

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM040517\
 Data File : BM009537.D
 Acq On : 04 Apr 2017 19:39
 Operator : SJ/MA
 Sample : I2493-17MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LSS-SP05-COMPMSD

Manual Integrations
 APPROVED

UGO
 4/5/2017 7:59:48 PM

Quant Time: Apr 05 04:44:22 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 31 12:10:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	113059	20.00	ng	-0.02
21) Naphthalene-d8	10.64	136	448194	20.00	ng	-0.02
38) Acenaphthene-d10	14.49	164	264868	20.00	ng	-0.02
63) Phenanthrene-d10	17.24	188	636968	20.00	ng	-0.01
75) Chrysene-d12	21.42	240	863302	20.00	ng	-0.01
86) Perylene-d12	23.74	264	728553	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	1055804	155.66	ng	-0.01
7) Phenol-d6	7.01	99	1411658	152.62	ng	-0.01
23) Nitrobenzene-d5	9.01	82	1171915	98.04	ng	-0.02
41) 2,4,6-Tribromophenol	15.98	330	599461	182.18	ng	-0.02
44) 2-Fluorobiphenyl	13.10	172	2211224	100.59	ng	-0.02
78) Terphenyl-d14	19.86	244	3654448	88.46	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.36	88	127677	37.43	ng	# 100
3) Pyridine	3.76	79	419607	43.53	ng	# 80
4) n-Nitrosodimethylamine	3.67	42	244753	47.70	ng	# 65
6) Aniline	7.18	93	294492	23.43	ng	91
8) 2-Chlorophenol	7.41	128	381489	55.47	ng	83
9) Benzaldehyde	6.99	77	190402	26.15	ng	85
10) Phenol	7.03	94	505267	50.41	ng	91
11) bis(2-Chloroethyl)ether	7.27	93	413126	50.29	ng	84
12) 1,3-Dichlorobenzene	7.73	146	402534	48.89	ng	# 90
13) 1,4-Dichlorobenzene	7.88	146	416973	49.01	ng	97
14) 1,2-Dichlorobenzene	8.19	146	402864	48.76	ng	# 91
15) Benzyl Alcohol	8.09	79	431524	53.36	ng	# 88
16) 2,2'-oxybis(1-Chloropropan	8.37	45	686276	48.06	ng	95
17) 2-Methylphenol	8.28	107	357029	54.44	ng	# 86
18) Hexachloroethane	8.92	117	175002	50.28	ng	90
19) n-Nitroso-di-n-propylamine	8.65	70	397861	53.36	ng	# 90
20) 3+4-Methylphenols	8.61	107	480155	53.85	ng	89
22) Acetophenone	8.67	105	654457	51.77	ng	# 89
24) Nitrobenzene	9.05	77	594324	50.75	ng	92
25) Isophorone	9.57	82	993364	54.47	ng	# 89
26) 2-Nitrophenol	9.76	139	201509	57.05	ng	# 52
27) 2,4-Dimethylphenol	9.81	122	369541	57.19	ng	# 84
28) bis(2-Chloroethoxy)methane	10.05	93	568275	50.33	ng	99
29) 2,4-Dichlorophenol	10.29	162	380131	56.28	ng	92
30) 1,2,4-Trichlorobenzene	10.50	180	429146	51.20	ng	95
31) Naphthalene	10.69	128	1199109	50.41	ng	99
32) Benzoic acid	9.93	122	182086	40.04	ng	# 73
33) 4-Chloroaniline	10.81	127	49720	5.42	ng	# 77
34) Hexachlorobutadiene	10.96	225	273342	47.63	ng	96
35) Caprolactam	11.59	113	100359m	48.66	ng	
36) 4-Chloro-3-methylphenol	11.92	107	447568	58.56	ng	# 81
37) 2-Methylnaphthalene	12.30	142	893276	54.40	ng	# 93
39) 1,2,4,5-Tetrachlorobenzene	12.67	216	492175	49.22	ng	# 100
40) Hexachlorocyclopentadiene	12.64	237	596420	103.91	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.91	196	328478	57.45	ng	98
43) 2,4,5-Trichlorophenol	12.98	196	350831	54.99	ng #	88
45) 1,1'-Biphenyl	13.32	154	1177507	53.42	ng	97
46) 2-Chloronaphthalene	13.36	162	922951	53.90	ng	94
47) 2-Nitroaniline	13.57	65	377279	59.81	ng #	72
48) Acenaphthylene	14.21	152	1483909	53.21	ng	99
49) Dimethylphthalate	13.94	163	1158815	56.98	ng	99
50) 2,6-Dinitrotoluene	14.07	165	252714	57.67	ng #	71
51) Acenaphthene	14.55	154	892973	53.07	ng	98
52) 3-Nitroaniline	14.40	138	104089	24.11	ng #	62
53) 2,4-Dinitrophenol	14.61	184	325400	108.04	ng	96
54) Dibenzofuran	14.89	168	1456900	58.55	ng	94
55) 4-Nitrophenol	14.70	139	422394	118.36	ng #	19
56) 2,4-Dinitrotoluene	14.86	165	370057	62.41	ng	96
57) Fluorene	15.54	166	1209439	54.08	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.12	232	334229	61.94	ng #	100
59) Diethylphthalate	15.31	149	1204787	58.40	ng	98
60) 4-Chlorophenyl-phenylether	15.53	204	635584	54.42	ng	100
61) 4-Nitroaniline	15.57	138	255806	54.55	ng #	31
62) Azobenzene	15.82	77	1515987	62.95	ng	85
64) 4,6-Dinitro-2-methylphenol	15.62	198	215888	53.80	ng	89
65) n-Nitrosodiphenylamine	15.75	169	964067	49.83	ng	99
66) 4-Bromophenyl-phenylether	16.43	248	390722	52.94	ng #	92
67) Hexachlorobenzene	16.54	284	424750	52.51	ng #	93
68) Atrazine	16.69	200	165342	22.72	ng	99
69) Pentachlorophenol	16.89	266	511525	103.52	ng	97
70) Phenanthrene	17.28	178	1967523	54.02	ng	99
71) Anthracene	17.37	178	2003939	54.34	ng	99
72) Carbazole	17.64	167	1597318	45.23	ng	98
73) Di-n-butylphthalate	18.19	149	2099252	58.19	ng #	98
74) Fluoranthene	19.29	202	2436080	56.47	ng	98
76) Benzidine	19.49	184	83658	3.25	ng #	95
77) Pyrene	19.66	202	2523172	51.53	ng	100
79) Butylbenzylphthalate	20.54	149	1032342	58.55	ng #	75
80) Benzo(a)anthracene	21.40	228	2729809	54.14	ng	99
81) 3,3'-Dichlorobenzidine	21.33	252	202680	10.73	ng #	96
82) Chrysene	21.46	228	2449340	50.88	ng	99
83) Bis(2-ethylhexyl)phthalate	21.31	149	1513105	57.51	ng	98
84) Di-n-octyl phthalate	22.21	149	2628363	60.03	ng #	91
85) Indeno(1,2,3-cd)pyrene	26.11	276	3010198	57.33	ng #	100
87) Benzo(b)fluoranthene	23.03	252	2747207	59.49	ng #	95
88) Benzo(k)fluoranthene	23.09	252	2719400	61.23	ng #	98
89) Benzo(a)pyrene	23.64	252	2607900	60.98	ng	100
90) Dibenzo(a,h)anthracene	26.11	278	2585973	61.73	ng #	96
91) Benzo(g,h,i)perylene	26.84	276	2409664	59.00	ng #	92

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

