

Data Path : Z:\HPCHEM1\BNA F\DATA\BF033115\
 Data File : BF078131.D
 Acq On : 1 Apr 2015 1:12
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Apr 01 08:39:38 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 01 08:27:38 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.05	152	43849	20.00	ng	0.00
21) Naphthalene-d8	8.62	136	176694	20.00	ng	0.00
38) Acenaphthene-d10	10.79	164	89792	20.00	ng	0.00
63) Phenanthrene-d10	12.63	188	175927	20.00	ng	0.00
75) Chrysene-d12	15.91	240	175523	20.00	ng	0.00
86) Perylene-d12	17.61	264	165317	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.32	112	200482	79.81	ng	0.00
7) Phenol-d6	6.64	99	249468	80.38	ng	0.00
23) Nitrobenzene-d5	7.74	82	226256	83.94	ng	0.00
41) 2,4,6-Tribromophenol	11.77	330	64734	75.35	ng	0.00
44) 2-Fluorobiphenyl	9.97	172	489422	80.10	ng	0.00
78) Terphenyl-d14	14.63	244	559545	77.87	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	1.89	88	40459	35.47	ng	# 100
3) Pyridine	2.53	79	109103	35.60	ng	96
4) n-Nitrosodimethylamine	2.48	42	43302	32.45	ng	96
6) Aniline	6.64	93	172568	38.87	ng	95
8) 2-Chlorophenol	6.78	128	115556	38.65	ng	89
9) Benzaldehyde	6.49	77	57924	45.55	ng	97
10) Phenol	6.65	94	150571	41.04	ng	# 78
11) bis(2-Chloroethyl)ether	6.74	93	105450	38.37	ng	91
12) 1,3-Dichlorobenzene	6.97	146	131745	39.61	ng	97
13) 1,4-Dichlorobenzene	7.07	146	134663	38.97	ng	99
14) 1,2-Dichlorobenzene	7.25	146	123194	38.89	ng	99
15) Benzyl Alcohol	7.24	79	96433	39.80	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.41	45	144850	39.11	ng	100
17) 2-Methylphenol	7.40	107	93371	38.36	ng	96
18) Hexachloroethane	7.66	117	46510	41.04	ng	94
19) n-Nitroso-di-n-propylamine	7.57	70	85154	39.66	ng	96
20) 3+4-Methylphenols	7.58	107	129242	40.46	ng	92
22) Acetophenone	7.56	105	162140	41.45	ng	# 92
24) Nitrobenzene	7.77	77	121267	41.76	ng	# 87
25) Isophorone	8.06	82	220172	40.45	ng	95
26) 2-Nitrophenol	8.16	139	55863	39.29	ng	# 83
27) 2,4-Dimethylphenol	8.24	122	105623	40.78	ng	97
28) bis(2-Chloroethoxy)methane	8.35	93	137457	41.60	ng	99
29) 2,4-Dichlorophenol	8.46	162	98934	40.07	ng	97
30) 1,2,4-Trichlorobenzene	8.56	180	110453	39.99	ng	99
31) Naphthalene	8.65	128	374895	41.31	ng	100
32) Benzoic acid	8.36	122	47409	33.79	ng	96
33) 4-Chloroaniline	8.73	127	143512	38.97	ng	# 92
34) Hexachlorobutadiene	8.81	225	65024	41.53	ng	99
35) Caprolactam	9.16	113	28583	39.61	ng	97
36) 4-Chloro-3-methylphenol	9.36	107	99532	40.07	ng	99
37) 2-Methylnaphthalene	9.50	142	240756	39.49	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.71	216	103465	38.63	ng	# 100
40) Hexachlorocyclopentadiene	9.70	237	45585	35.82	ng	99

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42) 2,4,6-Trichlorophenol	9.87	196	71026	38.29	nq	99
43) 2,4,5-Trichlorophenol	9.92	196	77777	38.81	nq	96
45) 1,1'-Biphenyl	10.09	154	290204	40.43	nq	98
46) 2-Chloronaphthalene	10.11	162	232919	39.93	nq	99
47) 2-Nitroaniline	10.25	65	60036	41.44	nq	88
48) Acenaphthylene	10.61	152	383215	39.51	nq	100
49) Dimethylphthalate	10.49	163	271092	40.71	nq	100
50) 2,6-Dinitrotoluene	10.56	165	59039	42.77	nq	# 85
51) Acenaphthene	10.83	154	220928	39.72	nq	100
52) 3-Nitroaniline	10.75	138	63430	39.56	nq	# 73
53) 2,4-Dinitrophenol	10.89	184	13863	32.64	nq	90
54) Dibenzofuran	11.05	168	322735	39.12	nq	100
55) 4-Nitrophenol	10.99	139	45938	38.68	nq	86
56) 2,4-Dinitrotoluene	11.05	165	73948	41.08	nq	# 84
57) Fluorene	11.47	166	274022	40.01	nq	99
58) 2,3,4,6-Tetrachlorophenol	11.21	232	58975	38.22	nq	# 100
59) Diethylphthalate	11.36	149	257940	39.75	nq	97
60) 4-Chlorophenyl-phenylether	11.48	204	122177	39.09	nq	99
61) 4-Nitroaniline	11.51	138	65752	41.15	nq	78
62) Azobenzene	11.68	77	282053	45.48	nq	99
64) 4,6-Dinitro-2-methylphenol	11.55	198	31078	38.36	nq	77
65) n-Nitrosodiphenylamine	11.63	169	236791	40.42	nq	99
66) 4-Bromophenyl-phenylether	12.09	248	71654	38.16	nq	96
67) Hexachlorobenzene	12.13	284	77974	37.80	nq	# 74
68) Atrazine	12.31	200	60475	36.84	nq	94
69) Pentachlorophenol	12.40	266	29987	29.77	nq	97
70) Phenanthrene	12.65	178	395178	39.13	nq	99
71) Anthracene	12.72	178	396920	39.78	nq	100
72) Carbazole	12.93	167	374074	39.41	nq	98
73) Di-n-butylphthalate	13.39	149	432308	40.62	nq	99
74) Fluoranthene	14.12	202	418274	37.55	nq	96
76) Benzidine	14.32	184	197842	45.64	nq	98
77) Pyrene	14.40	202	446667	40.47	nq	98
79) Butylbenzylphthalate	15.24	149	176650	40.59	nq	92
80) Benzo(a)anthracene	15.89	228	409829	39.39	nq	99
81) 3,3'-Dichlorobenzidine	15.87	252	128820	37.53	nq	97
82) Chrysene	15.93	228	375848	40.17	nq	99
83) Bis(2-ethylhexyl)phthalate	15.96	149	271866	41.24	nq	# 96
84) Di-n-octyl phthalate	16.75	149	454058	40.29	nq	100
85) Indeno(1,2,3-cd)pyrene	18.93	276	424467	36.35	nq	# 100
87) Benzo(b)fluoranthene	17.17	252	393934	36.50	nq	# 95
88) Benzo(k)fluoranthene	17.20	252	375626m	44.56	nq	
89) Benzo(a)pyrene	17.54	252	346567	38.49	nq	# 95
90) Dibenzo(a,h)anthracene	18.96	278	338235	37.87	nq	98
91) Benzo(g,h,i)perylene	19.32	276	344410	37.88	nq	100

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

