

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020916\
 Data File : BF084951.D
 Acq On : 9 Feb 2016 14:50
 Operator : SJ/IZ
 Sample : H1377-11MS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB03-2-3MS

Manual Integrations
 APPROVED

Sohil
 2/10/2016 1:44:52 PM

Quant Time: Feb 10 03:56:00 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 03 14:16:01 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.64	152	164305	20.00	ng	-0.05
21) Naphthalene-d8	7.93	136	669531	20.00	ng	-0.05
38) Acenaphthene-d10	9.68	164	316579	20.00	ng	-0.06
63) Phenanthrene-d10	11.15	188	563923	20.00	ng	-0.06
75) Chrysene-d12	13.78	240	404451	20.00	ng	-0.06
86) Perylene-d12	15.14	264	314531	20.00	ng	-0.07

System Monitoring Compounds						
5) 2-Fluorophenol	5.24	112	1494679	141.69	ng	-0.03
7) Phenol-d6	6.30	99	1899377	145.24	ng	-0.03
23) Nitrobenzene-d5	7.21	82	1181993	99.45	ng	-0.06
41) 2,4,6-Tribromophenol	10.46	330	428555	129.78	ng	-0.06
44) 2-Fluorobiphenyl	9.00	172	1581779	83.70	ng	-0.06
78) Terphenyl-d14	12.74	244	1737026	97.32	ng	-0.06

