

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040519\
 Data File : BF113621.D
 Acq On : 5 Apr 2019 17:52
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Apr 05 23:41:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 04 04:05:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	149018	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	586406	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	290732	20.00	ng	0.00
64) Phenanthrene-d10	11.20	188	580868	20.00	ng	0.00
76) Chrysene-d12	13.83	240	430696	20.00	ng	-0.01
87) Perylene-d12	15.23	264	336840	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	700427	80.15	ng	0.00
7) Phenol-d6	6.34	99	860833	79.92	ng	0.00
23) Nitrobenzene-d5	7.26	82	805767	80.49	ng	0.00
42) 2,4,6-Tribromophenol	10.51	330	277416	79.43	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	1422506	78.77	ng	0.00
79) Terphenyl-d14	12.78	244	1611629	76.19	ng	-0.01

Target Compounds					Qvalue
2) 1,4-Dioxane	2.33	88	157768	38.118 ng	98
3) Pyridine	3.03	79	435706	40.225 ng	97
4) n-Nitrosodimethylamine	2.98	42	186002	40.416 ng	92
6) Aniline	6.35	93	535730	41.035 ng	91
8) 2-Chlorophenol	6.47	128	408328	40.385 ng	99
9) Benzaldehyde	6.23	77	266343	41.588 ng	99
10) Phenol	6.36	94	501703	40.785 ng	85
11) bis(2-Chloroethyl)ether	6.43	93	379169	39.625 ng	99
12) 1,3-Dichlorobenzene	6.63	146	434119	39.870 ng	98
13) 1,4-Dichlorobenzene	6.70	146	438661	39.867 ng	98
14) 1,2-Dichlorobenzene	6.86	146	401086	40.384 ng	98
15) Benzyl Alcohol	6.84	79	307625	42.241 ng	98
16) 2,2'-oxybis(1-Chloropropan	6.97	45	578419	38.835 ng	84
17) 2-Methylphenol	6.96	107	314518	40.140 ng	95
18) Hexachloroethane	7.19	117	157276	39.270 ng	99
19) n-Nitroso-di-n-propylamine	7.11	70	266576	40.137 ng	98
20) 3+4-Methylphenols	7.12	107	400516	42.186 ng	93
22) Acetophenone	7.10	105	522324	40.555 ng	# 98
24) Nitrobenzene	7.27	77	411936	39.956 ng	97
25) Isophorone	7.52	82	771432	40.001 ng	98
26) 2-Nitrophenol	7.59	139	231580	41.735 ng	92
27) 2,4-Dimethylphenol	7.64	122	311721	39.296 ng	98
28) bis(2-Chloroethoxy)methane	7.72	93	484712	39.755 ng	99
29) 2,4-Dichlorophenol	7.84	162	352017	41.475 ng	100
30) 1,2,4-Trichlorobenzene	7.92	180	366564	39.771 ng	97
31) Naphthalene	7.99	128	1138698	40.077 ng	99
32) Benzoic acid	7.80	122	213113	34.694 ng	100
33) 4-Chloroaniline	8.05	127	483037	42.875 ng	99
34) Hexachlorobutadiene	8.11	225	225687	40.163 ng	99
35) Caprolactam	8.44	113	117713	43.691 ng	92
36) 4-Chloro-3-methylphenol	8.55	107	355530	41.729 ng	98
37) 2-Methylnaphthalene	8.68	142	753645	40.330 ng	98
38) 1-Methylnaphthalene	8.78	142	725165	40.244 ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.85	216	365999	38.926 ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.83	237	86533	35.274	ng	99
43) 2,4,6-Trichlorophenol	8.97	196	237210	39.967	ng	99
44) 2,4,5-Trichlorophenol	9.02	196	255246	40.058	ng	99
46) 1,1'-Biphenyl	9.15	154	942844	39.300	ng	99
47) 2-Chloronaphthalene	9.17	162	714911	39.358	ng	99
48) 2-Nitroaniline	9.27	65	230831	41.530	ng	97
49) Acenaphthylene	9.58	152	1154319	39.080	ng	100
50) Dimethylphthalate	9.45	163	844810	39.420	ng	99
51) 2,6-Dinitrotoluene	9.52	165	192100	40.665	ng	95
52) Acenaphthene	9.75	154	666131	39.397	ng	99
53) 3-Nitroaniline	9.69	138	223613	41.634	ng	94
54) 2,4-Dinitrophenol	9.80	184	84217	38.106	ng	92
55) Dibenzofuran	9.92	168	976566	39.374	ng	99
56) 4-Nitrophenol	9.87	139	124248	37.293	ng	91
57) 2,4-Dinitrotoluene	9.92	165	236513	41.397	ng	# 85
58) Fluorene	10.27	166	786958	40.117	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.05	232	223695	40.558	ng	97
60) Diethylphthalate	10.15	149	815178	39.452	ng	98
61) 4-Chlorophenyl-phenylether	10.26	204	386464	40.053	ng	93
62) 4-Nitroaniline	10.30	138	235507	43.121	ng	98
63) Azobenzene	10.42	77	767551	39.233	ng	97
65) 4,6-Dinitro-2-methylphenol	10.34	198	124461	40.700	ng	99
66) n-Nitrosodiphenylamine	10.38	169	701140	39.319	ng	100
67) 4-Bromophenyl-phenylether	10.74	248	256059	39.305	ng	94
68) Hexachlorobenzene	10.82	284	295364	39.568	ng	100
69) Atrazine	10.91	200	222381	39.565	ng	98
70) Pentachlorophenol	11.02	266	142618	40.546	ng	99
71) Phenanthrene	11.22	178	1145446	39.685	ng	100
72) Anthracene	11.27	178	1161456	39.550	ng	100
73) Carbazole	11.43	167	1121705	40.839	ng	100
74) Di-n-butylphthalate	11.76	149	1295128	39.639	ng	100
75) Fluoranthene	12.41	202	1262769	40.827	ng	98
77) Benzidine	12.53	184	617936	38.582	ng	99
78) Pyrene	12.63	202	1281495	39.104	ng	99
80) Butylbenzylphthalate	13.24	149	533669	40.006	ng	99
81) Benzo(a)anthracene	13.82	228	1020985	41.094	ng	100
82) 3,3'-Dichlorobenzidine	13.78	252	397279	41.515	ng	99
83) Chrysene	13.86	228	979838	38.905	ng	100
84) Bis(2-ethylhexyl)phthalate	13.80	149	661163	40.993	ng	99
85) Di-n-octyl phthalate	14.41	149	1022584	39.005	ng	100
86) Indeno(1,2,3-cd)pyrene	16.60	276	814939	35.046	ng	99
88) Benzo(b)fluoranthene	14.83	252	910505	44.159	ng	98
89) Benzo(k)fluoranthene	14.86	252	708250	38.274	ng	98
90) Benzo(a)pyrene	15.17	252	745504	40.689	ng	100
91) Dibenzo(a,h)anthracene	16.60	278	670329	39.123	ng	100
92) Benzo(g,h,i)perylene	17.00	276	683444	39.080	ng	100

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

