

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082218\
 Data File : BF108405.D
 Acq On : 23 Aug 2018 3:04
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 8/23/2018 3:43:35 PM

Quant Time: Aug 23 07:03:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 22 11:01:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	147088	20.00	ng	0.00
21) Naphthalene-d8	8.29	136	563516	20.00	ng	0.00
38) Acenaphthene-d10	10.05	164	264188	20.00	ng	0.00
63) Phenanthrene-d10	11.54	188	445613	20.00	ng	0.00
75) Chrysene-d12	14.19	240	305353	20.00	ng	0.00
86) Perylene-d12	15.75	264	271207	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	667496	82.16	ng	0.00
7) Phenol-d6	6.62	99	873579	80.92	ng	0.00
23) Nitrobenzene-d5	7.57	82	841542	83.33	ng	0.00
41) 2,4,6-Tribromophenol	10.84	330	190960	72.46	ng	0.00
44) 2-Fluorobiphenyl	9.36	172	1512520	91.92	ng	0.00
78) Terphenyl-d14	13.12	244	1088723	90.65	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.87	88	154842	37.092	ng	89
3) Pyridine	3.66	79	413178	38.422	ng	88
4) n-Nitrosodimethylamine	3.63	42	217958	36.020	ng	# 84
6) Aniline	6.67	93	590198	40.190	ng	98
8) 2-Chlorophenol	6.79	128	375464	41.033	ng	96
9) Benzaldehyde	6.56	77	325743	38.916	ng	97
10) Phenol	6.64	94	453178	40.580	ng	92
11) bis(2-Chloroethyl)ether	6.74	93	363014	40.208	ng	94
12) 1,3-Dichlorobenzene	6.94	146	437844	41.384	ng	98
13) 1,4-Dichlorobenzene	7.02	146	439991	41.391	ng	99
14) 1,2-Dichlorobenzene	7.18	146	401476	40.604	ng	96
15) Benzyl Alcohol	7.14	79	342669	37.703	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.27	45	688917	37.040	ng	99
17) 2-Methylphenol	7.25	107	326440	41.601	ng	95
18) Hexachloroethane	7.51	117	144434	38.165	ng	97
19) n-Nitroso-di-n-propylamine	7.41	70	283555	38.839	ng	94
20) 3+4-Methylphenols	7.40	107	396704	39.451	ng	88
22) Acetophenone	7.41	105	534934	40.598	ng	# 95
24) Nitrobenzene	7.59	77	422641	41.387	ng	95
25) Isophorone	7.82	82	756931	39.325	ng	98
26) 2-Nitrophenol	7.90	139	160110	41.517	ng	92
27) 2,4-Dimethylphenol	7.93	122	338526	39.626	ng	96
28) bis(2-Chloroethoxy)methane	8.02	93	452445	40.265	ng	99
29) 2,4-Dichlorophenol	8.14	162	320746	40.798	ng	98
30) 1,2,4-Trichlorobenzene	8.22	180	354969	40.952	ng	97
31) Naphthalene	8.31	128	1052244	41.059	ng	99
32) Benzoic acid	8.02	122	107147m	24.898	ng	
33) 4-Chloroaniline	8.35	127	450020	40.294	ng	96
34) Hexachlorobutadiene	8.42	225	231501	42.189	ng	99
35) Caprolactam	8.74	113	94512	38.362	ng	87
36) 4-Chloro-3-methylphenol	8.83	107	323700	38.167	ng	98
37) 2-Methylnaphthalene	9.00	142	722963	39.873	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.17	216	362790	44.717	ng	# 97
40) Hexachlorocyclopentadiene	9.15	237	33306	6.356	ng	99

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42) 2,4,6-Trichlorophenol	9.28	196	218518	43.404	nq	100
43) 2,4,5-Trichlorophenol	9.32	196	222884	40.217	nq	95
45) 1,1'-Biphenyl	9.47	154	860709	44.187	nq	98
46) 2-Chloronaphthalene	9.50	162	654929	42.541	nq	98
47) 2-Nitroaniline	9.59	65	224067	44.982	nq	90
48) Acenaphthylene	9.91	152	1028203	42.246	nq	99
49) Dimethylphthalate	9.76	163	804336	43.707	nq	99
50) 2,6-Dinitrotoluene	9.83	165	165769	43.054	nq	99
51) Acenaphthene	10.08	154	587481	39.409	nq	100
52) 3-Nitroaniline	10.00	138	184062	43.693	nq	91
53) 2,4-Dinitrophenol	10.11	184	2389	8.533	nq #	23
54) Dibenzofuran	10.25	168	885327	40.992	nq	96
55) 4-Nitrophenol	10.15	139	92014	28.844	nq	92
56) 2,4-Dinitrotoluene	10.24	165	209229	42.217	nq #	95
57) Fluorene	10.60	166	684825	40.056	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.37	232	155572	33.585	nq #	86
59) Diethylphthalate	10.46	149	774705	43.474	nq	96
60) 4-Chlorophenyl-phenylether	10.58	204	367464	41.248	nq	94
61) 4-Nitroaniline	10.62	138	172697	41.464	nq	92
62) Azobenzene	10.75	77	728908	39.607	nq	93
64) 4,6-Dinitro-2-methylphenol	10.64	198	6782	8.960	nq	92
65) n-Nitrosodiphenylamine	10.71	169	642391	48.783	nq	96
66) 4-Bromophenyl-phenylether	11.08	248	228331	47.150	nq	91
67) Hexachlorobenzene	11.15	284	222544	43.677	nq #	89
68) Atrazine	11.23	200	207213	46.916	nq	95
69) Pentachlorophenol	11.34	266	65186	26.729	nq	98
70) Phenanthrene	11.57	178	897801	42.716	nq	99
71) Anthracene	11.62	178	923162	42.854	nq	100
72) Carbazole	11.77	167	763741	39.904	nq	100
73) Di-n-butylphthalate	12.09	149	1083723	49.990	nq	99
74) Fluoranthene	12.76	202	818507	35.683	nq	96
76) Benzidine	12.88	184	433214	37.972	nq	99
77) Pyrene	13.00	202	789989	40.864	nq	99
79) Butylbenzylphthalate	13.59	149	358254	46.669	nq #	97
80) Benzo(a)anthracene	14.19	228	737604	42.091	nq	99
81) 3,3'-Dichlorobenzidine	14.14	252	266882	42.496	nq #	98
82) Chrysene	14.22	228	676856	42.130	nq	99
83) Bis(2-ethylhexyl)phthalate	14.15	149	474399	45.636	nq #	99
84) Di-n-octyl phthalate	14.77	149	801585	50.561	nq	97
85) Indeno(1,2,3-cd)pyrene	17.37	276	508384	36.520	nq #	90
87) Benzo(b)fluoranthene	15.28	252	717897m	44.299	nq	
88) Benzo(k)fluoranthene	15.31	252	624247	41.904	nq #	96
89) Benzo(a)pyrene	15.68	252	608119	41.905	nq	97
90) Dibenzo(a,h)anthracene	17.38	278	433958	36.142	nq #	93
91) Benzo(g,h,i)perylene	17.85	276	394856	33.397	nq #	89

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

