

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032317\
 Data File : BF093897.D
 Acq On : 24 Mar 2017 00:04
 Operator : SJ/MA
 Sample : I2367-17MSD
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-CC-DD-8-9-BASEMSD

Manual Integrations
 APPROVED

mohammad
 3/24/2017 4:10:55 PM

Quant Time: Mar 24 13:36:39 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 23 02:02:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.97	152	164386	20.00	ng	-0.01
21) Naphthalene-d8	7.48	136	664668	20.00	ng	-0.01
38) Acenaphthene-d10	9.63	164	298329	20.00	ng	-0.01
63) Phenanthrene-d10	11.45	188	500895	20.00	ng	-0.01
75) Chrysene-d12	14.71	240	292075	20.00	ng	0.00
86) Perylene-d12	16.34	264	282239	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	4.52	112	1114972	114.56	ng	0.01
7) Phenol-d6	5.57	99	1417262	117.71	ng	0.00
23) Nitrobenzene-d5	6.63	82	882558	75.78	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	291505	123.65	ng	0.00
44) 2-Fluorobiphenyl	8.81	172	1450847	74.18	ng	-0.01
78) Terphenyl-d14	13.43	244	983566	75.48	ng	-0.01

Target Compounds					Qvalue
2) 1,4-Dioxane	2.40	88	159070	34.02 ng	97
3) Pyridine	2.88	79	440748	34.59 ng	98
4) n-Nitrosodimethylamine	2.83	42	207797	39.63 ng	93
6) Aniline	5.60	93	376237	22.46 ng	96
8) 2-Chlorophenol	5.74	128	446290	41.72 ng	97
9) Benzaldehyde	5.47	77	129452	15.92 ng	99
10) Phenol	5.59	94	587181	42.86 ng	98
11) bis(2-Chloroethyl)ether	5.68	93	416305	38.88 ng	98
12) 1,3-Dichlorobenzene	5.91	146	467809	38.98 ng	98
13) 1,4-Dichlorobenzene	6.00	146	470984	38.75 ng	97
14) 1,2-Dichlorobenzene	6.17	146	440781	39.38 ng	96
15) Benzyl Alcohol	6.14	79	375794	42.64 ng	99
16) 2,2'-oxybis(1-Chloropropan	6.30	45	594517	36.03 ng	99
17) 2-Methylphenol	6.28	107	369060	40.28 ng	96
18) Hexachloroethane	6.57	117	163062	38.17 ng	# 86
19) n-Nitroso-di-n-propylamine	6.46	70	305275	39.45 ng	98
20) 3+4-Methylphenols	6.46	107	444893	40.09 ng	# 81
22) Acetophenone	6.45	105	609173	41.12 ng	# 92
24) Nitrobenzene	6.65	77	454427	39.56 ng	92
25) Isophorone	6.94	82	806258	39.40 ng	95
26) 2-Nitrophenol	7.02	139	247000	45.09 ng	95
27) 2,4-Dimethylphenol	7.09	122	391687	41.50 ng	96
28) bis(2-Chloroethoxy)methane	7.20	93	528579	39.17 ng	99
29) 2,4-Dichlorophenol	7.32	162	377890	42.82 ng	97
30) 1,2,4-Trichlorobenzene	7.42	180	364380	40.09 ng	98
31) Naphthalene	7.51	128	1235964	38.67 ng	100
33) 4-Chloroaniline	7.57	127	199999	15.44 ng	# 93
34) Hexachlorobutadiene	7.66	225	177293	38.03 ng	97
35) Caprolactam	8.02	113	109726m	38.99 ng	
36) 4-Chloro-3-methylphenol	8.19	107	391927	42.50 ng	91
37) 2-Methylnaphthalene	8.35	142	850428	39.92 ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.56	216	310356	38.03 ng	99
40) Hexachlorocyclopentadiene	8.55	237	262817	68.82 ng	100
42) 2,4,6-Trichlorophenol	8.70	196	254924	44.15 ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	8.75	196	252392	40.40	ng	94
45) 1,1'-Biphenyl	8.93	154	940175	39.07	ng	98
46) 2-Chloronaphthalene	8.95	162	739630	39.27	ng	93
47) 2-Nitroaniline	9.07	65	228163	40.83	ng	88
48) Acenaphthylene	9.46	152	1192398	38.09	ng	99
49) Dimethylphthalate	9.32	163	1054651	49.73	ng	99
50) 2,6-Dinitrotoluene	9.38	165	202534	41.66	ng	95
51) Acenaphthene	9.67	154	698741	38.25	ng	99
52) 3-Nitroaniline	9.57	138	154739	27.11	ng	96
53) 2,4-Dinitrophenol	9.70	184	14186	9.80	ng	# 74
54) Dibenzofuran	9.88	168	1020793	41.05	ng	99
55) 4-Nitrophenol	9.80	139	294173	65.80	ng	# 68
56) 2,4-Dinitrotoluene	9.87	165	254825	41.73	ng	# 81
57) Fluorene	10.30	166	761146	36.91	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.03	232	184925	43.14	ng	99
59) Diethylphthalate	10.19	149	859591	41.01	ng	100
60) 4-Chlorophenyl-phenylether	10.31	204	326307	37.14	ng	93
61) 4-Nitroaniline	10.33	138	191013	34.87	ng	96
62) Azobenzene	10.50	77	810386	40.87	ng	96
64) 4,6-Dinitro-2-methylphenol	10.37	198	84315	27.49	ng	# 76
65) n-Nitrosodiphenylamine	10.46	169	712505	41.13	ng	99
66) 4-Bromophenyl-phenylether	10.91	248	199350	40.47	ng	99
67) Hexachlorobenzene	10.98	284	203413	40.79	ng	# 87
68) Atrazine	11.13	200	207357	39.97	ng	97
69) Pentachlorophenol	11.23	266	205778	66.29	ng	97
70) Phenanthrene	11.49	178	1086930	39.01	ng	98
71) Anthracene	11.55	178	1103283	38.46	ng	98
72) Carbazole	11.75	167	986110	33.52	ng	98
73) Di-n-butylphthalate	12.20	149	1253438	40.97	ng	99
74) Fluoranthene	12.95	202	1000482	35.07	ng	98
76) Benzidine	13.11	184	295053	25.52	ng	# 94
77) Pyrene	13.22	202	966072	45.58	ng	99
79) Butylbenzylphthalate	14.04	149	427551	47.34	ng	# 86
80) Benzo(a)anthracene	14.69	228	658585	38.67	ng	99
81) 3,3'-Dichlorobenzidine	14.66	252	139183	22.86	ng	# 97
82) Chrysene	14.74	228	593329	37.54	ng	99
83) Bis(2-ethylhexyl)phthalate	14.76	149	556607	45.17	ng	# 98
84) Di-n-octyl phthalate	15.51	149	825643	40.11	ng	98
85) Indeno(1,2,3-cd)pyrene	17.74	276	679980	49.68	ng	99
87) Benzo(b)fluoranthene	15.93	252	637319	36.84	ng	98
88) Benzo(k)fluoranthene	15.96	252	568796m	33.60	ng	
89) Benzo(a)pyrene	16.29	252	596097	37.42	ng	100
90) Dibenzo(a,h)anthracene	17.76	278	569286	42.19	ng	98
91) Benzo(g,h,i)perylene	18.15	276	555696	40.87	ng	99

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

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