

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081217\
 Data File : BF097668.D
 Acq On : 13 Aug 2017 11:41
 Operator : SJ/JU
 Sample : I4736-08MSD
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-10-COMPMSD

Manual Integrations
 APPROVED

sohil
 8/14/2017 7:08:29 PM

Quant Time: Aug 14 08:05:14 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 11 15:55:57 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	118759	20.00	ng	0.00
21) Naphthalene-d8	9.53	136	481067	20.00	ng	0.00
38) Acenaphthene-d10	12.35	164	194403	20.00	ng	-0.02
63) Phenanthrene-d10	14.75	188	266443	20.00	ng	-0.01
75) Chrysene-d12	18.45	240	193083	20.00	ng	-0.01
86) Perylene-d12	20.12	264	166456	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.43	112	785084	105.08	ng	0.03
7) Phenol-d6	7.08	99	795761	85.86	ng	0.02
23) Nitrobenzene-d5	8.42	82	537117	70.01	ng	0.00
41) 2,4,6-Tribromophenol	13.65	330	103444	70.00	ng	-0.01
44) 2-Fluorobiphenyl	11.30	172	653604	50.74	ng	-0.02
78) Terphenyl-d14	17.10	244	384070	41.17	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.86	88	104956	32.370	ng	# 73
3) Pyridine	2.45	79	285118	30.149	ng	# 62
4) n-Nitrosodimethylamine	2.41	42	177311	32.314	ng	# 78
6) Aniline	7.00	93	228757	17.173	ng	99
8) 2-Chlorophenol	7.20	128	249369	29.115	ng	95
9) Benzaldehyde	6.82	77	102051	16.698	ng	93
10) Phenol	7.10	94	335353	33.046	ng	93
11) bis(2-Chloroethyl)ether	7.15	93	250538	30.780	ng	90
12) 1,3-Dichlorobenzene	7.40	146	251077	26.176	ng	98
13) 1,4-Dichlorobenzene	7.53	146	271164	27.774	ng	93
14) 1,2-Dichlorobenzene	7.76	146	270501	30.360	ng	98
15) Benzyl Alcohol	7.80	79	236030	36.854	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.00	45	576069	31.523	ng	91
17) 2-Methylphenol	8.05	107	217271	32.731	ng	99
18) Hexachloroethane	8.27	117	84265	25.227	ng	# 78
19) n-Nitroso-di-n-propylamine	8.23	70	200027	31.028	ng	# 81
20) 3+4-Methylphenols	8.31	107	273269	33.614	ng	# 59
22) Acetophenone	8.19	105	373656	32.878	ng	# 99
24) Nitrobenzene	8.45	77	304905	36.469	ng	98
25) Isophorone	8.85	82	487552	34.498	ng	98
26) 2-Nitrophenol	8.96	139	132617	31.360	ng	# 84
27) 2,4-Dimethylphenol	9.11	122	240374	34.086	ng	96
28) bis(2-Chloroethoxy)methane	9.22	93	320645	31.400	ng	99
29) 2,4-Dichlorophenol	9.40	162	194330	31.252	ng	95
30) 1,2,4-Trichlorobenzene	9.46	180	181717	29.476	ng	97
31) Naphthalene	9.56	128	705819	29.292	ng	100
33) 4-Chloroaniline	9.72	127	168145	16.845	ng	97
34) Hexachlorobutadiene	9.78	225	93492	27.762	ng	98
35) Caprolactam	10.30	113	61000	30.447	ng	# 41
36) 4-Chloro-3-methylphenol	10.59	107	216410	30.975	ng	90
37) 2-Methylnaphthalene	10.69	142	461709	30.453	ng	98
39) 1,2,4,5-Tetrachlorobenzene	10.97	216	153763	27.933	ng	99
40) Hexachlorocyclopentadiene	10.94	237	2973	14.065	ng	95
42) 2,4,6-Trichlorophenol	11.20	196	98154	27.063	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	11.30	196	89127	26.018	nq	99
45) 1,1'-Biphenyl	11.46	154	445455	26.708	nq	97
46) 2-Chloronaphthalene	11.47	162	332949	25.999	nq	# 92
47) 2-Nitroaniline	11.69	65	161397	32.173	nq	95
48) Acenaphthylene	12.12	152	589649	28.964	nq	98
49) Dimethylphthalate	12.00	163	589403	39.340	nq	99
50) 2,6-Dinitrotoluene	12.10	165	92629	29.939	nq	# 69
51) Acenaphthene	12.41	154	338915	27.779	nq	99
52) 3-Nitroaniline	12.37	138	92900	22.933	nq	95
54) Dibenzofuran	12.70	168	510263	29.875	nq	99
55) 4-Nitrophenol	12.80	139	58456	37.890	nq	# 58
56) 2,4-Dinitrotoluene	12.77	165	120280	29.127	nq	# 73
57) Fluorene	13.24	166	387155	27.562	nq	98
58) 2,3,4,6-Tetrachlorophenol	12.95	232	59973	25.492	nq	96
59) Diethylphthalate	13.15	149	448858	31.720	nq	97
60) 4-Chlorophenyl-phenylether	13.27	204	166867	27.827	nq	95
61) 4-Nitroaniline	13.37	138	95542	24.725	nq	92
62) Azobenzene	13.53	77	441314	31.347	nq	92
64) 4,6-Dinitro-2-methylphenol	13.46	198	6105	3.756	nq	# 1
65) n-Nitrosodiphenylamine	13.49	169	329810	33.496	nq	97
66) 4-Bromophenyl-phenylether	14.06	248	86045	31.128	nq	92
67) Hexachlorobenzene	14.14	284	84824	28.170	nq	# 86
68) Atrazine	14.40	200	93529	37.858	nq	95
69) Pentachlorophenol	14.53	266	5920m	16.038	nq	
70) Phenanthrene	14.79	178	485382	31.629	nq	99
71) Anthracene	14.87	178	461297	29.379	nq	97
72) Carbazole	15.17	167	441124	29.512	nq	99
73) Di-n-butylphthalate	15.74	149	556772	31.823	nq	99
74) Fluoranthene	16.55	202	402745	29.044	nq	99
76) Benzidine	16.78	184	229192	36.018	nq	# 93
77) Pyrene	16.86	202	412489	25.834	nq	99
79) Butylbenzylphthalate	17.77	149	209461	31.379	nq	# 73
80) Benzo(a)anthracene	18.44	228	342503	30.372	nq	99
81) 3,3'-Dichlorobenzidine	18.42	252	112485	28.921	nq	98
82) Chrysene	18.48	228	336625	30.012	nq	100
83) Bis(2-ethylhexyl)phthalate	18.52	149	314006	33.074	nq	98
84) Di-n-octyl phthalate	19.29	149	525050	31.667	nq	# 91
85) Indeno(1,2,3-cd)pyrene	21.33	276	253144	29.017	nq	95
87) Benzo(b)fluoranthene	19.71	252	325157m	32.531	nq	
88) Benzo(k)fluoranthene	19.74	252	277161m	27.532	nq	
89) Benzo(a)pyrene	20.06	252	281576	30.692	nq	97
90) Dibenzo(a,h)anthracene	21.33	278	208352	31.719	nq	# 93
91) Benzo(g,h,i)perylene	21.68	276	209330	29.049	nq	96

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

