

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091217\
 Data File : BF098454.D
 Acq On : 12 Sep 2017 17:03
 Operator : SJ/JU
 Sample : PB102265BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB102265BS

Manual Integrations
 APPROVED

Sohil
 9/18/2017 5:10:53 PM

Quant Time: Sep 13 01:34:00 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 02:14:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	221563	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	911590	20.00	ng	0.00
38) Acenaphthene-d10	9.70	164	393456	20.00	ng	0.00
63) Phenanthrene-d10	11.18	188	622725	20.00	ng	0.00
75) Chrysene-d12	13.82	240	390857	20.00	ng	0.00
86) Perylene-d12	15.21	264	345690	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.29	112	1845926	123.18	ng	0.01
7) Phenol-d6	6.32	99	2233324	117.38	ng	0.00
23) Nitrobenzene-d5	7.24	82	1430630	87.49	ng	0.00
41) 2,4,6-Tribromophenol	10.49	330	352271	134.15	ng	0.00
44) 2-Fluorobiphenyl	9.03	172	2215234	95.43	ng	0.00
78) Terphenyl-d14	12.76	244	1496729	82.66	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.38	88	260852	33.726	ng	90
3) Pyridine	3.06	79	768926	37.435	ng	87
4) n-Nitrosodimethylamine	3.03	42	383979	38.131	ng	100
6) Aniline	6.33	93	543174	22.756	ng	# 1
8) 2-Chlorophenol	6.46	128	666358	40.439	ng	96
9) Benzaldehyde	6.21	77	432185	34.748	ng	93
10) Phenol	6.33	94	834284	37.751	ng	79
11) bis(2-Chloroethyl)ether	6.41	93	663361	39.812	ng	95
12) 1,3-Dichlorobenzene	6.60	146	677284	40.835	ng	97
13) 1,4-Dichlorobenzene	6.68	146	682595	40.179	ng	95
14) 1,2-Dichlorobenzene	6.83	146	630249	40.719	ng	99
15) Benzyl Alcohol	6.82	79	523072	41.308	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.95	45	1063955	38.808	ng	54
17) 2-Methylphenol	6.94	107	557404	41.483	ng	# 89
18) Hexachloroethane	7.17	117	246238	37.847	ng	# 82
19) n-Nitroso-di-n-propylamine	7.09	70	475519	39.733	ng	99
20) 3+4-Methylphenols	7.10	107	719189	42.413	ng	# 65
22) Acetophenone	7.08	105	941186	39.947	ng	# 92
24) Nitrobenzene	7.26	77	736682	39.553	ng	95
25) Isophorone	7.50	82	1314798	39.742	ng	98
26) 2-Nitrophenol	7.57	139	339827	45.189	ng	# 81
27) 2,4-Dimethylphenol	7.62	122	592380	41.322	ng	95
28) bis(2-Chloroethoxy)methane	7.70	93	844108	41.814	ng	99
29) 2,4-Dichlorophenol	7.82	162	526447	44.155	ng	94
30) 1,2,4-Trichlorobenzene	7.89	180	565497	43.602	ng	99
31) Naphthalene	7.97	128	1871968	41.539	ng	99
32) Benzoic acid	7.77	122	304455	34.372	ng	99
33) 4-Chloroaniline	8.03	127	285034	15.563	ng	96
34) Hexachlorobutadiene	8.09	225	286210	42.987	ng	99
35) Caprolactam	8.41	113	183817m	44.709	ng	
36) 4-Chloro-3-methylphenol	8.53	107	597284	40.395	ng	# 83
37) 2-Methylnaphthalene	8.66	142	1198781	43.916	ng	100
39) 1,2,4,5-Tetrachlorobenzene	8.83	216	496425	40.505	ng	99
40) Hexachlorocyclopentadiene	8.81	237	143773	44.966	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.95	196	328651	45.653	ng	92
43) 2,4,5-Trichlorophenol	8.99	196	342877	45.569	ng	95
45) 1,1'-Biphenyl	9.13	154	1404421	39.879	ng	97
46) 2-Chloronaphthalene	9.15	162	1070440	42.561	ng	94
47) 2-Nitroaniline	9.26	65	417817	43.282	ng	97
48) Acenaphthylene	9.56	152	1511268	40.413	ng	99
49) Dimethylphthalate	9.43	163	1300282	42.693	ng	99
50) 2,6-Dinitrotoluene	9.50	165	274590	43.891	ng #	74
51) Acenaphthene	9.73	154	969734	40.795	ng	98
52) 3-Nitroaniline	9.66	138	176968	23.433	ng	98
53) 2,4-Dinitrophenol	9.78	184	38313	20.100	ng #	74
54) Dibenzofuran	9.90	168	1335615	42.650	ng	98
55) 4-Nitrophenol	9.85	139	350524	80.787	ng #	45
56) 2,4-Dinitrotoluene	9.90	165	308712	40.589	ng #	48
57) Fluorene	10.25	166	1020817	39.383	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.03	232	240273	47.515	ng	97
59) Diethylphthalate	10.13	149	1212338	41.855	ng	99
60) 4-Chlorophenyl-phenylether	10.24	204	493259	41.200	ng	91
61) 4-Nitroaniline	10.28	138	251958	33.036	ng	88
62) Azobenzene	10.40	77	1251646	40.093	ng	97
64) 4,6-Dinitro-2-methylphenol	10.32	198	38980	14.114	ng	78
65) n-Nitrosodiphenylamine	10.36	169	958957	47.426	ng	99
66) 4-Bromophenyl-phenylether	10.73	248	291424	49.308	ng	96
67) Hexachlorobenzene	10.79	284	268003	44.705	ng #	82
68) Atrazine	10.89	200	295264	47.432	ng	97
69) Pentachlorophenol	10.99	266	228970	85.737	ng	97
70) Phenanthrene	11.20	178	1426547	43.212	ng	99
71) Anthracene	11.26	178	1447495	42.707	ng	99
72) Carbazole	11.42	167	1243389	39.363	ng	99
73) Di-n-butylphthalate	11.74	149	1711782	45.229	ng	100
74) Fluoranthene	12.39	202	1277624	38.660	ng	98
76) Benzidine	12.52	184	832912	48.151	ng #	93
77) Pyrene	12.62	202	1293200	41.942	ng	99
79) Butylbenzylphthalate	13.24	149	612396	45.782	ng #	83
80) Benzo(a)anthracene	13.80	228	985442	43.004	ng	99
81) 3,3'-Dichlorobenzidine	13.77	252	371156	41.678	ng #	97
82) Chrysene	13.84	228	983024	42.439	ng	99
83) Bis(2-ethylhexyl)phthalate	13.80	149	773842	46.096	ng #	98
84) Di-n-octyl phthalate	14.40	149	1380617	47.934	ng	98
85) Indeno(1,2,3-cd)pyrene	16.56	276	709755	43.674	ng	98
87) Benzo(b)fluoranthene	14.82	252	955914	43.978	ng	98
88) Benzo(k)fluoranthene	14.84	252	866321	42.693	ng	98
89) Benzo(a)pyrene	15.16	252	848544	43.820	ng	99
90) Dibenzo(a,h)anthracene	16.56	278	594858	42.229	ng	96
91) Benzo(g,h,i)perylene	16.96	276	567425	40.049	ng	98

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

