

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030217\
 Data File : BF093334.D
 Acq On : 2 Mar 2017 21:28
 Operator : SJ/MA
 Sample : PB97296BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB97296BS

Manual Integrations
 APPROVED

mohammad
 3/3/2017 5:16:48 PM

Quant Time: Mar 03 01:18:00 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 00:22:20 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	153748	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	617684	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	293564	20.00	ng	-0.01
63) Phenanthrene-d10	11.33	188	534526	20.00	ng	-0.01
75) Chrysene-d12	13.97	240	407008	20.00	ng	-0.01
86) Perylene-d12	15.39	264	326616	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.40	112	1291506	144.11	ng	0.01
7) Phenol-d6	6.45	99	1553297	134.71	ng	0.00
23) Nitrobenzene-d5	7.37	82	978428	88.21	ng	-0.01
41) 2,4,6-Tribromophenol	10.65	330	273548	128.84	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1490031	91.95	ng	0.00
78) Terphenyl-d14	12.92	244	1327988	76.66	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.40	88	182990	37.79	ng	# 81
3) Pyridine	3.09	79	466901	37.92	ng	89
4) n-Nitrosodimethylamine	3.06	42	269101	46.54	ng	86
6) Aniline	6.46	93	388116	24.68	ng	# 78
8) 2-Chlorophenol	6.58	128	428830	43.37	ng	89
9) Benzaldehyde	6.34	77	131743	15.95	ng	86
10) Phenol	6.46	94	599842	44.46	ng	84
11) bis(2-Chloroethyl)ether	6.53	93	443902	41.22	ng	92
12) 1,3-Dichlorobenzene	6.73	146	462359m	39.93	ng	
13) 1,4-Dichlorobenzene	6.81	146	482611	41.18	ng	99
14) 1,2-Dichlorobenzene	6.97	146	458012	40.87	ng	96
15) Benzyl Alcohol	6.96	79	354721	41.55	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.08	45	765085	40.90	ng	91
17) 2-Methylphenol	7.07	107	364029	42.92	ng	95
18) Hexachloroethane	7.30	117	167035	41.75	ng	# 88
19) n-Nitroso-di-n-propylamine	7.22	70	323777	44.11	ng	97
20) 3+4-Methylphenols	7.23	107	495477	48.86	ng	# 76
22) Acetophenone	7.22	105	616697	43.73	ng	# 97
24) Nitrobenzene	7.39	77	526302	46.21	ng	99
25) Isophorone	7.63	82	926974	44.85	ng	94
26) 2-Nitrophenol	7.71	139	256816	47.43	ng	95
27) 2,4-Dimethylphenol	7.76	122	432179	45.45	ng	97
28) bis(2-Chloroethoxy)methane	7.85	93	606176	44.28	ng	99
29) 2,4-Dichlorophenol	7.96	162	380524	42.91	ng	97
30) 1,2,4-Trichlorobenzene	8.03	180	381563	39.93	ng	95
31) Naphthalene	8.11	128	1255295	40.09	ng	99
32) Benzoic acid	7.92	122	176299	35.01	ng	94
33) 4-Chloroaniline	8.17	127	244919	19.94	ng	95
34) Hexachlorobutadiene	8.22	225	199847	38.01	ng	98
35) Caprolactam	8.56	113	135535m	48.59	ng	
36) 4-Chloro-3-methylphenol	8.67	107	398858	43.14	ng	99
37) 2-Methylnaphthalene	8.81	142	841657	41.86	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.98	216	384279	46.63	ng	99
40) Hexachlorocyclopentadiene	8.96	237	216432	58.21	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.09	196	249890	46.15	nq	91
43) 2,4,5-Trichlorophenol	9.15	196	244445	41.20	nq	# 86
45) 1,1'-Biphenyl	9.28	154	1019627	47.97	nq	97
46) 2-Chloronaphthalene	9.30	162	764293	45.96	nq	97
47) 2-Nitroaniline	9.40	65	263675	47.16	nq	92
48) Acenaphthylene	9.71	152	1153191	40.14	nq	98
49) Dimethylphthalate	9.58	163	971740	49.14	nq	100
50) 2,6-Dinitrotoluene	9.65	165	218937	51.27	nq	92
51) Acenaphthene	9.88	154	673759	39.23	nq	97
52) 3-Nitroaniline	9.82	138	132867	27.19	nq	80
53) 2,4-Dinitrophenol	9.94	184	229631	104.19	nq	# 63
54) Dibenzofuran	10.05	168	1026088	44.67	nq	99
55) 4-Nitrophenol	10.01	139	220515	62.65	nq	95
56) 2,4-Dinitrotoluene	10.06	165	267743	47.09	nq	# 76
57) Fluorene	10.40	166	788287	44.60	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.19	232	207492	52.43	nq	92
59) Diethylphthalate	10.28	149	955150	51.26	nq	97
60) 4-Chlorophenyl-phenylether	10.40	204	392759	46.26	nq	# 75
61) 4-Nitroaniline	10.44	138	245501	49.05	nq	91
62) Azobenzene	10.56	77	1159172	60.21	nq	94
64) 4,6-Dinitro-2-methylphenol	10.48	198	150296	46.29	nq	80
65) n-Nitrosodiphenylamine	10.52	169	758393	44.41	nq	100
66) 4-Bromophenyl-phenylether	10.88	248	227279	40.70	nq	# 89
67) Hexachlorobenzene	10.96	284	228873	41.47	nq	# 89
68) Atrazine	11.05	200	231986	42.34	nq	99
69) Pentachlorophenol	11.15	266	180498	64.86	nq	96
70) Phenanthrene	11.37	178	1283918	41.93	nq	98
71) Anthracene	11.41	178	1189613	38.68	nq	99
72) Carbazole	11.57	167	1076266	36.61	nq	99
73) Di-n-butylphthalate	11.90	149	1471952	46.30	nq	# 95
74) Fluoranthene	12.54	202	1244095	39.38	nq	97
76) Benzidine	12.67	184	405782	23.40	nq	98
77) Pyrene	12.77	202	1263466	41.48	nq	99
79) Butylbenzylphthalate	13.39	149	583443	47.17	nq	98
80) Benzo(a)anthracene	13.96	228	1027592	41.28	nq	99
81) 3,3'-Dichlorobenzidine	13.93	252	235379	26.53	nq	# 95
82) Chrysene	14.00	228	969085	41.58	nq	99
83) Bis(2-ethylhexyl)phthalate	13.95	149	867884	51.31	nq	# 96
84) Di-n-octyl phthalate	14.56	149	1348441	48.95	nq	97
85) Indeno(1,2,3-cd)pyrene	16.79	276	708586	39.25	nq	# 94
87) Benzo(b)fluoranthene	14.99	252	857527m	39.89	nq	
88) Benzo(k)fluoranthene	15.03	252	846617m	45.61	nq	
89) Benzo(a)pyrene	15.33	252	781799	42.04	nq	# 96
90) Dibenzo(a,h)anthracene	16.79	278	600850	39.98	nq	# 94
91) Benzo(g,h,i)perylene	17.21	276	607509	40.01	nq	# 92

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

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