

Data Path : Z:\HPCHEM1\BNA F\DATA\BF090315\
 Data File : BF081411.D
 Acq On : 4 Sep 2015 9:26
 Operator : UM/IZ
 Sample : G3505-03MSD
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P005-BF001-01MSD

Manual Integrations
 APPROVED

MMDadoda
 9/4/2015 2:51:55 PM

Quant Time: Sep 04 11:54:26 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 01 17:35:35 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.36	152	55005	20.00	ng	-0.02
21) Naphthalene-d8	7.66	136	205102	20.00	ng	-0.02
38) Acenaphthene-d10	9.40	164	100225	20.00	ng	-0.01
63) Phenanthrene-d10	10.87	188	191648	20.00	ng	-0.01
75) Chrysene-d12	13.47	240	123691	20.00	ng	-0.02
86) Perylene-d12	14.78	264	91580	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	4.92	112	455918	128.60	ng	0.00
7) Phenol-d6	6.04	99	619520	132.02	ng	-0.01
23) Nitrobenzene-d5	6.95	82	401663	94.48	ng	-0.02
41) 2,4,6-Tribromophenol	10.19	330	128474	143.96	ng	-0.01
44) 2-Fluorobiphenyl	8.74	172	632323	80.85	ng	-0.02
78) Terphenyl-d14	12.45	244	509737	95.93	ng	-0.02

