

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101215\
 Data File : BF082234.D
 Acq On : 12 Oct 2015 22:16
 Operator : UM/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

MMdadoda
 10/13/2015 1:27:49 PM

Quant Time: Oct 13 02:54:36 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 13 02:05:16 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	76666	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	306046	20.00	ng	0.00
38) Acenaphthene-d10	9.87	164	150412	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	309187	20.00	ng	0.00
75) Chrysene-d12	14.01	240	285002	20.00	ng	0.00
86) Perylene-d12	15.44	264	229881	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	299111	78.74	ng	0.00
7) Phenol-d6	6.47	99	382810	77.44	ng	0.00
23) Nitrobenzene-d5	7.41	82	359977	78.99	ng	0.00
41) 2,4,6-Tribromophenol	10.67	330	117259	83.13	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	721411	79.54	ng	0.00
78) Terphenyl-d14	12.95	244	778664	72.40	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.37	88	72814	38.15	ng	95
3) Pyridine	3.10	79	173954	39.19	ng	97
4) n-Nitrosodimethylamine	3.05	42	76076	40.41	ng	99
6) Aniline	6.49	93	253449	39.77	ng	# 72
8) 2-Chlorophenol	6.62	128	182537	39.36	ng	92
9) Benzaldehyde	6.38	77	101738	40.30	ng	93
10) Phenol	6.48	94	201981	35.44	ng	99
11) bis(2-Chloroethyl)ether	6.57	93	176566	36.63	ng	94
12) 1,3-Dichlorobenzene	6.77	146	200658m	37.16	ng	
13) 1,4-Dichlorobenzene	6.85	146	206533	36.62	ng	96
14) 1,2-Dichlorobenzene	7.01	146	206237m	38.51	ng	
15) Benzyl Alcohol	6.98	79	143834	42.15	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.12	45	249016	37.10	ng	98
17) 2-Methylphenol	7.10	107	147401	40.20	ng	97
18) Hexachloroethane	7.35	117	69027	35.76	ng	# 73
19) n-Nitroso-di-n-propylamine	7.26	70	131448	37.86	ng	96
20) 3+4-Methylphenols	7.26	107	175702	40.21	ng	# 84
22) Acetophenone	7.25	105	233583	38.58	ng	98
24) Nitrobenzene	7.43	77	188873	38.29	ng	# 83
25) Isophorone	7.66	82	340201	36.99	ng	98
26) 2-Nitrophenol	7.74	139	98408	39.95	ng	# 86
27) 2,4-Dimethylphenol	7.78	122	166732	42.01	ng	96
28) bis(2-Chloroethoxy)methane	7.87	93	211698	39.56	ng	98
29) 2,4-Dichlorophenol	7.99	162	153079	38.59	ng	93
30) 1,2,4-Trichlorobenzene	8.06	180	168480	36.85	ng	96
31) Naphthalene	8.15	128	500732	37.74	ng	99
32) Benzoic acid	7.91	122	101136	37.97	ng	96
33) 4-Chloroaniline	8.19	127	216749	39.34	ng	# 95
34) Hexachlorobutadiene	8.25	225	100809	35.38	ng	99
35) Caprolactam	8.58	113	50172	43.58	ng	91
36) 4-Chloro-3-methylphenol	8.69	107	167344	43.14	ng	89
37) 2-Methylnaphthalene	8.83	142	362211	39.38	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.01	216	172610	38.58	ng	98
40) Hexachlorocyclopentadiene	8.98	237	63070	33.27	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.12	196	120438	40.53	nq	99
43) 2,4,5-Trichlorophenol	9.15	196	126676	39.35	nq	# 89
45) 1,1'-Biphenyl	9.30	154	470389	41.01	nq	96
46) 2-Chloronaphthalene	9.33	162	361012	39.98	nq	96
47) 2-Nitroaniline	9.43	65	110614	44.62	nq	# 84
48) Acenaphthylene	9.74	152	548150	38.97	nq	99
49) Dimethylphthalate	9.61	163	413985	39.09	nq	# 97
50) 2,6-Dinitrotoluene	9.67	165	89568	40.08	nq	# 82
51) Acenaphthene	9.91	154	333546	39.50	nq	98
52) 3-Nitroaniline	9.84	138	101601	40.25	nq	87
53) 2,4-Dinitrophenol	9.95	184	40639	36.34	nq	# 90
54) Dibenzofuran	10.08	168	448515	38.69	nq	96
55) 4-Nitrophenol	10.01	139	59746	42.12	nq	93
56) 2,4-Dinitrotoluene	10.08	165	120608	43.18	nq	89
57) Fluorene	10.42	166	363165	38.14	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.21	232	113407	43.12	nq	94
59) Diethylphthalate	10.30	149	402122	40.73	nq	98
60) 4-Chlorophenyl-phenylether	10.42	204	183776	42.14	nq	# 83
61) 4-Nitroaniline	10.46	138	108991	44.51	nq	97
62) Azobenzene	10.58	77	395869	40.97	nq	87
64) 4,6-Dinitro-2-methylphenol	10.48	198	64687	37.28	nq	# 67
65) n-Nitrosodiphenylamine	10.54	169	342445	36.99	nq	99
66) 4-Bromophenyl-phenylether	10.91	248	119829	38.73	nq	# 81
67) Hexachlorobenzene	10.97	284	126615	37.84	nq	# 81
68) Atrazine	11.07	200	126046	42.98	nq	95
69) Pentachlorophenol	11.17	266	66900	38.85	nq	97
70) Phenanthrene	11.39	178	592664	39.11	nq	98
71) Anthracene	11.44	178	623599	39.93	nq	100
72) Carbazole	11.60	167	569872	40.00	nq	99
73) Di-n-butylphthalate	11.93	149	672647	38.27	nq	# 97
74) Fluoranthene	12.57	202	651832	40.05	nq	96
76) Benzidine	12.71	184	312960	37.89	nq	100
77) Pyrene	12.80	202	668048	39.50	nq	99
79) Butylbenzylphthalate	13.43	149	309143	40.05	nq	# 83
80) Benzo(a)anthracene	14.00	228	573549	38.68	nq	99
81) 3,3'-Dichlorobenzidine	13.97	252	217126	38.58	nq	# 96
82) Chrysene	14.04	228	627622	40.17	nq	99
83) Bis(2-ethylhexyl)phthalate	13.98	149	385238	34.98	nq	# 95
84) Di-n-octyl phthalate	14.60	149	657586	34.77	nq	100
85) Indeno(1,2,3-cd)pyrene	16.85	276	505102	40.67	nq	99
87) Benzo(b)fluoranthene	15.03	252	506231	36.20	nq	# 97
88) Benzo(k)fluoranthene	15.05	252	479408m	40.78	nq	
89) Benzo(a)pyrene	15.38	252	464860	38.68	nq	# 96
90) Dibenzo(a,h)anthracene	16.87	278	413795	42.01	nq	97
91) Benzo(g,h,i)perylene	17.28	276	412224	43.08	nq	96

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

