

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111517\
 Data File : BF100602.D
 Acq On : 15 Nov 2017 22:29
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Nov 16 03:38:02 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 15 13:27:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	122822	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	467310	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	180027	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	262268	20.00	ng	0.00
75) Chrysene-d12	14.04	240	244252	20.00	ng	0.00
86) Perylene-d12	15.52	264	258895	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	605928	79.08	ng	0.00
7) Phenol-d6	6.51	99	735798	77.88	ng	0.00
23) Nitrobenzene-d5	7.45	82	652587	92.33	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	136342	74.32	ng	0.00
44) 2-Fluorobiphenyl	9.24	172	1094624	92.59	ng	0.00
78) Terphenyl-d14	12.99	244	723143	68.03	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.66	88	147612	39.532	ng	97
3) Pyridine	3.43	79	401096	39.373	ng	95
4) n-Nitrosodimethylamine	3.39	42	183612	37.123	ng	97
6) Aniline	6.55	93	495776	37.142	ng	100
8) 2-Chlorophenol	6.67	128	327745	40.434	ng	99
9) Benzaldehyde	6.44	77	241975	35.862	ng	94
10) Phenol	6.52	94	383203	38.634	ng	99
11) bis(2-Chloroethyl)ether	6.62	93	324597	38.961	ng	95
12) 1,3-Dichlorobenzene	6.83	146	374948	40.122	ng	98
13) 1,4-Dichlorobenzene	6.90	146	378289	39.994	ng	97
14) 1,2-Dichlorobenzene	7.06	146	344669	39.412	ng	97
15) Benzyl Alcohol	7.02	79	289862	39.309	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.16	45	431155	32.357	ng	97
17) 2-Methylphenol	7.13	107	273543	39.490	ng	99
18) Hexachloroethane	7.40	117	119582	38.344	ng	93
19) n-Nitroso-di-n-propylamine	7.30	70	241761	36.896	ng	95
20) 3+4-Methylphenols	7.29	107	331427	37.811	ng	90
22) Acetophenone	7.30	105	447899	39.306	ng	# 98
24) Nitrobenzene	7.47	77	334068	45.920	ng	98
25) Isophorone	7.70	82	577651	38.890	ng	100
26) 2-Nitrophenol	7.78	139	172724	50.342	ng	95
27) 2,4-Dimethylphenol	7.82	122	263162	39.523	ng	98
28) bis(2-Chloroethoxy)methane	7.92	93	394520	40.606	ng	98
29) 2,4-Dichlorophenol	8.02	162	267680	42.111	ng	97
30) 1,2,4-Trichlorobenzene	8.11	180	288846	42.009	ng	96
31) Naphthalene	8.19	128	897405	39.870	ng	100
32) Benzoic acid	7.92	122	184470	38.487	ng	97
33) 4-Chloroaniline	8.23	127	370649	39.561	ng	99
34) Hexachlorobutadiene	8.30	225	172929	42.300	ng	98
35) Caprolactam	8.61	113	75397	34.563	ng	94
36) 4-Chloro-3-methylphenol	8.71	107	254794	37.322	ng	97
37) 2-Methylnaphthalene	8.88	142	566713	37.925	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.05	216	294377	49.304	ng	# 97
40) Hexachlorocyclopentadiene	9.03	237	60934	21.684	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.15	196	176475	46.636	ng	99
43) 2,4,5-Trichlorophenol	9.19	196	181034	46.175	ng	93
45) 1,1'-Biphenyl	9.35	154	717400	46.074	ng	98
46) 2-Chloronaphthalene	9.37	162	514962	45.589	ng	96
47) 2-Nitroaniline	9.46	65	155448	46.656	ng	99
48) Acenaphthylene	9.79	152	789452	42.172	ng	99
49) Dimethylphthalate	9.64	163	608072	44.239	ng	100
50) 2,6-Dinitrotoluene	9.70	165	122903	45.172	ng	92
51) Acenaphthene	9.96	154	462813	40.651	ng	100
52) 3-Nitroaniline	9.87	138	137616	42.833	ng	92
53) 2,4-Dinitrophenol	9.97	184	20898	28.920	ng	# 14
54) Dibenzofuran	10.13	168	650138	40.744	ng	98
55) 4-Nitrophenol	10.02	139	87665	33.604	ng	88
56) 2,4-Dinitrotoluene	10.10	165	145301	42.332	ng	# 91
57) Fluorene	10.47	166	466224	39.884	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.24	232	118361	36.836	ng	# 87
59) Diethylphthalate	10.34	149	552379	40.158	ng	96
60) 4-Chlorophenyl-phenylether	10.46	204	241521	42.118	ng	93
61) 4-Nitroaniline	10.48	138	114626	37.626	ng	90
62) Azobenzene	10.62	77	469204	36.217	ng	93
64) 4,6-Dinitro-2-methylphenol	10.51	198	41340	34.214	ng	84
65) n-Nitrosodiphenylamine	10.57	169	458044	50.228	ng	96
66) 4-Bromophenyl-phenylether	10.95	248	138022	49.092	ng	# 82
67) Hexachlorobenzene	11.02	284	133661	45.456	ng	# 88
68) Atrazine	11.10	200	122525	44.386	ng	97
69) Pentachlorophenol	11.20	266	84668	40.156	ng	98
70) Phenanthrene	11.43	178	573641	41.542	ng	100
71) Anthracene	11.48	178	588400	41.999	ng	100
72) Carbazole	11.63	167	510811	39.516	ng	98
73) Di-n-butylphthalate	11.96	149	668173	44.169	ng	# 98
74) Fluoranthene	12.62	202	568782	38.865	ng	98
76) Benzidine	12.74	184	434500	44.419	ng	98
77) Pyrene	12.84	202	594307	34.134	ng	99
79) Butylbenzylphthalate	13.46	149	256592	36.137	ng	93
80) Benzo(a)anthracene	14.03	228	612478	41.640	ng	100
81) 3,3'-Dichlorobenzidine	13.99	252	235101	44.612	ng	98
82) Chrysene	14.07	228	578809	40.637	ng	99
83) Bis(2-ethylhexyl)phthalate	14.01	149	379258	40.610	ng	97
84) Di-n-octyl phthalate	14.63	149	717226	44.705	ng	99
85) Indeno(1,2,3-cd)pyrene	17.00	276	552946	47.995	ng	95
87) Benzo(b)fluoranthene	15.09	252	611499	39.868	ng	98
88) Benzo(k)fluoranthene	15.12	252	622053	42.923	ng	97
89) Benzo(a)pyrene	15.46	252	571106	42.000	ng	100
90) Dibenzo(a,h)anthracene	17.02	278	474009	42.625	ng	96
91) Benzo(g,h,i)perylene	17.45	276	440783	40.492	ng	94

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

