

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061620\
 Data File : BM026424.D
 Acq On : 17 Jun 2020 00:09
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Jun 17 01:20:23 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 16 10:23:36 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	103644	20.00	ng	0.00
21) Naphthalene-d8	10.16	136	419668	20.00	ng	0.00
39) Acenaphthene-d10	14.05	164	248341	20.00	ng	0.00
64) Phenanthrene-d10	16.81	188	525791	20.00	ng	0.00
76) Chrysene-d12	21.02	240	537520	20.00	ng	0.00
86) Perylene-d12	23.10	264	575580	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.04	112	505641	86.26	ng	0.00
7) Phenol-d6	6.61	99	698140	82.24	ng	0.00
23) Nitrobenzene-d5	8.55	82	710890	83.17	ng	0.00
42) 2,4,6-Tribromophenol	15.56	330	260747	86.68	ng	0.00
45) 2-Fluorobiphenyl	12.66	172	1438170	87.91	ng	0.00
79) Terphenyl-d14	19.46	244	2259302	81.59	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.02	88	113388	42.902	ng	98
3) Pyridine	3.40	79	322346	41.525	ng	99
4) n-Nitrosodimethylamine	3.32	42	149850	38.988	ng	96
6) Aniline	6.75	93	453331	40.715	ng	98
8) 2-Chlorophenol	6.97	128	285577	42.235	ng	99
9) Benzaldehyde	6.56	77	188176	34.766	ng	97
10) Phenol	6.63	94	369490	40.873	ng	99
11) bis(2-Chloroethyl)ether	6.85	93	289549	39.766	ng	97
12) 1,3-Dichlorobenzene	7.29	146	324818	43.390	ng	98
13) 1,4-Dichlorobenzene	7.43	146	325735	42.718	ng	99
14) 1,2-Dichlorobenzene	7.75	146	317020	42.561	ng	97
15) Benzyl Alcohol	7.65	79	277707	38.526	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.94	45	513364	37.751	ng	99
17) 2-Methylphenol	7.86	107	264891	40.091	ng	100
18) Hexachloroethane	8.46	117	117835	40.804	ng	98
19) n-Nitroso-di-n-propylamine	8.21	70	249579	37.730	ng	99
20) 3+4-Methylphenols	8.19	107	362536	40.258	ng	98
22) Acetophenone	8.22	105	449888	41.690	ng	# 98
24) Nitrobenzene	8.59	77	366174	40.925	ng	100
25) Isophorone	9.12	82	649576	40.453	ng	100
26) 2-Nitrophenol	9.29	139	151921	42.903	ng	97
27) 2,4-Dimethylphenol	9.37	122	233838	41.830	ng	98
28) bis(2-Chloroethoxy)methane	9.60	93	363361	39.885	ng	99
29) 2,4-Dichlorophenol	9.83	162	263873	43.022	ng	98
30) 1,2,4-Trichlorobenzene	10.03	180	294667	43.812	ng	99
31) Naphthalene	10.21	128	884599	41.828	ng	100
32) Benzoic acid	9.55	122	189337	42.752	ng	99
33) 4-Chloroaniline	10.33	127	375232	40.680	ng	99
34) Hexachlorobutadiene	10.50	225	190845	44.947	ng	97
35) Caprolactam	11.12	113	88157	40.911	ng	97
36) 4-Chloro-3-methylphenol	11.47	107	299227	40.037	ng	99
37) 2-Methylnaphthalene	11.83	142	619511	41.108	ng	100
38) 1-Methylnaphthalene	12.06	142	578118	40.495	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.22	216	317200	44.871	ng	99

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41) Hexachlorocyclopentadiene	12.19	237	65908	15.821	ng	96
43) 2,4,6-Trichlorophenol	12.47	196	212392	44.388	ng	98
44) 2,4,5-Trichlorophenol	12.55	196	246808	44.437	ng	98
46) 1,1'-Biphenyl	12.87	154	748864	43.011	ng	99
47) 2-Chloronaphthalene	12.91	162	611064	43.209	ng	99
48) 2-Nitroaniline	13.13	65	235093	41.945	ng	97
49) Acenaphthylene	13.77	152	968416	42.814	ng	99
50) Dimethylphthalate	13.53	163	753199	41.466	ng	100
51) 2,6-Dinitrotoluene	13.64	165	167203	41.949	ng	98
52) Acenaphthene	14.12	154	581838	42.848	ng	98
53) 3-Nitroaniline	13.97	138	186012	41.909	ng	96
54) 2,4-Dinitrophenol	14.20	184	42404	17.504	ng	92
55) Dibenzofuran	14.46	168	922341	42.156	ng	99
56) 4-Nitrophenol	14.31	139	143469	40.761	ng	98
57) 2,4-Dinitrotoluene	14.44	165	233696	42.168	ng	98
58) Fluorene	15.12	166	751952	42.313	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.70	232	201510	42.418	ng	100
60) Diethylphthalate	14.91	149	758871	41.017	ng	99
61) 4-Chlorophenyl-phenylether	15.12	204	401625	42.607	ng	99
62) 4-Nitroaniline	15.15	138	198507	41.937	ng	97
63) Azobenzene	15.41	77	827178	40.863	ng	98
65) 4,6-Dinitro-2-methylphenol	15.22	198	66290	20.733	ng	96
66) n-Nitrosodiphenylamine	15.33	169	643178	42.226	ng	99
67) 4-Bromophenyl-phenylether	16.01	248	251248	43.408	ng	99
68) Hexachlorobenzene	16.12	284	261029	43.219	ng	98
69) Atrazine	16.30	200	188405	41.403	ng	100
70) Pentachlorophenol	16.47	266	155261	42.688	ng	98
71) Phenanthrene	16.85	178	1146140	41.859	ng	100
72) Anthracene	16.94	178	1168517	42.761	ng	100
73) Carbazole	17.22	167	1098663	42.396	ng	99
74) Di-n-butylphthalate	17.80	149	1336963	43.747	ng	100
75) Fluoranthene	18.87	202	1354217	43.139	ng	100
77) Benzidine	19.08	184	523032	46.845	ng	99
78) Pyrene	19.24	202	1409701	39.692	ng	100
80) Butylbenzylphthalate	20.17	149	617016	42.961	ng	99
81) Benzo(a)anthracene	21.00	228	1372020	42.524	ng	100
82) 3,3'-Dichlorobenzidine	20.94	252	508947	50.986	ng	99
83) Chrysene	21.06	228	1330014	43.257	ng	100
84) Bis(2-ethylhexyl)phthalate	20.96	149	922688	45.056	ng	99
85) Di-n-octyl phthalate	21.81	149	1577312	49.929	ng	99
87) Indeno(1,2,3-cd)pyrene	25.15	276	1595462	47.918	ng	100
88) Benzo(b)fluoranthene	22.50	252	1404652	40.568	ng	99
89) Benzo(k)fluoranthene	22.53	252	1329300	39.517	ng	100
90) Benzo(a)pyrene	23.01	252	1306747	42.493	ng	100
91) Dibenzo(a,h)anthracene	25.16	278	1338461	48.152	ng	99
92) Benzo(g,h,i)perylene	25.77	276	1289410	48.747	ng	99

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

