

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031518\
 Data File : BF103682.D
 Acq On : 15 Mar 2018 21:36
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Mar 16 04:31:42 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 15 18:31:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	154300	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	645110	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	300184	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	529565	20.00	ng	0.00
75) Chrysene-d12	14.15	240	402756	20.00	ng	0.00
86) Perylene-d12	15.68	264	326235	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.58	112	818411	80.67	ng	0.00
7) Phenol-d6	6.59	99	999870	80.56	ng	0.00
23) Nitrobenzene-d5	7.53	82	859754	92.94	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	235673	91.31	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1453690	77.25	ng	0.00
78) Terphenyl-d14	13.09	244	1371991	79.07	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.83	88	201532	40.345	ng	92
3) Pyridine	3.60	79	549253	39.976	ng	91
4) n-Nitrosodimethylamine	3.56	42	201346	39.745	ng	94
6) Aniline	6.63	93	722373	40.761	ng	97
8) 2-Chlorophenol	6.75	128	436241	41.050	ng	97
9) Benzaldehyde	6.52	77	328815	40.327	ng	99
10) Phenol	6.60	94	530336	40.212	ng	99
11) bis(2-Chloroethyl)ether	6.70	93	437852	40.217	ng	94
12) 1,3-Dichlorobenzene	6.91	146	488778	40.382	ng	98
13) 1,4-Dichlorobenzene	6.99	146	485819	39.944	ng	99
14) 1,2-Dichlorobenzene	7.14	146	455887	40.393	ng	99
15) Benzyl Alcohol	7.10	79	394620	41.036	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.24	45	618768	39.959	ng	96
17) 2-Methylphenol	7.20	107	385000	40.682	ng	97
18) Hexachloroethane	7.48	117	180127	42.103	ng	93
19) n-Nitroso-di-n-propylamine	7.38	70	319148	40.124	ng	95
20) 3+4-Methylphenols	7.36	107	494617	41.183	ng	96
22) Acetophenone	7.37	105	649637	39.183	ng	96
24) Nitrobenzene	7.55	77	482031	44.384	ng	99
25) Isophorone	7.79	82	841509	39.296	ng	98
26) 2-Nitrophenol	7.86	139	196051	47.556	ng	99
27) 2,4-Dimethylphenol	7.89	122	376576	41.048	ng	98
28) bis(2-Chloroethoxy)methane	7.99	93	513579	39.605	ng	100
29) 2,4-Dichlorophenol	8.10	162	347475	41.631	ng	98
30) 1,2,4-Trichlorobenzene	8.19	180	386427	39.917	ng	99
31) Naphthalene	8.27	128	1283080	39.373	ng	99
32) Benzoic acid	8.00	122	301722	46.897	ng	97
33) 4-Chloroaniline	8.32	127	552699	40.427	ng	99
34) Hexachlorobutadiene	8.39	225	210054	39.576	ng	100
35) Caprolactam	8.69	113	118951	41.388	ng	90
36) 4-Chloro-3-methylphenol	8.79	107	382974	41.621	ng	99
37) 2-Methylnaphthalene	8.96	142	826825	40.010	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.13	216	348865	40.737	ng	99
40) Hexachlorocyclopentadiene	9.12	237	200021	45.264	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.23	196	239397	43.704	ng	98
43) 2,4,5-Trichlorophenol	9.27	196	270328	43.724	ng	98
45) 1,1'-Biphenyl	9.43	154	998609	39.895	ng	100
46) 2-Chloronaphthalene	9.46	162	793875	39.931	ng	99
47) 2-Nitroaniline	9.55	65	239324	46.298	ng	94
48) Acenaphthylene	9.87	152	1221153	39.610	ng	99
49) Dimethylphthalate	9.73	163	912232	39.831	ng	100
50) 2,6-Dinitrotoluene	9.79	165	194183	44.176	ng	90
51) Acenaphthene	10.05	154	743591	40.621	ng	100
52) 3-Nitroaniline	9.96	138	233072	44.221	ng	97
53) 2,4-Dinitrophenol	10.06	184	53259	52.942	ng	# 37
54) Dibenzofuran	10.22	168	1036265	39.636	ng	98
55) 4-Nitrophenol	10.10	139	177276	44.622	ng	95
56) 2,4-Dinitrotoluene	10.19	165	250914	45.130	ng	97
57) Fluorene	10.56	166	797856	39.328	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.33	232	200444	45.754	ng	97
59) Diethylphthalate	10.43	149	907237	39.911	ng	98
60) 4-Chlorophenyl-phenylether	10.55	204	369939	39.900	ng	94
61) 4-Nitroaniline	10.57	138	224187	44.192	ng	94
62) Azobenzene	10.71	77	907673	39.275	ng	97
64) 4,6-Dinitro-2-methylphenol	10.60	198	100799	50.477	ng	93
65) n-Nitrosodiphenylamine	10.67	169	728365	40.408	ng	100
66) 4-Bromophenyl-phenylether	11.04	248	223038	41.020	ng	96
67) Hexachlorobenzene	11.10	284	249707	40.238	ng	98
68) Atrazine	11.19	200	228507	40.980	ng	99
69) Pentachlorophenol	11.30	266	160593	45.012	ng	97
70) Phenanthrene	11.53	178	1150977	39.732	ng	99
71) Anthracene	11.58	178	1158958	39.391	ng	100
72) Carbazole	11.73	167	1131474	39.566	ng	99
73) Di-n-butylphthalate	12.05	149	1401760	40.852	ng	99
74) Fluoranthene	12.72	202	1171601	39.257	ng	100
76) Benzidine	12.84	184	710321	41.351	ng	98
77) Pyrene	12.95	202	1216790	41.138	ng	100
79) Butylbenzylphthalate	13.56	149	566094	43.198	ng	94
80) Benzo(a)anthracene	14.14	228	1016299	39.864	ng	99
81) 3,3'-Dichlorobenzidine	14.10	252	358279	42.013	ng	100
82) Chrysene	14.18	228	972717	40.025	ng	99
83) Bis(2-ethylhexyl)phthalate	14.12	149	810870	43.313	ng	99
84) Di-n-octyl phthalate	14.74	149	1191062	43.731	ng	98
85) Indeno(1,2,3-cd)pyrene	17.24	276	763369	35.968	ng	97
87) Benzo(b)fluoranthene	15.23	252	835916	40.557	ng	98
88) Benzo(k)fluoranthene	15.26	252	855446	43.177	ng	99
89) Benzo(a)pyrene	15.62	252	767540	41.058	ng	98
90) Dibenzo(a,h)anthracene	17.26	278	637389	39.376	ng	98
91) Benzo(g,h,i)perylene	17.72	276	614880	39.046	ng	96

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

