

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030317\
 Data File : BF093372.D
 Acq On : 3 Mar 2017 21:29
 Operator : SJ/MA
 Sample : I2086-04MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SRI-MW-4-20170302MS

Manual Integrations
 APPROVED

mohammad
 3/6/2017 4:23:27 PM

Quant Time: Mar 03 23:08:07 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 03 22:55:05 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	196616	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	819160	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	382255	20.00	ng	0.00
63) Phenanthrene-d10	11.31	188	669584	20.00	ng	0.00
75) Chrysene-d12	13.95	240	541004	20.00	ng	0.00
86) Perylene-d12	15.36	264	368415	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	784858	68.48	ng	0.00
7) Phenol-d6	6.43	99	658178	44.64	ng	-0.01
23) Nitrobenzene-d5	7.37	82	1373798	93.39	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	393796	142.44	ng	0.01
44) 2-Fluorobiphenyl	9.15	172	1874406	88.84	ng	0.00
78) Terphenyl-d14	12.89	244	1675685	72.78	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.35	88	102832	16.61	ng	# 81
3) Pyridine	3.07	79	234446	14.89	ng	89
4) n-Nitrosodimethylamine	3.00	42	139397	18.85	ng	# 85
6) Aniline	6.44	93	483591	24.04	ng	# 67
8) 2-Chlorophenol	6.57	128	467769	37.00	ng	85
9) Benzaldehyde	6.33	77	170623	16.15	ng	89
10) Phenol	6.45	94	290591	16.84	ng	80
11) bis(2-Chloroethyl)ether	6.52	93	583741	42.39	ng	92
12) 1,3-Dichlorobenzene	6.72	146	537411	36.29	ng	# 89
13) 1,4-Dichlorobenzene	6.80	146	550810	36.75	ng	99
14) 1,2-Dichlorobenzene	6.96	146	541795	37.80	ng	97
15) Benzyl Alcohol	6.94	79	314316	28.79	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.07	45	988077	41.30	ng	85
17) 2-Methylphenol	7.06	107	330143	30.44	ng	# 92
18) Hexachloroethane	7.30	117	176519	34.50	ng	# 84
19) n-Nitroso-di-n-propylamine	7.21	70	444337	47.34	ng	93
20) 3+4-Methylphenols	7.22	107	412447	31.80	ng	# 77
22) Acetophenone	7.21	105	849718	45.43	ng	# 95
24) Nitrobenzene	7.38	77	621997	41.18	ng	99
25) Isophorone	7.62	82	1256770	45.85	ng	95
26) 2-Nitrophenol	7.70	139	342356	47.68	ng	93
27) 2,4-Dimethylphenol	7.74	122	512901	40.68	ng	97
28) bis(2-Chloroethoxy)methane	7.84	93	829447	45.69	ng	99
29) 2,4-Dichlorophenol	7.95	162	482333	41.01	ng	97
30) 1,2,4-Trichlorobenzene	8.02	180	467573	36.89	ng	96
31) Naphthalene	8.10	128	1625931	39.16	ng	99
32) Benzoic acid	7.86	122	95297	18.51	ng	95
33) 4-Chloroaniline	8.16	127	420885	25.84	ng	98
34) Hexachlorobutadiene	8.21	225	240266	34.46	ng	98
35) Caprolactam	8.54	113	29212	7.90	ng	# 58
36) 4-Chloro-3-methylphenol	8.65	107	506264	41.29	ng	95
37) 2-Methylnaphthalene	8.78	142	1006412	37.75	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.96	216	509609	47.49	ng	100
40) Hexachlorocyclopentadiene	8.93	237	219894	45.42	ng	94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.07	196	325069	46.10	nq	93
43) 2,4,5-Trichlorophenol	9.13	196	363895	47.10	nq	# 83
45) 1,1'-Biphenyl	9.25	154	1406234	50.80	nq	98
46) 2-Chloronaphthalene	9.28	162	1033108	47.71	nq	98
47) 2-Nitroaniline	9.38	65	350612	48.16	nq	95
48) Acenaphthylene	9.69	152	1519558	40.62	nq	99
49) Dimethylphthalate	9.56	163	1413703	54.91	nq	100
50) 2,6-Dinitrotoluene	9.63	165	304866	54.83	nq	# 83
51) Acenaphthene	9.86	154	940903	42.07	nq	97
52) 3-Nitroaniline	9.79	138	169797	26.68	nq	96
53) 2,4-Dinitrophenol	9.92	184	271978	95.20	nq	# 78
54) Dibenzofuran	10.03	168	1284829	42.95	nq	97
55) 4-Nitrophenol	9.98	139	114210	24.92	nq	86
56) 2,4-Dinitrotoluene	10.03	165	303935	41.06	nq	# 73
57) Fluorene	10.37	166	1012223	43.98	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.16	232	258569	50.17	nq	99
59) Diethylphthalate	10.25	149	1116189	46.00	nq	100
60) 4-Chlorophenyl-phenylether	10.36	204	458960	41.51	nq	92
61) 4-Nitroaniline	10.41	138	264199	40.54	nq	93
62) Azobenzene	10.52	77	1310207	52.27	nq	97
64) 4,6-Dinitro-2-methylphenol	10.45	198	183378	45.16	nq	92
65) n-Nitrosodiphenylamine	10.49	169	912694	42.67	nq	99
66) 4-Bromophenyl-phenylether	10.85	248	307499	43.96	nq	92
67) Hexachlorobenzene	10.92	284	285903	41.36	nq	# 88
68) Atrazine	11.02	200	321877	46.89	nq	92
69) Pentachlorophenol	11.13	266	259941	73.48	nq	98
70) Phenanthrene	11.33	178	1475957	38.47	nq	99
71) Anthracene	11.39	178	1750998	45.45	nq	98
72) Carbazole	11.55	167	1582311	42.97	nq	98
73) Di-n-butylphthalate	11.87	149	1859620	46.69	nq	# 97
74) Fluoranthene	12.52	202	1627236	41.12	nq	98
76) Benzidine	12.65	184	613519	26.61	nq	96
77) Pyrene	12.75	202	1585128	39.15	nq	100
79) Butylbenzylphthalate	13.37	149	828295	50.38	nq	# 79
80) Benzo(a)anthracene	13.94	228	1371284	41.44	nq	100
81) 3,3'-Dichlorobenzidine	13.91	252	298427	25.30	nq	# 95
82) Chrysene	13.97	228	1024771	33.08	nq	99
83) Bis(2-ethylhexyl)phthalate	13.92	149	973600	43.30	nq	98
84) Di-n-octyl phthalate	14.53	149	1563624	42.71	nq	100
85) Indeno(1,2,3-cd)pyrene	16.74	276	984128	41.01	nq	# 94
87) Benzo(b)fluoranthene	14.97	252	1265428m	52.18	nq	
88) Benzo(k)fluoranthene	14.99	252	789568m	37.71	nq	
89) Benzo(a)pyrene	15.31	252	973697	46.42	nq	# 97
90) Dibenzo(a,h)anthracene	16.75	278	818924	48.31	nq	99
91) Benzo(g,h,i)perylene	17.15	276	870724	50.84	nq	# 89

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

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