

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100317\
 Data File : BF099026.D
 Acq On : 3 Oct 2017 12:48
 Operator : SJ/JU
 Sample : PB102810BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB102810BS

Manual Integrations
 APPROVED

Sohil
 10/4/2017 7:21:14 PM

Quant Time: Oct 03 18:41:51 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 03 17:12:12 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	136289	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	577808	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	263466	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	468264	20.00	ng	0.00
75) Chrysene-d12	14.04	240	326496	20.00	ng	0.00
86) Perylene-d12	15.52	264	267989	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.51	112	1080640	126.22	ng	0.02
7) Phenol-d6	6.51	99	1325564	125.51	ng	0.00
23) Nitrobenzene-d5	7.44	82	819783	90.99	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	291172	135.06	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1485026	94.10	ng	0.00
78) Terphenyl-d14	12.99	244	1266376	88.21	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.76	88	148065	36.205	ng	97
3) Pyridine	3.52	79	428666	38.746	ng	97
4) n-Nitrosodimethylamine	3.48	42	180785	38.535	ng	85
6) Aniline	6.54	93	267240	18.748	ng	100
8) 2-Chlorophenol	6.66	128	395948	40.839	ng	97
9) Benzaldehyde	6.43	77	198480	32.397	ng	95
10) Phenol	6.52	94	486876	41.220	ng	95
11) bis(2-Chloroethyl)ether	6.61	93	376010	40.948	ng	99
12) 1,3-Dichlorobenzene	6.82	146	416211	40.991	ng	99
13) 1,4-Dichlorobenzene	6.89	146	419011	40.559	ng	98
14) 1,2-Dichlorobenzene	7.05	146	392609	41.129	ng	98
15) Benzyl Alcohol	7.02	79	335057	43.245	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.15	45	573667	40.098	ng	98
17) 2-Methylphenol	7.13	107	342900	43.224	ng	98
18) Hexachloroethane	7.39	117	149474	40.253	ng	98
19) n-Nitroso-di-n-propylamine	7.29	70	263681	39.104	ng	98
20) 3+4-Methylphenols	7.28	107	404416	40.297	ng	# 74
22) Acetophenone	7.29	105	531184	37.944	ng	# 97
24) Nitrobenzene	7.46	77	407620	39.882	ng	98
25) Isophorone	7.70	82	767739	41.214	ng	98
26) 2-Nitrophenol	7.77	139	224279	45.633	ng	90
27) 2,4-Dimethylphenol	7.80	122	385562	41.540	ng	98
28) bis(2-Chloroethoxy)methane	7.90	93	490231	42.229	ng	99
29) 2,4-Dichlorophenol	8.02	162	348032	43.673	ng	96
30) 1,2,4-Trichlorobenzene	8.10	180	335765	42.127	ng	99
31) Naphthalene	8.18	128	1144218	42.164	ng	100
32) Benzoic acid	7.93	122	229925	30.157	ng	99
33) 4-Chloroaniline	8.22	127	102679	9.061	ng	96
34) Hexachlorobutadiene	8.29	225	164530	40.077	ng	99
35) Caprolactam	8.60	113	127159m	49.744	ng	
36) 4-Chloro-3-methylphenol	8.71	107	367279	41.004	ng	96
37) 2-Methylnaphthalene	8.87	142	799132	45.298	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.03	216	305908	38.933	ng	99
40) Hexachlorocyclopentadiene	9.02	237	291225	81.104	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.15	196	229496	41.229	nq	99
43) 2,4,5-Trichlorophenol	9.19	196	238032	42.837	nq	98
45) 1,1'-Biphenyl	9.33	154	916117	40.459	nq	99
46) 2-Chloronaphthalene	9.36	162	719561	42.324	nq	97
47) 2-Nitroaniline	9.46	65	219405	42.147	nq	96
48) Acenaphthylene	9.78	152	1111793	42.719	nq	100
49) Dimethylphthalate	9.63	163	859844	43.241	nq	100
50) 2,6-Dinitrotoluene	9.70	165	190446	43.992	nq	95
51) Acenaphthene	9.95	154	667566	40.493	nq	100
52) 3-Nitroaniline	9.86	138	88189	17.471	nq	99
53) 2,4-Dinitrophenol	9.98	184	169583	75.588	nq	89
54) Dibenzofuran	10.12	168	993193	45.247	nq	99
55) 4-Nitrophenol	10.03	139	333964	78.245	nq	88
56) 2,4-Dinitrotoluene	10.10	165	261328	46.513	nq	96
57) Fluorene	10.46	166	759138	43.028	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.24	232	179890	46.865	nq	98
59) Diethylphthalate	10.33	149	821518	42.280	nq	98
60) 4-Chlorophenyl-phenylether	10.45	204	336878	42.507	nq	99
61) 4-Nitroaniline	10.49	138	211665	43.464	nq	92
62) Azobenzene	10.62	77	791125	44.282	nq	93
64) 4,6-Dinitro-2-methylphenol	10.52	198	109949	38.592	nq	83
65) n-Nitrosodiphenylamine	10.57	169	684678	42.206	nq	99
66) 4-Bromophenyl-phenylether	10.95	248	200686	43.784	nq	# 89
67) Hexachlorobenzene	11.01	284	200416	40.940	nq	99
68) Atrazine	11.10	200	220653	45.209	nq	99
69) Pentachlorophenol	11.20	266	212598	69.780	nq	98
70) Phenanthrene	11.43	178	1065134	42.495	nq	99
71) Anthracene	11.48	178	1099667	42.878	nq	100
72) Carbazole	11.63	167	1048561	41.751	nq	99
73) Di-n-butylphthalate	11.95	149	1270054	42.014	nq	100
74) Fluoranthene	12.62	202	1087527	42.848	nq	99
76) Benzidine	12.73	184	278621	23.072	nq	99
77) Pyrene	12.85	202	1105818	44.417	nq	100
79) Butylbenzylphthalate	13.45	149	525459	44.908	nq	97
80) Benzo(a)anthracene	14.03	228	854445	43.219	nq	100
81) 3,3'-Dichlorobenzidine	13.99	252	102685	14.168	nq	98
82) Chrysene	14.07	228	802893	42.657	nq	99
83) Bis(2-ethylhexyl)phthalate	14.00	149	702999	43.634	nq	# 99
84) Di-n-octyl phthalate	14.62	149	1116092	43.021	nq	97
85) Indeno(1,2,3-cd)pyrene	17.00	276	678011	45.031	nq	96
87) Benzo(b)fluoranthene	15.08	252	749989m	45.517	nq	
88) Benzo(k)fluoranthene	15.12	252	664298	40.818	nq	99
89) Benzo(a)pyrene	15.46	252	668701	44.212	nq	98
90) Dibenzo(a,h)anthracene	17.02	278	558996	46.182	nq	96
91) Benzo(g,h,i)perylene	17.46	276	575480	48.098	nq	95

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

