

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082718\
 Data File : BM016629.D
 Acq On : 28 Aug 2018 05:51
 Operator : SJ/JU
 Sample : PB112404BS
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB112404BS

Quant Time: Aug 28 07:53:04 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 27 13:01:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	98654	20.00	ng	0.00
21) Naphthalene-d8	10.80	136	389791	20.00	ng	0.00
38) Acenaphthene-d10	14.64	164	295667	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	932322	20.00	ng	0.00
75) Chrysene-d12	21.53	240	1204533	20.00	ng	0.00
86) Perylene-d12	23.81	264	1123392	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.55	112	425389	84.50	ng	0.00
7) Phenol-d6	7.18	99	533871	80.14	ng	0.00
23) Nitrobenzene-d5	9.19	82	783872	91.82	ng	0.00
41) 2,4,6-Tribromophenol	16.14	330	726983	87.96	ng	0.00
44) 2-Fluorobiphenyl	13.26	172	2999277	97.51	ng	0.00
78) Terphenyl-d14	19.97	244	5938972	104.36	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.37	88	92988	33.433	ng	# 100
3) Pyridine	3.80	79	181516	31.368	ng	# 81
4) n-Nitrosodimethylamine	3.72	42	119709	46.435	ng	# 66
6) Aniline	7.34	93	169592	22.749	ng	90
8) 2-Chlorophenol	7.56	128	240152	40.089	ng	94
9) Benzaldehyde	7.16	77	139559	30.186	ng	# 72
10) Phenol	7.21	94	244236	37.071	ng	86
11) bis(2-Chloroethyl)ether	7.43	93	179737	36.381	ng	95
12) 1,3-Dichlorobenzene	7.88	146	280215	35.281	ng	# 82
13) 1,4-Dichlorobenzene	8.03	146	283267	34.753	ng	# 92
14) 1,2-Dichlorobenzene	8.35	146	289392	36.139	ng	95
15) Benzyl Alcohol	8.26	79	219453	35.472	ng	# 79
16) 2,2'-oxybis(1-Chloropropan	8.52	45	120598	35.451	ng	# 43
17) 2-Methylphenol	8.46	107	187190	39.236	ng	# 75
18) Hexachloroethane	9.06	117	168139	44.777	ng	# 50
19) n-Nitroso-di-n-propylamine	8.82	70	195902	37.507	ng	# 85
20) 3+4-Methylphenols	8.80	107	248525	37.675	ng	# 81
22) Acetophenone	8.85	105	375714	40.028	ng	# 98
24) Nitrobenzene	9.23	77	346127	40.579	ng	# 89
25) Isophorone	9.75	82	539078	47.247	ng	# 91
26) 2-Nitrophenol	9.94	139	134598	36.911	ng	# 26
27) 2,4-Dimethylphenol	9.99	122	226872	47.741	ng	# 76
28) bis(2-Chloroethoxy)methane	10.22	93	273846	40.207	ng	96
29) 2,4-Dichlorophenol	10.47	162	330786	46.509	ng	94
30) 1,2,4-Trichlorobenzene	10.67	180	519942	48.178	ng	# 95
31) Naphthalene	10.86	128	784735	42.255	ng	99
32) Benzoic acid	10.22	122	122369	30.776	ng	# 80
33) 4-Chloroaniline	11.00	127	96148	12.302	ng	# 87
34) Hexachlorobutadiene	11.10	225	546484	53.054	ng	97
35) Caprolactam	11.82	113	54812	32.117	ng	# 62
36) 4-Chloro-3-methylphenol	12.12	107	277764	40.624	ng	94
37) 2-Methylnaphthalene	12.46	142	669116	46.024	ng	# 85
39) 1,2,4,5-Tetrachlorobenzene	12.83	216	831096	49.771	ng	# 100
40) Hexachlorocyclopentadiene	12.78	237	959500	110.226	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.09	196	467511	46.819	nq	92
43) 2,4,5-Trichlorophenol	13.17	196	457364	46.890	nq #	93
45) 1,1'-Biphenyl	13.47	154	964150	37.189	nq	94
46) 2-Chloronaphthalene	13.52	162	805262	38.232	nq	95
47) 2-Nitroaniline	13.76	65	203719	36.312	nq #	82
48) Acenaphthylene	14.37	152	1190954	40.057	nq	98
49) Dimethylphthalate	14.11	163	1175298	41.375	nq #	98
50) 2,6-Dinitrotoluene	14.25	165	224025	40.434	nq #	70
51) Acenaphthene	14.71	154	706594	38.684	nq	95
52) 3-Nitroaniline	14.59	138	91700	18.190	nq #	68
53) 2,4-Dinitrophenol	14.82	184	174671	62.997	nq #	70
54) Dibenzofuran	15.04	168	1446821	41.378	nq	94
55) 4-Nitrophenol	14.91	139	193485	60.174	nq #	82
56) 2,4-Dinitrotoluene	15.04	165	355724	38.503	nq #	91
57) Fluorene	15.69	166	1193211	45.172	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.28	232	463582	49.449	nq #	100
59) Diethylphthalate	15.46	149	1118680	37.770	nq	94
60) 4-Chlorophenyl-phenylether	15.67	204	1028976	46.451	nq #	86
61) 4-Nitroaniline	15.76	138	138576	28.126	nq #	1
62) Azobenzene	15.97	77	1367200	43.339	nq	91
64) 4,6-Dinitro-2-methylphenol	15.82	198	213672	30.163	nq #	59
65) n-Nitrosodiphenylamine	15.90	169	1008117	38.436	nq	97
66) 4-Bromophenyl-phenylether	16.57	248	641718	44.796	nq	98
67) Hexachlorobenzene	16.69	284	716806	42.907	nq #	93
68) Atrazine	16.85	200	595216	49.726	nq	90
69) Pentachlorophenol	17.04	266	601555	68.054	nq	100
70) Phenanthrene	17.43	178	1996784	41.438	nq	98
71) Anthracene	17.52	178	2008491	41.975	nq	98
72) Carbazole	17.80	167	1371394	31.891	nq	99
73) Di-n-butylphthalate	18.32	149	1864638	35.567	nq #	97
74) Fluoranthene	19.43	202	2899506	41.038	nq	94
76) Benzidine	19.62	184	818887	35.351	nq	98
77) Pyrene	19.79	202	2926671	44.708	nq	98
79) Butylbenzylphthalate	20.66	149	861448	38.124	nq #	71
80) Benzo(a)anthracene	21.52	228	3031739	42.811	nq	99
81) 3,3'-Dichlorobenzidine	21.44	252	629425	19.968	nq #	95
82) Chrysene	21.57	228	2744719	40.621	nq	99
83) Bis(2-ethylhexyl)phthalate	21.40	149	1224901	36.465	nq #	95
84) Di-n-octyl phthalate	22.29	149	1970241	35.295	nq #	94
85) Indeno(1,2,3-cd)pyrene	26.12	276	3365106	34.631	nq #	92
87) Benzo(b)fluoranthene	23.12	252	3041643	46.969	nq #	91
88) Benzo(k)fluoranthene	23.17	252	2877845	43.676	nq #	93
89) Benzo(a)pyrene	23.71	252	2785880	44.213	nq #	94
90) Dibenzo(a,h)anthracene	26.11	278	2842019	40.246	nq #	87
91) Benzo(g,h,i)perylene	26.83	276	2588293	40.832	nq #	81

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

