

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101015\
 Data File : BF082175.D
 Acq On : 10 Oct 2015 14:34
 Operator : UM/IZ
 Sample : PB86035BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB86035BS

Manual Integrations
 APPROVED

MMDadoda
 10/12/2015 2:31:43 PM

Quant Time: Oct 12 02:21:24 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 12 01:46:21 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	87174	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	350375	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	181011	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	342151	20.00	ng	0.00
75) Chrysene-d12	14.01	240	322863	20.00	ng	0.00
86) Perylene-d12	15.44	264	260217	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.44	112	529503	122.59	ng	0.01
7) Phenol-d6	6.48	99	700029	124.54	ng	0.00
23) Nitrobenzene-d5	7.41	82	450570	86.36	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	199968	117.80	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	939647	86.09	ng	0.00
78) Terphenyl-d14	12.96	244	1030340	84.57	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.49	88	66600	30.69	ng	93
3) Pyridine	3.21	79	173028	34.28	ng	96
4) n-Nitrosodimethylamine	3.18	42	80331	37.53	ng	96
6) Aniline	6.50	93	244796	33.78	ng	98
8) 2-Chlorophenol	6.63	128	213385	40.47	ng	99
9) Benzaldehyde	6.39	77	105923	36.90	ng	95
10) Phenol	6.49	94	281359	43.42	ng	99
11) bis(2-Chloroethyl)ether	6.58	93	228549	41.70	ng	97
12) 1,3-Dichlorobenzene	6.78	146	243313	39.62	ng	99
13) 1,4-Dichlorobenzene	6.86	146	255070	39.77	ng	99
14) 1,2-Dichlorobenzene	7.01	146	229582	37.70	ng	98
15) Benzyl Alcohol	6.98	79	164183	42.31	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.12	45	300674	39.40	ng	100
17) 2-Methylphenol	7.10	107	165844	39.77	ng	98
18) Hexachloroethane	7.35	117	87917	40.05	ng	96
19) n-Nitroso-di-n-propylamine	7.26	70	145839	36.95	ng	# 88
20) 3+4-Methylphenols	7.26	107	214456	43.16	ng	94
22) Acetophenone	7.26	105	287804	41.52	ng	# 97
24) Nitrobenzene	7.43	77	220943	39.12	ng	94
25) Isophorone	7.67	82	416224	39.53	ng	98
26) 2-Nitrophenol	7.75	139	122170	43.32	ng	93
27) 2,4-Dimethylphenol	7.78	122	192048	42.27	ng	99
28) bis(2-Chloroethoxy)methane	7.89	93	250343	40.86	ng	98
29) 2,4-Dichlorophenol	7.99	162	207843	45.76	ng	98
30) 1,2,4-Trichlorobenzene	8.07	180	211094	40.33	ng	98
31) Naphthalene	8.15	128	627656	41.33	ng	100
32) Benzoic acid	7.92	122	128409	42.11	ng	98
33) 4-Chloroaniline	8.21	127	172145	27.29	ng	95
34) Hexachlorobutadiene	8.26	225	132785	40.71	ng	99
35) Caprolactam	8.57	113	58728m	44.56	ng	
36) 4-Chloro-3-methylphenol	8.69	107	190213	42.83	ng	99
37) 2-Methylnaphthalene	8.85	142	437976	41.60	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.01	216	214982	39.93	ng	98
40) Hexachlorocyclopentadiene	8.99	237	233883	85.24	ng	99

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42) 2,4,6-Trichlorophenol	9.12	196	139068	38.89	ng	100
43) 2,4,5-Trichlorophenol	9.17	196	155001	40.01	ng	99
45) 1,1'-Biphenyl	9.30	154	513407	37.20	ng	98
46) 2-Chloronaphthalene	9.34	162	421293	38.77	ng	97
47) 2-Nitroaniline	9.43	65	119064	39.91	ng	97
48) Acenaphthylene	9.75	152	687347	40.61	ng	99
49) Dimethylphthalate	9.61	163	498948	39.14	ng	100
50) 2,6-Dinitrotoluene	9.68	165	112589	41.86	ng	# 86
51) Acenaphthene	9.92	154	408067	40.16	ng	99
52) 3-Nitroaniline	9.85	138	76834	25.29	ng	95
53) 2,4-Dinitrophenol	9.95	184	126444	83.80	ng	95
54) Dibenzofuran	10.09	168	602632	43.20	ng	99
55) 4-Nitrophenol	10.01	139	147806	72.98	ng	86
56) 2,4-Dinitrotoluene	10.08	165	147703	43.94	ng	98
57) Fluorene	10.43	166	485502	42.37	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.21	232	133601	42.21	ng	97
59) Diethylphthalate	10.31	149	473881	39.88	ng	99
60) 4-Chlorophenyl-phenylether	10.42	204	207018	39.44	ng	99
61) 4-Nitroaniline	10.47	138	118694	40.28	ng	95
62) Azobenzene	10.58	77	466932	40.15	ng	99
64) 4,6-Dinitro-2-methylphenol	10.49	198	84605	44.06	ng	90
65) n-Nitrosodiphenylamine	10.55	169	422344	41.23	ng	99
66) 4-Bromophenyl-phenylether	10.91	248	140118	40.92	ng	99
67) Hexachlorobenzene	10.98	284	152907	41.29	ng	98
68) Atrazine	11.07	200	122448	37.73	ng	98
69) Pentachlorophenol	11.18	266	178162	93.49	ng	98
70) Phenanthrene	11.39	178	690315	41.16	ng	100
71) Anthracene	11.45	178	760080	43.98	ng	99
72) Carbazole	11.60	167	626551	39.74	ng	100
73) Di-n-butylphthalate	11.93	149	804160	41.35	ng	99
74) Fluoranthene	12.58	202	773631	42.95	ng	99
76) Benzidine	12.71	184	358320	38.30	ng	98
77) Pyrene	12.81	202	776168	40.51	ng	100
79) Butylbenzylphthalate	13.43	149	333035	38.09	ng	98
80) Benzo(a)anthracene	14.00	228	670545	39.92	ng	99
81) 3,3'-Dichlorobenzidine	13.97	252	187576	29.42	ng	100
82) Chrysene	14.04	228	609563	34.44	ng	100
83) Bis(2-ethylhexyl)phthalate	13.99	149	496262	39.78	ng	100
84) Di-n-octyl phthalate	14.60	149	809233	37.77	ng	100
85) Indeno(1,2,3-cd)pyrene	16.86	276	548614	39.00	ng	100
87) Benzo(b)fluoranthene	15.03	252	717790m	45.34	ng	
88) Benzo(k)fluoranthene	15.06	252	481758m	36.20	ng	
89) Benzo(a)pyrene	15.38	252	565869	41.59	ng	100
90) Dibenzo(a,h)anthracene	16.87	278	462686	41.50	ng	100
91) Benzo(g,h,i)perylene	17.28	276	446014	41.18	ng	99

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

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