

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112017\
 Data File : BF100741.D
 Acq On : 20 Nov 2017 20:20
 Operator : SJ/JU
 Sample : I6521-01MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-(0-12)MSD

Manual Integrations
 APPROVED

SOHIL
 11/21/2017 2:15:33 PM

Quant Time: Nov 21 04:41:13 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 17 15:22:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	95809	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	380681	20.00	ng	-0.01
38) Acenaphthene-d10	9.91	164	179814	20.00	ng	-0.01
63) Phenanthrene-d10	11.40	188	348378	20.00	ng	0.00
75) Chrysene-d12	14.04	240	227263	20.00	ng	0.00
86) Perylene-d12	15.50	264	191547	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	700749	117.24	ng	0.00
7) Phenol-d6	6.50	99	882104	119.68	ng	0.00
23) Nitrobenzene-d5	7.44	82	542065	94.14	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	254110	138.68	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1041617	88.21	ng	0.00
78) Terphenyl-d14	12.98	244	997125	100.81	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.71	88	90939	31.221	ng	97
3) Pyridine	3.48	79	247309	31.121	ng	95
4) n-Nitrosodimethylamine	3.43	42	134666	34.904	ng	98
6) Aniline	6.55	93	283271	27.205	ng	99
8) 2-Chlorophenol	6.66	128	289808	45.834	ng	94
9) Benzaldehyde	6.42	77	101551	19.294	ng	96
10) Phenol	6.52	94	375995	48.595	ng	98
11) bis(2-Chloroethyl)ether	6.61	93	264014	40.624	ng	98
12) 1,3-Dichlorobenzene	6.82	146	310662	42.615	ng	97
13) 1,4-Dichlorobenzene	6.89	146	318182	43.124	ng	96
14) 1,2-Dichlorobenzene	7.05	146	300672	44.075	ng	96
15) Benzyl Alcohol	7.02	79	254084	44.172	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.15	45	464743	44.711	ng	98
17) 2-Methylphenol	7.12	107	252859	46.796	ng	99
18) Hexachloroethane	7.39	117	104397	42.914	ng	91
19) n-Nitroso-di-n-propylamine	7.29	70	201243	39.372	ng	99
20) 3+4-Methylphenols	7.27	107	305248	44.642	ng	# 89
22) Acetophenone	7.29	105	386610	41.648	ng	# 98
24) Nitrobenzene	7.46	77	300429	50.693	ng	98
25) Isophorone	7.70	82	555199	45.884	ng	99
26) 2-Nitrophenol	7.77	139	165492	58.263	ng	95
27) 2,4-Dimethylphenol	7.80	122	276931	51.055	ng	99
28) bis(2-Chloroethoxy)methane	7.90	93	353288	44.636	ng	98
29) 2,4-Dichlorophenol	8.01	162	269404	52.027	ng	97
30) 1,2,4-Trichlorobenzene	8.09	180	263429	47.031	ng	98
31) Naphthalene	8.18	128	835543	45.569	ng	99
32) Benzoic acid	7.87	122	27196m	14.640	ng	
33) 4-Chloroaniline	8.22	127	118599	15.539	ng	97
34) Hexachlorobutadiene	8.29	225	149935	45.021	ng	97
35) Caprolactam	8.60	113	85451m	48.086	ng	
36) 4-Chloro-3-methylphenol	8.70	107	274658	49.387	ng	99
37) 2-Methylnaphthalene	8.87	142	572124	46.999	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.03	216	266393	44.670	ng	97
40) Hexachlorocyclopentadiene	9.02	237	259002	92.279	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.15	196	201240	53.244	nq	98
43) 2,4,5-Trichlorophenol	9.19	196	209005	53.373	nq	94
45) 1,1'-Biphenyl	9.33	154	680540	43.759	nq	99
46) 2-Chloronaphthalene	9.36	162	544891	48.295	nq	97
47) 2-Nitroaniline	9.46	65	172453	51.513	nq	95
48) Acenaphthylene	9.77	152	827034	44.232	nq	100
49) Dimethylphthalate	9.64	163	711590	51.831	nq	98
50) 2,6-Dinitrotoluene	9.70	165	148708	54.721	nq	88
51) Acenaphthene	9.95	154	496310	43.645	nq	99
52) 3-Nitroaniline	9.86	138	106315	33.130	nq	96
53) 2,4-Dinitrophenol	9.97	184	99738	82.480	nq #	12
54) Dibenzofuran	10.12	168	739682	46.410	nq	96
55) 4-Nitrophenol	10.03	139	246977	94.783	nq	94
56) 2,4-Dinitrotoluene	10.10	165	193853	56.544	nq #	90
57) Fluorene	10.46	166	584136	50.030	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.23	232	169800	52.907	nq #	85
59) Diethylphthalate	10.33	149	608803	44.312	nq	96
60) 4-Chlorophenyl-phenylether	10.45	204	268216	46.828	nq	92
61) 4-Nitroaniline	10.49	138	159228	52.328	nq	88
62) Azobenzene	10.61	77	544288	42.063	nq	95
64) 4,6-Dinitro-2-methylphenol	10.51	198	99701	55.130	nq	85
65) n-Nitrosodiphenylamine	10.57	169	536477	44.288	nq	96
66) 4-Bromophenyl-phenylether	10.94	248	176414	47.238	nq #	90
67) Hexachlorobenzene	11.01	284	183684	47.027	nq #	85
68) Atrazine	11.10	200	173531	47.325	nq	95
69) Pentachlorophenol	11.20	266	219585	78.402	nq	98
70) Phenanthrene	11.42	178	860649	46.921	nq	98
71) Anthracene	11.47	178	911944	49.004	nq	99
72) Carbazole	11.63	167	801416	46.673	nq	99
73) Di-n-butylphthalate	11.95	149	932343	46.398	nq	99
74) Fluoranthene	12.61	202	1044386	53.724	nq	97
76) Benzidine	12.73	184	211425	23.230	nq	98
77) Pyrene	12.85	202	1024918	63.266	nq	99
79) Butylbenzylphthalate	13.45	149	357715	54.144	nq	94
80) Benzo(a)anthracene	14.03	228	743421	54.321	nq	99
81) 3,3'-Dichlorobenzidine	13.99	252	126203	25.738	nq #	96
82) Chrysene	14.06	228	661259	49.896	nq	100
83) Bis(2-ethylhexyl)phthalate	14.00	149	471929	54.311	nq	99
84) Di-n-octyl phthalate	14.62	149	670619	44.925	nq	99
85) Indeno(1,2,3-cd)pyrene	16.99	276	646642	60.324	nq	95
87) Benzo(b)fluoranthene	15.07	252	592369	52.200	nq	99
88) Benzo(k)fluoranthene	15.10	252	486541	45.376	nq	98
89) Benzo(a)pyrene	15.44	252	534989	53.177	nq	100
90) Dibenzo(a,h)anthracene	17.01	278	503630	61.212	nq	97
91) Benzo(g,h,i)perylene	17.44	276	573107	71.159	nq	94

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

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