

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082118\
 Data File : BM016485.D
 Acq On : 21 Aug 2018 14:34
 Operator : SJ/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

Sohil
 8/22/2018 6:06:52 PM

Quant Time: Aug 21 15:15:16 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	147168	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	601489	20.00	ng	0.00
38) Acenaphthene-d10	14.66	164	397187	20.00	ng	0.00
63) Phenanthrene-d10	17.40	188	1186828	20.00	ng	0.00
75) Chrysene-d12	21.55	240	1782278	20.00	ng	0.00
86) Perylene-d12	23.84	264	1910839	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.56	112	783294m	89.45	ng	0.00
7) Phenol-d6	7.20	99	1026882	87.24	ng	0.00
23) Nitrobenzene-d5	9.21	82	1348232	93.35	ng	0.00
41) 2,4,6-Tribromophenol	16.16	330	1141296	190.25	ng	0.00
44) 2-Fluorobiphenyl	13.28	172	4370655	114.42	ng	0.00
78) Terphenyl-d14	20.00	244	9665525	81.97	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.38	88	208859	47.626	ng	# 100
3) Pyridine	3.80	79	442161m	50.081	ng	
4) n-Nitrosodimethylamine	3.73	42	197351m	19.400	ng	
6) Aniline	7.36	93	560221	45.995	ng	91
8) 2-Chlorophenol	7.59	128	463277	42.910	ng	96
9) Benzaldehyde	7.17	77	333364m	43.074	ng	
10) Phenol	7.23	94	494735	43.391	ng	99
11) bis(2-Chloroethyl)ether	7.45	93	374280	43.240	ng	95
12) 1,3-Dichlorobenzene	7.90	146	600652	48.542	ng	# 89
13) 1,4-Dichlorobenzene	8.06	146	600709	48.023	ng	98
14) 1,2-Dichlorobenzene	8.37	146	610904	49.715	ng	94
15) Benzyl Alcohol	8.29	79	445348	41.685	ng	# 81
16) 2,2'-oxybis(1-Chloropropan	8.54	45	259402	23.807	ng	# 52
17) 2-Methylphenol	8.48	107	352987	42.844	ng	# 77
18) Hexachloroethane	9.09	117	280487	43.543	ng	# 57
19) n-Nitroso-di-n-propylamine	8.84	70	396282	44.345	ng	# 88
20) 3+4-Methylphenols	8.82	107	502497	44.382	ng	82
22) Acetophenone	8.87	105	712525	44.169	ng	# 95
24) Nitrobenzene	9.26	77	662130	47.559	ng	94
25) Isophorone	9.77	82	917734	47.642	ng	# 94
26) 2-Nitrophenol	9.96	139	287843m	49.149	ng	
27) 2,4-Dimethylphenol	10.01	122	381871m	48.775	ng	
28) bis(2-Chloroethoxy)methane	10.25	93	537071	48.198	ng	99
29) 2,4-Dichlorophenol	10.49	162	571527	53.998	ng	97
30) 1,2,4-Trichlorobenzene	10.69	180	851633	68.042	ng	95
31) Naphthalene	10.88	128	1419957	46.547	ng	99
32) Benzoic acid	10.22	122	310579m	48.839	ng	
33) 4-Chloroaniline	11.01	127	609688	46.632	ng	# 88
34) Hexachlorobutadiene	11.13	225	791483	73.361	ng	97
35) Caprolactam	11.84	113	135058m	49.773	ng	
36) 4-Chloro-3-methylphenol	12.13	107	521503	43.208	ng	89
37) 2-Methylnaphthalene	12.49	142	1140505	50.396	ng	# 88
39) 1,2,4,5-Tetrachlorobenzene	12.86	216	1156948	71.153	ng	# 100
40) Hexachlorocyclopentadiene	12.81	237	439353	87.076	ng	95

