

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100716\
 Data File : BF090479.D
 Acq On : 8 Oct 2016 10:34
 Operator : UM/SJ
 Sample : H5134-04MS
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MJSB-16(6.5-8.5)MS

Manual Integrations
 APPROVED

sohil
 10/10/2016 6:51:05 PM

Quant Time: Oct 10 03:51:20 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF100516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 05 14:57:00 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	219651	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	560380	20.00	ng	0.01
38) Acenaphthene-d10	10.10	164	180018	20.00	ng	0.05
63) Phenanthrene-d10	11.57	188	401533	20.00	ng	0.02
75) Chrysene-d12	14.20	240	470997	20.00	ng	0.00
86) Perylene-d12	15.70	264	443658	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	1313229	89.23	ng	0.00
7) Phenol-d6	6.63	99	1907375	98.95	ng	0.01
23) Nitrobenzene-d5	7.56	82	952784	82.70	ng	-0.01
41) 2,4,6-Tribromophenol	10.89	330	216093	147.63	ng	0.03
44) 2-Fluorobiphenyl	9.40	172	1054219	97.57	ng	0.02
78) Terphenyl-d14	13.14	244	1281209	61.16	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.69	88	176593	24.34	ng	# 85
3) Pyridine	3.47	79	566847	28.02	ng	# 76
4) n-Nitrosodimethylamine	3.40	42	216425	25.93	ng	# 46
6) Aniline	6.66	93	346040	13.00	ng	97
8) 2-Chlorophenol	6.78	128	509250	31.44	ng	95
9) Benzaldehyde	6.53	77	213596	21.92	ng	92
10) Phenol	6.65	94	789287	35.22	ng	94
11) bis(2-Chloroethyl)ether	6.73	93	519825	30.29	ng	96
12) 1,3-Dichlorobenzene	6.94	146	521124	32.28	ng	98
13) 1,4-Dichlorobenzene	7.01	146	494673	30.17	ng	98
14) 1,2-Dichlorobenzene	7.17	146	467645m	29.60	ng	
15) Benzyl Alcohol	7.14	79	537367	36.67	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.27	45	477500	16.41	ng	# 46
17) 2-Methylphenol	7.25	107	402535	30.22	ng	# 81
18) Hexachloroethane	7.58	117	2340034	361.98	ng	# 23
19) n-Nitroso-di-n-propylamine	7.41	70	394290m	27.90	ng	
20) 3+4-Methylphenols	7.41	107	493211	28.61	ng	# 72
22) Acetophenone	7.41	105	630483	47.50	ng	# 89
24) Nitrobenzene	7.58	77	541627m	47.36	ng	
25) Isophorone	7.82	82	1163893	55.26	ng	# 44
26) 2-Nitrophenol	7.90	139	146385	29.74	ng	# 1
27) 2,4-Dimethylphenol	7.95	122	395167	41.28	ng	89
28) bis(2-Chloroethoxy)methane	8.03	93	501611	40.27	ng	# 15
29) 2,4-Dichlorophenol	8.15	162	445394	61.19	ng	85
30) 1,2,4-Trichlorobenzene	8.23	180	401251	53.49	ng	96
31) Naphthalene	8.33	128	1752962	64.84	ng	# 77
33) 4-Chloroaniline	8.36	127	659865m	59.05	ng	
34) Hexachlorobutadiene	8.44	225	209220	49.92	ng	99
35) Caprolactam	8.75	113	555602	227.04	ng	# 65
36) 4-Chloro-3-methylphenol	8.86	107	311969	34.23	ng	93
37) 2-Methylnaphthalene	9.05	142	814691	49.78	ng	93
39) 1,2,4,5-Tetrachlorobenzene	9.22	216	232891	49.07	ng	# 73
40) Hexachlorocyclopentadiene	9.19	237	26412	10.15	ng	94
42) 2,4,6-Trichlorophenol	9.32	196	184451m	56.32	ng	

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	9.35	196	72038m	23.26	ng	
45) 1,1'-Biphenyl	9.50	154	493351	36.98	ng	98
46) 2-Chloronaphthalene	9.53	162	396578	40.24	ng	# 84
47) 2-Nitroaniline	9.61	65	160356	41.23	ng	92
48) Acenaphthylene	9.96	152	734773	45.18	ng	# 71
49) Dimethylphthalate	9.79	163	618187	53.62	ng	# 91
50) 2,6-Dinitrotoluene	9.87	165	157360	61.48	ng	# 90
51) Acenaphthene	10.12	154	594258	58.61	ng	# 80
52) 3-Nitroaniline	10.02	138	129760m	43.40	ng	
54) Dibenzofuran	10.30	168	658717	50.19	ng	# 43
55) 4-Nitrophenol	10.18	139	306737	117.25	ng	# 51
57) Fluorene	10.65	166	849282	77.00	ng	# 87
58) 2,3,4,6-Tetrachlorophenol	10.43	232	84131m	36.38	ng	
59) Diethylphthalate	10.50	149	347731m	29.15	ng	
60) 4-Chlorophenyl-phenylether	10.63	204	178348	35.56	ng	# 66
61) 4-Nitroaniline	10.63	138	91731	30.45	ng	92
62) Azobenzene	10.79	77	469221	34.06	ng	# 26
66) 4-Bromophenyl-phenylether	11.11	248	158957	40.39	ng	# 59
67) Hexachlorobenzene	11.23	284	105604	24.09	ng	# 75
68) Atrazine	11.27	200	181444	54.69	ng	77
69) Pentachlorophenol	11.39	266	147145	56.40	ng	99
70) Phenanthrene	11.61	178	2898058	140.95	ng	95
71) Anthracene	11.65	178	1193038m	56.82	ng	
72) Carbazole	11.80	167	656267	33.64	ng	# 87
73) Di-n-butylphthalate	12.12	149	876160	36.91	ng	98
74) Fluoranthene	12.78	202	2132288	105.73	ng	95
76) Benzidine	12.89	184	259733	20.97	ng	# 85
77) Pyrene	13.01	202	2458514	78.28	ng	98
79) Butylbenzylphthalate	13.61	149	499178	30.98	ng	98
80) Benzo(a)anthracene	14.19	228	2082048	78.79	ng	99
81) 3,3'-Dichlorobenzidine	14.14	252	274347	28.63	ng	# 98
82) Chrysene	14.22	228	1841401	75.71	ng	99
83) Bis(2-ethylhexyl)phthalate	14.17	149	833394	36.38	ng	98
84) Di-n-octyl phthalate	14.78	149	1219258	35.90	ng	# 91
85) Indeno(1,2,3-cd)pyrene	17.23	276	1361743	78.53	ng	97
87) Benzo(b)fluoranthene	15.26	252	2701712m	95.60	ng	
88) Benzo(k)fluoranthene	15.29	252	1145503m	43.15	ng	
89) Benzo(a)pyrene	15.64	252	2005355	82.33	ng	97
90) Dibenzo(a,h)anthracene	17.24	278	766046	36.97	ng	98
91) Benzo(g,h,i)perylene	17.69	276	1107604	55.70	ng	99

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

