

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092019\
 Data File : BM022739.D
 Acq On : 20 Sep 2019 23:21
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Sep 21 02:15:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM092019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 20 16:34:01 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	206842	20.00	ng	0.00
21) Naphthalene-d8	10.62	136	838177	20.00	ng	0.00
39) Acenaphthene-d10	14.46	164	481789	20.00	ng	0.00
64) Phenanthrene-d10	17.20	188	922971	20.00	ng	0.00
76) Chrysene-d12	21.37	240	878128	20.00	ng	-0.01
87) Perylene-d12	23.64	264	889688	20.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.40	112	1021959	78.36	ng	0.00
7) Phenol-d6	6.99	99	1331467	79.49	ng	0.00
23) Nitrobenzene-d5	8.98	82	1190182	78.82	ng	0.00
42) 2,4,6-Tribromophenol	15.94	330	473411	78.71	ng	0.00
45) 2-Fluorobiphenyl	13.08	172	2749084	78.57	ng	0.00
79) Terphenyl-d14	19.82	244	3800650	83.81	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.32	88	202340	38.557	ng	# 100
3) Pyridine	3.72	79	571900	41.344	ng	99
4) n-Nitrosodimethylamine	3.62	42	200728	39.862	ng	95
6) Aniline	7.15	93	798072	39.734	ng	100
8) 2-Chlorophenol	7.39	128	521860	39.273	ng	99
9) Benzaldehyde	6.96	77	433993	45.501	ng	99
10) Phenol	7.02	94	623629	39.544	ng	99
11) bis(2-Chloroethyl)ether	7.25	93	497204	40.129	ng	99
12) 1,3-Dichlorobenzene	7.72	146	621605	39.140	ng	99
13) 1,4-Dichlorobenzene	7.86	146	625800	38.922	ng	99
14) 1,2-Dichlorobenzene	8.17	146	597875	39.004	ng	100
15) Benzyl Alcohol	8.06	79	418645	39.941	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.35	45	723239	40.183	ng	98
17) 2-Methylphenol	8.26	107	418628	40.104	ng	99
18) Hexachloroethane	8.90	117	228731	38.787	ng	97
19) n-Nitroso-di-n-propylamine	8.63	70	382617	40.904	ng	100
20) 3+4-Methylphenols	8.59	107	557283	40.083	ng	99
22) Acetophenone	8.65	105	885167	42.783	ng	99
24) Nitrobenzene	9.02	77	573111	39.670	ng	100
25) Isophorone	9.55	82	993631	40.021	ng	100
26) 2-Nitrophenol	9.73	139	238234	38.198	ng	96
27) 2,4-Dimethylphenol	9.79	122	414281	39.678	ng	100
28) bis(2-Chloroethoxy)methane	10.03	93	681337	39.658	ng	100
29) 2,4-Dichlorophenol	10.26	162	464750	39.370	ng	99
30) 1,2,4-Trichlorobenzene	10.48	180	560761	38.610	ng	99
31) Naphthalene	10.67	128	1615312	39.004	ng	100
32) Benzoic acid	9.93	122	307339	43.273	ng	96
33) 4-Chloroaniline	10.77	127	710558	39.516	ng	99
34) Hexachlorobutadiene	10.96	225	332495	38.405	ng	100
35) Caprolactam	11.55	113	139069	39.738	ng	94
36) 4-Chloro-3-methylphenol	11.89	107	458092	39.775	ng	100
37) 2-Methylnaphthalene	12.28	142	1133169	39.308	ng	99
38) 1-Methylnaphthalene	12.50	142	1070010	39.103	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.65	216	724046	44.352	ng	100

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41) Hexachlorocyclopentadiene	12.64	237	340871	38.270	nq	99
43) 2,4,6-Trichlorophenol	12.89	196	367418	38.995	nq	99
44) 2,4,5-Trichlorophenol	12.96	196	380797	38.783	nq	100
46) 1,1'-Biphenyl	13.29	154	1730309	43.394	nq	100
47) 2-Chloronaphthalene	13.33	162	1160299	39.041	nq	100
48) 2-Nitroaniline	13.53	65	312274	40.351	nq	95
49) Acenaphthylene	14.18	152	1721990	39.126	nq	100
50) Dimethylphthalate	13.92	163	1356007	38.946	nq	99
51) 2,6-Dinitrotoluene	14.03	165	285369	39.183	nq	99
52) Acenaphthene	14.52	154	1129805	39.239	nq	100
53) 3-Nitroaniline	14.36	138	330217	39.385	nq	100
54) 2,4-Dinitrophenol	14.56	184	126508	36.224	nq	99
55) Dibenzofuran	14.86	168	1623906	38.632	nq	100
56) 4-Nitrophenol	14.66	139	253253	39.298	nq	99
57) 2,4-Dinitrotoluene	14.81	165	374870	39.290	nq	95
58) Fluorene	15.50	166	1332274	39.003	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.08	232	331128	38.669	nq	99
60) Diethylphthalate	15.28	149	1326332	39.177	nq	100
61) 4-Chlorophenyl-phenylether	15.50	204	687205	39.154	nq	98
62) 4-Nitroaniline	15.52	138	347116	39.489	nq	99
63) Azobenzene	15.79	77	1227972	40.166	nq	99
65) 4,6-Dinitro-2-methylphenol	15.57	198	165215	37.329	nq	99
66) n-Nitrosodiphenylamine	15.71	169	1124239	39.871	nq	100
67) 4-Bromophenyl-phenylether	16.39	248	403957	39.350	nq	99
68) Hexachlorobenzene	16.51	284	471610	39.394	nq	99
69) Atrazine	16.66	200	364711	39.788	nq	99
70) Pentachlorophenol	16.84	266	252369	39.528	nq	99
71) Phenanthrene	17.24	178	2043251	39.233	nq	99
72) Anthracene	17.33	178	2006185	39.747	nq	99
73) Carbazole	17.59	167	1851722	39.425	nq	100
74) Di-n-butylphthalate	18.16	149	2140015	40.183	nq	100
75) Fluoranthene	19.25	202	2280882	38.812	nq	99
77) Benzidine	19.43	184	1018390	42.274	nq	99
78) Pyrene	19.61	202	2359360	40.998	nq	100
80) Butylbenzylphthalate	20.51	149	892534	40.809	nq	95
81) Benzo(a)anthracene	21.36	228	2261048	39.728	nq	100
82) 3,3'-Dichlorobenzidine	21.29	252	804109	39.841	nq	99
83) Chrysene	21.40	228	2178853	39.489	nq	100
84) Bis(2-ethylhexyl)phthalate	21.29	149	1340723	41.372	nq	100
85) Di-n-octyl phthalate	22.17	149	2062519	39.934	nq	100
86) Indeno(1,2,3-cd)pyrene	25.95	276	2403641	36.499	nq	# 100
88) Benzo(b)fluoranthene	22.96	252	2171789	40.317	nq	100
89) Benzo(k)fluoranthene	23.00	252	2125205	39.761	nq	100
90) Benzo(a)pyrene	23.54	252	1972914	39.747	nq	100
91) Dibenzo(a,h)anthracene	25.96	278	2004299	38.603	nq	99
92) Benzo(g,h,i)perylene	26.66	276	1866641	37.852	nq	100

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

