol	6.80	94	364683	33.142	ng	95
11) bis(2-Chloroethyl)ether	7.02	93	273288	31.461	ng	94
12) 1,3-Dichlorobenzene	7.47	146	343967	34.129	ng	96
13) 1,4-Dichlorobenzene	7.62	146	345610	33.456	ng	99
14) 1,2-Dichlorobenzene	7.92	146	335523	33.496	ng	98
15) Benzyl Alcohol	7.83	79	252733	30.272	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.10	45	309398	23.394	ng	91
17) 2-Methylphenol	8.03	107	250606	33.808	ng	96
18) Hexachloroethane	8.63	117	109102	29.965	ng	98
19) n-Nitroso-di-n-propylamine	8.38	70	216687	28.787	ng	91
20) 3+4-Methylphenols	8.36	107	330549	32.095	ng	97
22) Acetophenone	8.40	105	434277	34.554	ng	# 92
24) Nitrobenzene	8.77	77	316336	31.964	ng	96
25) Isophorone	9.30	82	565619	37.543	ng	98
26) 2-Nitrophenol	9.48	139	145261	31.865	ng	94
27) 2,4-Dimethylphenol	9.55	122	271522	41.329	ng	98
28) bis(2-Chloroethoxy)methane	9.78	93	351860	34.483	ng	98
29) 2,4-Dichlorophenol	10.01	162	267211	35.038	ng	98
30) 1,2,4-Trichlorobenzene	10.22	180	299490	33.764	ng	97
31) Naphthalene	10.40	128	912980	36.875	ng	99
32) Benzoic acid	9.70	122	167301	28.112	ng	95
33) 4-Chloroaniline	10.53	127	98921	9.123	ng	97
34) Hexachlorobutadiene	10.67	225	186230	33.819	ng	98
35) Caprolactam	11.32	113	82682m	29.440	ng	
36) 4-Chloro-3-methylphenol	11.66	107	283770	32.228	ng	98
37) 2-Methylnaphthalene	12.02	142	648088	37.036	ng	98
38) 1-Methylnaphthalene	12.24	142	619192	36.088	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.39	216	328115	34.922	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.36	237	220310	45.377	ng	99
43) 2,4,6-Trichlorophenol	12.65	196	213484	36.275	ng	96
44) 2,4,5-Trichlorophenol	12.72	196	224525	36.132	ng	99
46) 1,1'-Biphenyl	13.05	154	797717	33.354	ng	99
47) 2-Chloronaphthalene	13.09	162	621439	35.007	ng	99
48) 2-Nitroaniline	13.32	65	175679	32.000	ng	88
49) Acenaphthylene	13.95	152	1000315	37.505	ng	99
50) Dimethylphthalate	13.70	163	731429	33.764	ng	98
51) 2,6-Dinitrotoluene	13.82	165	152602	31.635	ng	93
52) Acenaphthene	14.29	154	575564	35.163	ng	99
53) 3-Nitroaniline	14.16	138	98498	18.755	ng	96
54) 2,4-Dinitrophenol	14.37	184	94738	36.428	ng	95
55) Dibenzofuran	14.63	168	915062	34.209	ng	98
56) 4-Nitrophenol	14.48	139	262195	60.670	ng	91
57) 2,4-Dinitrotoluene	14.62	165	200693	30.864	ng	96
58) Fluorene	15.29	166	737889	36.033	ng	97
59) 2,3,4,6-Tetrachlorophenol	14.86	232	206432	36.118	ng	99
60) Diethylphthalate	15.07	149	736180	31.452	ng	99
61) 4-Chlorophenyl-phenylether	15.29	204	367881	31.619	ng	100
62) 4-Nitroaniline	15.33	138	159606	32.246	ng	94
63) Azobenzene	15.58	77	677283	31.301	ng	94
65) 4,6-Dinitro-2-methylphenol	15.39	198	69614	18.009	ng	93
66) n-Nitrosodiphenylamine	15.50	169	635645	35.431	ng	99
67) 4-Bromophenyl-phenylether	16.18	248	231849	35.132	ng	98
68) Hexachlorobenzene	16.29	284	254780	34.246	ng	92
69) Atrazine	16.47	200	220147	33.363	ng	95
70) Pentachlorophenol	16.64	266	308516	59.184	ng	99
71) Phenanthrene	17.03	178	1140058	37.652	ng	99
72) Anthracene	17.12	178	1176984	39.947	ng	99
73) Carbazole	17.40	167	1055133	35.440	ng	100
74) Di-n-butylphthalate	17.96	149	1234939	37.261	ng	99
75) Fluoranthene	19.06	202	1328079	38.766	ng	99
77) Benzidine	19.26	184	1214158	60.827	ng	99
78) Pyrene	19.42	202	1381117	36.517	ng	99
80) Butylbenzylphthalate	20.33	149	577018	37.563	ng	98
81) Benzo(a)anthracene	21.18	228	1337536	37.875	ng	99
82) 3,3'-Dichlorobenzidine	21.12	252	516507	37.896	ng	97
83) Chrysene	21.23	228	1273857	37.139	ng	99
84) Bis(2-ethylhexyl)phthalate	21.12	149	824728	36.317	ng	99
85) Di-n-octyl phthalate	21.98	149	1449936	39.873	ng	# 95
86) Indeno(1,2,3-cd)pyrene	25.56	276	1772202	64.657	ng	100
88) Benzo(b)fluoranthene	22.73	252	1424695	35.055	ng	99
89) Benzo(k)fluoranthene	22.77	252	1375006	33.473	ng	99
90) Benzo(a)pyrene	23.29	252	1413466	38.143	ng	99
91) Dibenzo(a,h)anthracene	25.57	278	1502503	48.228	ng	99
92) Benzo(g,h,i)perylene	26.22	276	1506369	51.418	ng	99

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

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