

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092217\
 Data File : BF098704.D
 Acq On : 22 Sep 2017 18:56
 Operator : SJ/JU
 Sample : PB102532BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB102532BS

Manual Integrations
 APPROVED

Sohil
 9/25/2017 6:21:12 PM

Quant Time: Sep 23 01:51:26 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 10:38:37 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	151073	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	641842	20.00	ng	0.00
38) Acenaphthene-d10	9.97	164	293018	20.00	ng	0.00
63) Phenanthrene-d10	11.46	188	526201	20.00	ng	0.00
75) Chrysene-d12	14.09	240	419491	20.00	ng	0.00
86) Perylene-d12	15.59	264	358004	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.56	112	1216617	122.29	ng	0.02
7) Phenol-d6	6.56	99	1469072	121.94	ng	0.01
23) Nitrobenzene-d5	7.49	82	893504	94.43	ng	0.00
41) 2,4,6-Tribromophenol	10.76	330	366927	136.93	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	1595444	86.64	ng	0.00
78) Terphenyl-d14	13.04	244	1502149	76.89	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.82	88	159163	33.870	ng	98
3) Pyridine	3.59	79	475808	38.099	ng	99
4) n-Nitrosodimethylamine	3.54	42	214676	39.229	ng	97
6) Aniline	6.60	93	413652	25.919	ng	99
8) 2-Chlorophenol	6.72	128	419222	38.792	ng	98
9) Benzaldehyde	6.48	77	122372	18.109	ng	97
10) Phenol	6.57	94	509014	38.281	ng	97
11) bis(2-Chloroethyl)ether	6.66	93	409362	39.962	ng	99
12) 1,3-Dichlorobenzene	6.87	146	440955	38.629	ng	98
13) 1,4-Dichlorobenzene	6.95	146	450846	39.023	ng	100
14) 1,2-Dichlorobenzene	7.10	146	420546	39.031	ng	99
15) Benzyl Alcohol	7.07	79	367608	42.287	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.20	45	642701	39.371	ng	100
17) 2-Methylphenol	7.17	107	364578	41.401	ng	97
18) Hexachloroethane	7.44	117	160640	40.389	ng	94
19) n-Nitroso-di-n-propylamine	7.34	70	286291	38.360	ng	98
20) 3+4-Methylphenols	7.33	107	446852	39.357	ng	# 85
22) Acetophenone	7.34	105	580054	36.533	ng	# 96
24) Nitrobenzene	7.51	77	440437	40.301	ng	100
25) Isophorone	7.75	82	794880	38.518	ng	99
26) 2-Nitrophenol	7.82	139	223684	46.414	ng	99
27) 2,4-Dimethylphenol	7.86	122	395862	39.060	ng	99
28) bis(2-Chloroethoxy)methane	7.96	93	511563	39.837	ng	100
29) 2,4-Dichlorophenol	8.06	162	350253	40.176	ng	98
30) 1,2,4-Trichlorobenzene	8.15	180	348306	39.611	ng	100
31) Naphthalene	8.23	128	1194919	39.225	ng	100
32) Benzoic acid	7.98	122	279008	42.417	ng	97
33) 4-Chloroaniline	8.27	127	224516	17.727	ng	99
34) Hexachlorobutadiene	8.35	225	184924	39.036	ng	97
35) Caprolactam	8.65	113	116572m	42.881	ng	
36) 4-Chloro-3-methylphenol	8.76	107	382698	38.457	ng	94
37) 2-Methylnaphthalene	8.92	142	824432	42.148	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.09	216	304726	35.163	ng	99
40) Hexachlorocyclopentadiene	9.07	237	383968	101.959	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.20	196	241154	41.292	nq	98
43) 2,4,5-Trichlorophenol	9.24	196	246750	41.204	nq	95
45) 1,1'-Biphenyl	9.39	154	930599	36.967	nq	99
46) 2-Chloronaphthalene	9.42	162	739929	39.197	nq	97
47) 2-Nitroaniline	9.50	65	238012	44.906	nq	95
48) Acenaphthylene	9.83	152	1131526	38.972	nq	99
49) Dimethylphthalate	9.69	163	864529	39.343	nq	99
50) 2,6-Dinitrotoluene	9.75	165	192167	44.318	nq	98
51) Acenaphthene	10.00	154	763261	42.352	nq	97
52) 3-Nitroaniline	9.92	138	163175	31.568	nq	97
53) 2,4-Dinitrophenol	10.02	184	146163	77.644	nq	98
54) Dibenzofuran	10.17	168	1023100	41.575	nq	99
55) 4-Nitrophenol	10.07	139	374189	83.190	nq	94
56) 2,4-Dinitrotoluene	10.15	165	264250	45.364	nq	96
57) Fluorene	10.52	166	784384	39.656	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.29	232	197957	46.463	nq	99
59) Diethylphthalate	10.39	149	855646	39.735	nq	99
60) 4-Chlorophenyl-phenylether	10.50	204	345416	39.863	nq	99
61) 4-Nitroaniline	10.53	138	222509	44.884	nq	97
62) Azobenzene	10.67	77	840837	41.751	nq	96
64) 4,6-Dinitro-2-methylphenol	10.56	198	108312	42.528	nq	93
65) n-Nitrosodiphenylamine	10.63	169	708901	39.139	nq	99
66) 4-Bromophenyl-phenylether	11.00	248	215673	40.592	nq	91
67) Hexachlorobenzene	11.06	284	230856	38.131	nq	93
68) Atrazine	11.15	200	223479	41.018	nq	98
69) Pentachlorophenol	11.26	266	277099	80.077	nq	99
70) Phenanthrene	11.48	178	1136487	39.814	nq	99
71) Anthracene	11.53	178	1181318	40.587	nq	100
72) Carbazole	11.68	167	1103196	38.790	nq	99
73) Di-n-butylphthalate	12.01	149	1326442	40.155	nq	99
74) Fluoranthene	12.67	202	1179742	41.471	nq	99
76) Benzidine	12.79	184	349603	20.524	nq	98
77) Pyrene	12.90	202	1211438	38.898	nq	100
79) Butylbenzylphthalate	13.51	149	607256	47.132	nq	95
80) Benzo(a)anthracene	14.08	228	1021710	40.435	nq	100
81) 3,3'-Dichlorobenzidine	14.04	252	293407	32.539	nq	97
82) Chrysene	14.12	228	958206	39.268	nq	99
83) Bis(2-ethylhexyl)phthalate	14.06	149	775506	43.514	nq	98
84) Di-n-octyl phthalate	14.68	149	1239391	46.306	nq	99
85) Indeno(1,2,3-cd)pyrene	17.10	276	803895	42.469	nq	100
87) Benzo(b)fluoranthene	15.14	252	903084	43.496	nq	98
88) Benzo(k)fluoranthene	15.18	252	764637	36.853	nq	98
89) Benzo(a)pyrene	15.52	252	771383	40.910	nq	# 96
90) Dibenzo(a,h)anthracene	17.12	278	665346	41.723	nq	99
91) Benzo(g,h,i)perylene	17.56	276	674966	42.106	nq	98

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

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