

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071417\
 Data File : BF096780.D
 Acq On : 14 Jul 2017 19:14
 Operator : SJ/JU
 Sample : PB100644BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB100644BS

Manual Integrations
 APPROVED

mohammad
 7/17/2017 3:56:32 PM

Quant Time: Jul 15 00:36:52 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	6.63	152	104395	20.00	ng	0.00
21) Naphthalene-d8	7.91	136	418347	20.00	ng	-0.01
38) Acenaphthene-d10	9.66	164	173111	20.00	ng	-0.01
63) Phenanthrene-d10	11.13	188	310977	20.00	ng	-0.01
75) Chrysene-d12	13.77	240	211650	20.00	ng	0.00
86) Perylene-d12	15.14	264	178467	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.24	112	896142	129.07	ng	0.00
7) Phenol-d6	6.28	99	953133	114.40	ng	0.00
23) Nitrobenzene-d5	7.19	82	654081	96.84	ng	-0.01
41) 2,4,6-Tribromophenol	10.45	330	207162	140.20	ng	0.00
44) 2-Fluorobiphenyl	8.99	172	1019533	93.05	ng	-0.01
78) Terphenyl-d14	12.72	244	913646	74.86	ng	-0.01

Target Compounds					Qvalue	
2) 1,4-Dioxane	2.29	88	110925	31.049	ng	# 76
3) Pyridine	2.96	79	279524	37.229	ng	# 61
4) n-Nitrosodimethylamine	2.90	42	177271	42.220	ng	# 77
6) Aniline	6.29	93	350647	34.166	ng	# 23
8) 2-Chlorophenol	6.42	128	294501	39.318	ng	100
9) Benzaldehyde	6.17	77	143346	29.766	ng	100
10) Phenol	6.29	94	354435	37.630	ng	78
11) bis(2-Chloroethyl)ether	6.37	93	278152	39.271	ng	92
12) 1,3-Dichlorobenzene	6.57	146	339037	42.582	ng	99
13) 1,4-Dichlorobenzene	6.65	146	352841	43.760	ng	97
14) 1,2-Dichlorobenzene	6.80	146	324661	44.269	ng	96
15) Benzyl Alcohol	6.77	79	262939	48.203	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.92	45	660976	45.218	ng	93
17) 2-Methylphenol	6.89	107	267821	45.981	ng	99
18) Hexachloroethane	7.14	117	122165	42.728	ng	# 82
19) n-Nitroso-di-n-propylamine	7.05	70	213752	42.583	ng	# 85
20) 3+4-Methylphenols	7.05	107	316117	44.725	ng	# 70
22) Acetophenone	7.04	105	433413	46.353	ng	# 97
24) Nitrobenzene	7.21	77	319991	45.810	ng	94
25) Isophorone	7.46	82	537609	42.789	ng	95
26) 2-Nitrophenol	7.53	139	186647	48.055	ng	# 88
27) 2,4-Dimethylphenol	7.57	122	306693	47.997	ng	97
28) bis(2-Chloroethoxy)methane	7.67	93	398724	46.963	ng	100
29) 2,4-Dichlorophenol	7.77	162	276006	47.719	ng	98
30) 1,2,4-Trichlorobenzene	7.85	180	235745	42.868	ng	97
31) Naphthalene	7.93	128	945102	47.689	ng	99
32) Benzoic acid	7.72	122	231713	56.737	ng	93
33) 4-Chloroaniline	7.98	127	313690	39.951	ng	97
34) Hexachlorobutadiene	8.06	225	132343	45.080	ng	99
35) Caprolactam	8.36	113	92357m	49.834	ng	
36) 4-Chloro-3-methylphenol	8.47	107	257670	41.431	ng	89
37) 2-Methylnaphthalene	8.62	142	566915	44.870	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.79	216	196782	42.051	ng	99
40) Hexachlorocyclopentadiene	8.78	237	184509	103.286	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.90	196	155626	45.085	ng	100
43) 2,4,5-Trichlorophenol	8.95	196	167828	48.209	ng	95
45) 1,1'-Biphenyl	9.09	154	638488	43.005	ng	98
46) 2-Chloronaphthalene	9.11	162	497602	45.516	ng	94
47) 2-Nitroaniline	9.20	65	176507	46.970	ng	93
48) Acenaphthylene	9.52	152	795816	45.687	ng	99
49) Dimethylphthalate	9.39	163	603589	46.306	ng	99
50) 2,6-Dinitrotoluene	9.45	165	131575	46.603	ng #	86
51) Acenaphthene	9.69	154	459257	44.631	ng	98
52) 3-Nitroaniline	9.62	138	125854	37.629	ng	92
53) 2,4-Dinitrophenol	9.73	184	145063	101.766	ng #	75
54) Dibenzofuran	9.87	168	687170	47.654	ng	98
55) 4-Nitrophenol	9.79	139	237026	97.003	ng #	67
56) 2,4-Dinitrotoluene	9.86	165	172906	47.659	ng #	76
57) Fluorene	10.21	166	503160	47.220	ng	97
58) 2,3,4,6-Tetrachlorophenol	9.99	232	130733	54.148	ng #	97
59) Diethylphthalate	10.09	149	580924	45.781	ng	99
60) 4-Chlorophenyl-phenylether	10.20	204	209630	48.534	ng	92
61) 4-Nitroaniline	10.23	138	165682	52.752	ng	93
62) Azobenzene	10.36	77	551027	49.224	ng	92
64) 4,6-Dinitro-2-methylphenol	10.27	198	96894	49.626	ng #	75
65) n-Nitrosodiphenylamine	10.32	169	484529	43.588	ng	98
66) 4-Bromophenyl-phenylether	10.69	248	142086	44.752	ng	97
67) Hexachlorobenzene	10.76	284	151375	42.808	ng #	92
68) Atrazine	10.85	200	149102	47.080	ng	93
69) Pentachlorophenol	10.95	266	154175	90.251	ng	99
70) Phenanthrene	11.16	178	774784	45.821	ng	99
71) Anthracene	11.22	178	806974	47.202	ng	99
72) Carbazole	11.37	167	766628	46.306	ng	99
73) Di-n-butylphthalate	11.71	149	937611	45.475	ng	99
74) Fluoranthene	12.34	202	846556	52.306	ng	99
76) Benzidine	12.47	184	469743	67.747	ng #	92
77) Pyrene	12.57	202	878208	45.720	ng	100
79) Butylbenzylphthalate	13.20	149	426265	93.412	ng #	83
80) Benzo(a)anthracene	13.75	228	627172	47.558	ng	100
81) 3,3'-Dichlorobenzidine	13.72	252	183403	38.675	ng #	94
82) Chrysene	13.79	228	623492	48.026	ng	99
83) Bis(2-ethylhexyl)phthalate	13.76	149	487989	43.906	ng	99
84) Di-n-octyl phthalate	14.37	149	737044	40.117	ng #	95
85) Indeno(1,2,3-cd)pyrene	16.45	276	470388	39.024	ng	99
87) Benzo(b)fluoranthene	14.76	252	480573	44.648	ng	97
88) Benzo(k)fluoranthene	14.79	252	443785m	43.540	ng	
89) Benzo(a)pyrene	15.09	252	456430	46.234	ng	99
90) Dibenzo(a,h)anthracene	16.46	278	385767	42.467	ng	96
91) Benzo(g,h,i)perylene	16.84	276	414389	42.600	ng	99

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

