

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021517\
 Data File : BF092968.D
 Acq On : 15 Feb 2017 23:09
 Operator : SJ/MA
 Sample : I1816-01MSD
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 H-1MSD

Manual Integrations
 APPROVED

mohammad
 2/16/2017 11:37:57 AM

Quant Time: Feb 16 07:34:09 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 13 18:04:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	163571	20.00	ng	-0.01
21) Naphthalene-d8	8.17	136	657702	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	355450	20.00	ng	-0.01
63) Phenanthrene-d10	11.39	188	517729	20.00	ng	-0.01
75) Chrysene-d12	14.03	240	370719	20.00	ng	-0.02
86) Perylene-d12	15.46	264	323005	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.48	112	1339682	140.51	ng	0.00
7) Phenol-d6	6.52	99	1713947	139.72	ng	-0.01
23) Nitrobenzene-d5	7.45	82	1080826	91.51	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	310280	120.69	ng	-0.01
44) 2-Fluorobiphenyl	9.24	172	1888074	96.23	ng	0.00
78) Terphenyl-d14	12.98	244	1438081	91.15	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.52	88	205320	39.85	ng	# 82
3) Pyridine	3.25	79	532776	40.67	ng	86
4) n-Nitrosodimethylamine	3.22	42	307654	50.02	ng	90
6) Aniline	6.54	93	229032	13.69	ng	# 52
8) 2-Chlorophenol	6.66	128	510156	48.50	ng	92
9) Benzaldehyde	6.42	77	133873	15.23	ng	97
10) Phenol	6.53	94	681799	47.50	ng	# 79
11) bis(2-Chloroethyl)ether	6.61	93	525144	45.84	ng	93
12) 1,3-Dichlorobenzene	6.82	146	587987	47.73	ng	98
13) 1,4-Dichlorobenzene	6.89	146	566329	45.42	ng	98
14) 1,2-Dichlorobenzene	7.05	146	570070	47.81	ng	98
15) Benzyl Alcohol	7.02	79	438815	48.32	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.16	45	979989	49.24	ng	95
17) 2-Methylphenol	7.14	107	457232	50.67	ng	97
18) Hexachloroethane	7.39	117	206485	48.51	ng	95
19) n-Nitroso-di-n-propylamine	7.30	70	356419	45.64	ng	96
20) 3+4-Methylphenols	7.30	107	547586	50.75	ng	95
22) Acetophenone	7.29	105	711296	47.37	ng	95
24) Nitrobenzene	7.47	77	631329	52.06	ng	# 86
25) Isophorone	7.71	82	1108237	50.36	ng	99
26) 2-Nitrophenol	7.78	139	290578	50.41	ng	94
27) 2,4-Dimethylphenol	7.82	122	489268	48.33	ng	96
28) bis(2-Chloroethoxy)methane	7.92	93	667006	45.76	ng	99
29) 2,4-Dichlorophenol	8.03	162	487052	51.58	ng	97
30) 1,2,4-Trichlorobenzene	8.11	180	494185	48.57	ng	98
31) Naphthalene	8.19	128	1539718	46.18	ng	99
32) Benzoic acid	7.96	122	234846	41.99	ng	91
33) 4-Chloroaniline	8.24	127	71015	5.43	ng	98
34) Hexachlorobutadiene	8.30	225	261512	46.71	ng	98
35) Caprolactam	8.61	113	145113m	48.86	ng	
36) 4-Chloro-3-methylphenol	8.73	107	470852	47.83	ng	97
37) 2-Methylnaphthalene	8.88	142	1030804	48.15	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.04	216	423948	42.49	ng	99
40) Hexachlorocyclopentadiene	9.02	237	289672	64.35	ng	94

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42) 2,4,6-Trichlorophenol	9.16	196	320206	48.84	ng	94
43) 2,4,5-Trichlorophenol	9.21	196	327530	45.59	ng #	82
45) 1,1'-Biphenyl	9.33	154	1127943	43.82	ng	96
46) 2-Chloronaphthalene	9.37	162	982477	48.80	ng	94
47) 2-Nitroaniline	9.46	65	273312	40.37	ng	98
48) Acenaphthylene	9.78	152	1486910	42.75	ng	99
49) Dimethylphthalate	9.64	163	1578668	65.94	ng	98
50) 2,6-Dinitrotoluene	9.71	165	267259	51.69	ng #	82
51) Acenaphthene	9.95	154	922563	44.36	ng	96
52) 3-Nitroaniline	9.87	138	87985	14.87	ng	88
53) 2,4-Dinitrophenol	9.98	184	175803	67.62	ng #	92
54) Dibenzofuran	10.12	168	1264527	45.46	ng	95
55) 4-Nitrophenol	10.05	139	404305	94.87	ng #	65
56) 2,4-Dinitrotoluene	10.11	165	323636	47.02	ng	93
57) Fluorene	10.46	166	975972	45.61	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.24	232	237017	49.46	ng	97
59) Diethylphthalate	10.34	149	1017615	45.10	ng	98
60) 4-Chlorophenyl-phenylether	10.45	204	441911	42.99	ng #	86
61) 4-Nitroaniline	10.49	138	222784	36.76	ng	93
62) Azobenzene	10.61	77	1063968m	45.64	ng	
64) 4,6-Dinitro-2-methylphenol	10.52	198	137184	43.78	ng	85
65) n-Nitrosodiphenylamine	10.58	169	875180	52.92	ng	100
66) 4-Bromophenyl-phenylether	10.94	248	275651	50.96	ng #	80
67) Hexachlorobenzene	11.00	284	251997	47.14	ng #	77
68) Atrazine	11.10	200	285467	53.79	ng	95
69) Pentachlorophenol	11.21	266	295967	104.81	ng	97
70) Phenanthrene	11.43	178	1433959	48.34	ng	98
71) Anthracene	11.47	178	1296348	43.52	ng	98
72) Carbazole	11.63	167	1331297	46.76	ng	98
73) Di-n-butylphthalate	11.95	149	1544273	50.15	ng	98
74) Fluoranthene	12.60	202	1257230	41.09	ng	98
76) Benzidine	12.73	184	314546	19.91	ng	96
77) Pyrene	12.83	202	1253492	45.18	ng	99
79) Butylbenzylphthalate	13.45	149	596247	52.92	ng	94
80) Benzo(a)anthracene	14.02	228	936342	41.30	ng	99
81) 3,3'-Dichlorobenzidine	13.99	252	189461	23.44	ng #	95
82) Chrysene	14.05	228	885126m	41.69	ng	
83) Bis(2-ethylhexyl)phthalate	14.01	149	847079	54.98	ng #	97
84) Di-n-octyl phthalate	14.61	149	1292825	51.53	ng	99
85) Indeno(1,2,3-cd)pyrene	16.88	276	952924	57.95	ng #	94
87) Benzo(h)fluoranthene	15.05	252	899091m	42.29	ng	
88) Benzo(k)fluoranthene	15.08	252	918440m	50.03	ng	
89) Benzo(a)pyrene	15.40	252	861856	46.87	ng	97
90) Dibenzo(a,h)anthracene	16.90	278	792303	53.31	ng	98
91) Benzo(g,h,i)perylene	17.31	276	799003	53.21	ng #	92

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

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