

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM081318\
 Data File : BM016337.D
 Acq On : 13 Aug 2018 23:27
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Aug 14 00:05:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM072318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 10 18:44:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	29096	20.00	ng	-0.02
21) Naphthalene-d8	10.93	136	118595	20.00	ng	-0.02
38) Acenaphthene-d10	14.76	164	59369	20.00	ng	-0.01
63) Phenanthrene-d10	17.49	188	122152	20.00	ng	-0.02
75) Chrysene-d12	21.62	240	150735	20.00	ng	-0.02
86) Perylene-d12	23.95	264	157706	20.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.65	112	143325	81.83	ng	-0.01
7) Phenol-d6	7.29	99	173542	75.87	ng	-0.01
23) Nitrobenzene-d5	9.31	82	223922	87.82	ng	-0.01
41) 2,4,6-Tribromophenol	16.25	330	54430	72.29	ng	-0.02
44) 2-Fluorobiphenyl	13.37	172	455921	93.44	ng	-0.02
78) Terphenyl-d14	20.07	244	591743	82.18	ng	-0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.45	88	37424	45.601	ng	# 100
3) Pyridine	3.89	79	72179	36.509	ng	# 65
4) n-Nitrosodimethylamine	3.80	42	74218	48.130	ng	# 28
6) Aniline	7.45	93	94566	35.905	ng	# 83
8) 2-Chlorophenol	7.68	128	82262	39.516	ng	97
9) Benzaldehyde	7.27	77	59955	37.305	ng	84
10) Phenol	7.32	94	85804	37.513	ng	95
11) bis(2-Chloroethyl)ether	7.55	93	67829	37.537	ng	85
12) 1,3-Dichlorobenzene	8.00	146	96548	39.622	ng	# 87
13) 1,4-Dichlorobenzene	8.16	146	98069	39.683	ng	# 95
14) 1,2-Dichlorobenzene	8.48	146	94649	39.141	ng	# 92
15) Benzyl Alcohol	8.38	79	72694	39.911	ng	# 80
16) 2,2'-oxybis(1-Chloropropan	8.65	45	88376	31.956	ng	94
17) 2-Methylphenol	8.58	107	60358	38.607	ng	# 77
18) Hexachloroethane	9.19	117	47045	44.039	ng	98
19) n-Nitroso-di-n-propylamine	8.93	70	62851	37.592	ng	# 64
20) 3+4-Methylphenols	8.91	107	80341	35.718	ng	# 84
22) Acetophenone	8.96	105	117952	39.499	ng	# 79
24) Nitrobenzene	9.35	77	105242	41.005	ng	94
25) Isophorone	9.87	82	136123	39.029	ng	# 94
26) 2-Nitrophenol	10.06	139	41702	38.923	ng	# 42
27) 2,4-Dimethylphenol	10.11	122	57165	38.297	ng	# 76
28) bis(2-Chloroethoxy)methane	10.35	93	82986	36.608	ng	98
29) 2,4-Dichlorophenol	10.59	162	71531	40.328	ng	95
30) 1,2,4-Trichlorobenzene	10.80	180	89043	41.317	ng	# 96
31) Naphthalene	10.99	128	230473	38.398	ng	99
32) Benzoic acid	10.30	122	38535	33.866	ng	# 76
33) 4-Chloroaniline	11.12	127	92795	37.886	ng	# 86
34) Hexachlorobutadiene	11.23	225	73647	49.266	ng	97
35) Caprolactam	11.93	113	17003	34.610	ng	# 84
36) 4-Chloro-3-methylphenol	12.23	107	76062	38.988	ng	87
37) 2-Methylnaphthalene	12.59	142	155716	38.728	ng	# 86
39) 1,2,4,5-Tetrachlorobenzene	12.95	216	95015	47.009	ng	# 100
40) Hexachlorocyclopentadiene	12.90	237	16735	22.923	ng	98

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM081318\
 Data File : BM016337.D
 Acq On : 13 Aug 2018 23:27
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Aug 14 00:05:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM072318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 10 18:44:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.20	196	55792	44.119	ng	98
43) 2,4,5-Trichlorophenol	13.28	196	55319	42.410	ng #	79
45) 1,1'-Biphenyl	13.59	154	225229	42.037	ng	99
46) 2-Chloronaphthalene	13.64	162	173194	41.561	ng #	88
47) 2-Nitroaniline	13.86	65	58042	42.129	ng #	74
48) Acenaphthylene	14.48	152	247236	40.527	ng	98
49) Dimethylphthalate	14.22	163	216822	41.731	ng	99
50) 2,6-Dinitrotoluene	14.36	165	42482	40.637	ng #	49
51) Acenaphthene	14.82	154	148277	39.520	ng	96
52) 3-Nitroaniline	14.69	138	40843	36.172	ng #	71
53) 2,4-Dinitrophenol	14.92	184	12166	31.148	ng #	59
54) Dibenzofuran	15.15	168	249483	40.351	ng #	77
55) 4-Nitrophenol	15.02	139	23769	29.364	ng #	83
56) 2,4-Dinitrotoluene	15.15	165	61244	40.972	ng #	56
57) Fluorene	15.80	166	184503	40.157	ng	100
58) 2,3,4,6-Tetrachlorophenol	15.39	232	47264	42.155	ng #	100
59) Diethylphthalate	15.56	149	238663	42.617	ng	95
60) 4-Chlorophenyl-phenylether	15.79	204	110990	43.395	ng #	81
61) 4-Nitroaniline	15.86	138	38849	35.860	ng #	1
62) Azobenzene	16.08	77	268564	45.536	ng #	73
64) 4,6-Dinitro-2-methylphenol	15.92	198	26010	35.081	ng	84
65) n-Nitrosodiphenylamine	16.00	169	168046	41.779	ng	96
66) 4-Bromophenyl-phenylether	16.67	248	62061	44.868	ng #	72
67) Hexachlorobenzene	16.80	284	60993	40.042	ng #	56
68) Atrazine	16.95	200	60465	41.943	ng	96
69) Pentachlorophenol	17.15	266	33987	36.033	ng	96
70) Phenanthrene	17.53	178	263909	38.590	ng	98
71) Anthracene	17.63	178	262565	39.504	ng	97
72) Carbazole	17.90	167	265141	38.506	ng #	95
73) Di-n-butylphthalate	18.42	149	368525	42.951	ng #	96
74) Fluoranthene	19.53	202	306164	40.134	ng	97
76) Benzidine	19.72	184	141399	33.569	ng	98
77) Pyrene	19.89	202	326924	36.192	ng	98
79) Butylbenzylphthalate	20.74	149	176916	38.553	ng #	76
80) Benzo(a)anthracene	21.60	228	373707	40.253	ng	98
81) 3,3'-Dichlorobenzidine	21.53	252	145134	38.812	ng	96
82) Chrysene	21.66	228	353948	39.744	ng	99
83) Bis(2-ethylhexyl)phthalate	21.49	149	258405	38.855	ng	90
84) Di-n-octyl phthalate	22.39	149	446891	39.982	ng #	88
85) Indeno(1,2,3-cd)pyrene	26.33	276	455348	43.087	ng	98
87) Benzo(b)fluoranthene	23.24	252	371240	39.981	ng #	93
88) Benzo(k)fluoranthene	23.29	252	373189	39.658	ng #	96
89) Benzo(a)pyrene	23.85	252	362935	40.659	ng	97
90) Dibenzo(a,h)anthracene	26.33	278	386460	41.787	ng #	93
91) Benzo(g,h,i)perylene	27.06	276	352881	40.343	ng #	90

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