Target Compounds						Qvalue
2) 1,4-Dioxane	2.17	88	147244	31.87	ng	# 73
3) Pyridine	2.82	79	405178	33.66	ng	# 74
4) n-Nitrosodimethylamine	2.77	42	257283	41.32	ng	# 80
6) Aniline	6.30	93	435672	27.23	ng	# 58
8) 2-Chlorophenol	6.43	128	471953	39.53	ng	# 90
9) Benzaldehyde	6.18	77	140696	18.22	ng	# 96
10) Phenol	6.32	94	674305	46.03	ng	# 85
11) bis(2-Chloroethyl)ether	6.38	93	489174	42.82	ng	# 88
12) 1,3-Dichlorobenzene	6.58	146	492615	39.05	ng	# 98
13) 1,4-Dichlorobenzene	6.65	146	475879	37.73	ng	# 97
14) 1,2-Dichlorobenzene	6.81	146	489224	41.65	ng	# 98
15) Benzyl Alcohol	6.80	79	423621	46.91	ng	# 92
16) 2,2'-oxybis(1-Chloropropan	6.93	45	783665	40.78	ng	# 89
17) 2-Methylphenol	6.92	107	407348	43.14	ng	# 88
18) Hexachloroethane	7.15	117	197023	40.57	ng	# 87
19) n-Nitroso-di-n-propylamine	7.07	70	356772	43.52	ng	# 76
20) 3+4-Methylphenols	7.07	107	501123	42.75	ng	# 83
22) Acetophenone	7.06	105	696540	39.47	ng	# 98
24) Nitrobenzene	7.23	77	547286	43.00	ng	# 99
25) Isophorone	7.47	82	935485	40.48	ng	# 99
26) 2-Nitrophenol	7.55	139	277827	44.26	ng	# 94
27) 2,4-Dimethylphenol	7.60	122	408392	37.40	ng	# 96
28) bis(2-Chloroethoxy)methane	7.69	93	573501	41.83	ng	# 99
29) 2,4-Dichlorophenol	7.79	162	383378	41.04	ng	# 96
30) 1,2,4-Trichlorobenzene	7.87	180	436662	41.38	ng	# 96
31) Naphthalene	7.95	128	1398764	41.65	ng	# 99
32) Benzoic acid	7.70	122	65044	9.31	ng	# 96
33) 4-Chloroaniline	8.01	127	263508	18.97	ng	# 89
34) Hexachlorobutadiene	8.06	225	231268	38.01	ng	# 99
35) Caprolactam	8.38	113	127940m	44.43	ng	# 95
36) 4-Chloro-3-methylphenol	8.50	107	480111	45.05	ng	# 97
37) 2-Methylnaphthalene	8.64	142	920654	43.80	ng	# 99
39) 1,2,4,5-Tetrachlorobenzene	8.81	216	407891	40.57	ng	# 99
40) Hexachlorocyclopentadiene	8.80	237	328742	75.72	ng	# 98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.92	196	270636	41.78	nq	94
43) 2,4,5-Trichlorophenol	8.97	196	274263	41.67	nq	# 86
45) 1,1'-Biphenyl	9.10	154	1194948	42.26	nq	97
46) 2-Chloronaphthalene	9.13	162	877828	42.16	nq	# 88
47) 2-Nitroaniline	9.23	65	315169	47.80	nq	# 83
48) Acenaphthylene	9.54	152	1339101	42.38	nq	99
49) Dimethylphthalate	9.41	163	1189046	49.31	nq	# 90
50) 2,6-Dinitrotoluene	9.47	165	203613	38.66	nq	# 43
51) Acenaphthene	9.71	154	852169	41.45	nq	98
52) 3-Nitroaniline	9.64	138	184288	31.13	nq	# 84
53) 2,4-Dinitrophenol	9.76	184	173139	71.85	nq	90
54) Dibenzofuran	9.88	168	1172066	46.65	nq	96
55) 4-Nitrophenol	9.82	139	273317	62.83	nq	# 79
56) 2,4-Dinitrotoluene	9.88	165	302280	44.99	nq	91
57) Fluorene	10.22	166	908451	43.30	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.01	232	226013	45.11	nq	89
59) Diethylphthalate	10.11	149	932614	39.61	nq	99
60) 4-Chlorophenyl-phenylether	10.21	204	368227	41.15	nq	92
61) 4-Nitroaniline	10.26	138	256978	44.55	nq	90
62) Azobenzene	10.37	77	967981	42.76	nq	88
64) 4,6-Dinitro-2-methylphenol	10.28	198	143084	41.11	nq	# 51
65) n-Nitrosodiphenylamine	10.34	169	765180	42.73	nq	99
66) 4-Bromophenyl-phenylether	10.70	248	275096	44.17	nq	# 93
67) Hexachlorobenzene	10.76	284	265572	39.13	nq	# 83
68) Atrazine	10.88	200	283007	50.05	nq	97
69) Pentachlorophenol	10.97	266	233510	73.21	nq	98
70) Phenanthrene	11.17	178	1217823	38.94	nq	97
71) Anthracene	11.23	178	1396133	44.30	nq	99
72) Carbazole	11.39	167	1264404	46.01	nq	98
73) Di-n-butylphthalate	11.72	149	1362327	38.94	nq	# 96
74) Fluoranthene	12.35	202	1237275	39.08	nq	99
76) Benzidine	12.49	184	359218	25.70	nq	97
77) Pyrene	12.58	202	1243276	40.15	nq	98
79) Butylbenzylphthalate	13.21	149	589917	41.99	nq	# 83
80) Benzo(a)anthracene	13.77	228	1096218	43.67	nq	99
81) 3,3'-Dichlorobenzidine	13.73	252	165271	18.59	nq	# 95
82) Chrysene	13.80	228	822237	33.78	nq	100
83) Bis(2-ethylhexyl)phthalate	13.78	149	784458	44.87	nq	# 97
84) Di-n-octyl phthalate	14.39	149	1148685	39.83	nq	# 88
85) Indeno(1,2,3-cd)pyrene	16.40	276	805353	42.03	nq	98
87) Benzo(b)fluoranthene	14.77	252	1021093m	48.32	nq	
88) Benzo(k)fluoranthene	14.80	252	647788m	37.07	nq	
89) Benzo(a)pyrene	15.08	252	750849	41.70	nq	# 97
90) Dibenzo(a,h)anthracene	16.41	278	666670	43.98	nq	# 95
91) Benzo(g,h,i)perylene	16.77	276	687282	45.01	nq	97

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

