

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121115\
 Data File : BF083702.D
 Acq On : 11 Dec 2015 17:10
 Operator : UM/IZ
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

MMdadoda
 12/14/2015 6:22:22 PM

Quant Time: Dec 11 18:20:22 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 11 17:53:49 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	128908	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	524222	20.00	ng	0.00
38) Acenaphthene-d10	9.96	164	230021	20.00	ng	0.00
63) Phenanthrene-d10	11.45	188	443627	20.00	ng	0.00
75) Chrysene-d12	14.08	240	303800	20.00	ng	0.00
86) Perylene-d12	15.51	264	249041	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.53	112	888014	104.48	ng	0.00
7) Phenol-d6	6.56	99	1068411	94.08	ng	0.01
23) Nitrobenzene-d5	7.49	82	894982	79.66	ng	0.01
41) 2,4,6-Tribromophenol	10.75	330	254934	124.40	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	1413417	86.48	ng	0.00
78) Terphenyl-d14	13.03	244	1312365	101.64	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.56	88	207142	49.33	ng	# 87
3) Pyridine	3.30	79	557756	54.68	ng	89
4) n-Nitrosodimethylamine	3.24	42	226197	42.12	ng	# 75
6) Aniline	6.59	93	725304	51.83	ng	# 82
8) 2-Chlorophenol	6.70	128	459143	49.40	ng	99
9) Benzaldehyde	6.46	77	261913	43.41	ng	97
10) Phenol	6.57	94	640276	46.55	ng	# 71
11) bis(2-Chloroethyl)ether	6.66	93	418649	41.37	ng	95
12) 1,3-Dichlorobenzene	6.86	146	491592m	47.46	ng	
13) 1,4-Dichlorobenzene	6.94	146	520245	48.95	ng	99
14) 1,2-Dichlorobenzene	7.10	146	466646	47.31	ng	94
15) Benzyl Alcohol	7.06	79	370278	45.83	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.21	45	584473	31.23	ng	67
17) 2-Methylphenol	7.17	107	362976	48.00	ng	# 86
18) Hexachloroethane	7.45	117	179959	47.95	ng	93
19) n-Nitroso-di-n-propylamine	7.34	70	314169	39.63	ng	91
20) 3+4-Methylphenols	7.33	107	468490	47.05	ng	# 60
22) Acetophenone	7.33	105	562670	38.26	ng	# 92
24) Nitrobenzene	7.50	77	422653	34.04	ng	94
25) Isophorone	7.76	82	856660	39.29	ng	# 88
26) 2-Nitrophenol	7.82	139	202430	41.06	ng	# 85
27) 2,4-Dimethylphenol	7.86	122	403763	46.81	ng	99
28) bis(2-Chloroethoxy)methane	7.96	93	521847	43.10	ng	# 96
29) 2,4-Dichlorophenol	8.06	162	371272	47.33	ng	99
30) 1,2,4-Trichlorobenzene	8.16	180	403863	46.61	ng	98
31) Naphthalene	8.24	128	1307381	47.09	ng	98
32) Benzoic acid	7.97	122	255810	52.66	ng	94
33) 4-Chloroaniline	8.28	127	532781	52.62	ng	# 83
34) Hexachlorobutadiene	8.36	225	222195	43.89	ng	97
35) Caprolactam	8.65	113	127548	54.18	ng	# 62
36) 4-Chloro-3-methylphenol	8.75	107	383088	43.47	ng	95
37) 2-Methylnaphthalene	8.92	142	755090	43.30	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.09	216	378485	50.12	ng	# 100
40) Hexachlorocyclopentadiene	9.08	237	177880	173.80	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.20	196	255655	56.90	nq	99
43) 2,4,5-Trichlorophenol	9.23	196	232343	52.23	nq #	91
45) 1,1'-Biphenyl	9.39	154	1054668	52.75	nq	96
46) 2-Chloronaphthalene	9.41	162	795000	52.67	nq	96
47) 2-Nitroaniline	9.49	65	194653	35.76	nq	93
48) Acenaphthylene	9.82	152	1169964	47.01	nq	99
49) Dimethylphthalate	9.69	163	853401	46.61	nq	100
50) 2,6-Dinitrotoluene	9.74	165	201584	53.32	nq #	74
51) Acenaphthene	10.00	154	695962	48.84	nq	99
52) 3-Nitroaniline	9.90	138	193454	48.15	nq	87
53) 2,4-Dinitrophenol	10.01	184	55404	34.38	nq #	1
54) Dibenzofuran	10.17	168	941820	47.66	nq	91
55) 4-Nitrophenol	10.05	139	160850	90.93	nq	93
56) 2,4-Dinitrotoluene	10.14	165	266173	53.09	nq #	67
57) Fluorene	10.51	166	733200	42.51	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.28	232	188171	57.13	nq #	100
59) Diethylphthalate	10.38	149	825927	46.39	nq	99
60) 4-Chlorophenyl-phenylether	10.51	204	367798	44.80	nq #	85
61) 4-Nitroaniline	10.52	138	215568	59.06	nq	97
62) Azobenzene	10.66	77	734069	37.47	nq	97
64) 4,6-Dinitro-2-methylphenol	10.54	198	94396	35.44	nq	92
65) n-Nitrosodiphenylamine	10.62	169	720810	49.23	nq	98
66) 4-Bromophenyl-phenylether	10.99	248	221570	44.78	nq #	84
67) Hexachlorobenzene	11.06	284	245419	46.75	nq	98
68) Atrazine	11.14	200	213774	49.07	nq	96
69) Pentachlorophenol	11.24	266	154299	134.24	nq	99
70) Phenanthrene	11.47	178	1212263	47.56	nq	98
71) Anthracene	11.52	178	1155327	43.76	nq	98
72) Carbazole	11.66	167	1050020	46.37	nq #	95
73) Di-n-butylphthalate	12.01	149	1206673	42.37	nq	99
74) Fluoranthene	12.65	202	1226631	47.04	nq	94
76) Benzidine	12.77	184	616124	65.20	nq	99
77) Pyrene	12.88	202	1216303	55.61	nq	98
79) Butylbenzylphthalate	13.51	149	436445	44.96	nq #	80
80) Benzo(a)anthracene	14.07	228	903277	50.80	nq	99
81) 3,3'-Dichlorobenzidine	14.02	252	288831	48.08	nq #	97
82) Chrysene	14.10	228	858938	48.19	nq	100
83) Bis(2-ethylhexyl)phthalate	14.07	149	450717	33.45	nq	98
84) Di-n-octyl phthalate	14.67	149	670256	32.55	nq #	100
85) Indeno(1,2,3-cd)pyrene	16.95	276	921182	76.56	nq #	100
87) Benzo(b)fluoranthene	15.09	252	703649m	49.62	nq	
88) Benzo(k)fluoranthene	15.13	252	760314m	48.66	nq	
89) Benzo(a)pyrene	15.45	252	730337	52.78	nq	97
90) Dibenzo(a,h)anthracene	16.97	278	765224	72.32	nq	97
91) Benzo(g,h,i)perylene	17.38	276	791474	76.72	nq	97

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

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