

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
 Data File : BF099212.D
 Acq On : 8 Oct 2017 1:02
 Operator : SJ/JU
 Sample : PB103045BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB103045BS

Manual Integrations
 APPROVED

Sohil
 10/9/2017 6:35:03 PM

Quant Time: Oct 08 05:21:07 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Oct 08 03:54:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	121817	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	476173	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	185267	20.00	ng	0.00
63) Phenanthrene-d10	11.39	188	285812	20.00	ng	0.00
75) Chrysene-d12	14.03	240	208055	20.00	ng	0.00
86) Perylene-d12	15.51	264	165705	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	896013	117.09	ng	0.02
7) Phenol-d6	6.50	99	1064657	112.78	ng	0.00
23) Nitrobenzene-d5	7.43	82	618149	83.25	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	164222	108.33	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1072370	96.63	ng	0.00
78) Terphenyl-d14	12.97	244	751140	82.11	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.71	88	120266	32.901	ng	96
3) Pyridine	3.49	79	348061	35.198	ng	95
4) n-Nitrosodimethylamine	3.44	42	138046	32.921	ng	81
6) Aniline	6.53	93	391914	30.761	ng	98
8) 2-Chlorophenol	6.66	128	318045	36.701	ng	95
9) Benzaldehyde	6.42	77	130369	23.808	ng	90
10) Phenol	6.52	94	393909	37.311	ng	99
11) bis(2-Chloroethyl)ether	6.60	93	306564	37.352	ng	98
12) 1,3-Dichlorobenzene	6.81	146	342738	37.765	ng	98
13) 1,4-Dichlorobenzene	6.89	146	347251	37.606	ng	97
14) 1,2-Dichlorobenzene	7.04	146	322703	37.822	ng	98
15) Benzyl Alcohol	7.01	79	241744	34.908	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.14	45	443823	34.708	ng	96
17) 2-Methylphenol	7.12	107	261853	36.929	ng	97
18) Hexachloroethane	7.37	117	95250	28.698	ng	94
19) n-Nitroso-di-n-propylamine	7.28	70	201754	33.475	ng	95
20) 3+4-Methylphenols	7.27	107	307175	34.244	ng	# 76
22) Acetophenone	7.28	105	413602	35.851	ng	# 97
24) Nitrobenzene	7.45	77	308923	36.677	ng	97
25) Isophorone	7.69	82	571059	37.199	ng	99
26) 2-Nitrophenol	7.77	139	131895	32.564	ng	# 84
27) 2,4-Dimethylphenol	7.80	122	275470	36.013	ng	95
28) bis(2-Chloroethoxy)methane	7.90	93	375599	39.261	ng	100
29) 2,4-Dichlorophenol	8.01	162	249606	38.007	ng	97
30) 1,2,4-Trichlorobenzene	8.09	180	259928	39.573	ng	98
31) Naphthalene	8.17	128	847556	37.898	ng	100
32) Benzoic acid	7.92	122	112277	17.869	ng	95
33) 4-Chloroaniline	8.22	127	263137	28.175	ng	96
34) Hexachlorobutadiene	8.28	225	130153	38.471	ng	99
35) Caprolactam	8.60	113	79469m	37.723	ng	
36) 4-Chloro-3-methylphenol	8.70	107	247189	33.487	ng	92
37) 2-Methylnaphthalene	8.86	142	559939	38.514	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.03	216	220886	39.978	ng	98
40) Hexachlorocyclopentadiene	9.01	237	15831	6.270	ng	99

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42) 2,4,6-Trichlorophenol	9.14	196	148932	38.049	ng	98
43) 2,4,5-Trichlorophenol	9.19	196	150043	38.400	ng	# 93
45) 1,1'-Biphenyl	9.33	154	629998	39.566	ng	98
46) 2-Chloronaphthalene	9.35	162	485657	40.624	ng	97
47) 2-Nitroaniline	9.45	65	146856	40.118	ng	90
48) Acenaphthylene	9.77	152	707643	38.667	ng	100
49) Dimethylphthalate	9.62	163	584236	41.782	ng	100
50) 2,6-Dinitrotoluene	9.69	165	108346	35.591	ng	# 80
51) Acenaphthene	9.94	154	417319	35.998	ng	98
52) 3-Nitroaniline	9.86	138	150579	42.422	ng	# 89
53) 2,4-Dinitrophenol	9.97	184	10902	12.205	ng	95
54) Dibenzofuran	10.11	168	619930	40.163	ng	99
55) 4-Nitrophenol	10.02	139	164395	54.773	ng	90
56) 2,4-Dinitrotoluene	10.09	165	125449	31.753	ng	96
57) Fluorene	10.45	166	465035	37.484	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.23	232	102343	37.916	ng	98
59) Diethylphthalate	10.32	149	543738	39.795	ng	98
60) 4-Chlorophenyl-phenylether	10.45	204	217289	38.990	ng	93
61) 4-Nitroaniline	10.47	138	143461	41.893	ng	89
62) Azobenzene	10.60	77	456525	36.339	ng	93
64) 4,6-Dinitro-2-methylphenol	10.50	198	11151	6.412	ng	85
65) n-Nitrosodiphenylamine	10.56	169	416332	42.048	ng	100
66) 4-Bromophenyl-phenylether	10.93	248	118980	42.529	ng	94
67) Hexachlorobenzene	11.00	284	115427	38.630	ng	94
68) Atrazine	11.09	200	123218	41.362	ng	99
69) Pentachlorophenol	11.20	266	103883	55.863	ng	99
70) Phenanthrene	11.42	178	617691	40.375	ng	98
71) Anthracene	11.47	178	637531	40.728	ng	99
72) Carbazole	11.63	167	591042	38.557	ng	98
73) Di-n-butylphthalate	11.95	149	755951	40.971	ng	99
74) Fluoranthene	12.60	202	615191	39.711	ng	99
76) Benzidine	12.73	184	519268	67.479	ng	98
77) Pyrene	12.84	202	628760	39.632	ng	99
79) Butylbenzylphthalate	13.44	149	304530	40.843	ng	92
80) Benzo(a)anthracene	14.02	228	506741	40.223	ng	99
81) 3,3'-Dichlorobenzidine	13.98	252	193117	41.815	ng	98
82) Chrysene	14.06	228	472581	39.401	ng	99
83) Bis(2-ethylhexyl)phthalate	14.00	149	419542	40.864	ng	99
84) Di-n-octyl phthalate	14.61	149	706634	42.744	ng	95
85) Indeno(1,2,3-cd)pyrene	17.00	276	354829	36.982	ng	95
87) Benzo(b)fluoranthene	15.07	252	432839	42.484	ng	# 97
88) Benzo(k)fluoranthene	15.10	252	384438	38.203	ng	98
89) Benzo(a)pyrene	15.44	252	377168	40.330	ng	# 96
90) Dibenzo(a,h)anthracene	17.01	278	299232	39.981	ng	96
91) Benzo(g,h,i)perylene	17.45	276	285391	38.576	ng	96

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

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