

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
 Data File : BF101651.D
 Acq On : 22 Dec 2017 14:16
 Operator : SJ/JU
 Sample : PB105201BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB105201BS

Manual Integrations
 APPROVED

Sohil
 12/26/2017 1:37:14 PM

Quant Time: Dec 22 16:13:28 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 22 16:07:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	140479	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	568253	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	257840	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	496219	20.00	ng	0.00
75) Chrysene-d12	13.97	240	367376	20.00	ng	0.00
86) Perylene-d12	15.43	264	303371	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	991270	105.11	ng	0.01
7) Phenol-d6	6.44	99	1150144	97.99	ng	0.00
23) Nitrobenzene-d5	7.36	82	749560	73.43	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	280054	112.72	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1270378	76.43	ng	0.00
78) Terphenyl-d14	12.90	244	1256610	82.46	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.59	88	151151	31.192	ng	95
3) Pyridine	3.35	79	376479	27.962	ng	96
4) n-Nitrosodimethylamine	3.30	42	197141	31.802	ng	98
6) Aniline	6.46	93	187820	11.515	ng	# 73
8) 2-Chlorophenol	6.58	128	332904	33.557	ng	99
9) Benzaldehyde	6.35	77	183083	21.122	ng	97
10) Phenol	6.45	94	422575	31.589	ng	94
11) bis(2-Chloroethyl)ether	6.53	93	350964	33.447	ng	99
12) 1,3-Dichlorobenzene	6.73	146	366407	32.725	ng	99
13) 1,4-Dichlorobenzene	6.80	146	371474	32.489	ng	99
14) 1,2-Dichlorobenzene	6.96	146	337892	32.831	ng	98
15) Benzyl Alcohol	6.94	79	254671	31.524	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.06	45	569181	35.442	ng	# 47
17) 2-Methylphenol	7.05	107	275386	30.178	ng	# 90
18) Hexachloroethane	7.29	117	134635	33.868	ng	96
19) n-Nitroso-di-n-propylamine	7.20	70	246430	31.971	ng	93
20) 3+4-Methylphenols	7.21	107	368885	38.147	ng	# 79
22) Acetophenone	7.20	105	475666	36.685	ng	# 92
24) Nitrobenzene	7.37	77	378411	33.494	ng	96
25) Isophorone	7.61	82	690338	33.710	ng	96
26) 2-Nitrophenol	7.69	139	183879	37.883	ng	96
27) 2,4-Dimethylphenol	7.73	122	305339	37.029	ng	97
28) bis(2-Chloroethoxy)methane	7.82	93	435456	33.475	ng	98
29) 2,4-Dichlorophenol	7.95	162	286466	35.164	ng	98
30) 1,2,4-Trichlorobenzene	8.01	180	288862	33.600	ng	98
31) Naphthalene	8.09	128	947022	33.103	ng	99
32) Benzoic acid	7.89	122	216400	40.449	ng	96
33) 4-Chloroaniline	8.16	127	84510	6.797	ng	98
34) Hexachlorobutadiene	8.20	225	163540	32.890	ng	97
35) Caprolactam	8.53	113	100470m	35.517	ng	
36) 4-Chloro-3-methylphenol	8.65	107	302014	33.729	ng	100
37) 2-Methylnaphthalene	8.78	142	635382	34.090	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.95	216	300518	34.521	ng	98
40) Hexachlorocyclopentadiene	8.93	237	79023	53.887	ng	99

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42) 2,4,6-Trichlorophenol	9.07	196	188239	35.918	ng	99
43) 2,4,5-Trichlorophenol	9.12	196	188321	32.818	ng	97
45) 1,1'-Biphenyl	9.25	154	783634	34.270	ng	98
46) 2-Chloronaphthalene	9.27	162	587171	33.559	ng	98
47) 2-Nitroaniline	9.39	65	208017	34.416	ng	94
48) Acenaphthylene	9.69	152	871834	31.548	ng	99
49) Dimethylphthalate	9.55	163	648998	31.293	ng	99
50) 2,6-Dinitrotoluene	9.63	165	145234	34.455	ng	91
51) Acenaphthene	9.86	154	526502	31.711	ng	99
52) 3-Nitroaniline	9.80	138	89368	17.005	ng	98
53) 2,4-Dinitrophenol	9.93	184	112997	77.358	ng #	20
54) Dibenzofuran	10.03	168	748976	35.497	ng	99
55) 4-Nitrophenol	9.98	139	140218	61.189	ng #	85
56) 2,4-Dinitrotoluene	10.04	165	184981	38.950	ng #	87
57) Fluorene	10.38	166	621335	34.810	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.16	232	153392	37.206	ng	89
59) Diethylphthalate	10.25	149	676758	32.951	ng	98
60) 4-Chlorophenyl-phenylether	10.36	204	299713	35.036	ng	94
61) 4-Nitroaniline	10.42	138	144146	27.288	ng	98
62) Azobenzene	10.53	77	700269	32.914	ng	96
64) 4,6-Dinitro-2-methylphenol	10.46	198	84530	33.381	ng	90
65) n-Nitrosodiphenylamine	10.49	169	572167	31.227	ng	95
66) 4-Bromophenyl-phenylether	10.85	248	188422	33.794	ng	93
67) Hexachlorobenzene	10.93	284	195644	33.154	ng	93
68) Atrazine	11.02	200	186392	34.117	ng	97
69) Pentachlorophenol	11.13	266	149194	66.242	ng	98
70) Phenanthrene	11.34	178	899017	32.578	ng	99
71) Anthracene	11.40	178	923046	32.746	ng	99
72) Carbazole	11.56	167	817757	30.986	ng	99
73) Di-n-butylphthalate	11.87	149	1057279	33.007	ng	99
74) Fluoranthene	12.53	202	945047	32.627	ng	100
77) Pyrene	12.76	202	985598	35.747	ng	100
79) Butylbenzylphthalate	13.37	149	446214	35.767	ng	96
80) Benzo(a)anthracene	13.96	228	797525	32.876	ng	100
81) 3,3'-Dichlorobenzidine	13.92	252	174284	22.034	ng #	98
82) Chrysene	13.99	228	760140	33.188	ng	99
83) Bis(2-ethylhexyl)phthalate	13.92	149	571408	36.161	ng	99
84) Di-n-octyl phthalate	14.53	149	929152	32.971	ng #	95
85) Indeno(1,2,3-cd)pyrene	16.90	276	621069	30.450	ng	96
87) Benzo(b)fluoranthene	15.00	252	622724	33.677	ng	98
88) Benzo(k)fluoranthene	15.03	252	613302	34.113	ng	98
89) Benzo(a)pyrene	15.37	252	584281	34.276	ng	99
90) Dibenzo(a,h)anthracene	16.90	278	522272	35.685	ng	95
91) Benzo(g,h,i)perylene	17.34	276	549141	37.892	ng	95

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

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