

Data Path : Z:\HPCHEM1\BNA F\DATA\BF123015\
 Data File : BF084165.D
 Acq On : 30 Dec 2015 23:59
 Operator : UM/SJ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Sohil
 12/31/2015 6:41:07 PM

Quant Time: Dec 31 03:49:27 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 31 03:17:24 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	359256	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	1516286	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	680062	20.00	ng	-0.01
63) Phenanthrene-d10	11.43	188	1319197	20.00	ng	0.00
75) Chrysene-d12	14.07	240	1136892	20.00	ng	0.00
86) Perylene-d12	15.49	264	854729	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	460934	21.51	ng	-0.01
7) Phenol-d6	6.51	99	605973	22.73	ng	-0.01
23) Nitrobenzene-d5	7.46	82	512650	19.50	ng	-0.01
41) 2,4,6-Tribromophenol	10.73	330	150488	20.74	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	1017898	20.96	ng	0.00
78) Terphenyl-d14	13.01	244	970545	15.98	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.53	88	106987	11.90	ng	98
3) Pyridine	3.28	79	278664	12.24	ng	98
4) n-Nitrosodimethylamine	3.20	42	140197	13.52	ng	# 88
6) Aniline	6.56	93	424833	12.39	ng	99
8) 2-Chlorophenol	6.68	128	261950	10.08	ng	# 82
9) Benzaldehyde	6.44	77	220727	15.21	ng	93
10) Phenol	6.52	94	359133	11.85	ng	90
11) bis(2-Chloroethyl)ether	6.62	93	238718	10.87	ng	95
12) 1,3-Dichlorobenzene	6.84	146	286017m	10.14	ng	
13) 1,4-Dichlorobenzene	6.92	146	262905	9.17	ng	95
14) 1,2-Dichlorobenzene	7.07	146	277454	10.58	ng	94
15) Benzyl Alcohol	7.04	79	257047	12.41	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.18	45	490946	20.02	ng	99
17) 2-Methylphenol	7.14	107	202841	9.92	ng	97
18) Hexachloroethane	7.41	117	101793	9.67	ng	# 71
19) n-Nitroso-di-n-propylamine	7.31	70	205633	12.42	ng	93
20) 3+4-Methylphenols	7.30	107	290605	11.18	ng	91
22) Acetophenone	7.31	105	390184	10.38	ng	# 70
24) Nitrobenzene	7.48	77	255107	9.38	ng	# 79
25) Isophorone	7.72	82	535485	10.99	ng	99
26) 2-Nitrophenol	7.80	139	96505	7.00	ng	# 82
27) 2,4-Dimethylphenol	7.84	122	249805m	10.83	ng	
28) bis(2-Chloroethoxy)methane	7.94	93	331354	11.60	ng	99
29) 2,4-Dichlorophenol	8.03	162	203649	9.32	ng	93
30) 1,2,4-Trichlorobenzene	8.13	180	221096	9.36	ng	# 95
31) Naphthalene	8.21	128	736397	9.29	ng	94
32) Benzoic acid	7.89	122	120308m	9.65	ng	
33) 4-Chloroaniline	8.25	127	318952	9.79	ng	# 93
34) Hexachlorobutadiene	8.33	225	121037	8.90	ng	95
35) Caprolactam	8.59	113	71906	11.64	ng	93
36) 4-Chloro-3-methylphenol	8.73	107	242789	9.92	ng	98
37) 2-Methylnaphthalene	8.90	142	523235	10.37	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.07	216	200305	9.43	ng	99
40) Hexachlorocyclopentadiene	9.06	237	124480	21.65	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.17	196	152493m	10.90	nq	
43) 2,4,5-Trichlorophenol	9.21	196	156921m	11.22	nq	
45) 1,1'-Biphenyl	9.37	154	642289	10.79	nq	92
46) 2-Chloronaphthalene	9.39	162	451151	9.91	nq	# 88
47) 2-Nitroaniline	9.47	65	104178	8.34	nq	# 79
48) Acenaphthylene	9.80	152	773630	10.24	nq	98
49) Dimethylphthalate	9.66	163	540658	9.70	nq	98
50) 2,6-Dinitrotoluene	9.72	165	101110	8.65	nq	# 84
51) Acenaphthene	9.97	154	490732	10.97	nq	99
52) 3-Nitroaniline	9.88	138	117585	9.38	nq	# 96
53) 2,4-Dinitrophenol	9.98	184	24502	6.92	nq	# 9
54) Dibenzofuran	10.14	168	670410	10.86	nq	96
55) 4-Nitrophenol	10.03	139	98350	13.54	nq	# 83
56) 2,4-Dinitrotoluene	10.12	165	113119	7.20	nq	# 88
57) Fluorene	10.49	166	520670	10.18	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.26	232	140217	13.72	nq	94
59) Diethylphthalate	10.36	149	569472	9.98	nq	96
60) 4-Chlorophenyl-phenylether	10.49	204	230486	9.69	nq	# 86
61) 4-Nitroaniline	10.49	138	108224	9.02	nq	90
62) Azobenzene	10.64	77	520252	10.68	nq	93
64) 4,6-Dinitro-2-methylphenol	10.52	198	41228	5.59	nq	88
65) n-Nitrosodiphenylamine	10.60	169	426928	9.34	nq	97
66) 4-Bromophenyl-phenylether	10.97	248	143051	9.37	nq	# 73
67) Hexachlorobenzene	11.04	284	161028	9.72	nq	# 77
68) Atrazine	11.13	200	138395	9.72	nq	91
69) Pentachlorophenol	11.22	266	90592	16.20	nq	97
70) Phenanthrene	11.45	178	810185	10.52	nq	98
71) Anthracene	11.49	178	773167	9.62	nq	95
72) Carbazole	11.65	167	720814	10.49	nq	# 94
73) Di-n-butylphthalate	12.00	149	871289	9.85	nq	# 95
74) Fluoranthene	12.64	202	805740	9.95	nq	92
76) Benzidine	12.75	184	455987	11.50	nq	99
77) Pyrene	12.87	202	823343	8.69	nq	96
79) Butylbenzylphthalate	13.49	149	335857	8.32	nq	# 81
80) Benzo(a)anthracene	14.05	228	645955	9.32	nq	95
81) 3,3'-Dichlorobenzidine	14.01	252	228520	10.57	nq	# 96
82) Chrysene	14.09	228	622315	10.03	nq	97
83) Bis(2-ethylhexyl)phthalate	14.05	149	437786	8.61	nq	# 99
84) Di-n-octyl phthalate	14.66	149	699224	9.54	nq	100
85) Indeno(1,2,3-cd)pyrene	16.91	276	408296	6.57	nq	100
87) Benzo(b)fluoranthene	15.08	252	548177	8.64	nq	# 96
88) Benzo(k)fluoranthene	15.11	252	516826m	12.07	nq	
89) Benzo(a)pyrene	15.44	252	490043	9.41	nq	# 96
90) Dibenzo(a,h)anthracene	16.93	278	338918	6.83	nq	99
91) Benzo(g,h,i)perylene	17.33	276	333125	6.56	nq	96

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

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