

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072415\
 Data File : BF080632.D
 Acq On : 24 Jul 2015 18:23
 Operator : TP/UM
 Sample : SSTDICC050
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

apatel
 7/27/2015 4:21:45 PM

Quant Time: Jul 24 23:13:03 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 24 22:49:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	50950	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	173156	20.00	ng	0.01
38) Acenaphthene-d10	9.97	164	90910	20.00	ng	0.00
63) Phenanthrene-d10	11.46	188	162233	20.00	ng	0.00
75) Chrysene-d12	14.23	240	142386	20.00	ng	0.07
86) Perylene-d12	15.85	264	128412	20.00	ng	0.11

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	287207	97.62	ng	0.00
7) Phenol-d6	6.54	99	355008	99.37	ng	0.00
23) Nitrobenzene-d5	7.49	82	364213	118.03	ng	0.00
41) 2,4,6-Tribromophenol	10.76	330	85698	116.28	ng	0.00
44) 2-Fluorobiphenyl	9.30	172	604320	98.14	ng	0.00
78) Terphenyl-d14	13.07	244	682329	101.47	ng	0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.45	88	60816	44.77	ng	98
3) Pyridine	3.22	79	167841	46.31	ng	93
4) n-Nitrosodimethylamine	3.14	42	102098	44.90	ng	99
6) Aniline	6.58	93	243827	47.38	ng	98
8) 2-Chlorophenol	6.70	128	154537	49.08	ng	97
9) Benzaldehyde	6.48	77	108188	41.97	ng	95
10) Phenol	6.56	94	221026	51.61	ng	97
11) bis(2-Chloroethyl)ether	6.66	93	160174	52.88	ng	99
12) 1,3-Dichlorobenzene	6.88	146	175509m	44.66	ng	
13) 1,4-Dichlorobenzene	6.94	146	178376	46.04	ng	97
14) 1,2-Dichlorobenzene	7.10	146	180334	50.01	ng	99
15) Benzyl Alcohol	7.07	79	143124	46.00	ng	# 86
16) 2,2'-oxybis(1-Chloropropan	7.21	45	264985	44.86	ng	99
17) 2-Methylphenol	7.17	107	118239	47.53	ng	98
18) Hexachloroethane	7.45	117	66504	48.81	ng	96
19) n-Nitroso-di-n-propylamine	7.33	70	115490	47.88	ng	97
20) 3+4-Methylphenols	7.33	107	168171	49.06	ng	# 83
22) Acetophenone	7.33	105	214995	51.35	ng	95
24) Nitrobenzene	7.52	77	165507	51.45	ng	# 76
25) Isophorone	7.76	82	297635	52.96	ng	# 86
26) 2-Nitrophenol	7.84	139	64121	54.61	ng	94
27) 2,4-Dimethylphenol	7.87	122	124645	49.21	ng	87
28) bis(2-Chloroethoxy)methane	7.96	93	170623	54.07	ng	98
29) 2,4-Dichlorophenol	8.06	162	117333	54.64	ng	98
30) 1,2,4-Trichlorobenzene	8.16	180	136329	49.81	ng	96
31) Naphthalene	8.24	128	418450	48.71	ng	100
32) Benzoic acid	7.93	122	72552	63.69	ng	98
33) 4-Chloroaniline	8.28	127	177674	52.10	ng	98
34) Hexachlorobutadiene	8.36	225	96545	58.85	ng	99
35) Caprolactam	8.62	113	33970	59.12	ng	# 71
36) 4-Chloro-3-methylphenol	8.76	107	129135	52.44	ng	# 74
37) 2-Methylnaphthalene	8.93	142	304939	60.94	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.09	216	128772	43.02	ng	100
40) Hexachlorocyclopentadiene	9.09	237	87047	54.30	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.21	196	96808m	55.67	ng	
43) 2,4,5-Trichlorophenol	9.24	196	102622m	58.47	ng	
45) 1,1'-Biphenyl	9.39	154	317633	44.00	ng	98
46) 2-Chloronaphthalene	9.42	162	272835	46.58	ng	# 85
47) 2-Nitroaniline	9.50	65	82103	45.80	ng	96
48) Acenaphthylene	9.84	152	454702	51.21	ng	100
49) Dimethylphthalate	9.69	163	311532	47.32	ng	98
50) 2,6-Dinitrotoluene	9.76	165	63560	52.05	ng	# 41
51) Acenaphthene	10.01	154	281346	52.70	ng	99
52) 3-Nitroaniline	9.92	138	74010	53.66	ng	95
53) 2,4-Dinitrophenol	10.02	184	27254	74.61	ng	# 84
54) Dibenzofuran	10.18	168	389948	50.91	ng	98
55) 4-Nitrophenol	10.06	139	59015	59.12	ng	88
56) 2,4-Dinitrotoluene	10.16	165	82240	53.60	ng	# 43
57) Fluorene	10.52	166	309020	50.49	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.29	232	81379	59.79	ng	100
59) Diethylphthalate	10.40	149	346076	54.71	ng	99
60) 4-Chlorophenyl-phenylether	10.52	204	136911	49.63	ng	98
61) 4-Nitroaniline	10.52	138	67572	49.09	ng	98
62) Azobenzene	10.67	77	303728	47.15	ng	98
64) 4,6-Dinitro-2-methylphenol	10.56	198	42917	74.30	ng	84
65) n-Nitrosodiphenylamine	10.64	169	255473	48.83	ng	100
66) 4-Bromophenyl-phenylether	11.00	248	76832	46.93	ng	# 85
67) Hexachlorobenzene	11.07	284	93203	49.89	ng	# 87
68) Atrazine	11.16	200	76994	53.86	ng	90
69) Pentachlorophenol	11.26	266	49694	56.66	ng	96
70) Phenanthrene	11.48	178	438370	48.40	ng	98
71) Anthracene	11.54	178	445428	52.02	ng	99
72) Carbazole	11.69	167	399501	53.61	ng	99
73) Di-n-butylphthalate	12.03	149	490529	56.63	ng	99
74) Fluoranthene	12.68	202	450226	55.32	ng	96
76) Benzidine	12.81	184	246657	57.31	ng	98
77) Pyrene	12.92	202	453558	48.26	ng	97
79) Butylbenzylphthalate	13.59	149	208513	63.98	ng	94
80) Benzo(a)anthracene	14.20	228	428897	53.53	ng	99
81) 3,3'-Dichlorobenzidine	14.17	252	142743	61.40	ng	97
82) Chrysene	14.25	228	408291	51.45	ng	99
83) Bis(2-ethylhexyl)phthalate	14.21	149	301894	62.27	ng	# 98
84) Di-n-octyl phthalate	14.96	149	499857	77.64	ng	99
85) Indeno(1,2,3-cd)pyrene	17.35	276	478071	73.92	ng	100
87) Benzo(b)fluoranthene	15.41	252	377381	49.16	ng	98
88) Benzo(k)fluoranthene	15.45	252	385584m	50.69	ng	
89) Benzo(a)pyrene	15.79	252	365886	53.52	ng	# 95
90) Dibenzo(a,h)anthracene	17.37	278	392048	59.16	ng	99
91) Benzo(g,h,i)perylene	17.79	276	398491	59.32	ng	97

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

