

Data Path : Z:\HPCHEM1\BNA F\DATA\BF123115\
 Data File : BF084211.D
 Acq On : 31 Dec 2015 22:49
 Operator : UM/SJ
 Sample : PB87590BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB87590BS

Manual Integrations
 APPROVED

UMANGI
 1/4/2016 4:29:44 PM

Quant Time: Jan 02 00:13:30 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 31 14:15:57 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	287430	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	1217353	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	545184	20.00	ng	-0.01
63) Phenanthrene-d10	11.42	188	999951	20.00	ng	0.00
75) Chrysene-d12	14.07	240	734274	20.00	ng	0.00
86) Perylene-d12	15.51	264	535158	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	1757620	98.91	ng	0.01
7) Phenol-d6	6.53	99	2428853	108.83	ng	0.00
23) Nitrobenzene-d5	7.47	82	1643674	75.15	ng	0.00
41) 2,4,6-Tribromophenol	10.74	330	631114	114.66	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	2347675	75.01	ng	0.00
78) Terphenyl-d14	13.01	244	2212431	67.26	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.61	88	261832	31.10	ng	# 74
3) Pyridine	3.34	79	663352	28.26	ng	# 72
4) n-Nitrosodimethylamine	3.29	42	406615	33.75	ng	# 77
6) Aniline	6.56	93	724529	22.19	ng	# 81
8) 2-Chlorophenol	6.68	128	694822	34.46	ng	83
9) Benzaldehyde	6.44	77	255839	17.61	ng	94
10) Phenol	6.54	94	1058182	41.70	ng	87
11) bis(2-Chloroethyl)ether	6.64	93	673889	34.28	ng	# 82
12) 1,3-Dichlorobenzene	6.84	146	747220m	35.93	ng	
13) 1,4-Dichlorobenzene	6.92	146	763512m	36.61	ng	
14) 1,2-Dichlorobenzene	7.07	146	666517	33.37	ng	97
15) Benzyl Alcohol	7.05	79	687918	36.64	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.18	45	1233430	34.46	ng	84
17) 2-Methylphenol	7.15	107	581323	34.40	ng	94
18) Hexachloroethane	7.41	117	290648	34.85	ng	91
19) n-Nitroso-di-n-propylamine	7.32	70	555636	36.78	ng	# 77
20) 3+4-Methylphenols	7.31	107	763270	38.43	ng	# 81
22) Acetophenone	7.31	105	955688	32.65	ng	# 84
24) Nitrobenzene	7.48	77	725661	32.71	ng	94
25) Isophorone	7.72	82	1410167	33.71	ng	97
26) 2-Nitrophenol	7.80	139	374350	37.08	ng	# 85
27) 2,4-Dimethylphenol	7.84	122	691945	33.86	ng	96
28) bis(2-Chloroethoxy)methane	7.94	93	891719	34.90	ng	99
29) 2,4-Dichlorophenol	8.04	162	596197	35.02	ng	97
30) 1,2,4-Trichlorobenzene	8.13	180	639110	36.36	ng	# 94
31) Naphthalene	8.21	128	2101174	38.04	ng	99
32) Benzoic acid	7.95	122	424031	34.30	ng	90
33) 4-Chloroaniline	8.26	127	385901	15.24	ng	# 93
34) Hexachlorobutadiene	8.33	225	347072	34.35	ng	99
35) Caprolactam	8.62	113	210134m	36.62	ng	
36) 4-Chloro-3-methylphenol	8.74	107	708519	37.16	ng	95
37) 2-Methylnaphthalene	8.90	142	1305041	32.92	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.07	216	591431	37.74	ng	99
40) Hexachlorocyclopentadiene	9.06	237	702000	74.20	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.17	196	414880	35.38	nq	96
43) 2,4,5-Trichlorophenol	9.22	196	437165	38.89	nq	# 91
45) 1,1'-Biphenyl	9.37	154	1686054	38.91	nq	100
46) 2-Chloronaphthalene	9.39	162	1244507	36.57	nq	94
47) 2-Nitroaniline	9.48	65	419403	37.34	nq	98
48) Acenaphthylene	9.80	152	1877709	37.34	nq	97
49) Dimethylphthalate	9.66	163	1410592	36.19	nq	# 89
50) 2,6-Dinitrotoluene	9.73	165	325262	38.05	nq	# 87
51) Acenaphthene	9.97	154	1365925	40.50	nq	98
52) 3-Nitroaniline	9.89	138	234808	22.17	nq	87
53) 2,4-Dinitrophenol	10.00	184	273661	75.37	nq	# 77
54) Dibenzofuran	10.16	168	1714764	38.85	nq	97
55) 4-Nitrophenol	10.05	139	570190	74.16	nq	# 81
56) 2,4-Dinitrotoluene	10.13	165	466520	43.12	nq	# 74
57) Fluorene	10.49	166	1232937	36.02	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.27	232	355251	37.02	nq	91
59) Diethylphthalate	10.37	149	1551751	40.38	nq	97
60) 4-Chlorophenyl-phenylether	10.49	204	647218	45.50	nq	99
61) 4-Nitroaniline	10.51	138	368998	39.08	nq	86
62) Azobenzene	10.65	77	1686339	40.01	nq	93
64) 4,6-Dinitro-2-methylphenol	10.53	198	226994	37.19	nq	80
65) n-Nitrosodiphenylamine	10.60	169	1121242	35.07	nq	99
66) 4-Bromophenyl-phenylether	10.98	248	413874	38.45	nq	98
67) Hexachlorobenzene	11.04	284	413647	34.15	nq	# 82
68) Atrazine	11.13	200	378131	36.14	nq	97
69) Pentachlorophenol	11.23	266	473505	75.39	nq	99
70) Phenanthrene	11.45	178	1971677	37.70	nq	97
71) Anthracene	11.50	178	2068072	40.33	nq	97
72) Carbazole	11.65	167	1640726	32.73	nq	98
73) Di-n-butylphthalate	12.00	149	2291621	42.37	nq	# 96
74) Fluoranthene	12.64	202	1909383	34.94	nq	98
76) Benzidine	12.76	184	506340	19.23	nq	98
77) Pyrene	12.86	202	1969477	34.18	nq	97
79) Butylbenzylphthalate	13.49	149	979006	39.26	nq	# 92
80) Benzo(a)anthracene	14.05	228	1453993	33.72	nq	98
81) 3,3'-Dichlorobenzidine	14.02	252	348708	23.17	nq	# 99
82) Chrysene	14.09	228	1374349	33.17	nq	97
83) Bis(2-ethylhexyl)phthalate	14.05	149	1129742	35.86	nq	98
84) Di-n-octyl phthalate	14.66	149	1800658	33.86	nq	# 86
85) Indeno(1,2,3-cd)pyrene	16.92	276	1251803	38.12	nq	99
87) Benzo(b)fluoranthene	15.08	252	1509881m	43.13	nq	
88) Benzo(k)fluoranthene	15.12	252	886352m	30.96	nq	
89) Benzo(a)pyrene	15.44	252	1148260	38.67	nq	# 95
90) Dibenzo(a,h)anthracene	16.95	278	1036071	40.55	nq	# 95
91) Benzo(g,h,i)perylene	17.35	276	1056193	39.92	nq	98

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