Target Compounds						Qvalue
2) 1,4-Dioxane	1.75	88	56177	35.28	ng	# 89
4) n-Nitrosodimethylamine	2.25	42	104274	42.75	ng	# 74
6) Aniline	6.03	93	66710	10.07	ng	# 55
8) 2-Chlorophenol	6.16	128	175152	44.83	ng	92
9) Benzaldehyde	5.91	77	24672	8.39	ng	# 83
10) Phenol	6.06	94	203189	37.34	ng	# 49
11) bis(2-Chloroethyl)ether	6.12	93	175475	44.69	ng	90
12) 1,3-Dichlorobenzene	6.31	146	185611	44.47	ng	# 92
13) 1,4-Dichlorobenzene	6.39	146	184391	43.39	ng	# 90
14) 1,2-Dichlorobenzene	6.54	146	153675	40.16	ng	# 86
15) Benzyl Alcohol	6.54	79	157217	40.84	ng	# 72
16) 2,2'-oxybis(1-Chloropropan	6.67	45	332199	44.14	ng	# 9
17) 2-Methylphenol	6.67	107	144297	46.29	ng	# 75
18) Hexachloroethane	6.88	117	70607	42.70	ng	91
19) n-Nitroso-di-n-propylamine	6.81	70	136059	42.60	ng	# 86
20) 3+4-Methylphenols	6.83	107	186160	44.29	ng	88
22) Acetophenone	6.80	105	261062	46.60	ng	# 78
24) Nitrobenzene	6.97	77	227784	51.60	ng	# 67
25) Isophorone	7.21	82	338049	42.75	ng	# 81
26) 2-Nitrophenol	7.29	139	93539	48.61	ng	# 66
27) 2,4-Dimethylphenol	7.35	122	131362	39.53	ng	# 73
28) bis(2-Chloroethoxy)methane	7.44	93	206108	45.13	ng	# 92
29) 2,4-Dichlorophenol	7.54	162	139459	48.39	ng	97
30) 1,2,4-Trichlorobenzene	7.61	180	152199	48.61	ng	99
31) Naphthalene	7.68	128	471071	43.07	ng	98
32) Benzoic acid	7.44	122	6190	5.01	ng	# 59
33) 4-Chloroaniline	7.75	127	54387	12.19	ng	# 86
34) Hexachlorobutadiene	7.80	225	85098	44.66	ng	96
35) Caprolactam	8.12	113	26859m	30.39	ng	
36) 4-Chloro-3-methylphenol	8.25	107	150977	46.80	ng	# 69
37) 2-Methylnaphthalene	8.38	142	335596	47.15	ng	# 90
39) 1,2,4,5-Tetrachlorobenzene	8.55	216	131565	41.51	ng	98
40) Hexachlorocyclopentadiene	8.52	237	102383	65.82	ng	94
42) 2,4,6-Trichlorophenol	8.66	196	90446	41.35	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	8.71	196	93805	42.03	nq	# 83
45) 1,1'-Biphenyl	8.84	154	403820	43.79	nq	94
46) 2-Chloronaphthalene	8.86	162	281726	42.31	nq	# 88
47) 2-Nitroaniline	8.97	65	127065	46.82	nq	# 76
48) Acenaphthylene	9.27	152	496486	42.03	nq	98
49) Dimethylphthalate	9.16	163	483497	62.09	nq	# 97
50) 2,6-Dinitrotoluene	9.21	165	65446	37.03	nq	# 1
51) Acenaphthene	9.44	154	300289	44.68	nq	90
52) 3-Nitroaniline	9.38	138	58170	28.91	nq	# 87
53) 2,4-Dinitrophenol	9.48	184	29343	39.42	nq	# 19
54) Dibenzofuran	9.61	168	448707	45.73	nq	# 92
55) 4-Nitrophenol	9.58	139	113604	89.87	nq	# 36
56) 2,4-Dinitrotoluene	9.61	165	88330	39.25	nq	# 24
57) Fluorene	9.94	166	307855	41.34	nq	97
58) 2,3,4,6-Tetrachlorophenol	9.74	232	81323	47.73	nq	97
59) Diethylphthalate	9.85	149	323019	39.44	nq	95
60) 4-Chlorophenyl-phenylether	9.95	204	159084	45.84	nq	96
61) 4-Nitroaniline	9.99	138	75038	36.91	nq	# 27
62) Azobenzene	10.10	77	383058	42.06	nq	82
64) 4,6-Dinitro-2-methylphenol	10.02	198	36682	34.22	nq	# 77
65) n-Nitrosodiphenylamine	10.07	169	263923	37.85	nq	91
66) 4-Bromophenyl-phenylether	10.43	248	88065	45.45	nq	94
67) Hexachlorobenzene	10.49	284	85840	42.37	nq	# 77
68) Atrazine	10.62	200	86725	42.97	nq	81
69) Pentachlorophenol	10.68	266	77532	78.76	nq	98
70) Phenanthrene	10.89	178	458162	43.00	nq	97
71) Anthracene	10.95	178	443377	40.50	nq	99
72) Carbazole	11.11	167	451378	43.94	nq	96
73) Di-n-butylphthalate	11.46	149	551618	45.48	nq	# 99
74) Fluoranthene	12.06	202	483515	41.92	nq	92
77) Pyrene	12.29	202	505908	55.22	nq	99
79) Butylbenzylphthalate	12.94	149	210004	52.70	nq	92
80) Benzo(a)anthracene	13.46	228	335300	42.57	nq	97
81) 3,3'-Dichlorobenzidine	13.44	252	57937	21.80	nq	# 95
82) Chrysene	13.50	228	267723	37.92	nq	96
83) Bis(2-ethylhexyl)phthalate	13.50	149	250931	47.72	nq	# 89
84) Di-n-octyl phthalate	14.10	149	350840	40.52	nq	# 96
85) Indeno(1,2,3-cd)pyrene	15.86	276	272439	52.59	nq	# 87
87) Benzo(b)fluoranthene	14.46	252	323077m	49.41	nq	
88) Benzo(k)fluoranthene	14.48	252	214203m	39.48	nq	
89) Benzo(a)pyrene	14.73	252	248999	44.94	nq	# 87
90) Dibenzo(a,h)anthracene	15.87	278	222176	51.90	nq	# 81
91) Benzo(g,h,i)perylene	16.18	276	224585	54.96	nq	# 75

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

