

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082118\
 Data File : BM016482.D
 Acq On : 21 Aug 2018 12:45
 Operator : SJ/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Sohil
 8/22/2018 6:06:44 PM

Quant Time: Aug 21 15:02:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 21 14:43:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	123335	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	503976	20.00	ng	0.00
38) Acenaphthene-d10	14.66	164	335689	20.00	ng	0.00
63) Phenanthrene-d10	17.40	188	1067328	20.00	ng	0.00
75) Chrysene-d12	21.54	240	1521279	20.00	ng	0.00
86) Perylene-d12	23.84	264	1815038	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.57	112	133233m	18.15	ng	0.01
7) Phenol-d6	7.19	99	176730	17.92	ng	0.00
23) Nitrobenzene-d5	9.21	82	234167	19.35	ng	0.00
41) 2,4,6-Tribromophenol	16.16	330	171482	33.82	ng	0.00
44) 2-Fluorobiphenyl	13.28	172	656723	20.34	ng	0.00
78) Terphenyl-d14	19.99	244	1467855	14.58	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.38	88	39931m	10.865	ng	
3) Pyridine	3.83	79	75318m	10.179	ng	
4) n-Nitrosodimethylamine	3.75	42	32511m	3.813	ng	
6) Aniline	7.36	93	105217	10.308	ng	# 90
8) 2-Chlorophenol	7.59	128	81885m	9.050	ng	
9) Benzaldehyde	7.19	77	72247m	11.139	ng	
10) Phenol	7.22	94	86152	9.016	ng	93
11) bis(2-Chloroethyl)ether	7.45	93	63815	8.797	ng	94
12) 1,3-Dichlorobenzene	7.91	146	109042	10.515	ng	# 91
13) 1,4-Dichlorobenzene	8.06	146	114349	10.908	ng	93
14) 1,2-Dichlorobenzene	8.37	146	106658	10.357	ng	97
15) Benzyl Alcohol	8.29	79	91127m	10.178	ng	
16) 2,2'-oxybis(1-Chloropropan	8.56	45	46077	5.046	ng	# 47
17) 2-Methylphenol	8.48	107	64261	9.307	ng	# 73
18) Hexachloroethane	9.08	117	52657	9.754	ng	# 57
19) n-Nitroso-di-n-propylamine	8.83	70	65281	8.717	ng	# 86
20) 3+4-Methylphenols	8.83	107	84974m	8.955	ng	
22) Acetophenone	8.87	105	129390	9.573	ng	# 97
24) Nitrobenzene	9.26	77	123325m	10.572	ng	
25) Isophorone	9.76	82	153002m	9.480	ng	
26) 2-Nitrophenol	9.98	139	43190m	8.801	ng	
27) 2,4-Dimethylphenol	10.02	122	70979m	10.820	ng	
28) bis(2-Chloroethoxy)methane	10.25	93	92760	9.935	ng	# 87
29) 2,4-Dichlorophenol	10.51	162	94789m	10.688	ng	
30) 1,2,4-Trichlorobenzene	10.69	180	137750	13.135	ng	94
31) Naphthalene	10.88	128	250702	9.808	ng	96
32) Benzoic acid	10.16	122	47223m	8.863	ng	
33) 4-Chloroaniline	11.02	127	106875m	9.756	ng	
34) Hexachlorobutadiene	11.13	225	134922	14.925	ng	90
35) Caprolactam	11.82	113	20841m	9.167	ng	
36) 4-Chloro-3-methylphenol	12.14	107	88211m	8.723	ng	
37) 2-Methylnaphthalene	12.49	142	198344	10.460	ng	# 88
39) 1,2,4,5-Tetrachlorobenzene	12.86	216	201503	14.663	ng	# 100
40) Hexachlorocyclopentadiene	12.80	237	35315	8.281	ng	92

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.11	196	113433	13.207	ng	93
43) 2,4,5-Trichlorophenol	13.20	196	103925	12.402	ng	# 90
45) 1,1'-Biphenyl	13.49	154	301263	9.696	ng	96
46) 2-Chloronaphthalene	13.55	162	247851	10.225	ng	96
47) 2-Nitroaniline	13.79	65	65974m	7.849	ng	
48) Acenaphthylene	14.39	152	335900	9.216	ng	99
49) Dimethylphthalate	14.12	163	338131	10.271	ng	# 98
50) 2,6-Dinitrotoluene	14.27	165	63789	10.209	ng	# 67
51) Acenaphthene	14.73	154	213031	9.685	ng	93
52) 3-Nitroaniline	14.62	138	55138m	8.957	ng	
53) 2,4-Dinitrophenol	14.86	184	26792m	9.831	ng	
54) Dibenzofuran	15.06	168	389092	10.039	ng	# 77
55) 4-Nitrophenol	14.96	139	29215m	7.265	ng	
56) 2,4-Dinitrotoluene	15.07	165	93699	8.971	ng	98
57) Fluorene	15.72	166	285526	9.643	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.30	232	106555	13.572	ng	# 100
59) Diethylphthalate	15.47	149	350678	9.388	ng	99
60) 4-Chlorophenyl-phenylether	15.70	204	248141	13.246	ng	95
61) 4-Nitroaniline	15.77	138	56582m	9.506	ng	
62) Azobenzene	15.99	77	390082	9.055	ng	91
64) 4,6-Dinitro-2-methylphenol	15.84	198	72509m	9.742	ng	
65) n-Nitrosodiphenylamine	15.92	169	288201	7.980	ng	98
66) 4-Bromophenyl-phenylether	16.59	248	167235	11.650	ng	97
67) Hexachlorobenzene	16.71	284	184872	13.123	ng	# 93
68) Atrazine	16.86	200	154049	11.635	ng	93
69) Pentachlorophenol	17.06	266	92292	11.442	ng	94
70) Phenanthrene	17.45	178	529867	8.762	ng	99
71) Anthracene	17.54	178	524327	8.734	ng	98
72) Carbazole	17.82	167	468918	8.070	ng	99
73) Di-n-butylphthalate	18.34	149	586226	7.030	ng	# 95
74) Fluoranthene	19.44	202	776116	10.624	ng	95
76) Benzidine	19.64	184	351707	10.504	ng	97
77) Pyrene	19.81	202	825555	8.224	ng	99
79) Butylbenzylphthalate	20.67	149	278110	5.911	ng	# 82
80) Benzo(a)anthracene	21.53	228	928005	9.428	ng	99
81) 3,3'-Dichlorobenzidine	21.46	252	406047	11.588	ng	# 97
82) Chrysene	21.59	228	876241	9.473	ng	100
83) Bis(2-ethylhexyl)phthalate	21.42	149	420020	6.258	ng	# 88
84) Di-n-octyl phthalate	22.31	149	714904	7.101	ng	# 92
85) Indeno(1,2,3-cd)pyrene	26.16	276	1291050	17.064	ng	# 94
87) Benzo(b)fluoranthene	23.14	252	1023919	8.517	ng	# 92
88) Benzo(k)fluoranthene	23.19	252	1054716	8.778	ng	# 94
89) Benzo(a)pyrene	23.74	252	1005619	9.240	ng	# 95
90) Dibenzo(a,h)anthracene	26.16	278	1093672	11.207	ng	# 87
91) Benzo(g,h,i)perylene	26.88	276	1025066	11.687	ng	# 82

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

