

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011019\
 Data File : BM018531.D
 Acq On : 11 Jan 2019 07:34
 Operator : JU/SJ
 Sample : PB116287BSD
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB116287BSD

Manual Integrations
 APPROVED

Sohil
 1/11/2019 4:41:16 PM

Quant Time: Jan 11 10:29:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 10 05:57:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	491532	20.00	ng	0.00
21) Naphthalene-d8	10.55	136	2100418	20.00	ng	0.00
39) Acenaphthene-d10	14.41	164	1233879	20.00	ng	0.00
64) Phenanthrene-d10	17.16	188	2783571	20.00	ng	0.00
76) Chrysene-d12	21.35	240	2530865	20.00	ng	0.00
87) Perylene-d12	23.64	264	2238508	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.35	112	1935623	72.71	ng	0.00
7) Phenol-d6	6.93	99	2546882	75.33	ng	0.00
23) Nitrobenzene-d5	8.93	82	2290956	63.23	ng	0.00
42) 2,4,6-Tribromophenol	15.90	330	1021679	65.03	ng	0.00
45) 2-Fluorobiphenyl	13.02	172	6091278	64.42	ng	0.00
79) Terphenyl-d14	19.79	244	7857447	65.45	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.29	88	247873	22.846	ng	96
3) Pyridine	3.68	79	722291	26.703	ng	95
4) n-Nitrosodimethylamine	3.60	42	376014	33.617	ng	# 94
6) Aniline	7.10	93	616767	16.353	ng	99
8) 2-Chlorophenol	7.33	128	1123997	36.153	ng	94
9) Benzaldehyde	6.91	77	310999	13.338	ng	98
10) Phenol	6.96	94	1242392	36.161	ng	98
11) bis(2-Chloroethyl)ether	7.19	93	850767	31.515	ng	95
12) 1,3-Dichlorobenzene	7.65	146	1113716	30.080	ng	99
13) 1,4-Dichlorobenzene	7.79	146	1138806	30.347	ng	99
14) 1,2-Dichlorobenzene	8.11	146	1102025	30.970	ng	98
15) Benzyl Alcohol	8.00	79	881440	36.095	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.29	45	1083647	31.411	ng	94
17) 2-Methylphenol	8.20	107	875528	37.542	ng	99
18) Hexachloroethane	8.83	117	397275	29.691	ng	94
19) n-Nitroso-di-n-propylamine	8.56	70	716607	32.569	ng	94
20) 3+4-Methylphenols	8.53	107	1210891	37.612	ng	97
22) Acetophenone	8.59	105	1464283	28.181	ng	# 98
24) Nitrobenzene	8.97	77	1084643	30.425	ng	97
25) Isophorone	9.49	82	1988418	36.242	ng	98
26) 2-Nitrophenol	9.67	139	518787	30.095	ng	94
27) 2,4-Dimethylphenol	9.73	122	1021460	39.533	ng	98
28) bis(2-Chloroethoxy)methane	9.97	93	1229048	32.427	ng	99
29) 2,4-Dichlorophenol	10.20	162	1104214	35.890	ng	99
30) 1,2,4-Trichlorobenzene	10.41	180	1132437	29.654	ng	99
31) Naphthalene	10.60	128	3290065	32.524	ng	100
32) Benzoic acid	9.86	122	516118m	32.151	ng	
33) 4-Chloroaniline	10.72	127	448542	11.259	ng	98
34) Hexachlorobutadiene	10.87	225	679630	28.171	ng	98
35) Caprolactam	11.52	113	328437	31.398	ng	89
36) 4-Chloro-3-methylphenol	11.85	107	1136061	35.196	ng	100
37) 2-Methylnaphthalene	12.22	142	2503054	35.022	ng	99
38) 1-Methylnaphthalene	12.44	142	2377604	34.382	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.58	216	1299147	30.109	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.55	237	1121855	47.374	ng	100
43) 2,4,6-Trichlorophenol	12.83	196	903439	34.480	ng	100
44) 2,4,5-Trichlorophenol	12.91	196	961470	34.890	ng	96
46) 1,1'-Biphenyl	13.24	154	3230084	29.981	ng	98
47) 2-Chloronaphthalene	13.28	162	2512681	30.076	ng	100
48) 2-Nitroaniline	13.50	65	692633	33.698	ng	93
49) Acenaphthylene	14.13	152	4120019	33.840	ng	100
50) Dimethylphthalate	13.87	163	3265883	32.229	ng	100
51) 2,6-Dinitrotoluene	14.00	165	701656	32.688	ng	88
52) Acenaphthene	14.47	154	2365653	31.756	ng	98
53) 3-Nitroaniline	14.33	138	506161	23.390	ng	95
54) 2,4-Dinitrophenol	14.53	184	200915	22.935	ng	# 91
55) Dibenzofuran	14.81	168	3867045	31.417	ng	98
56) 4-Nitrophenol	14.64	139	1010671	54.061	ng	94
57) 2,4-Dinitrotoluene	14.78	165	936759	30.844	ng	100
58) Fluorene	15.46	166	3121345	34.042	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.03	232	817665	33.475	ng	96
60) Diethylphthalate	15.23	149	3375179	32.993	ng	99
61) 4-Chlorophenyl-phenylether	15.45	204	1613819	30.538	ng	99
62) 4-Nitroaniline	15.50	138	792512	33.575	ng	98
63) Azobenzene	15.74	77	2656380	31.849	ng	97
65) 4,6-Dinitro-2-methylphenol	15.54	198	181088	14.453	ng	90
66) n-Nitrosodiphenylamine	15.67	169	2773975	32.127	ng	99
67) 4-Bromophenyl-phenylether	16.35	248	982157	31.530	ng	97
68) Hexachlorobenzene	16.45	284	1076385	30.867	ng	100
69) Atrazine	16.63	200	1016110	32.552	ng	99
70) Pentachlorophenol	16.80	266	1073026	46.984	ng	98
71) Phenanthrene	17.20	178	4962371	33.517	ng	99
72) Anthracene	17.29	178	5013798	35.216	ng	100
73) Carbazole	17.57	167	4535382	31.006	ng	99
74) Di-n-butylphthalate	18.12	149	5625348	35.109	ng	100
75) Fluoranthene	19.22	202	5705076	33.422	ng	97
77) Benzidine	19.42	184	1528613	24.461	ng	100
78) Pyrene	19.59	202	5734647	38.227	ng	99
80) Butylbenzylphthalate	20.49	149	2570589	40.157	ng	97
81) Benzo(a)anthracene	21.34	228	5126273	34.382	ng	99
82) 3,3'-Dichlorobenzidine	21.27	252	1157922	20.538	ng	100
83) Chrysene	21.39	228	4727879	32.838	ng	99
84) Bis(2-ethylhexyl)phthalate	21.26	149	3628685	39.256	ng	99
85) Di-n-octyl phthalate	22.15	149	6022761	38.739	ng	# 94
86) Indeno(1,2,3-cd)pyrene	25.96	276	4744569	28.578	ng	96
88) Benzo(b)fluoranthene	22.94	252	4488502	35.319	ng	99
89) Benzo(k)fluoranthene	22.99	252	4358811	34.034	ng	98
90) Benzo(a)pyrene	23.54	252	4203200	34.547	ng	98
91) Dibenzo(a,h)anthracene	25.97	278	4010325	32.531	ng	98
92) Benzo(g,h,i)perylene	26.67	276	3913927	32.626	ng	97

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

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