

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070916\
 Data File : BF088887.D
 Acq On : 9 Jul 2016 22:58
 Operator : UM/SJ
 Sample : H3813-31MSD 25X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-WW001-01MSD

Manual Integrations

sohil
 7/11/2016 8:07:45 PM

Quant Time: Jul 11 15:55:29 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 08 18:51:20 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	89900m	20.00	ng	-0.01
21) Naphthalene-d8	8.01	136	363092	20.00	ng	-0.01
38) Acenaphthene-d10	9.76	164	173677	20.00	ng	-0.01
63) Phenanthrene-d10	11.23	188	342516	20.00	ng	-0.01
75) Chrysene-d12	13.85	240	184116	20.00	ng	-0.02
86) Perylene-d12	15.23	264	172079	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	13403	2.23	ng	-0.02
7) Phenol-d6	6.34	99	11719	1.51	ng	-0.03
23) Nitrobenzene-d5	7.28	82	22032	3.18	ng	-0.02
41) 2,4,6-Tribromophenol	10.54	330	8601	5.33	ng	-0.02
44) 2-Fluorobiphenyl	9.08	172	40960	3.73	ng	-0.01
78) Terphenyl-d14	12.81	244	35681	4.32	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.24	88	1536	0.52	ng	# 71
3) Pyridine	2.99	79	566	0.07	ng	# 23
4) n-Nitrosodimethylamine	2.87	42	210	0.06	ng	# 86
6) Aniline	6.38	93	2815	0.25	ng	99
8) 2-Chlorophenol	6.50	128	3980	0.62	ng	86
9) Benzaldehyde	6.26	77	2869	0.55	ng	97
10) Phenol	6.36	94	2385	0.27	ng	# 59
11) bis(2-Chloroethyl)ether	6.44	93	4449	0.67	ng	88
12) 1,3-Dichlorobenzene	6.66	146	4694m	0.66	ng	
13) 1,4-Dichlorobenzene	6.73	146	4754m	0.68	ng	
14) 1,2-Dichlorobenzene	6.89	146	5026	0.76	ng	# 92
15) Benzyl Alcohol	6.86	79	2935	0.51	ng	92
16) 2,2'-oxybis(1-Chloropropan	7.00	45	5417	0.57	ng	88
17) 2-Methylphenol	6.98	107	2724	0.52	ng	96
18) Hexachloroethane	7.23	117	1865	0.68	ng	90
19) n-Nitroso-di-n-propylamine	7.13	70	3716	0.71	ng	88
20) 3+4-Methylphenols	7.13	107	3611	0.57	ng	# 60
22) Acetophenone	7.13	105	6885	0.78	ng	# 92
24) Nitrobenzene	7.30	77	4992	0.71	ng	93
25) Isophorone	7.54	82	10076	0.79	ng	97
26) 2-Nitrophenol	7.62	139	2070	0.72	ng	# 87
27) 2,4-Dimethylphenol	7.67	122	3912	0.66	ng	96
28) bis(2-Chloroethoxy)methane	7.76	93	6681	0.80	ng	# 97
29) 2,4-Dichlorophenol	7.86	162	3235	0.67	ng	92
30) 1,2,4-Trichlorobenzene	7.95	180	4156	0.79	ng	97
31) Naphthalene	8.02	128	15116	0.84	ng	99
33) 4-Chloroaniline	8.08	127	3502	0.44	ng	# 95
34) Hexachlorobutadiene	8.15	225	2082	0.73	ng	98
36) 4-Chloro-3-methylphenol	8.56	107	4194	0.74	ng	95
37) 2-Methylnaphthalene	8.72	142	10469	0.91	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.88	216	3774	0.65	ng	# 1
40) Hexachlorocyclopentadiene	8.88	237	900	0.32	ng	94
42) 2,4,6-Trichlorophenol	8.99	196	2698	0.71	ng	98
43) 2,4,5-Trichlorophenol	9.04	196	2542	0.78	ng	# 97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
45) 1,1'-Biphenyl	9.18	154	11824	0.82	nq	98
46) 2-Chloronaphthalene	9.20	162	9788	0.87	nq	98
47) 2-Nitroaniline	9.29	65	2498	0.62	nq	90
48) Acenaphthylene	9.61	152	14891	0.83	nq	98
49) Dimethylphthalate	9.47	163	15699	1.27	nq	100
50) 2,6-Dinitrotoluene	9.54	165	1928	0.70	nq	97
51) Acenaphthene	9.78	154	10106	0.92	nq	98
52) 3-Nitroaniline	9.70	138	1933	0.55	nq	# 95
54) Dibenzofuran	9.95	168	13973	0.97	nq	# 90
55) 4-Nitrophenol	9.87	139	1777	0.67	nq	# 82
56) 2,4-Dinitrotoluene	9.94	165	2446	0.68	nq	# 91
57) Fluorene	10.30	166	11570	1.04	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.08	232	1941	0.77	nq	# 53
59) Diethylphthalate	10.17	149	11059	0.89	nq	95
60) 4-Chlorophenyl-phenylether	10.30	204	5332	1.03	nq	95
61) 4-Nitroaniline	10.30	138	1987	0.59	nq	# 77
62) Azobenzene	10.46	77	11357	0.80	nq	87
65) n-Nitrosodiphenylamine	10.41	169	9973	0.86	nq	99
66) 4-Bromophenyl-phenylether	10.78	248	2806	0.81	nq	# 66
67) Hexachlorobenzene	10.84	284	3028	0.79	nq	# 88
68) Atrazine	10.95	200	3247	0.90	nq	90
69) Pentachlorophenol	11.04	266	3594	1.59	nq	93
70) Phenanthrene	11.26	178	14275	0.76	nq	98
71) Anthracene	11.30	178	16633	0.89	nq	97
72) Carbazole	11.46	167	13807	0.79	nq	98
73) Di-n-butylphthalate	11.80	149	19764	0.97	nq	100
74) Fluoranthene	12.43	202	14956	0.79	nq	99
77) Pvrene	12.66	202	13918	0.89	nq	100
79) Butylbenzylphthalate	13.29	149	7464	1.07	nq	# 86
80) Benzo(a)anthracene	13.84	228	9874	0.83	nq	98
81) 3,3'-Dichlorobenzidine	13.82	252	1161	0.26	nq	# 94
82) Chrysene	13.87	228	9202	0.78	nq	95
83) Bis(2-ethylhexyl)phthalate	13.85	149	10453	1.31	nq	# 98
84) Di-n-octyl phthalate	14.47	149	15115	1.12	nq	# 100
85) Indeno(1,2,3-cd)pvrene	16.51	276	8585	0.95	nq	# 100
87) Benzo(b)fluoranthene	14.85	252	8273m	0.77	nq	
88) Benzo(k)fluoranthene	14.87	252	7198m	0.77	nq	
89) Benzo(a)pyrene	15.17	252	7406	0.81	nq	# 83
90) Dibenzo(a,h)anthracene	16.54	278	7130	0.93	nq	# 95
91) Benzo(g,h,i)perylene	16.90	276	7578	0.96	nq	96

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

