

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071917\
 Data File : BF096874.D
 Acq On : 19 Jul 2017 17:11
 Operator : SJ/JU
 Sample : I4272-01MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HDD-P1LWC2MS

Manual Integrations
 APPROVED

Sohil
 7/20/2017 3:04:24 PM

Quant Time: Jul 19 19:02:34 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF071317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 18 13:33:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.59	152	93624	20.00	ng	-0.01
21) Naphthalene-d8	7.87	136	419476	20.00	ng	-0.02
38) Acenaphthene-d10	9.62	164	201461	20.00	ng	-0.01
63) Phenanthrene-d10	11.10	188	316764	20.00	ng	0.00
75) Chrysene-d12	13.72	240	191700	20.00	ng	-0.01
86) Perylene-d12	15.09	264	174483	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.20	112	884158	141.99	ng	0.00
7) Phenol-d6	6.25	99	1074229	143.76	ng	0.00
23) Nitrobenzene-d5	7.16	82	616949	91.10	ng	-0.01
41) 2,4,6-Tribromophenol	10.41	330	221599	128.86	ng	-0.01
44) 2-Fluorobiphenyl	8.95	172	1090660	85.53	ng	-0.01
78) Terphenyl-d14	12.68	244	871510	78.84	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.23	88	115925	36.182	ng	# 75
3) Pyridine	2.88	79	325053	48.274	ng	# 61
4) n-Nitrosodimethylamine	2.83	42	202822	53.863	ng	# 80
6) Aniline	6.25	93	381587	41.458	ng	# 47
8) 2-Chlorophenol	6.38	128	325222	48.415	ng	# 99
9) Benzaldehyde	6.13	77	126879	29.378	ng	# 95
10) Phenol	6.26	94	462253	54.724	ng	# 97
11) bis(2-Chloroethyl)ether	6.33	93	322013	50.694	ng	# 91
12) 1,3-Dichlorobenzene	6.53	146	324550	45.452	ng	# 99
13) 1,4-Dichlorobenzene	6.60	146	329736	45.600	ng	# 94
14) 1,2-Dichlorobenzene	6.76	146	302444	45.984	ng	# 98
15) Benzyl Alcohol	6.74	79	249910	51.085	ng	# 98
16) 2,2'-oxybis(1-Chloropropan	6.87	45	602000	45.922	ng	# 97
17) 2-Methylphenol	6.86	107	256889	49.178	ng	# 96
18) Hexachloroethane	7.10	117	120813	47.117	ng	# 81
19) n-Nitroso-di-n-propylamine	7.02	70	209844	46.614	ng	# 85
20) 3+4-Methylphenols	7.02	107	321139	50.663	ng	# 70
22) Acetophenone	7.00	105	436665	46.575	ng	# 98
24) Nitrobenzene	7.17	77	323016	46.118	ng	# 93
25) Isophorone	7.42	82	591914	46.984	ng	# 93
26) 2-Nitrophenol	7.49	139	186235	47.820	ng	# 90
27) 2,4-Dimethylphenol	7.55	122	303603	47.386	ng	# 96
28) bis(2-Chloroethoxy)methane	7.63	93	400405	47.034	ng	# 100
29) 2,4-Dichlorophenol	7.74	162	273843	47.218	ng	# 98
30) 1,2,4-Trichlorobenzene	7.82	180	252942	45.871	ng	# 98
31) Naphthalene	7.89	128	923711	46.484	ng	# 99
32) Benzoic acid	7.65	122	56032	13.683	ng	# 92
33) 4-Chloroaniline	7.95	127	302094	38.371	ng	# 97
34) Hexachlorobutadiene	8.02	225	130947	44.484	ng	# 95
35) Caprolactam	8.33	113	93833m	50.495	ng	#