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082118\
 Data File : BM016485.D
 Acq On : 21 Aug 2018 14:34
 Operator : SJ/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

Sohil
 8/22/2018 6:06:52 PM

Quant Time: Aug 21 15:15:16 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)
42) 2,4,6-Trichlorophenol	13.10	196	722264	71.071	nq	97
43) 2,4,5-Trichlorophenol	13.19	196	689734	69.566	nq #	94
45) 1,1'-Biphenyl	13.50	154	1806841	49.149	nq	98
46) 2-Chloronaphthalene	13.55	162	1462414	50.991	nq	96
47) 2-Nitroaniline	13.77	65	389830	39.196	nq	85
48) Acenaphthylene	14.39	152	2080697	48.248	nq	99
49) Dimethylphthalate	14.13	163	1975243	50.712	nq #	98
50) 2,6-Dinitrotoluene	14.27	165	404234	54.679	nq #	76
51) Acenaphthene	14.73	154	1257980	48.337	nq	95
52) 3-Nitroaniline	14.61	138	345633m	47.451	nq	
53) 2,4-Dinitrophenol	14.83	184	205289	63.666	nq #	75
54) Dibenzofuran	15.07	168	2474122	53.953	nq	93
55) 4-Nitrophenol	14.92	139	223232	46.920	nq #	85
56) 2,4-Dinitrotoluene	15.06	165	632294	51.166	nq #	93
57) Fluorene	15.72	166	1854104	52.922	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.30	232	649810	69.951	nq #	100
59) Diethylphthalate	15.48	149	2043753	46.241	nq	98
60) 4-Chlorophenyl-phenylether	15.70	204	1543090	69.620	nq	92
61) 4-Nitroaniline	15.77	138	340757	48.386	nq #	1
62) Azobenzene	15.99	77	2180495	42.780	nq	92
64) 4,6-Dinitro-2-methylphenol	15.83	198	464477	56.122	nq #	62
65) n-Nitrosodiphenylamine	15.92	169	1744784	43.445	nq	98
66) 4-Bromophenyl-phenylether	16.59	248	938361	58.784	nq	95
67) Hexachlorobenzene	16.71	284	1105486	70.571	nq	93
68) Atrazine	16.87	200	770112	52.309	nq	90
69) Pentachlorophenol	17.06	266	592225	66.027	nq	98
70) Phenanthrene	17.44	178	3209212	47.726	nq	100
71) Anthracene	17.54	178	3206068	48.029	nq	99
72) Carbazole	17.82	167	2859735	44.263	nq	98
73) Di-n-butylphthalate	18.34	149	3575152	38.558	nq #	97
74) Fluoranthene	19.44	202	4725922	58.180	nq	95
76) Benzidine	19.64	184	1718480	43.809	nq	98
77) Pyrene	19.81	202	5031676	42.782	nq	99
79) Butylbenzylphthalate	20.67	149	1774854	32.198	nq #	82
80) Benzo(a)anthracene	21.53	228	5371920	46.586	nq	99
81) 3,3'-Dichlorobenzidine	21.46	252	2410517	58.719	nq #	95
82) Chrysene	21.59	228	5099495	47.057	nq	99
83) Bis(2-ethylhexyl)phthalate	21.43	149	2647408	33.668	nq #	94
84) Di-n-octyl phthalate	22.31	149	4365415	37.013	nq #	92
85) Indeno(1,2,3-cd)pyrene	26.17	276	7284353	82.178	nq #	94
87) Benzo(b)fluoranthene	23.15	252	5708537	45.101	nq #	91
88) Benzo(k)fluoranthene	23.19	252	5828639	46.079	nq #	94
89) Benzo(a)pyrene	23.74	252	5543259	48.382	nq #	95
90) Dibenzo(a,h)anthracene	26.17	278	6263164	60.960	nq #	88
91) Benzo(g,h,i)perylene	26.88	276	5490306	59.456	nq #	82

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

