

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111617\
 Data File : BF100622.D
 Acq On : 16 Nov 2017 11:19
 Operator : SJ/JU
 Sample : PB104248BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB104248BS

Manual Integrations
 APPROVED

SOHIL
 11/17/2017 1:24:31 PM

Quant Time: Nov 16 13:05:43 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 16 12:15:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	102291	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	416029	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	196548	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	389516	20.00	ng	0.00
75) Chrysene-d12	14.04	240	296076	20.00	ng	0.00
86) Perylene-d12	15.51	264	221061	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.51	112	667960	104.68	ng	0.01
7) Phenol-d6	6.51	99	810406	102.99	ng	0.00
23) Nitrobenzene-d5	7.45	82	512246	81.40	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	232004	115.84	ng	0.00
44) 2-Fluorobiphenyl	9.24	172	1018111	78.88	ng	0.00
78) Terphenyl-d14	12.99	244	1018973	79.08	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.76	88	94059	30.246	ng	96
3) Pyridine	3.52	79	258048	30.415	ng	96
4) n-Nitrosodimethylamine	3.47	42	120737	29.310	ng	98
6) Aniline	6.55	93	146033	13.136	ng	99
8) 2-Chlorophenol	6.67	128	245101	36.307	ng	97
9) Benzaldehyde	6.43	77	100164	17.824	ng	93
10) Phenol	6.52	94	292431	35.400	ng	99
11) bis(2-Chloroethyl)ether	6.62	93	229436	33.066	ng	95
12) 1,3-Dichlorobenzene	6.82	146	278635	35.800	ng	98
13) 1,4-Dichlorobenzene	6.90	146	279655	35.501	ng	98
14) 1,2-Dichlorobenzene	7.05	146	265651	36.473	ng	98
15) Benzyl Alcohol	7.02	79	210993	34.356	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.16	45	415072	37.402	ng	99
17) 2-Methylphenol	7.13	107	212056	36.758	ng	99
18) Hexachloroethane	7.39	117	96179	37.030	ng	98
19) n-Nitroso-di-n-propylamine	7.29	70	170804	31.299	ng	98
20) 3+4-Methylphenols	7.28	107	256135	35.086	ng	# 87
22) Acetophenone	7.29	105	332798	32.805	ng	# 99
24) Nitrobenzene	7.46	77	257067	39.691	ng	99
25) Isophorone	7.70	82	470818	35.605	ng	100
26) 2-Nitrophenol	7.78	139	137733	45.653	ng	92
27) 2,4-Dimethylphenol	7.81	122	237755	40.108	ng	98
28) bis(2-Chloroethoxy)methane	7.91	93	302849	35.013	ng	99
29) 2,4-Dichlorophenol	8.02	162	219835	38.847	ng	98
30) 1,2,4-Trichlorobenzene	8.10	180	229506	37.493	ng	96
31) Naphthalene	8.19	128	709263	35.396	ng	99
32) Benzoic acid	7.90	122	53583	18.870	ng	98
33) 4-Chloroaniline	8.23	127	73712	8.837	ng	98
34) Hexachlorobutadiene	8.30	225	130240	35.784	ng	99
35) Caprolactam	8.60	113	62374m	32.118	ng	
36) 4-Chloro-3-methylphenol	8.71	107	224012	36.858	ng	94
37) 2-Methylnaphthalene	8.87	142	493601	37.103	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.04	216	224488	34.439	ng	98
40) Hexachlorocyclopentadiene	9.03	237	228912	74.614	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.15	196	163575	39.594	nq	98
43) 2,4,5-Trichlorophenol	9.19	196	167829	39.209	nq	# 91
45) 1,1'-Biphenyl	9.34	154	590765	34.752	nq	98
46) 2-Chloronaphthalene	9.37	162	463327	37.570	nq	95
47) 2-Nitroaniline	9.46	65	143400	39.854	nq	93
48) Acenaphthylene	9.78	152	718782	35.169	nq	99
49) Dimethylphthalate	9.64	163	536338	35.740	nq	99
50) 2,6-Dinitrotoluene	9.70	165	123533	41.587	nq	97
51) Acenaphthene	9.95	154	449133	36.134	nq	99
52) 3-Nitroaniline	9.87	138	60595	17.275	nq	# 90
53) 2,4-Dinitrophenol	9.97	184	63084	57.085	nq	# 16
54) Dibenzofuran	10.13	168	631544	36.252	nq	97
55) 4-Nitrophenol	10.03	139	203357	71.398	nq	95
56) 2,4-Dinitrotoluene	10.10	165	163590	43.654	nq	93
57) Fluorene	10.47	166	481558	37.733	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.24	232	142601	40.649	nq	# 84
59) Diethylphthalate	10.34	149	533834	35.547	nq	96
60) 4-Chlorophenyl-phenylether	10.46	204	237864	37.993	nq	91
61) 4-Nitroaniline	10.49	138	125841	37.835	nq	85
62) Azobenzene	10.62	77	491502	34.750	nq	97
64) 4,6-Dinitro-2-methylphenol	10.51	198	59709	33.509	nq	83
65) n-Nitrosodiphenylamine	10.57	169	467369	34.508	nq	94
66) 4-Bromophenyl-phenylether	10.95	248	152284	36.470	nq	90
67) Hexachlorobenzene	11.02	284	162565	37.225	nq	# 85
68) Atrazine	11.10	200	152318	37.153	nq	97
69) Pentachlorophenol	11.20	266	164484	52.526	nq	98
70) Phenanthrene	11.43	178	738998	36.034	nq	99
71) Anthracene	11.48	178	761452	36.596	nq	99
72) Carbazole	11.63	167	684566	35.657	nq	99
73) Di-n-butylphthalate	11.96	149	858588	38.215	nq	# 98
74) Fluoranthene	12.62	202	794044	36.533	nq	98
76) Benzidine	12.73	184	206135	17.385	nq	98
77) Pyrene	12.84	202	814026	38.569	nq	99
79) Butylbenzylphthalate	13.46	149	354874	41.230	nq	94
80) Benzo(a)anthracene	14.03	228	661542	37.104	nq	99
81) 3,3'-Dichlorobenzidine	13.99	252	110624	17.317	nq	# 95
82) Chrysene	14.07	228	622003	36.026	nq	99
83) Bis(2-ethylhexyl)phthalate	14.02	149	479893	42.392	nq	99
84) Di-n-octyl phthalate	14.62	149	707576	36.384	nq	99
85) Indeno(1,2,3-cd)pyrene	17.00	276	510349	36.544	nq	95
87) Benzo(b)fluoranthene	15.08	252	513491	39.208	nq	100
88) Benzo(k)fluoranthene	15.11	252	448567	36.249	nq	98
89) Benzo(a)pyrene	15.45	252	443996	38.240	nq	99
90) Dibenzo(a,h)anthracene	17.02	278	414977	43.703	nq	97
91) Benzo(g,h,i)perylene	17.45	276	438511	47.178	nq	# 93

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

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