

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061520\
 Data File : BM026398.D
 Acq On : 15 Jun 2020 19:08
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Jun 15 19:48:27 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 15 12:22:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	92734	20.00	ng	0.00
21) Naphthalene-d8	10.16	136	380103	20.00	ng	0.00
39) Acenaphthene-d10	14.06	164	228126	20.00	ng	0.00
64) Phenanthrene-d10	16.81	188	491201	20.00	ng	0.00
76) Chrysene-d12	21.02	240	517616	20.00	ng	0.00
86) Perylene-d12	23.11	264	537306	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.04	112	453497	86.46	ng	0.00
7) Phenol-d6	6.61	99	635629	83.69	ng	0.00
23) Nitrobenzene-d5	8.55	82	656792	84.84	ng	0.00
42) 2,4,6-Tribromophenol	15.56	330	239195	86.56	ng	0.00
45) 2-Fluorobiphenyl	12.67	172	1325110	88.17	ng	0.00
79) Terphenyl-d14	19.47	244	2125152	79.70	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.02	88	104571	44.221	ng	99
3) Pyridine	3.40	79	295089	42.485	ng	99
4) n-Nitrosodimethylamine	3.32	42	140966	40.991	ng	97
6) Aniline	6.75	93	412023	41.359	ng	99
8) 2-Chlorophenol	6.97	128	253525	41.906	ng	97
9) Benzaldehyde	6.56	77	171070	35.324	ng	98
10) Phenol	6.64	94	336326	41.581	ng	98
11) bis(2-Chloroethyl)ether	6.85	93	263862	40.501	ng	97
12) 1,3-Dichlorobenzene	7.29	146	287466	42.918	ng	98
13) 1,4-Dichlorobenzene	7.44	146	291253	42.690	ng	99
14) 1,2-Dichlorobenzene	7.75	146	283263	42.503	ng	98
15) Benzyl Alcohol	7.66	79	265235	41.125	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.95	45	491972	40.434	ng	99
17) 2-Methylphenol	7.86	107	245303	41.495	ng	98
18) Hexachloroethane	8.46	117	109819	42.503	ng	99
19) n-Nitroso-di-n-propylamine	8.22	70	238297	40.263	ng	97
20) 3+4-Methylphenols	8.19	107	332996	41.328	ng	98
22) Acetophenone	8.22	105	414162	42.374	ng	# 97
24) Nitrobenzene	8.59	77	343337	42.367	ng	99
25) Isophorone	9.12	82	615394	42.314	ng	99
26) 2-Nitrophenol	9.29	139	141433	44.098	ng	95
27) 2,4-Dimethylphenol	9.37	122	215232	42.510	ng	94
28) bis(2-Chloroethoxy)methane	9.60	93	344626	41.766	ng	99
29) 2,4-Dichlorophenol	9.83	162	238019	42.846	ng	97
30) 1,2,4-Trichlorobenzene	10.03	180	262673	43.121	ng	99
31) Naphthalene	10.21	128	800244	41.778	ng	99
32) Benzoic acid	9.54	122	145407	37.036	ng	98
33) 4-Chloroaniline	10.33	127	349067	41.783	ng	99
34) Hexachlorobutadiene	10.50	225	172611	44.885	ng	98
35) Caprolactam	11.12	113	83331	42.697	ng	96
36) 4-Chloro-3-methylphenol	11.48	107	283129	41.826	ng	99
37) 2-Methylnaphthalene	11.83	142	563471	41.281	ng	100
38) 1-Methylnaphthalene	12.06	142	539445	41.719	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.22	216	287397	44.257	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.20	237	148454	38.793	ng	100
43) 2,4,6-Trichlorophenol	12.47	196	193374	43.995	ng	99
44) 2,4,5-Trichlorophenol	12.55	196	226963	44.485	ng	99
46) 1,1'-Biphenyl	12.88	154	693714	43.374	ng	99
47) 2-Chloronaphthalene	12.92	162	563187	43.353	ng	100
48) 2-Nitroaniline	13.13	65	225182	43.737	ng	98
49) Acenaphthylene	13.77	152	893194	42.988	ng	100
50) Dimethylphthalate	13.53	163	718875	43.083	ng	99
51) 2,6-Dinitrotoluene	13.65	165	158203	43.208	ng	96
52) Acenaphthene	14.12	154	540141	43.302	ng	99
53) 3-Nitroaniline	13.98	138	174485	42.795	ng	96
54) 2,4-Dinitrophenol	14.20	184	88592	39.811	ng	98
55) Dibenzofuran	14.46	168	860468	42.813	ng	99
56) 4-Nitrophenol	14.32	139	129235	39.970	ng	98
57) 2,4-Dinitrotoluene	14.45	165	219856	43.186	ng	97
58) Fluorene	15.12	166	700174	42.891	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.70	232	184886	42.367	ng	100
60) Diethylphthalate	14.92	149	730470	42.980	ng	99
61) 4-Chlorophenyl-phenylether	15.12	204	373075	43.086	ng	98
62) 4-Nitroaniline	15.15	138	186041	42.786	ng	99
63) Azobenzene	15.42	77	793142	42.654	ng	99
65) 4,6-Dinitro-2-methylphenol	15.22	198	124998	41.848	ng	96
66) n-Nitrosodiphenylamine	15.34	169	604712	42.496	ng	100
67) 4-Bromophenyl-phenylether	16.02	248	233159	43.119	ng	97
68) Hexachlorobenzene	16.13	284	244584	43.347	ng	97
69) Atrazine	16.31	200	187177	44.029	ng	99
70) Pentachlorophenol	16.47	266	134859	39.689	ng	98
71) Phenanthrene	16.85	178	1082743	42.328	ng	99
72) Anthracene	16.94	178	1088436	42.635	ng	100
73) Carbazole	17.22	167	1044991	43.164	ng	99
74) Di-n-butylphthalate	17.81	149	1256167	43.998	ng	100
75) Fluoranthene	18.88	202	1285999	43.851	ng	99
77) Benzidine	19.08	184	477451	44.407	ng	99
78) Pyrene	19.24	202	1333369	38.987	ng	100
80) Butylbenzylphthalate	20.18	149	584159	42.238	ng	95
81) Benzo(a)anthracene	21.00	228	1312372	42.240	ng	100
82) 3,3'-Dichlorobenzidine	20.95	252	468143	48.701	ng	99
83) Chrysene	21.06	228	1259056	42.524	ng	99
84) Bis(2-ethylhexyl)phthalate	20.97	149	859313	43.575	ng	99
85) Di-n-octyl phthalate	21.82	149	1488105	48.917	ng	100
87) Indeno(1,2,3-cd)pyrene	25.16	276	1504446	48.404	ng	100
88) Benzo(b)fluoranthene	22.50	252	1349534	41.753	ng	99
89) Benzo(k)fluoranthene	22.54	252	1244818	39.642	ng	99
90) Benzo(a)pyrene	23.02	252	1218235	42.437	ng	99
91) Dibenzo(a,h)anthracene	25.16	278	1258410	48.497	ng	99
92) Benzo(g,h,i)perylene	25.77	276	1206037	48.843	ng	99

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

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