

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052319\
 Data File : BM020520.D
 Acq On : 23 May 2019 17:44
 Operator : JU/SJ
 Sample : PB119565BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB119565BS

Manual Integrations
 APPROVED

Sohil
 5/24/2019 5:01:38 PM

Quant Time: May 24 03:12:42 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 23 13:59:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	53675	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	188280	20.00	ng	0.00
39) Acenaphthene-d10	14.33	164	115677	20.00	ng	0.00
64) Phenanthrene-d10	17.09	188	286074	20.00	ng	0.00
76) Chrysene-d12	21.28	240	307511	20.00	ng	0.00
87) Perylene-d12	23.53	264	351888	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.25	112	397682	121.11	ng	0.00
7) Phenol-d6	6.85	99	515520	114.92	ng	0.00
23) Nitrobenzene-d5	8.84	82	287490	60.28	ng	0.00
42) 2,4,6-Tribromophenol	15.83	330	239375	133.32	ng	0.00
45) 2-Fluorobiphenyl	12.94	172	589751	62.73	ng	0.00
79) Terphenyl-d14	19.72	244	1101845	62.50	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.17	88	73283	45.504	ng	# 83
3) Pyridine	3.58	79	150894	36.635	ng	# 88
4) n-Nitrosodimethylamine	3.50	42	35904	27.184	ng	# 51
6) Aniline	7.01	93	142246	25.190	ng	# 94
8) 2-Chlorophenol	7.23	128	129931	36.340	ng	# 94
9) Benzaldehyde	6.83	77	53890	15.872	ng	# 91
10) Phenol	6.87	94	173773	35.983	ng	# 85
11) bis(2-Chloroethyl)ether	7.10	93	121253	34.545	ng	# 94
12) 1,3-Dichlorobenzene	7.56	146	148516	35.701	ng	# 93
13) 1,4-Dichlorobenzene	7.70	146	150311	35.768	ng	# 97
14) 1,2-Dichlorobenzene	8.02	146	141719	35.397	ng	# 96
15) Benzyl Alcohol	7.92	79	123126	28.464	ng	# 98
16) 2,2'-oxybis(1-Chloropropan	8.20	45	79462	27.302	ng	# 91
17) 2-Methylphenol	8.12	107	105134	33.818	ng	# 95
18) Hexachloroethane	8.73	117	56578	32.289	ng	# 87
19) n-Nitroso-di-n-propylamine	8.48	70	108539	28.280	ng	# 94
20) 3+4-Methylphenols	8.45	107	139710	33.814	ng	# 96
22) Acetophenone	8.50	105	189415	36.810	ng	# 95
24) Nitrobenzene	8.88	77	167536	33.882	ng	# 95
25) Isophorone	9.40	82	281816	34.267	ng	# 98
26) 2-Nitrophenol	9.59	139	61362	41.106	ng	# 91
27) 2,4-Dimethylphenol	9.65	122	105081	42.282	ng	# 93
28) bis(2-Chloroethoxy)methane	9.89	93	151836	33.899	ng	# 98
29) 2,4-Dichlorophenol	10.12	162	122868	39.207	ng	# 98
30) 1,2,4-Trichlorobenzene	10.32	180	150325	38.984	ng	# 98
31) Naphthalene	10.51	128	352226	36.050	ng	# 99
32) Benzoic acid	9.79	122	65192m	38.401	ng	# 96
33) 4-Chloroaniline	10.64	127	77046	19.159	ng	# 96
34) Hexachlorobutadiene	10.77	225	105770	38.775	ng	# 96
35) Caprolactam	11.45	113	29300	31.268	ng	# 97
36) 4-Chloro-3-methylphenol	11.76	107	122387	33.233	ng	# 92
37) 2-Methylnaphthalene	12.13	142	254918	35.788	ng	# 97
38) 1-Methylnaphthalene	12.35	142	243068	34.897	ng	# 98
40) 1,2,4,5-Tetrachlorobenzene	12.50	216	183860	40.625	ng	# 96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.46	237	172161	64.606	nq	97
43) 2,4,6-Trichlorophenol	12.75	196	108718	42.268	nq	97
44) 2,4,5-Trichlorophenol	12.83	196	117294	40.821	nq	96
46) 1,1'-Biphenyl	13.16	154	339079	35.613	nq	97
47) 2-Chloronaphthalene	13.20	162	277697	36.529	nq	97
48) 2-Nitroaniline	13.43	65	73716	27.237	nq	88
49) Acenaphthylene	14.05	152	415249	34.131	nq	99
50) Dimethylphthalate	13.79	163	341103	34.694	nq	99
51) 2,6-Dinitrotoluene	13.93	165	72976	35.241	nq	90
52) Acenaphthene	14.39	154	242206	34.716	nq	98
53) 3-Nitroaniline	14.26	138	38252	19.903	nq	97
54) 2,4-Dinitrophenol	14.48	184	80520	77.880	nq	# 85
55) Dibenzofuran	14.73	168	427571	36.333	nq	96
56) 4-Nitrophenol	14.57	139	103137	62.897	nq	95
57) 2,4-Dinitrotoluene	14.72	165	101765	36.167	nq	97
58) Fluorene	15.38	166	336785	34.846	nq	100
59) 2,3,4,6-Tetrachlorophenol	14.96	232	115292	42.223	nq	# 91
60) Diethylphthalate	15.16	149	326846	32.997	nq	98
61) 4-Chlorophenyl-phenylether	15.37	204	209849	38.320	nq	93
62) 4-Nitroaniline	15.43	138	52527m	24.765	nq	
63) Azobenzene	15.67	77	356308	30.492	nq	97
65) 4,6-Dinitro-2-methylphenol	15.49	198	53654	38.787	nq	91
66) n-Nitrosodiphenylamine	15.60	169	296169	35.183	nq	100
67) 4-Bromophenyl-phenylether	16.27	248	133524	38.063	nq	# 91
68) Hexachlorobenzene	16.38	284	157235	38.950	nq	# 93
69) Atrazine	16.55	200	109190	33.192	nq	98
70) Pentachlorophenol	16.73	266	178633	77.340	nq	98
71) Phenanthrene	17.13	178	558649	35.377	nq	100
72) Anthracene	17.22	178	558407	35.367	nq	99
73) Carbazole	17.50	167	470379	33.018	nq	99
74) Di-n-butylphthalate	18.04	149	524756	30.607	nq	98
75) Fluoranthene	19.15	202	717169	35.894	nq	99
77) Benzidine	19.35	184	214975	21.866	nq	98
78) Pyrene	19.52	202	730302	35.373	nq	99
80) Butylbenzylphthalate	20.41	149	234083	31.178	nq	99
81) Benzo(a)anthracene	21.26	228	729523	33.828	nq	98
82) 3,3'-Dichlorobenzidine	21.20	252	199301	24.214	nq	# 97
83) Chrysene	21.31	228	729719	34.890	nq	99
84) Bis(2-ethylhexyl)phthalate	21.18	149	353096	30.308	nq	98
85) Di-n-octyl phthalate	22.06	149	628992	32.483	nq	# 96
86) Indeno(1,2,3-cd)pyrene	25.81	276	995058	39.998	nq	# 100
88) Benzo(b)fluoranthene	22.85	252	831725m	35.793	nq	
89) Benzo(k)fluoranthene	22.89	252	760597m	35.883	nq	
90) Benzo(a)pyrene	23.43	252	780109	36.960	nq	98
91) Dibenzo(a,h)anthracene	25.81	278	844766	38.486	nq	97
92) Benzo(g,h,i)perylene	26.50	276	838096	40.217	nq	97

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

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