

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082218\
 Data File : BM016541.D
 Acq On : 23 Aug 2018 04:48
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

Sohil
 8/23/2018 6:39:45 PM

Quant Time: Aug 23 07:51:20 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 16:31:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	123560	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	518456	20.00	ng	0.00
38) Acenaphthene-d10	14.66	164	435362	20.00	ng	-0.01
63) Phenanthrene-d10	17.40	188	1439873	20.00	ng	0.00
75) Chrysene-d12	21.54	240	2233507	20.00	ng	0.00
86) Perylene-d12	23.83	264	2319478	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.55	112	452687	71.79	ng	0.00
7) Phenol-d6	7.19	99	625523	74.97	ng	-0.01
23) Nitrobenzene-d5	9.20	82	963679	84.87	ng	0.00
41) 2,4,6-Tribromophenol	16.15	330	995923	81.83	ng	0.00
44) 2-Fluorobiphenyl	13.27	172	3930809	86.79	ng	0.00
78) Terphenyl-d14	19.99	244	9717096	92.08	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.37	88	133644	38.365	ng	# 100
3) Pyridine	3.80	79	251806	34.743	ng	# 82
4) n-Nitrosodimethylamine	3.73	42	130700	40.479	ng	# 63
6) Aniline	7.35	93	347966	37.268	ng	# 89
8) 2-Chlorophenol	7.58	128	290916	38.774	ng	# 95
9) Benzaldehyde	7.17	77	233552	40.333	ng	# 82
10) Phenol	7.22	94	306743	37.174	ng	# 92
11) bis(2-Chloroethyl)ether	7.45	93	225936	36.514	ng	# 93
12) 1,3-Dichlorobenzene	7.90	146	383095	38.512	ng	# 85
13) 1,4-Dichlorobenzene	8.05	146	387008	37.910	ng	# 96
14) 1,2-Dichlorobenzene	8.36	146	384334	38.321	ng	# 93
15) Benzyl Alcohol	8.28	79	284028	36.656	ng	# 75
16) 2,2'-oxybis(1-Chloropropan	8.55	45	150137	35.238	ng	# 43
17) 2-Methylphenol	8.47	107	223784	37.451	ng	# 75
18) Hexachloroethane	9.08	117	203203	43.207	ng	# 39
19) n-Nitroso-di-n-propylamine	8.83	70	257536	39.368	ng	# 87
20) 3+4-Methylphenols	8.80	107	309390	37.448	ng	# 79
22) Acetophenone	8.86	105	468563	37.532	ng	# 97
24) Nitrobenzene	9.25	77	434824	38.326	ng	# 91
25) Isophorone	9.76	82	588840	38.801	ng	# 95
26) 2-Nitrophenol	9.95	139	179028	36.911	ng	# 44
27) 2,4-Dimethylphenol	10.00	122	230135	36.409	ng	# 66
28) bis(2-Chloroethoxy)methane	10.24	93	332737	36.729	ng	# 98
29) 2,4-Dichlorophenol	10.48	162	427771	45.219	ng	# 89
30) 1,2,4-Trichlorobenzene	10.68	180	710284	49.482	ng	# 96
31) Naphthalene	10.87	128	948730	38.408	ng	# 97
32) Benzoic acid	10.22	122	172305	32.581	ng	# 73
33) 4-Chloroaniline	11.00	127	404327	38.894	ng	# 86
34) Hexachlorobutadiene	11.12	225	697174	50.886	ng	# 96
35) Caprolactam	11.84	113	84475	37.214	ng	# 55
36) 4-Chloro-3-methylphenol	12.13	107	351077	38.603	ng	# 90
37) 2-Methylnaphthalene	12.47	142	779593	40.315	ng	# 87
39) 1,2,4,5-Tetrachlorobenzene	12.85	216	1041884	42.374	ng	# 100
40) Hexachlorocyclopentadiene	12.80	237	402894	44.758	ng	# 98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.10	196	632463	43.015	nq	96
43) 2,4,5-Trichlorophenol	13.18	196	633918m	44.137	nq	
45) 1,1'-Biphenyl	13.49	154	1326976	34.761	nq	98
46) 2-Chloronaphthalene	13.54	162	1128293	36.380	nq	98
47) 2-Nitroaniline	13.77	65	269862	32.668	nq	87
48) Acenaphthylene	14.38	152	1550210	35.410	nq	99
49) Dimethylphthalate	14.12	163	1586232	37.924	nq	# 96
50) 2,6-Dinitrotoluene	14.26	165	309284	37.911	nq	# 76
51) Acenaphthene	14.72	154	947786	35.239	nq	96
52) 3-Nitroaniline	14.60	138	235757	31.760	nq	# 78
53) 2,4-Dinitrophenol	14.83	184	124989	30.614	nq	# 66
54) Dibenzofuran	15.06	168	2042283	39.666	nq	95
55) 4-Nitrophenol	14.92	139	145808	30.796	nq	# 84
56) 2,4-Dinitrotoluene	15.06	165	518190	38.091	nq	# 96
57) Fluorene	15.70	166	1529226	39.317	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.29	232	607855	44.034	nq	# 100
59) Diethylphthalate	15.47	149	1563114	35.842	nq	# 93
60) 4-Chlorophenyl-phenylether	15.69	204	1425126	43.692	nq	# 84
61) 4-Nitroaniline	15.77	138	244169	33.656	nq	# 1
62) Azobenzene	15.99	77	1694996	36.490	nq	89
64) 4,6-Dinitro-2-methylphenol	15.83	198	374596	34.239	nq	# 56
65) n-Nitrosodiphenylamine	15.92	169	1456873	35.966	nq	99
66) 4-Bromophenyl-phenylether	16.59	248	884788	39.992	nq	96
67) Hexachlorobenzene	16.70	284	1028690	39.870	nq	93
68) Atrazine	16.86	200	758025	41.005	nq	95
69) Pentachlorophenol	17.05	266	535122	39.199	nq	97
70) Phenanthrene	17.44	178	2839205	38.151	nq	100
71) Anthracene	17.53	178	2791390	37.773	nq	99
72) Carbazole	17.81	167	2365245	35.615	nq	99
73) Di-n-butylphthalate	18.33	149	2779319	34.327	nq	# 96
74) Fluoranthene	19.44	202	4354542	39.907	nq	94
76) Benzidine	19.63	184	1734759	40.388	nq	99
77) Pyrene	19.80	202	4626197	38.113	nq	99
79) Butylbenzylphthalate	20.67	149	1406078	33.559	nq	# 80
80) Benzo(a)anthracene	21.53	228	5063732	38.563	nq	99
81) 3,3'-Dichlorobenzidine	21.46	252	2184188	37.368	nq	# 97
82) Chrysene	21.58	228	4681944	37.369	nq	99
83) Bis(2-ethylhexyl)phthalate	21.42	149	2056761	33.021	nq	# 97
84) Di-n-octyl phthalate	22.30	149	3410336	32.947	nq	# 92
85) Indeno(1,2,3-cd)pyrene	26.15	276	6353572	35.263	nq	# 92
87) Benzo(b)fluoranthene	23.14	252	5460195	40.837	nq	# 90
88) Benzo(k)fluoranthene	23.18	252	5169772	38.000	nq	# 93
89) Benzo(a)pyrene	23.73	252	4992294	38.373	nq	# 95
90) Dibenzo(a,h)anthracene	26.15	278	5418029	37.160	nq	# 87
91) Benzo(g,h,i)perylene	26.86	276	4770739	36.451	nq	# 81

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

