

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042318\
 Data File : BF104699.D
 Acq On : 23 Apr 2018 10:31
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 4/24/2018 3:48:20 PM

Quant Time: Apr 23 15:46:21 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 23 15:38:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	118507	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	485757	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	227915	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	424847	20.00	ng	0.00
75) Chrysene-d12	13.99	240	333580	20.00	ng	0.00
86) Perylene-d12	15.45	264	288378	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	571307	73.71	ng	0.00
7) Phenol-d6	6.45	99	690194	72.48	ng	0.00
23) Nitrobenzene-d5	7.37	82	635734	85.46	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	191624	81.05	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1087060	75.35	ng	0.00
78) Terphenyl-d14	12.92	244	1104266	75.47	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.50	88	133060	36.845	ng	# 88
3) Pyridine	3.25	79	356805	35.141	ng	87
4) n-Nitrosodimethylamine	3.21	42	113939	32.102	ng	# 75
6) Aniline	6.47	93	465047	35.151	ng	99
8) 2-Chlorophenol	6.59	128	308794	37.749	ng	96
9) Benzaldehyde	6.36	77	168746	33.085	ng	97
10) Phenol	6.46	94	383230	37.929	ng	88
11) bis(2-Chloroethyl)ether	6.55	93	295009	36.588	ng	95
12) 1,3-Dichlorobenzene	6.75	146	348173	37.570	ng	100
13) 1,4-Dichlorobenzene	6.82	146	354515	38.376	ng	98
14) 1,2-Dichlorobenzene	6.97	146	329066	38.439	ng	98
15) Benzyl Alcohol	6.95	79	258758	35.776	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.08	45	371930	34.348	ng	85
17) 2-Methylphenol	7.07	107	269219	37.861	ng	96
18) Hexachloroethane	7.32	117	128284	38.948	ng	98
19) n-Nitroso-di-n-propylamine	7.22	70	204024	32.787	ng	94
20) 3+4-Methylphenols	7.22	107	321064	36.406	ng	# 68
22) Acetophenone	7.22	105	434271	36.313	ng	97
24) Nitrobenzene	7.39	77	334021	40.669	ng	99
25) Isophorone	7.63	82	562160	35.736	ng	98
26) 2-Nitrophenol	7.71	139	173897	52.009	ng	91
27) 2,4-Dimethylphenol	7.75	122	244604	35.232	ng	99
28) bis(2-Chloroethoxy)methane	7.84	93	353724	36.777	ng	99
29) 2,4-Dichlorophenol	7.95	162	252794	38.162	ng	99
30) 1,2,4-Trichlorobenzene	8.03	180	277815	38.215	ng	98
31) Naphthalene	8.12	128	913727	37.776	ng	100
32) Benzoic acid	7.87	122	185483m	33.484	ng	
33) 4-Chloroaniline	8.16	127	394153	38.036	ng	97
34) Hexachlorobutadiene	8.22	225	155378	39.020	ng	99
35) Caprolactam	8.55	113	87060m	37.674	ng	
36) 4-Chloro-3-methylphenol	8.65	107	273075	38.445	ng	98
37) 2-Methylnaphthalene	8.80	142	592469	37.867	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	260953	39.998	ng	99
40) Hexachlorocyclopentadiene	8.95	237	150526	41.212	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.09	196	172131	38.006	nq	97
43) 2,4,5-Trichlorophenol	9.13	196	193038	38.089	nq	97
45) 1,1'-Biphenyl	9.27	154	731167	38.188	nq	100
46) 2-Chloronaphthalene	9.29	162	577503	38.186	nq	99
47) 2-Nitroaniline	9.39	65	175703	46.980	nq	96
48) Acenaphthylene	9.71	152	885073	37.301	nq	99
49) Dimethylphthalate	9.57	163	697895	38.830	nq	100
50) 2,6-Dinitrotoluene	9.64	165	149687	45.010	nq #	86
51) Acenaphthene	9.88	154	516461	35.674	nq	100
52) 3-Nitroaniline	9.81	138	176398	45.307	nq	93
53) 2,4-Dinitrophenol	9.92	184	62447	54.165	nq #	20
54) Dibenzofuran	10.05	168	738549	36.606	nq	97
55) 4-Nitrophenol	9.97	139	113081	36.974	nq	92
56) 2,4-Dinitrotoluene	10.05	165	196401	45.022	nq #	92
57) Fluorene	10.40	166	598440	40.751	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.17	232	144191	38.135	nq	94
59) Diethylphthalate	10.27	149	661667	36.965	nq	97
60) 4-Chlorophenyl-phenylether	10.39	204	269446	41.199	nq	100
61) 4-Nitroaniline	10.42	138	177802	46.706	nq	97
62) Azobenzene	10.55	77	608004	35.553	nq	92
64) 4,6-Dinitro-2-methylphenol	10.46	198	103115	49.555	nq #	69
65) n-Nitrosodiphenylamine	10.51	169	538969	36.589	nq	99
66) 4-Bromophenyl-phenylether	10.87	248	170570	37.745	nq	98
67) Hexachlorobenzene	10.95	284	194648	38.280	nq #	90
68) Atrazine	11.03	200	165310	37.179	nq	100
69) Pentachlorophenol	11.15	266	117409	37.323	nq	96
70) Phenanthrene	11.36	178	864947	36.558	nq	99
71) Anthracene	11.42	178	887106	36.886	nq	99
72) Carbazole	11.57	167	864818	36.799	nq	100
73) Di-n-butylphthalate	11.89	149	1034107	36.022	nq	100
74) Fluoranthene	12.55	202	898628	36.353	nq	99
76) Benzidine	12.67	184	385186m	32.122	nq	
77) Pyrene	12.78	202	933765	37.764	nq	100
79) Butylbenzylphthalate	13.39	149	436471	37.469	nq	97
80) Benzo(a)anthracene	13.97	228	796933	39.990	nq	99
81) 3,3'-Dichlorobenzidine	13.93	252	281667	37.907	nq	99
82) Chrysene	14.01	228	764090	37.328	nq	99
83) Bis(2-ethylhexyl)phthalate	13.95	149	579474	38.675	nq	100
84) Di-n-octyl phthalate	14.57	149	879195	35.117	nq #	95
85) Indeno(1,2,3-cd)pyrene	16.91	276	601891	34.614	nq	98
87) Benzo(b)fluoranthene	15.03	252	677934	37.222	nq	98
88) Benzo(k)fluoranthene	15.06	252	677911	39.425	nq	99
89) Benzo(a)pyrene	15.39	252	621365	38.129	nq	98
90) Dibenzo(a,h)anthracene	16.93	278	497486	37.169	nq	98
91) Benzo(g,h,i)perylene	17.35	276	484793	37.386	nq	98

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

