

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071417\
 Data File : BF096784.D
 Acq On : 14 Jul 2017 21:07
 Operator : SJ/JU
 Sample : PB100646BSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB100646BSD

Manual Integrations
 APPROVED

mohammad
 7/17/2017 3:56:58 PM

Quant Time: Jul 15 00:49:08 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 13 14:49:34 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.62	152	107424	20.00	ng	-0.01
21) Naphthalene-d8	7.91	136	448840	20.00	ng	-0.01
38) Acenaphthene-d10	9.66	164	201538	20.00	ng	-0.01
63) Phenanthrene-d10	11.13	188	353859	20.00	ng	-0.01
75) Chrysene-d12	13.77	240	233419	20.00	ng	0.00
86) Perylene-d12	15.14	264	206027	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.23	112	524826	73.46	ng	0.00
7) Phenol-d6	6.27	99	681482	79.49	ng	-0.01
23) Nitrobenzene-d5	7.19	82	581690	80.27	ng	-0.01
41) 2,4,6-Tribromophenol	10.45	330	143528	83.43	ng	-0.01
44) 2-Fluorobiphenyl	8.99	172	1027425	80.54	ng	-0.01
78) Terphenyl-d14	12.72	244	867698	64.47	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.29	88	96840	26.342	ng	# 75
3) Pyridine	2.95	79	253816	32.852	ng	# 67
4) n-Nitrosodimethylamine	2.89	42	158444	36.672	ng	# 79
6) Aniline	6.29	93	316832	30.001	ng	# 41
8) 2-Chlorophenol	6.41	128	302790	39.285	ng	# 95
9) Benzaldehyde	6.16	77	154018	31.080	ng	# 98
10) Phenol	6.28	94	372053	38.387	ng	# 74
11) bis(2-Chloroethyl)ether	6.36	93	288708	39.612	ng	# 90
12) 1,3-Dichlorobenzene	6.56	146	312438	38.135	ng	# 98
13) 1,4-Dichlorobenzene	6.64	146	314731	37.933	ng	# 94
14) 1,2-Dichlorobenzene	6.80	146	295280	39.128	ng	# 96
15) Benzyl Alcohol	6.77	79	241348	42.997	ng	# 99
16) 2,2'-oxybis(1-Chloropropan	6.91	45	620217	41.234	ng	# 91
17) 2-Methylphenol	6.89	107	245703	40.994	ng	# 99
18) Hexachloroethane	7.14	117	111398	37.864	ng	# 86
19) n-Nitroso-di-n-propylamine	7.05	70	200310	38.780	ng	# 83
20) 3+4-Methylphenols	7.05	107	299528	41.183	ng	# 79
22) Acetophenone	7.04	105	400342	39.908	ng	# 95
24) Nitrobenzene	7.21	77	300450	40.090	ng	# 87
25) Isophorone	7.45	82	555787	41.230	ng	# 96
26) 2-Nitrophenol	7.53	139	172246	41.334	ng	# 90
27) 2,4-Dimethylphenol	7.57	122	277605	40.494	ng	# 95
28) bis(2-Chloroethoxy)methane	7.66	93	380160	41.735	ng	# 99
29) 2,4-Dichlorophenol	7.77	162	257242	41.454	ng	# 94
30) 1,2,4-Trichlorobenzene	7.85	180	233169	39.519	ng	# 98
31) Naphthalene	7.93	128	853327	40.133	ng	# 100
32) Benzoic acid	7.71	122	199825	45.605	ng	# 94
33) 4-Chloroaniline	7.98	127	191596	22.744	ng	# 95
34) Hexachlorobutadiene	8.06	225	121747	38.653	ng	# 99
35) Caprolactam	8.35	113	91469m	46.002	ng	#
36) 4-Chloro-3-methylphenol	8.47	107	279511	41.890	ng	# 87
37) 2-Methylnaphthalene	8.62	142	580682	42.837	ng	# 98
39) 1,2,4,5-Tetrachlorobenzene	8.79	216	201053	36.904	ng	# 99
40) Hexachlorocyclopentadiene	8.78	237	172206	83.791	ng	# 99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.90	196	162313	40.390	ng	99
43) 2,4,5-Trichlorophenol	8.95	196	171369	42.283	ng	97
45) 1,1'-Biphenyl	9.09	154	660950	38.239	ng	97
46) 2-Chloronaphthalene	9.11	162	516463	40.578	ng	95
47) 2-Nitroaniline	9.20	65	189600	43.337	ng	94
48) Acenaphthylene	9.52	152	819052	40.388	ng	99
49) Dimethylphthalate	9.39	163	628755	41.433	ng	99
50) 2,6-Dinitrotoluene	9.45	165	138361	42.094	ng	92
51) Acenaphthene	9.69	154	471395	39.349	ng	98
52) 3-Nitroaniline	9.61	138	107509	27.610	ng	91
53) 2,4-Dinitrophenol	9.73	184	144428	88.006	ng #	75
54) Dibenzofuran	9.86	168	712543	42.444	ng	98
55) 4-Nitrophenol	9.79	139	243295	85.524	ng #	67
56) 2,4-Dinitrotoluene	9.86	165	183088	43.348	ng #	79
57) Fluorene	10.20	166	524971	42.318	ng	98
58) 2,3,4,6-Tetrachlorophenol	9.99	232	134143	47.724	ng #	98
59) Diethylphthalate	10.09	149	601090	40.688	ng	99
60) 4-Chlorophenyl-phenylether	10.20	204	213399	40.410	ng	98
61) 4-Nitroaniline	10.23	138	172502	47.176	ng	90
62) Azobenzene	10.36	77	585695	44.941	ng	92
64) 4,6-Dinitro-2-methylphenol	10.26	198	97276	43.784	ng	89
65) n-Nitrosodiphenylamine	10.32	169	493301	38.999	ng	99
66) 4-Bromophenyl-phenylether	10.69	248	145169	40.182	ng	96
67) Hexachlorobenzene	10.76	284	154842	38.482	ng #	93
68) Atrazine	10.85	200	154545	42.885	ng	95
69) Pentachlorophenol	10.95	266	156864	80.697	ng	95
70) Phenanthrene	11.16	178	807426	41.965	ng	99
71) Anthracene	11.22	178	828192	42.573	ng	99
72) Carbazole	11.37	167	816423	43.338	ng	99
73) Di-n-butylphthalate	11.71	149	951040	40.536	ng	98
74) Fluoranthene	12.34	202	799926	43.435	ng	99
76) Benzidine	12.46	184	354356	46.340	ng #	93
77) Pyrene	12.57	202	845200	39.898	ng	99
79) Butylbenzylphthalate	13.20	149	417323	82.924	ng #	81
80) Benzo(a)anthracene	13.75	228	645426	44.378	ng	99
81) 3,3'-Dichlorobenzidine	13.72	252	148728	28.438	ng #	96
82) Chrysene	13.79	228	636489	44.455	ng	99
83) Bis(2-ethylhexyl)phthalate	13.76	149	487350	39.759	ng	99
84) Di-n-octyl phthalate	14.37	149	888281	43.840	ng #	94
85) Indeno(1,2,3-cd)pyrene	16.44	276	442807	34.033	ng	100
87) Benzo(b)fluoranthene	14.76	252	534453	43.012	ng	98
88) Benzo(k)fluoranthene	14.79	252	504107m	42.843	ng	
89) Benzo(a)pyrene	15.09	252	500844	43.947	ng	100
90) Dibenzo(a,h)anthracene	16.46	278	363879	34.699	ng	98
91) Benzo(g,h,i)perylene	16.84	276	385857	34.361	ng	97

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

