

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031119\
 Data File : BM019097.D
 Acq On : 11 Mar 2019 17:28
 Operator : JU/SJ
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM031119

Manual Integrations
 APPROVED

Sohil
 3/12/2019 4:45:14 PM

Quant Time: Mar 11 18:29:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 11 17:30:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	201371	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	927397	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	547836	20.00	ng	0.00
64) Phenanthrene-d10	17.07	188	1245534	20.00	ng	0.00
76) Chrysene-d12	21.28	240	1387463	20.00	ng	0.00
87) Perylene-d12	23.52	264	1412362	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.26	112	982585	79.02	ng	0.00
7) Phenol-d6	6.85	99	1475135	81.11	ng	0.00
23) Nitrobenzene-d5	8.84	82	1579089	80.49	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	681560	84.26	ng	0.00
45) 2-Fluorobiphenyl	12.93	172	3392390	80.55	ng	0.00
79) Terphenyl-d14	19.72	244	5557750	79.52	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.22	88	205681	37.184	ng	100
3) Pyridine	3.62	79	546643	36.324	ng	99
4) n-Nitrosodimethylamine	3.55	42	177735	37.990	ng	98
6) Aniline	7.01	93	946327	40.300	ng	99
8) 2-Chlorophenol	7.23	128	596465	39.918	ng	99
9) Benzaldehyde	6.83	77	469818	41.067	ng	100
10) Phenol	6.87	94	791287	40.369	ng	100
11) bis(2-Chloroethyl)ether	7.11	93	599907	40.109	ng	99
12) 1,3-Dichlorobenzene	7.55	146	626448	39.616	ng	99
13) 1,4-Dichlorobenzene	7.70	146	632070	39.207	ng	98
14) 1,2-Dichlorobenzene	8.02	146	613247	39.098	ng	99
15) Benzyl Alcohol	7.92	79	616258	40.940	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.20	45	593951	38.891	ng	99
17) 2-Methylphenol	8.12	107	517846	40.542	ng	99
18) Hexachloroethane	8.73	117	238942	39.887	ng	97
19) n-Nitroso-di-n-propylamine	8.48	70	611648	40.992	ng	98
20) 3+4-Methylphenols	8.45	107	708414	40.859	ng	99
22) Acetophenone	8.50	105	1035456	40.401	ng	# 99
24) Nitrobenzene	8.88	77	826390	40.501	ng	98
25) Isophorone	9.40	82	1529521	40.857	ng	100
26) 2-Nitrophenol	9.58	139	356701	42.175	ng	96
27) 2,4-Dimethylphenol	9.64	122	513182	40.814	ng	100
28) bis(2-Chloroethoxy)methane	9.88	93	899535	40.580	ng	99
29) 2,4-Dichlorophenol	10.10	162	581359	41.748	ng	99
30) 1,2,4-Trichlorobenzene	10.32	180	625403	39.829	ng	100
31) Naphthalene	10.50	128	1944830	39.787	ng	100
32) Benzoic acid	9.81	122	426926m	40.092	ng	
33) 4-Chloroaniline	10.63	127	821655	41.006	ng	99
34) Hexachlorobutadiene	10.77	225	359662	39.914	ng	98
35) Caprolactam	11.45	113	213516m	43.042	ng	
36) 4-Chloro-3-methylphenol	11.76	107	701853	40.847	ng	98
37) 2-Methylnaphthalene	12.12	142	1420460	39.990	ng	99
38) 1-Methylnaphthalene	12.35	142	1404011	40.102	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.49	216	703458	40.593	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.46	237	326118	41.546	ng	100
43) 2,4,6-Trichlorophenol	12.75	196	466848	41.686	ng	99
44) 2,4,5-Trichlorophenol	12.82	196	510548	42.633	ng	99
46) 1,1'-Biphenyl	13.15	154	1921889	40.171	ng	99
47) 2-Chloronaphthalene	13.19	162	1447147	40.523	ng	99
48) 2-Nitroaniline	13.42	65	514452	42.954	ng	98
49) Acenaphthylene	14.05	152	2463083	40.791	ng	100
50) Dimethylphthalate	13.79	163	1909903	40.753	ng	99
51) 2,6-Dinitrotoluene	13.92	165	426070	42.057	ng	96
52) Acenaphthene	14.39	154	1397333	40.273	ng	100
53) 3-Nitroaniline	14.26	138	438471	40.865	ng	98
54) 2,4-Dinitrophenol	14.46	184	250365	42.554	ng	98
55) Dibenzofuran	14.72	168	2249932	40.210	ng	99
56) 4-Nitrophenol	14.56	139	370872	42.334	ng	99
57) 2,4-Dinitrotoluene	14.71	165	602638	40.980	ng	98
58) Fluorene	15.37	166	1876584	40.687	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.95	232	461073	41.357	ng	100
60) Diethylphthalate	15.16	149	1943131	40.939	ng	99
61) 4-Chlorophenyl-phenylether	15.37	204	905193	40.409	ng	99
62) 4-Nitroaniline	15.43	138	450177	41.632	ng	99
63) Azobenzene	15.66	77	2058918	41.223	ng	99
65) 4,6-Dinitro-2-methylphenol	15.47	198	333682	40.883	ng	96
66) n-Nitrosodiphenylamine	15.59	169	1588687	40.064	ng	99
67) 4-Bromophenyl-phenylether	16.27	248	551470	40.150	ng	98
68) Hexachlorobenzene	16.37	284	657396	40.258	ng	99
69) Atrazine	16.55	200	552960	41.425	ng	99
70) Pentachlorophenol	16.72	266	398223	40.477	ng	100
71) Phenanthrene	17.12	178	2912465	39.894	ng	100
72) Anthracene	17.21	178	2876087	40.242	ng	99
73) Carbazole	17.49	167	2747205	40.727	ng	99
74) Di-n-butylphthalate	18.04	149	3353653	41.114	ng	99
75) Fluoranthene	19.14	202	3377799	40.322	ng	99
77) Benzidine	19.34	184	1876386	47.685	ng	99
78) Pyrene	19.51	202	3517615	40.070	ng	100
80) Butylbenzylphthalate	20.41	149	1583684	41.419	ng	98
81) Benzo(a)anthracene	21.26	228	3461994	39.773	ng	100
82) 3,3'-Dichlorobenzidine	21.20	252	1385528	41.319	ng	99
83) Chrysene	21.32	228	3365894	39.970	ng	100
84) Bis(2-ethylhexyl)phthalate	21.18	149	2306810	41.469	ng	99
85) Di-n-octyl phthalate	22.06	149	3956142	42.235	ng	100
86) Indeno(1,2,3-cd)pyrene	25.79	276	3859491	42.576	ng	100
88) Benzo(b)fluoranthene	22.84	252	3710920	40.549	ng	100
89) Benzo(k)fluoranthene	22.89	252	3326836	39.522	ng	99
90) Benzo(a)pyrene	23.42	252	3312912	40.282	ng	100
91) Dibenzo(a,h)anthracene	25.80	278	3258784	41.421	ng	100
92) Benzo(g,h,i)perylene	26.49	276	2950643	40.850	ng	100

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

