

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010815\
 Data File : BF076783.D
 Acq On : 8 Jan 2015 19:14
 Operator : IZ / TP
 Sample : PB81245BS
 Misc : S
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB81245BS

Manual Integrations
 APPROVED

TEJASKUMAR
 1/9/2015 5:22:52 PM

Quant Time: Jan 09 11:56:25 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121714.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 08 09:08:31 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	35917	20.00	ng	0.00
21) Naphthalene-d8	11.04	136	162073	20.00	ng	0.00
38) Acenaphthene-d10	15.18	164	86548	20.00	ng	-0.01
63) Phenanthrene-d10	17.90	188	154284	20.00	ng	0.00
75) Chrysene-d12	21.40	240	140307	20.00	ng	0.00
86) Perylene-d12	23.12	264	124924	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.21	112	277835	131.14	ng	0.00
7) Phenol-d6	7.52	99	359196	138.82	ng	0.00
23) Nitrobenzene-d5	9.43	82	220850	81.04	ng	0.00
41) 2,4,6-Tribromophenol	16.81	330	93299	127.55	ng	0.00
44) 2-Fluorobiphenyl	13.69	172	467298	84.02	ng	-0.01
78) Terphenyl-d14	20.12	244	493541	77.59	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.41	88	22869	25.25	ng	# 100
3) Pyridine	1.92	79	67268	29.51	ng	97
4) n-Nitrosodimethylamine	1.86	42	41096	37.26	ng	98
6) Aniline	7.42	93	66090	17.89	ng	# 83
8) 2-Chlorophenol	7.66	128	89547	36.91	ng	98
9) Benzaldehyde	7.14	77	49941	25.37	ng	97
10) Phenol	7.54	94	104142	33.16	ng	79
11) bis(2-Chloroethyl)ether	7.66	93	80997	35.83	ng	# 75
12) 1,3-Dichlorobenzene	7.98	146	92775	32.94	ng	# 96
13) 1,4-Dichlorobenzene	8.17	146	97316	34.11	ng	98
14) 1,2-Dichlorobenzene	8.48	146	91182	33.70	ng	96
15) Benzyl Alcohol	8.55	79	75358	33.63	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.90	45	101915	31.12	ng	# 31
17) 2-Methylphenol	8.89	107	70524	35.37	ng	# 93
18) Hexachloroethane	9.24	117	32991	32.36	ng	# 88
19) n-Nitroso-di-n-propylamine	9.18	70	64064	33.90	ng	99
20) 3+4-Methylphenols	9.28	107	94162	34.98	ng	97
22) Acetophenone	9.11	105	132194	34.03	ng	# 96
24) Nitrobenzene	9.46	77	98224	34.76	ng	98
25) Isophorone	10.06	82	172226	32.79	ng	# 92
26) 2-Nitrophenol	10.21	139	46687	33.83	ng	# 85
27) 2,4-Dimethylphenol	10.47	122	82015	36.72	ng	93
28) bis(2-Chloroethoxy)methane	10.69	93	103653	33.62	ng	98
29) 2,4-Dichlorophenol	10.80	162	75081	34.20	ng	92
30) 1,2,4-Trichlorobenzene	10.95	180	76328	32.25	ng	95
31) Naphthalene	11.09	128	261768	32.60	ng	98
32) Benzoic acid	10.79	122	42231m	33.12	ng	
33) 4-Chloroaniline	11.33	127	59117	18.04	ng	95
34) Hexachlorobutadiene	11.45	225	42994	31.25	ng	94
35) Caprolactam	12.15	113	23269m	36.76	ng	
36) 4-Chloro-3-methylphenol	12.61	107	85259	34.90	ng	91
37) 2-Methylnaphthalene	12.75	142	190289	33.97	ng	96
39) 1,2,4,5-Tetrachlorobenzene	13.15	216	74432	34.96	ng	# 81
40) Hexachlorocyclopentadiene	13.13	237	79356	65.71	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.49	196	55810	36.13	ng	92
43) 2,4,5-Trichlorophenol	13.58	196	53203	32.01	ng	96
45) 1,1'-Biphenyl	13.90	154	229289	35.66	ng	96
46) 2-Chloronaphthalene	13.88	162	181222	35.54	ng	97
47) 2-Nitroaniline	14.21	65	59857	38.58	ng	97
48) Acenaphthylene	14.84	152	294369	33.94	ng	98
49) Dimethylphthalate	14.77	163	220112	35.95	ng	99
50) 2,6-Dinitrotoluene	14.86	165	46187	36.82	ng	# 70
51) Acenaphthene	15.26	154	168070	34.25	ng	99
52) 3-Nitroaniline	15.21	138	37285	26.18	ng	93
53) 2,4-Dinitrophenol	15.49	184	25415	44.24	ng	88
54) Dibenzofuran	15.67	168	264670	37.29	ng	99
55) 4-Nitrophenol	15.77	139	71986	65.92	ng	# 74
56) 2,4-Dinitrotoluene	15.77	165	62990	37.68	ng	# 55
57) Fluorene	16.37	166	219226	36.77	ng	97
58) 2,3,4,6-Tetrachlorophenol	15.99	232	43933	35.55	ng	# 100
59) Diethylphthalate	16.35	149	229972	36.78	ng	98
60) 4-Chlorophenyl-phenylether	16.45	204	92556	35.87	ng	# 83
61) 4-Nitroaniline	16.49	138	49478	33.05	ng	96
62) Azobenzene	16.72	77	228976	37.16	ng	88
64) 4,6-Dinitro-2-methylphenol	16.56	198	23860	24.72	ng	# 79
65) n-Nitrosodiphenylamine	16.68	169	183932	33.94	ng	97
66) 4-Bromophenyl-phenylether	17.27	248	55231	36.75	ng	90
67) Hexachlorobenzene	17.29	284	55337	31.55	ng	# 86
68) Atrazine	17.63	200	67413	41.63	ng	97
69) Pentachlorophenol	17.65	266	42466	49.46	ng	96
70) Phenanthrene	17.93	178	314524	33.79	ng	99
71) Anthracene	18.00	178	317356	34.75	ng	99
72) Carbazole	18.29	167	298581	33.71	ng	100
73) Di-n-butylphthalate	18.86	149	366164	33.65	ng	99
74) Fluoranthene	19.57	202	326127	34.31	ng	96
76) Benzidine	19.80	184	176873	30.25	ng	99
77) Pyrene	19.85	202	338686	34.50	ng	98
79) Butylbenzylphthalate	20.77	149	161119	34.33	ng	# 83
80) Benzo(a)anthracene	21.39	228	287587	32.92	ng	99
81) 3,3'-Dichlorobenzidine	21.39	252	59111	20.24	ng	98
82) Chrysene	21.42	228	279952	35.01	ng	99
83) Bis(2-ethylhexyl)phthalate	21.51	149	228623	32.20	ng	97
84) Di-n-octyl phthalate	22.31	149	393829	32.49	ng	99
85) Indeno(1,2,3-cd)pyrene	24.33	276	279107m	31.81	ng	
87) Benzo(b)fluoranthene	22.68	252	277136m	33.63	ng	
88) Benzo(k)fluoranthene	22.71	252	249973	32.57	ng	# 96
89) Benzo(a)pyrene	23.07	252	257343	34.99	ng	# 97
90) Dibenzo(a,h)anthracene	24.36	278	230983	31.69	ng	# 95
91) Benzo(g,h,i)perylene	24.64	276	228263	32.19	ng	98

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

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