

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080315\
 Data File : BF080815.D
 Acq On : 3 Aug 2015 22:32
 Operator : TP
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

MMDadoda
 8/5/2015 2:28:03 PM

Quant Time: Aug 04 02:49:40 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 04 01:58:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	191830	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	697156	20.00	ng	0.00
38) Acenaphthene-d10	9.87	164	306130	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	512674	20.00	ng	0.01
75) Chrysene-d12	13.99	240	354924	20.00	ng	0.00
86) Perylene-d12	15.39	264	289176	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	1190083	103.43	ng	0.00
7) Phenol-d6	6.48	99	1686754	108.98	ng	0.01
23) Nitrobenzene-d5	7.41	82	1737633	123.00	ng	0.01
41) 2,4,6-Tribromophenol	10.66	330	321557	109.62	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	2209292	109.09	ng	0.00
78) Terphenyl-d14	12.93	244	1844041	103.54	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	2.49	88	360402	61.84	ng	99
3) Pyridine	3.20	79	981551	63.76	ng	97
4) n-Nitrosodimethylamine	3.16	42	539105	63.05	ng	99
6) Aniline	6.51	93	1270899	57.43	ng	98
8) 2-Chlorophenol	6.62	128	700635	54.96	ng	82
9) Benzaldehyde	6.38	77	477671	57.05	ng	96
10) Phenol	6.49	94	936909	51.83	ng	89
11) bis(2-Chloroethyl)ether	6.58	93	674788	52.06	ng	89
12) 1,3-Dichlorobenzene	6.78	146	790659m	55.33	ng	
13) 1,4-Dichlorobenzene	6.85	146	743338	52.20	ng	98
14) 1,2-Dichlorobenzene	7.01	146	769732m	58.44	ng	
15) Benzyl Alcohol	6.99	79	660991	62.89	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.13	45	1457460	56.87	ng	95
17) 2-Methylphenol	7.09	107	599673	53.36	ng	94
18) Hexachloroethane	7.36	117	299573	57.07	ng	# 71
19) n-Nitroso-di-n-propylamine	7.26	70	551374	55.07	ng	98
20) 3+4-Methylphenols	7.25	107	672388	51.51	ng	93
22) Acetophenone	7.25	105	886148	52.35	ng	# 81
24) Nitrobenzene	7.42	77	776424	53.85	ng	# 85
25) Isophorone	7.66	82	1449030	57.35	ng	96
26) 2-Nitrophenol	7.73	139	347855	56.44	ng	91
27) 2,4-Dimethylphenol	7.78	122	639901	54.49	ng	94
28) bis(2-Chloroethoxy)methane	7.88	93	915951	60.00	ng	98
29) 2,4-Dichlorophenol	7.98	162	580974	59.12	ng	93
30) 1,2,4-Trichlorobenzene	8.06	180	620109	57.81	ng	96
31) Naphthalene	8.14	128	1908109	55.71	ng	98
32) Benzoic acid	7.92	122	447622	62.93	ng	97
33) 4-Chloroaniline	8.19	127	741266	51.44	ng	94
34) Hexachlorobutadiene	8.26	225	362043	55.02	ng	97
35) Caprolactam	8.58	113	154313	56.69	ng	# 58
36) 4-Chloro-3-methylphenol	8.68	107	611998	58.48	ng	86
37) 2-Methylnaphthalene	8.83	142	1098578	51.18	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.00	216	605486	62.28	ng	99
40) Hexachlorocyclopentadiene	8.99	237	377831	63.53	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.10	196	402631	58.33	ng	99
43) 2,4,5-Trichlorophenol	9.15	196	413713	61.24	ng	# 87
45) 1,1'-Biphenyl	9.30	154	1356450	58.54	ng	96
46) 2-Chloronaphthalene	9.32	162	1101365	57.27	ng	95
47) 2-Nitroaniline	9.41	65	390185	54.93	ng	88
48) Acenaphthylene	9.73	152	1612902	53.79	ng	99
49) Dimethylphthalate	9.61	163	1263832	61.79	ng	99
50) 2,6-Dinitrotoluene	9.66	165	298317	67.73	ng	# 76
51) Acenaphthene	9.90	154	994701	55.05	ng	100
52) 3-Nitroaniline	9.82	138	297177	58.14	ng	84
53) 2,4-Dinitrophenol	9.93	184	140340	62.27	ng	# 58
54) Dibenzofuran	10.08	168	1262741	52.80	ng	95
55) 4-Nitrophenol	9.98	139	229792	62.89	ng	96
56) 2,4-Dinitrotoluene	10.06	165	369001	66.26	ng	# 78
57) Fluorene	10.42	166	950221	51.21	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.19	232	262289	54.00	ng	98
59) Diethylphthalate	10.30	149	1172945	60.05	ng	96
60) 4-Chlorophenyl-phenylether	10.42	204	540382	59.38	ng	# 78
61) 4-Nitroaniline	10.44	138	265693	56.16	ng	96
62) Azobenzene	10.57	77	1326664	58.63	ng	96
64) 4,6-Dinitro-2-methylphenol	10.46	198	181052	57.29	ng	# 32
65) n-Nitrosodiphenylamine	10.53	169	929084	51.77	ng	99
66) 4-Bromophenyl-phenylether	10.90	248	301056	53.44	ng	# 89
67) Hexachlorobenzene	10.97	284	381403	59.81	ng	# 84
68) Atrazine	11.06	200	174993	42.76	ng	99
69) Pentachlorophenol	11.15	266	199222	58.53	ng	97
70) Phenanthrene	11.38	178	1527336	57.28	ng	100
71) Anthracene	11.42	178	1417715	53.86	ng	100
72) Carbazole	11.58	167	1297506	56.93	ng	97
73) Di-n-butylphthalate	11.92	149	1515281	55.52	ng	100
74) Fluoranthene	12.56	202	1262027	49.42	ng	99
76) Benzidine	12.68	184	374139	27.95	ng	98
77) Pyrene	12.78	202	1418253	53.91	ng	100
79) Butylbenzylphthalate	13.41	149	633275	61.12	ng	91
80) Benzo(a)anthracene	13.97	228	1088345	53.71	ng	98
81) 3,3'-Dichlorobenzidine	13.94	252	408051	56.75	ng	# 97
82) Chrysene	14.01	228	1040483	52.83	ng	100
83) Bis(2-ethylhexyl)phthalate	13.97	149	750777	58.71	ng	100
84) Di-n-octyl phthalate	14.58	149	1181179	58.40	ng	98
85) Indeno(1,2,3-cd)pyrene	16.77	276	1180490	65.35	ng	100
87) Benzo(b)fluoranthene	14.99	252	881505m	52.07	ng	
88) Benzo(k)fluoranthene	15.03	252	973704m	66.01	ng	
89) Benzo(a)pyrene	15.33	252	848732	58.79	ng	98
90) Dibenzo(a,h)anthracene	16.80	278	945660	67.62	ng	98
91) Benzo(g,h,i)perylene	17.20	276	1015605	69.65	ng	98

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

