

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100815\
 Data File : BF082122.D
 Acq On : 8 Oct 2015 11:12
 Operator : UM/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 09 05:03:34 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 08 02:46:05 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	100240	20.00	ng	-0.01
21) Naphthalene-d8	8.15	136	401659	20.00	ng	-0.01
38) Acenaphthene-d10	9.91	164	203557	20.00	ng	-0.01
63) Phenanthrene-d10	11.39	188	401195	20.00	ng	-0.01
75) Chrysene-d12	14.05	240	367424	20.00	ng	-0.01
86) Perylene-d12	15.49	264	302564	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.45	112	405054	73.35	ng	-0.01
7) Phenol-d6	6.49	99	516693	74.36	ng	-0.01
23) Nitrobenzene-d5	7.44	82	491902	81.64	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	147174	71.39	ng	-0.02
44) 2-Fluorobiphenyl	9.23	172	1014400	83.01	ng	-0.01
78) Terphenyl-d14	12.98	244	1004048	89.59	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.42	88	92312	38.44	ng	87
3) Pyridine	3.15	79	224565	35.92	ng	# 84
4) n-Nitrosodimethylamine	3.12	42	93584	32.96	ng	96
6) Aniline	6.53	93	338822	36.30	ng	# 87
8) 2-Chlorophenol	6.65	128	231401	37.95	ng	94
9) Benzaldehyde	6.41	77	143076	31.44	ng	# 79
10) Phenol	6.51	94	273452	35.08	ng	# 71
11) bis(2-Chloroethyl)ether	6.61	93	251254	41.57	ng	92
12) 1,3-Dichlorobenzene	6.80	146	271032m	37.63	ng	
13) 1,4-Dichlorobenzene	6.88	146	292557m	38.90	ng	
14) 1,2-Dichlorobenzene	7.04	146	259185	38.67	ng	98
15) Benzyl Alcohol	7.01	79	178139	35.52	ng	# 76
16) 2,2'-oxybis(1-Chloropropan	7.14	45	317353	37.14	ng	82
17) 2-Methylphenol	7.12	107	189536	38.34	ng	# 86
18) Hexachloroethane	7.37	117	93499	39.71	ng	90
19) n-Nitroso-di-n-propylamine	7.28	70	160366	37.97	ng	# 78
20) 3+4-Methylphenols	7.28	107	234232	37.51	ng	# 59
22) Acetophenone	7.28	105	309218	37.65	ng	# 77
24) Nitrobenzene	7.45	77	238126	36.40	ng	# 67
25) Isophorone	7.69	82	465468	38.98	ng	# 89
26) 2-Nitrophenol	7.77	139	138074	43.96	ng	# 73
27) 2,4-Dimethylphenol	7.81	122	204911	37.09	ng	# 81
28) bis(2-Chloroethoxy)methane	7.91	93	292200	41.20	ng	# 95
29) 2,4-Dichlorophenol	8.01	162	220816	41.22	ng	99
30) 1,2,4-Trichlorobenzene	8.09	180	231590	38.72	ng	97
31) Naphthalene	8.17	128	642302	36.09	ng	97
32) Benzoic acid	7.93	122	110648	35.70	ng	# 70
33) 4-Chloroaniline	8.23	127	303675	39.46	ng	# 91
34) Hexachlorobutadiene	8.29	225	137536	38.88	ng	97
35) Caprolactam	8.61	113	59595	40.18	ng	# 82
36) 4-Chloro-3-methylphenol	8.71	107	218563	41.72	ng	83
37) 2-Methylnaphthalene	8.87	142	488944	40.25	ng	# 91
39) 1,2,4,5-Tetrachlorobenzene	9.03	216	227529	36.41	ng	99
40) Hexachlorocyclopentadiene	9.02	237	111905	34.77	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.14	196	154190	35.99	nq	100
43) 2,4,5-Trichlorophenol	9.19	196	177180	40.22	nq	# 94
45) 1,1'-Biphenyl	9.34	154	602688	38.38	nq	94
46) 2-Chloronaphthalene	9.36	162	488666	39.63	nq	96
47) 2-Nitroaniline	9.46	65	138957	38.39	nq	93
48) Acenaphthylene	9.77	152	730125	36.99	nq	96
49) Dimethylphthalate	9.63	163	517426	36.83	nq	# 97
50) 2,6-Dinitrotoluene	9.70	165	131678	43.17	nq	# 86
51) Acenaphthene	9.94	154	450805	38.47	nq	90
52) 3-Nitroaniline	9.87	138	150784	42.72	nq	# 80
53) 2,4-Dinitrophenol	9.98	184	51176	47.75	nq	# 78
54) Dibenzofuran	10.11	168	621479	37.52	nq	# 87
55) 4-Nitrophenol	10.03	139	80590	31.60	nq	# 76
56) 2,4-Dinitrotoluene	10.10	165	161907	41.63	nq	# 83
57) Fluorene	10.46	166	497021	41.81	nq	92
58) 2,3,4,6-Tetrachlorophenol	10.23	232	140194	40.12	nq	95
59) Diethylphthalate	10.33	149	532981	40.61	nq	99
60) 4-Chlorophenyl-phenylether	10.45	204	229573	37.86	nq	# 80
61) 4-Nitroaniline	10.49	138	146824	41.49	nq	# 54
62) Azobenzene	10.61	77	502030	39.19	nq	88
64) 4,6-Dinitro-2-methylphenol	10.51	198	88607	48.54	nq	# 64
65) n-Nitrosodiphenylamine	10.57	169	451064	37.16	nq	91
66) 4-Bromophenyl-phenylether	10.94	248	156021	38.57	nq	# 89
67) Hexachlorobenzene	11.01	284	173315	37.96	nq	# 78
68) Atrazine	11.10	200	139396	37.01	nq	83
69) Pentachlorophenol	11.20	266	94862	36.47	nq	98
70) Phenanthrene	11.42	178	742904	38.23	nq	96
71) Anthracene	11.47	178	841428	41.26	nq	97
72) Carbazole	11.63	167	755808	40.95	nq	97
73) Di-n-butylphthalate	11.96	149	829401	44.10	nq	# 98
74) Fluoranthene	12.61	202	797622	37.11	nq	92
76) Benzidine	12.74	184	451283	40.76	nq	98
77) Pyrene	12.84	202	820569	37.38	nq	96
79) Butylbenzylphthalate	13.45	149	371603	44.99	nq	# 80
80) Benzo(a)anthracene	14.02	228	746625	37.74	nq	96
81) 3,3'-Dichlorobenzidine	14.00	252	281174	42.08	nq	# 91
82) Chrysene	14.07	228	809717	49.35	nq	94
83) Bis(2-ethylhexyl)phthalate	14.01	149	518700	50.06	nq	# 92
84) Di-n-octyl phthalate	14.63	149	882256	52.33	nq	# 92
85) Indeno(1,2,3-cd)pyrene	16.90	276	615586	39.77	nq	# 88
87) Benzo(b)fluoranthene	15.06	252	611984	35.43	nq	# 87
88) Benzo(k)fluoranthene	15.10	252	710957m	44.90	nq	
89) Benzo(a)pyrene	15.42	252	600507	40.08	nq	# 87
90) Dibenzo(a,h)anthracene	16.93	278	505180	40.18	nq	# 81
91) Benzo(g,h,i)perylene	17.34	276	496522	38.54	nq	# 77

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

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