36) 4-Chloro-3-methylphenol	8.45	107	296928	47.615	ng	# 89
37) 2-Methylnaphthalene	8.59	142	635263	50.144	ng	# 98
39) 1,2,4,5-Tetrachlorobenzene	8.76	216	222309	40.821	ng	# 99
40) Hexachlorocyclopentadiene	8.75	237	173764	84.534	ng	# 98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.87	196	172762	43.006	nq	98
43) 2,4,5-Trichlorophenol	8.92	196	183637	45.327	nq	95
45) 1,1'-Biphenyl	9.05	154	737790	42.701	nq	98
46) 2-Chloronaphthalene	9.07	162	580257	45.608	nq	# 93
47) 2-Nitroaniline	9.17	65	223582	51.124	nq	96
48) Acenaphthylene	9.49	152	906614	44.723	nq	99
49) Dimethylphthalate	9.36	163	831442	54.811	nq	99
50) 2,6-Dinitrotoluene	9.42	165	157018	47.788	nq	# 83
51) Acenaphthene	9.66	154	523613	43.725	nq	98
52) 3-Nitroaniline	9.58	138	144984	37.248	nq	94
53) 2,4-Dinitrophenol	9.69	184	69562	45.896	nq	# 76
54) Dibenzofuran	9.83	168	745868	44.446	nq	98
55) 4-Nitrophenol	9.76	139	242556	85.297	nq	# 64
56) 2,4-Dinitrotoluene	9.82	165	184394	43.673	nq	# 69
57) Fluorene	10.17	166	539034	43.468	nq	98
58) 2,3,4,6-Tetrachlorophenol	9.95	232	132936	47.312	nq	97
59) Diethylphthalate	10.06	149	619636	41.960	nq	100
60) 4-Chlorophenyl-phenylether	10.16	204	223534	43.009	nq	96
61) 4-Nitroaniline	10.19	138	172503	47.195	nq	92
62) Azobenzene	10.32	77	600227	46.074	nq	93
64) 4,6-Dinitro-2-methylphenol	10.23	198	83199	41.833	nq	# 76
65) n-Nitrosodiphenylamine	10.29	169	519472	45.877	nq	98
66) 4-Bromophenyl-phenylether	10.65	248	158671	49.063	nq	95
67) Hexachlorobenzene	10.72	284	168801	46.864	nq	# 87
68) Atrazine	10.82	200	154679	47.948	nq	94
69) Pentachlorophenol	10.92	266	153608	88.276	nq	95
70) Phenanthrene	11.12	178	813812	47.250	nq	99
71) Anthracene	11.17	178	824775	47.362	nq	99
72) Carbazole	11.33	167	796892	47.255	nq	98
73) Di-n-butylphthalate	11.67	149	959264	45.675	nq	98
74) Fluoranthene	12.30	202	807681	48.992	nq	99
76) Benzidine	12.43	184	395766	63.018	nq	# 93
77) Pyrene	12.53	202	816947	46.957	nq	99
79) Butylbenzylphthalate	13.16	149	389216	47.085	nq	# 80
80) Benzo(a)anthracene	13.71	228	573743	48.035	nq	99
81) 3,3'-Dichlorobenzidine	13.68	252	143741	33.465	nq	# 95
82) Chrysene	13.75	228	562099	47.803	nq	99
83) Bis(2-ethylhexyl)phthalate	13.72	149	481279	47.809	nq	98
84) Di-n-octyl phthalate	14.33	149	809559	48.650	nq	# 95
85) Indeno(1,2,3-cd)pyrene	16.38	276	503842	45.147	nq	100
87) Benzo(b)fluoranthene	14.72	252	472374	44.888	nq	99
88) Benzo(k)fluoranthene	14.74	252	464139m	46.577	nq	
89) Benzo(a)pyrene	15.04	252	462185	47.886	nq	100
90) Dibenzo(a,h)anthracene	16.39	278	409590	46.119	nq	99
91) Benzo(g,h,i)perylene	16.76	276	441142	46.386	nq	97

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

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