

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071717\
 Data File : BF096822.D
 Acq On : 17 Jul 2017 16:31
 Operator : SJ/JU
 Sample : I4196-01MS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-PS-01-012MS

Manual Integrations
 APPROVED

mohammad
 7/20/2017 8:38:19 AM

Quant Time: Jul 17 18:56:05 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF071317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 17 18:17:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.61	152	110839	20.00	ng	0.00
21) Naphthalene-d8	7.89	136	522957	20.00	ng	0.00
38) Acenaphthene-d10	9.64	164	194998	20.00	ng	0.00
63) Phenanthrene-d10	11.12	188	297730	20.00	ng	0.00
75) Chrysene-d12	13.74	240	199225	20.00	ng	0.00
86) Perylene-d12	15.11	264	204827	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.23	112	661647	89.76	ng	0.02
7) Phenol-d6	6.26	99	814473	92.07	ng	0.00
23) Nitrobenzene-d5	7.17	82	523762	62.04	ng	0.00
41) 2,4,6-Tribromophenol	10.43	330	148352	89.13	ng	0.00
44) 2-Fluorobiphenyl	8.97	172	917475	74.34	ng	0.00
78) Terphenyl-d14	12.70	244	515292	44.86	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.27	88	96987	25.569	ng	# 75
3) Pyridine	2.93	79	238814	29.958	ng	# 63
4) n-Nitrosodimethylamine	2.87	42	151363	33.954	ng	# 81
6) Aniline	6.27	93	277764	25.491	ng	# 19
8) 2-Chlorophenol	6.40	128	247116	31.074	ng	# 98
9) Benzaldehyde	6.15	77	134937	26.391	ng	# 99
10) Phenol	6.28	94	331837	33.183	ng	# 94
11) bis(2-Chloroethyl)ether	6.35	93	265617	35.321	ng	# 90
12) 1,3-Dichlorobenzene	6.55	146	264747	31.319	ng	# 98
13) 1,4-Dichlorobenzene	6.63	146	269018	31.425	ng	# 95
14) 1,2-Dichlorobenzene	6.78	146	253493	32.556	ng	# 98
15) Benzyl Alcohol	6.76	79	195745	33.798	ng	# 100
16) 2,2'-oxybis(1-Chloropropan	6.90	45	580044	37.375	ng	# 93
17) 2-Methylphenol	6.88	107	221761	35.860	ng	# 98
18) Hexachloroethane	7.12	117	98957	32.599	ng	# 78
19) n-Nitroso-di-n-propylamine	7.03	70	185941	34.889	ng	# 81
20) 3+4-Methylphenols	7.04	107	273430	36.436	ng	# 65
22) Acetophenone	7.03	105	377730	32.317	ng	# 99
24) Nitrobenzene	7.19	77	263608	30.189	ng	# 95
25) Isophorone	7.44	82	456794	29.084	ng	# 95
26) 2-Nitrophenol	7.52	139	146773	30.230	ng	# 92
27) 2,4-Dimethylphenol	7.56	122	254989	31.923	ng	# 96
28) bis(2-Chloroethoxy)methane	7.65	93	314391	29.623	ng	# 100
29) 2,4-Dichlorophenol	7.76	162	212405	29.377	ng	# 98
30) 1,2,4-Trichlorobenzene	7.84	180	208833	30.378	ng	# 98
31) Naphthalene	7.92	128	786912	31.764	ng	# 100
32) Benzoic acid	7.66	122	28861	5.653	ng	# 96
33) 4-Chloroaniline	7.97	127	231387	23.574	ng	# 95
34) Hexachlorobutadiene	8.04	225	104344	28.433	ng	# 97
35) Caprolactam	8.33	113	74644m	32.220	ng	# 91
