

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061116\
 Data File : BF087931.D
 Acq On : 12 Jun 2016 6:20
 Operator : UM/SJ
 Sample : H3501-25MS
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW1MS

Manual Integrations
 APPROVED

umangi
 6/13/2016 7:44:46 PM

Quant Time: Jun 13 12:15:51 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 13 01:40:32 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	105046	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	431870	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	213031	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	381575	20.00	ng	0.00
75) Chrysene-d12	13.97	240	273415	20.00	ng	0.00
86) Perylene-d12	15.37	264	235975	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.39	112	462760	75.31	ng	0.00
7) Phenol-d6	6.44	99	401799	53.96	ng	0.00
23) Nitrobenzene-d5	7.37	82	718140	97.10	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	252666	112.33	ng	0.00
44) 2-Fluorobiphenyl	9.18	172	1221245	90.41	ng	0.00
78) Terphenyl-d14	12.91	244	1007947	83.89	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.39	88	52999	17.75	ng	# 28
3) Pyridine	3.11	79	115078	16.32	ng	96
4) n-Nitrosodimethylamine	3.03	42	66294	22.20	ng	86
6) Aniline	6.47	93	193987	18.30	ng	98
8) 2-Chlorophenol	6.59	128	320344	44.05	ng	84
9) Benzaldehyde	6.34	77	26644	4.56	ng	94
10) Phenol	6.46	94	168057	20.50	ng	# 67
11) bis(2-Chloroethyl)ether	6.54	93	326322	49.97	ng	95
12) 1,3-Dichlorobenzene	6.74	146	342401	43.88	ng	97
13) 1,4-Dichlorobenzene	6.82	146	363490	46.24	ng	98
14) 1,2-Dichlorobenzene	6.97	146	312249m	43.68	ng	
15) Benzyl Alcohol	6.95	79	193042	33.13	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.10	45	448314	50.82	ng	89
17) 2-Methylphenol	7.07	107	222947	37.80	ng	99
18) Hexachloroethane	7.31	117	120602	43.24	ng	# 83
19) n-Nitroso-di-n-propylamine	7.23	70	273569	57.42	ng	# 87
20) 3+4-Methylphenols	7.22	107	231349	33.62	ng	# 39
22) Acetophenone	7.22	105	523722	54.26	ng	97
24) Nitrobenzene	7.39	77	396335	55.32	ng	92
25) Isophorone	7.63	82	649355	47.40	ng	99
26) 2-Nitrophenol	7.71	139	219524	57.20	ng	# 75
27) 2,4-Dimethylphenol	7.76	122	348666	49.35	ng	98
28) bis(2-Chloroethoxy)methane	7.85	93	472529	55.61	ng	98
29) 2,4-Dichlorophenol	7.95	162	293092	49.68	ng	99
30) 1,2,4-Trichlorobenzene	8.03	180	292440	45.52	ng	100
31) Naphthalene	8.11	128	995734	46.59	ng	100
32) Benzoic acid	7.87	122	103823	26.68	ng	97
33) 4-Chloroaniline	8.17	127	139993	14.67	ng	# 91
34) Hexachlorobutadiene	8.24	225	159466	47.07	ng	98
35) Caprolactam	8.58	113	21623	11.07	ng	95
36) 4-Chloro-3-methylphenol	8.65	107	289331	45.86	ng	97
37) 2-Methylnaphthalene	8.81	142	750523	53.28	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	274603	52.46	ng	# 1
40) Hexachlorocyclopentadiene	8.96	237	268313	78.08	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.08	196	216510	50.82	nq	97
43) 2,4,5-Trichlorophenol	9.14	196	231136	52.15	nq	97
45) 1,1'-Biphenyl	9.28	154	899667	58.07	nq	95
46) 2-Chloronaphthalene	9.30	162	709497	51.47	nq	# 92
47) 2-Nitroaniline	9.39	65	191158	47.01	nq	92
48) Acenaphthylene	9.71	152	1078225	49.17	nq	99
49) Dimethylphthalate	9.59	163	874150	56.65	nq	99
50) 2,6-Dinitrotoluene	9.64	165	214557	60.71	nq	93
51) Acenaphthene	9.88	154	778452	56.91	nq	99
52) 3-Nitroaniline	9.80	138	96844	23.75	nq	93
53) 2,4-Dinitrophenol	9.92	184	221183	105.63	nq	# 70
54) Dibenzofuran	10.06	168	940261	49.91	nq	# 91
55) 4-Nitrophenol	9.98	139	117854	37.24	nq	93
56) 2,4-Dinitrotoluene	10.04	165	273411	60.20	nq	96
57) Fluorene	10.40	166	732441	59.04	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.17	232	166499	47.80	nq	# 54
59) Diethylphthalate	10.27	149	730009	46.73	nq	99
60) 4-Chlorophenyl-phenylether	10.39	204	272810	43.55	nq	95
61) 4-Nitroaniline	10.42	138	120749	31.22	nq	90
62) Azobenzene	10.55	77	758367	49.10	nq	98
64) 4,6-Dinitro-2-methylphenol	10.44	198	120071	50.30	nq	95
65) n-Nitrosodiphenylamine	10.51	169	669199	52.85	nq	100
66) 4-Bromophenyl-phenylether	10.88	248	204993	50.08	nq	# 90
67) Hexachlorobenzene	10.95	284	227665	53.53	nq	# 72
68) Atrazine	11.05	200	211548	55.18	nq	94
69) Pentachlorophenol	11.14	266	238868	106.54	nq	98
70) Phenanthrene	11.36	178	1122069	56.04	nq	99
71) Anthracene	11.41	178	928508	45.97	nq	100
72) Carbazole	11.57	167	1026373	54.30	nq	100
73) Di-n-butylphthalate	11.90	149	1079637	45.12	nq	99
74) Fluoranthene	12.54	202	982235	46.48	nq	98
77) Pyrene	12.77	202	996321	49.11	nq	100
79) Butylbenzylphthalate	13.39	149	533003	59.42	nq	94
80) Benzo(a)anthracene	13.95	228	723448	46.43	nq	100
81) 3,3'-Dichlorobenzidine	13.93	252	10026	2.39	nq	# 84
82) Chrysene	13.99	228	751901	51.30	nq	99
83) Bis(2-ethylhexyl)phthalate	13.95	149	564069	51.66	nq	# 99
84) Di-n-octyl phthalate	14.56	149	1063522	59.94	nq	# 100
85) Indeno(1,2,3-cd)pyrene	16.75	276	781710	57.40	nq	# 100
87) Benzo(b)fluoranthene	14.97	252	960561m	58.40	nq	
88) Benzo(k)fluoranthene	15.01	252	520449m	41.95	nq	
89) Benzo(a)pyrene	15.33	252	720513	54.15	nq	# 92
90) Dibenzo(a,h)anthracene	16.77	278	639113	51.71	nq	98
91) Benzo(g,h,i)perylene	17.17	276	598134	47.05	nq	97

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

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