

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070316\
 Data File : BF088651.D
 Acq On : 3 Jul 2016 16:02
 Operator : UM/SJ
 Sample : H3837-08MSD
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C1.2-6MSD

Manual Integrations

APPROVED
 umangi
 7/5/2016 7:51:47 PM

Quant Time: Jul 04 05:54:10 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 30 18:01:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	111490	20.00	ng	-0.02
21) Naphthalene-d8	8.07	136	439910	20.00	ng	-0.03
38) Acenaphthene-d10	9.82	164	204324	20.00	ng	-0.03
63) Phenanthrene-d10	11.30	188	422082	20.00	ng	-0.02
75) Chrysene-d12	13.92	240	282983	20.00	ng	-0.03
86) Perylene-d12	15.31	264	264738	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.39	112	946848	127.28	ng	-0.01
7) Phenol-d6	6.42	99	1146076	119.23	ng	-0.02
23) Nitrobenzene-d5	7.35	82	767607	91.44	ng	-0.03
41) 2,4,6-Tribromophenol	10.60	330	276633	145.80	ng	-0.03
44) 2-Fluorobiphenyl	9.15	172	1338216	103.45	ng	-0.02
78) Terphenyl-d14	12.88	244	1134222	89.27	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.39	88	115126	31.22	ng	# 34
3) Pyridine	3.10	79	43117	4.03	ng	89
4) n-Nitrosodimethylamine	3.02	42	82879	20.27	ng	87
6) Aniline	6.44	93	122331	8.73	ng	# 87
8) 2-Chlorophenol	6.57	128	194293	24.30	ng	96
9) Benzaldehyde	6.33	77	123639	18.99	ng	99
10) Phenol	6.43	94	178517	16.27	ng	# 62
11) bis(2-Chloroethyl)ether	6.52	93	185729	22.70	ng	90
12) 1,3-Dichlorobenzene	6.73	146	177749m	20.20	ng	
13) 1,4-Dichlorobenzene	6.80	146	185449	21.24	ng	94
14) 1,2-Dichlorobenzene	6.96	146	183357	22.43	ng	95
15) Benzyl Alcohol	6.92	79	167460	23.67	ng	94
16) 2,2'-oxybis(1-Chloropropan	7.07	45	242956	20.51	ng	91
17) 2-Methylphenol	7.04	107	151995	23.30	ng	97
18) Hexachloroethane	7.30	117	74927	22.18	ng	94
19) n-Nitroso-di-n-propylamine	7.20	70	133482	20.67	ng	94
20) 3+4-Methylphenols	7.20	107	196677	25.12	ng	# 83
22) Acetophenone	7.20	105	273489	25.61	ng	# 90
24) Nitrobenzene	7.37	77	232416	27.46	ng	91
25) Isophorone	7.61	82	396290	25.49	ng	99
26) 2-Nitrophenol	7.69	139	92272	26.59	ng	# 75
27) 2,4-Dimethylphenol	7.72	122	159388	22.05	ng	88
28) bis(2-Chloroethoxy)methane	7.83	93	255548	25.25	ng	# 96
29) 2,4-Dichlorophenol	7.93	162	155733	26.66	ng	95
30) 1,2,4-Trichlorobenzene	8.01	180	151887	23.69	ng	100
31) Naphthalene	8.09	128	537495	24.78	ng	99
32) Benzoic acid	7.79	122	21605	4.12	ng	93
33) 4-Chloroaniline	8.14	127	104557	10.84	ng	97
34) Hexachlorobutadiene	8.22	225	84989	24.76	ng	100
35) Caprolactam	8.54	113	49678m	25.04	ng	
36) 4-Chloro-3-methylphenol	8.63	107	176672	25.57	ng	92
37) 2-Methylnaphthalene	8.79	142	339003	24.28	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.95	216	131104	19.29	ng	# 1
40) Hexachlorocyclopentadiene	8.94	237	237386	71.17	ng	99

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42) 2,4,6-Trichlorophenol	9.06	196	110742	24.72	nq	99
43) 2,4,5-Trichlorophenol	9.10	196	110324	28.62	nq #	90
45) 1,1'-Biphenyl	9.24	154	421582	24.92	nq	99
46) 2-Chloronaphthalene	9.27	162	329171	24.78	nq	99
47) 2-Nitroaniline	9.36	65	113536	24.09	nq	98
48) Acenaphthylene	9.68	152	518775	24.55	nq	100
49) Dimethylphthalate	9.55	163	557727	38.31	nq	98
50) 2,6-Dinitrotoluene	9.61	165	90096	27.93	nq	93
51) Acenaphthene	9.85	154	352513	27.21	nq	98
52) 3-Nitroaniline	9.77	138	81268	19.82	nq	95
53) 2,4-Dinitrophenol	9.87	184	76286	53.55	nq #	85
54) Dibenzofuran	10.02	168	440286	25.97	nq #	88
55) 4-Nitrophenol	9.94	139	253400	81.31	nq #	80
56) 2,4-Dinitrotoluene	10.01	165	121169	28.73	nq	98
57) Fluorene	10.36	166	356951	27.32	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.15	232	83601	28.03	nq #	52
59) Diethylphthalate	10.25	149	421560	28.86	nq	99
60) 4-Chlorophenyl-phenylether	10.36	204	161789	26.65	nq	92
61) 4-Nitroaniline	10.38	138	80701	20.45	nq	83
62) Azobenzene	10.52	77	467356	28.05	nq	92
64) 4,6-Dinitro-2-methylphenol	10.41	198	52688	23.34	nq	90
65) n-Nitrosodiphenylamine	10.48	169	334147	23.39	nq	100
66) 4-Bromophenyl-phenylether	10.84	248	95338	22.41	nq #	75
67) Hexachlorobenzene	10.91	284	111553	23.69	nq #	85
68) Atrazine	11.00	200	95693	21.50	nq	91
69) Pentachlorophenol	11.11	266	177725	63.78	nq	98
70) Phenanthrene	11.32	178	573663	24.86	nq	98
71) Anthracene	11.37	178	532284	23.20	nq	99
72) Carbazole	11.53	167	525710	24.35	nq	97
73) Di-n-butylphthalate	11.86	149	616103	24.53	nq	99
74) Fluoranthene	12.50	202	560287	23.95	nq	99
76) Benzidine	12.63	184	44800	3.13	nq	98
77) Pyrene	12.73	202	588676	24.53	nq	99
79) Butylbenzylphthalate	13.36	149	272190	25.43	nq	97
80) Benzo(a)anthracene	13.91	228	447566	24.56	nq	99
81) 3,3'-Dichlorobenzidine	13.87	252	98953	14.44	nq #	93
82) Chrysene	13.95	228	452385	25.09	nq	98
83) Bis(2-ethylhexyl)phthalate	13.92	149	349200	28.58	nq	99
84) Di-n-octyl phthalate	14.53	149	581628	28.02	nq #	100
85) Indeno(1,2,3-cd)pyrene	16.64	276	346543	24.99	nq #	100
87) Benzo(b)fluoranthene	14.93	252	428066m	25.78	nq	
88) Benzo(k)fluoranthene	14.95	252	335763m	23.22	nq	
89) Benzo(a)pyrene	15.26	252	354411	25.21	nq #	96
90) Dibenzo(a,h)anthracene	16.66	278	286103	24.37	nq	98
91) Benzo(g,h,i)perylene	17.04	276	287790	23.69	nq	99

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