36) 4-Chloro-3-methylphenol	8.46	107	237029	30.489	ng	# 98
37) 2-Methylnaphthalene	8.60	142	544881	34.499	ng	# 98
39) 1,2,4,5-Tetrachlorobenzene	8.77	216	190603	36.159	ng	# 100
40) Hexachlorocyclopentadiene	8.76	237	54265	30.650	ng	# 99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.89	196	130698	33.614	nq	99
43) 2,4,5-Trichlorophenol	8.93	196	146927	37.468	nq	94
45) 1,1'-Biphenyl	9.07	154	586942	35.096	nq	98
46) 2-Chloronaphthalene	9.09	162	457136	37.121	nq	94
47) 2-Nitroaniline	9.19	65	163951	38.732	nq	96
48) Acenaphthylene	9.50	152	628976	32.056	nq	99
49) Dimethylphthalate	9.37	163	640421	43.617	nq	100
50) 2,6-Dinitrotoluene	9.43	165	101370	31.875	nq	# 87
51) Acenaphthene	9.67	154	361730	31.208	nq	99
52) 3-Nitroaniline	9.60	138	104015	27.608	nq	86
53) 2,4-Dinitrophenol	9.71	184	16172	16.145	nq	# 69
54) Dibenzofuran	9.85	168	548147	33.747	nq	99
55) 4-Nitrophenol	9.77	139	154468	56.120	nq	# 69
56) 2,4-Dinitrotoluene	9.83	165	125217	30.640	nq	# 75
57) Fluorene	10.19	166	390483	32.532	nq	99
58) 2,3,4,6-Tetrachlorophenol	9.97	232	96309	35.413	nq	95
59) Diethylphthalate	10.07	149	474771	33.216	nq	99
60) 4-Chlorophenyl-phenylether	10.18	204	172231	32.077	nq	97
61) 4-Nitroaniline	10.20	138	108395	30.638	nq	92
62) Azobenzene	10.34	77	418633	33.199	nq	92
64) 4,6-Dinitro-2-methylphenol	10.24	198	21154	11.316	nq	# 75
65) n-Nitrosodiphenylamine	10.30	169	374527	35.191	nq	99
66) 4-Bromophenyl-phenylether	10.67	248	108499	35.694	nq	95
67) Hexachlorobenzene	10.73	284	108282	31.984	nq	# 86
68) Atrazine	10.83	200	106383	35.085	nq	96
69) Pentachlorophenol	10.93	266	85349	52.185	nq	99
70) Phenanthrene	11.14	178	536100	33.116	nq	99
71) Anthracene	11.19	178	542575	33.149	nq	99
72) Carbazole	11.34	167	498831	31.471	nq	99
73) Di-n-butylphthalate	11.69	149	696811	35.299	nq	99
74) Fluoranthene	12.32	202	459693	29.667	nq	97
76) Benzidine	12.44	184	283064	43.370	nq	# 92
77) Pyrene	12.54	202	454519	25.138	nq	99
79) Butylbenzylphthalate	13.17	149	238083	55.428	nq	# 86
80) Benzo(a)anthracene	13.73	228	387736	31.236	nq	98
81) 3,3'-Dichlorobenzidine	13.69	252	138028	30.921	nq	# 95
82) Chrysene	13.76	228	387635	31.721	nq	98
83) Bis(2-ethylhexyl)phthalate	13.74	149	332578	31.789	nq	100
84) Di-n-octyl phthalate	14.34	149	593733	34.332	nq	# 93
85) Indeno(1,2,3-cd)pyrene	16.40	276	293619	27.430	nq	100
87) Benzo(b)fluoranthene	14.73	252	374652	30.328	nq	99
88) Benzo(k)fluoranthene	14.76	252	369479m	31.585	nq	
89) Benzo(a)pyrene	15.06	252	349651	30.860	nq	99
90) Dibenzo(a,h)anthracene	16.42	278	248780	23.862	nq	98
91) Benzo(g,h,i)perylene	16.79	276	236414	21.176	nq	96

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

