

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062116\
 Data File : BF088241.D
 Acq On : 22 Jun 2016 5:54
 Operator : UM/SJ
 Sample : H3597-02MS
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EDD-3EMS

Manual Integrations
 APPROVED

sohil
 6/22/2016 6:08:13 PM

Quant Time: Jun 22 12:41:26 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 22 01:34:10 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	86880m	20.00	ng	0.00
21) Naphthalene-d8	7.94	136	342005	20.00	ng	0.00
38) Acenaphthene-d10	9.69	164	158188	20.00	ng	0.00
63) Phenanthrene-d10	11.16	188	295963	20.00	ng	0.00
75) Chrysene-d12	13.79	240	210325	20.00	ng	0.00
86) Perylene-d12	15.17	264	175689	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.24	112	438549	86.29	ng	0.00
7) Phenol-d6	6.31	99	543477	88.24	ng	0.00
23) Nitrobenzene-d5	7.22	82	332722	56.81	ng	-0.01
41) 2,4,6-Tribromophenol	10.48	330	127199	76.15	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	672689	65.54	ng	0.00
78) Terphenyl-d14	12.74	244	551276	59.64	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.17	88	68274	27.65	ng	# 33
3) Pyridine	2.82	79	177838	30.50	ng	94
4) n-Nitrosodimethylamine	2.78	42	70099	28.39	ng	# 78
6) Aniline	6.40	93	188754	21.53	ng	# 57
8) 2-Chlorophenol	6.43	128	192357	31.98	ng	99
10) Phenol	6.32	94	227393	35.22	ng	73
11) bis(2-Chloroethyl)ether	6.40	93	188754	34.95	ng	# 83
12) 1,3-Dichlorobenzene	6.59	146	197513m	30.60	ng	
13) 1,4-Dichlorobenzene	6.67	146	198639m	30.55	ng	
14) 1,2-Dichlorobenzene	6.82	146	200855	33.97	ng	95
15) Benzyl Alcohol	6.81	79	129918	26.96	ng	99
16) 2,2'-oxybis(1-Chloropropan	6.94	45	198517	27.21	ng	# 31
17) 2-Methylphenol	6.92	107	152002	31.16	ng	98
18) Hexachloroethane	7.15	117	72879m	31.59	ng	
19) n-Nitroso-di-n-propylamine	7.07	70	116118	29.47	ng	89
20) 3+4-Methylphenols	7.08	107	211832	37.22	ng	# 82
22) Acetophenone	7.07	105	263852	34.52	ng	97
24) Nitrobenzene	7.24	77	201109	35.45	ng	100
25) Isophorone	7.48	82	353748	32.61	ng	96
26) 2-Nitrophenol	7.56	139	95676	31.48	ng	# 74
27) 2,4-Dimethylphenol	7.61	122	179243	32.03	ng	95
28) bis(2-Chloroethoxy)methane	7.70	93	236438	35.14	ng	98
29) 2,4-Dichlorophenol	7.80	162	168480	36.06	ng	98
30) 1,2,4-Trichlorobenzene	7.88	180	168265	33.08	ng	97
31) Naphthalene	7.95	128	553582	32.71	ng	99
32) Benzoic acid	7.61	122	179243	58.17	ng	# 13
33) 4-Chloroaniline	8.02	127	105495	13.96	ng	# 93
34) Hexachlorobutadiene	8.08	225	86916	32.40	ng	99
35) Caprolactam	8.39	113	51672	33.39	ng	# 79
36) 4-Chloro-3-methylphenol	8.51	107	167679	33.56	ng	94
37) 2-Methylnaphthalene	8.65	142	370978	33.26	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.82	216	150921	33.74	ng	# 1
40) Hexachlorocyclopentadiene	8.80	237	51939	20.36	ng	99
42) 2,4,6-Trichlorophenol	8.94	196	82740	26.16	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	8.99	196	87218	26.50	nq	97
45) 1,1'-Biphenyl	9.12	154	441440	35.93	nq	95
46) 2-Chloronaphthalene	9.14	162	347675	33.96	nq	93
47) 2-Nitroaniline	9.24	65	100600	33.31	nq	# 80
48) Acenaphthylene	9.55	152	532933	32.73	nq	99
49) Dimethylphthalate	9.43	163	664592	58.00	nq	100
50) 2,6-Dinitrotoluene	9.48	165	86492	32.96	nq	# 62
51) Acenaphthene	9.72	154	312708	30.79	nq	100
52) 3-Nitroaniline	9.66	138	72858	24.06	nq	91
53) 2,4-Dinitrophenol	9.84	184	3531	6.07	nq	# 1
54) Dibenzofuran	9.90	168	456804	32.66	nq	95
55) 4-Nitrophenol	9.84	139	59093	25.14	nq	98
56) 2,4-Dinitrotoluene	9.90	165	108717	32.24	nq	# 86
57) Fluorene	10.24	166	358835	37.11	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.02	232	64254	24.84	nq	# 60
59) Diethylphthalate	10.12	149	405814	34.99	nq	98
60) 4-Chlorophenyl-phenylether	10.23	204	156955	33.74	nq	98
61) 4-Nitroaniline	10.27	138	105257	36.65	nq	91
62) Azobenzene	10.39	77	388612	33.88	nq	97
64) 4,6-Dinitro-2-methylphenol	10.31	198	18456	11.70	nq	# 68
65) n-Nitrosodiphenylamine	10.35	169	342939	34.92	nq	100
66) 4-Bromophenyl-phenylether	10.72	248	102001	32.12	nq	95
67) Hexachlorobenzene	10.78	284	102033	30.93	nq	98
68) Atrazine	10.88	200	87658	29.48	nq	90
69) Pentachlorophenol	10.98	266	44748	25.73	nq	97
70) Phenanthrene	11.19	178	513866	33.09	nq	100
71) Anthracene	11.24	178	542439	34.63	nq	98
72) Carbazole	11.40	167	520179	35.48	nq	98
73) Di-n-butylphthalate	11.74	149	611330	32.94	nq	99
74) Fluoranthene	12.38	202	561199	34.24	nq	96
76) Benzidine	12.50	184	189733	32.58	nq	98
77) Pyrene	12.59	202	553654	35.47	nq	99
79) Butylbenzylphthalate	13.22	149	249487	36.16	nq	90
80) Benzo(a)anthracene	13.78	228	381869	31.86	nq	100
81) 3,3'-Dichlorobenzidine	13.75	252	86687	26.90	nq	# 94
82) Chrysene	13.82	228	417382	37.02	nq	98
83) Bis(2-ethylhexyl)phthalate	13.78	149	326500	38.87	nq	99
84) Di-n-octyl phthalate	14.39	149	519643	38.07	nq	# 100
85) Indeno(1,2,3-cd)pyrene	16.46	276	353214	33.72	nq	# 100
87) Benzo(b)fluoranthene	14.79	252	334101m	27.28	nq	
88) Benzo(k)fluoranthene	14.82	252	360780m	39.06	nq	
89) Benzo(a)pyrene	15.12	252	346099	34.93	nq	# 94
90) Dibenzo(a,h)anthracene	16.46	278	294474	32.00	nq	99
91) Benzo(g,h,i)perylene	16.83	276	312254	32.99	nq	99

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

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