

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030416\
 Data File : BF085237.D
 Acq On : 5 Mar 2016 7:16
 Operator : SJ/IZ
 Sample : H1683-04MS
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-16-COMP(0.5-5.0)MS

Manual Integrations
 APPROVED

UMANGI
 3/7/2016 3:30:53 PM

Quant Time: Mar 07 12:24:35 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 04 00:54:58 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	163878	20.00	ng	-0.01
21) Naphthalene-d8	8.26	136	606704	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	261331	20.00	ng	-0.01
63) Phenanthrene-d10	11.50	188	397362	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	271787	20.00	ng	-0.01
86) Perylene-d12	15.60	264	276594	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	1416253	144.96	ng	0.01
7) Phenol-d6	6.61	99	1700540	132.86	ng	0.00
23) Nitrobenzene-d5	7.54	82	1157930	102.13	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	357499	159.88	ng	0.00
44) 2-Fluorobiphenyl	9.34	172	1470636	85.58	ng	-0.01
78) Terphenyl-d14	13.09	244	1034437	88.35	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.72	88	208966	41.83	ng	99
3) Pyridine	3.49	79	604114	48.31	ng	100
4) n-Nitrosodimethylamine	3.44	42	274412	51.28	ng	100
6) Aniline	6.64	93	523672	29.19	ng	96
8) 2-Chlorophenol	6.76	128	487761	41.47	ng	94
9) Benzaldehyde	6.52	77	179865	27.40	ng	98
10) Phenol	6.62	94	735387m	53.64	ng	
11) bis(2-Chloroethyl)ether	6.71	93	523547	45.98	ng	99
12) 1,3-Dichlorobenzene	6.92	146	561891	44.50	ng	98
13) 1,4-Dichlorobenzene	7.00	146	547827m	43.29	ng	
14) 1,2-Dichlorobenzene	7.14	146	511787	44.57	ng	99
15) Benzyl Alcohol	7.11	79	428985	45.66	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.25	45	643725	39.22	ng	98
17) 2-Methylphenol	7.22	107	455197	47.43	ng	96
18) Hexachloroethane	7.49	117	214929	46.04	ng	97
19) n-Nitroso-di-n-propylamine	7.40	70	380316	45.59	ng	96
20) 3+4-Methylphenols	7.37	107	509316	44.80	ng	98
22) Acetophenone	7.38	105	684413	46.22	ng	# 98
24) Nitrobenzene	7.56	77	524501	45.54	ng	99
25) Isophorone	7.80	82	1031227	49.08	ng	98
26) 2-Nitrophenol	7.88	139	290324	51.80	ng	# 89
27) 2,4-Dimethylphenol	7.91	122	516535	50.43	ng	94
28) bis(2-Chloroethoxy)methane	8.00	93	579915	49.56	ng	99
29) 2,4-Dichlorophenol	8.12	162	429407	51.98	ng	95
30) 1,2,4-Trichlorobenzene	8.20	180	423412	47.41	ng	98
31) Naphthalene	8.29	128	1484914	49.78	ng	98
32) Benzoic acid	7.98	122	75667	11.12	ng	# 49
33) 4-Chloroaniline	8.32	127	230886	19.13	ng	98
34) Hexachlorobutadiene	8.40	225	238870	51.40	ng	99
35) Caprolactam	8.69	113	173417	64.28	ng	# 87
36) 4-Chloro-3-methylphenol	8.81	107	468744	50.02	ng	95
37) 2-Methylnaphthalene	8.97	142	890762	46.53	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.14	216	376856	50.42	ng	97
40) Hexachlorocyclopentadiene	9.13	237	226348	56.15	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.25	196	259740	46.69	nq	96
43) 2,4,5-Trichlorophenol	9.29	196	261122	49.50	nq	99
45) 1,1'-Biphenyl	9.44	154	1099926	49.27	nq	95
46) 2-Chloronaphthalene	9.46	162	825793	48.52	nq	94
47) 2-Nitroaniline	9.56	65	290951	55.17	nq	89
48) Acenaphthylene	9.88	152	1295478	48.91	nq	98
49) Dimethylphthalate	9.74	163	1033121	51.51	nq	99
50) 2,6-Dinitrotoluene	9.80	165	197512	46.03	nq	# 81
51) Acenaphthene	10.05	154	838550	52.14	nq	99
52) 3-Nitroaniline	9.97	138	208584	40.63	nq	# 98
53) 2,4-Dinitrophenol	10.07	184	62867	34.00	nq	# 48
54) Dibenzofuran	10.23	168	1110426	52.70	nq	92
55) 4-Nitrophenol	10.13	139	397262	98.47	nq	94
56) 2,4-Dinitrotoluene	10.21	165	303679	55.30	nq	# 66
57) Fluorene	10.57	166	855843	51.71	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.34	232	206026	55.60	nq	# 97
59) Diethylphthalate	10.44	149	897509	46.11	nq	96
60) 4-Chlorophenyl-phenylether	10.56	204	408911	50.78	nq	# 85
61) 4-Nitroaniline	10.57	138	176177	36.69	nq	99
62) Azobenzene	10.72	77	890355	46.43	nq	83
64) 4,6-Dinitro-2-methylphenol	10.61	198	68825	28.93	nq	# 49
65) n-Nitrosodiphenylamine	10.68	169	802277	64.91	nq	96
66) 4-Bromophenyl-phenylether	11.05	248	213423	55.31	nq	# 81
67) Hexachlorobenzene	11.12	284	225404	54.76	nq	# 91
68) Atrazine	11.20	200	225057	65.48	nq	97
69) Pentachlorophenol	11.30	266	216305	94.67	nq	99
70) Phenanthrene	11.53	178	1477465	70.95	nq	99
71) Anthracene	11.58	178	1135740	51.37	nq	99
72) Carbazole	11.73	167	850812	43.89	nq	98
73) Di-n-butylphthalate	12.06	149	1197722	48.88	nq	99
74) Fluoranthene	12.72	202	1906365	91.22	nq	97
76) Benzidine	12.84	184	345361	42.70	nq	98
77) Pyrene	12.95	202	1762551	86.16	nq	99
79) Butylbenzylphthalate	13.56	149	459099	49.46	nq	# 78
80) Benzo(a)anthracene	14.13	228	1203217	73.95	nq	100
81) 3,3'-Dichlorobenzidine	14.08	252	119662	23.31	nq	98
82) Chrysene	14.16	228	844949	56.22	nq	98
83) Bis(2-ethylhexyl)phthalate	14.12	149	593893	48.62	nq	# 95
84) Di-n-octyl phthalate	14.72	149	949400	51.04	nq	98
85) Indeno(1,2,3-cd)pyrene	17.09	276	973591	83.00	nq	97
87) Benzo(b)fluoranthene	15.18	252	1476094m	80.07	nq	
88) Benzo(k)fluoranthene	15.21	252	626400m	42.12	nq	
89) Benzo(a)pyrene	15.56	252	1056417	69.15	nq	98
90) Dibenzo(a,h)anthracene	17.11	278	757466m	58.08	nq	
91) Benzo(g,h,i)perylene	17.55	276	793623	60.52	nq	96

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

