

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081717\
 Data File : BF097786.D
 Acq On : 17 Aug 2017 15:16
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 8/18/2017 7:02:20 PM

Quant Time: Aug 18 02:04:48 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 15 13:25:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	151994	20.00	ng	-0.01
21) Naphthalene-d8	8.06	136	594485	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	257636	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	485874	20.00	ng	0.00
75) Chrysene-d12	13.93	240	356714	20.00	ng	0.00
86) Perylene-d12	15.36	264	269352	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.38	112	826765	82.74	ng	0.00
7) Phenol-d6	6.41	99	1135347	83.62	ng	0.00
23) Nitrobenzene-d5	7.35	82	990356	90.50	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	209257	93.61	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	1411769	80.51	ng	0.00
78) Terphenyl-d14	12.89	244	1373803	80.31	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.48	88	229242	41.136	ng	88
3) Pyridine	3.20	79	642480	41.703	ng	86
4) n-Nitrosodimethylamine	3.16	42	232443	39.212	ng	# 75
6) Aniline	6.44	93	775674	40.669	ng	98
8) 2-Chlorophenol	6.56	128	441232	41.870	ng	98
9) Benzaldehyde	6.33	77	320253	37.240	ng	91
10) Phenol	6.42	94	669826	42.183	ng	95
11) bis(2-Chloroethyl)ether	6.52	93	486949	40.630	ng	97
12) 1,3-Dichlorobenzene	6.72	146	468625	41.342	ng	# 93
13) 1,4-Dichlorobenzene	6.79	146	470643	40.824	ng	# 92
14) 1,2-Dichlorobenzene	6.95	146	436336	41.047	ng	98
15) Benzyl Alcohol	6.92	79	416305	40.955	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.06	45	628575	38.981	ng	91
17) 2-Methylphenol	7.03	107	380427	40.965	ng	98
18) Hexachloroethane	7.29	117	187554	41.495	ng	# 85
19) n-Nitroso-di-n-propylamine	7.20	70	363890	39.152	ng	90
20) 3+4-Methylphenols	7.19	107	458358	40.410	ng	98
22) Acetophenone	7.19	105	618247	39.367	ng	# 87
24) Nitrobenzene	7.36	77	537953	44.150	ng	92
25) Isophorone	7.60	82	909702	39.978	ng	97
26) 2-Nitrophenol	7.67	139	213085	48.241	ng	# 76
27) 2,4-Dimethylphenol	7.72	122	394300	41.549	ng	93
28) bis(2-Chloroethoxy)methane	7.82	93	608869	40.700	ng	99
29) 2,4-Dichlorophenol	7.92	162	337315	42.819	ng	92
30) 1,2,4-Trichlorobenzene	8.00	180	362852	41.550	ng	100
31) Naphthalene	8.09	128	1262099	40.894	ng	100
32) Benzoic acid	7.85	122	301943	44.246	ng	88
33) 4-Chloroaniline	8.13	127	537874	41.324	ng	99
34) Hexachlorobutadiene	8.20	225	212889	41.753	ng	99
35) Caprolactam	8.52	113	115872	43.267	ng	86
36) 4-Chloro-3-methylphenol	8.62	107	419885	41.485	ng	# 80
37) 2-Methylnaphthalene	8.77	142	773162	40.861	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.94	216	328781	41.582	ng	98
40) Hexachlorocyclopentadiene	8.93	237	173416	45.654	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.05	196	231523	43.614	nq	96
43) 2,4,5-Trichlorophenol	9.09	196	229939	43.662	nq	95
45) 1,1'-Biphenyl	9.24	154	963705	41.019	nq	96
46) 2-Chloronaphthalene	9.26	162	710417	40.925	nq	# 93
47) 2-Nitroaniline	9.36	65	276741	47.554	nq	95
48) Acenaphthylene	9.68	152	1123605	41.218	nq	99
49) Dimethylphthalate	9.55	163	766171	40.684	nq	99
50) 2,6-Dinitrotoluene	9.60	165	169032	44.966	nq	# 74
51) Acenaphthene	9.85	154	715671	41.627	nq	99
52) 3-Nitroaniline	9.77	138	221809	45.734	nq	82
53) 2,4-Dinitrophenol	9.87	184	81027m	49.710	nq	
54) Dibenzofuran	10.02	168	944245	40.979	nq	98
55) 4-Nitrophenol	9.93	139	171799	44.405	nq	# 51
56) 2,4-Dinitrotoluene	10.00	165	220937	46.833	nq	# 62
57) Fluorene	10.36	166	732698	41.129	nq	96
58) 2,3,4,6-Tetrachlorophenol	10.14	232	178632	43.811	nq	91
59) Diethylphthalate	10.24	149	777327	39.470	nq	99
60) 4-Chlorophenyl-phenylether	10.36	204	336312	40.636	nq	88
61) 4-Nitroaniline	10.39	138	213566m	46.331	nq	
62) Azobenzene	10.52	77	900451	38.492	nq	92
64) 4,6-Dinitro-2-methylphenol	10.41	198	109926	48.009	nq	85
65) n-Nitrosodiphenylamine	10.47	169	674373	40.759	nq	99
66) 4-Bromophenyl-phenylether	10.85	248	212526	42.122	nq	96
67) Hexachlorobenzene	10.91	284	218030	42.396	nq	93
68) Atrazine	11.00	200	170921	38.953	nq	98
69) Pentachlorophenol	11.10	266	134681	44.916	nq	99
70) Phenanthrene	11.32	178	1040009	40.169	nq	100
71) Anthracene	11.37	178	1074333	40.911	nq	100
72) Carbazole	11.53	167	1019658	40.906	nq	99
73) Di-n-butylphthalate	11.87	149	1192319	41.527	nq	99
74) Fluoranthene	12.51	202	1114311	41.234	nq	98
76) Benzidine	12.63	184	424154	36.265	nq	94
77) Pyrene	12.73	202	1150554	39.436	nq	98
79) Butylbenzylphthalate	13.36	149	470697	42.080	nq	# 79
80) Benzo(a)anthracene	13.92	228	840937	39.334	nq	99
81) 3,3'-Dichlorobenzidine	13.89	252	312933	43.377	nq	# 97
82) Chrysene	13.96	228	844451	39.706	nq	100
83) Bis(2-ethylhexyl)phthalate	13.92	149	535482	43.201	nq	# 99
84) Di-n-octyl phthalate	14.53	149	921612	42.733	nq	99
85) Indeno(1,2,3-cd)pyrene	16.76	276	616226	39.388	nq	95
87) Benzo(b)fluoranthene	14.95	252	709486	41.788	nq	99
88) Benzo(k)fluoranthene	14.98	252	644007	40.374	nq	# 93
89) Benzo(a)pyrene	15.30	252	622407	41.410	nq	98
90) Dibenzo(a,h)anthracene	16.77	278	527692	43.125	nq	98
91) Benzo(g,h,i)perylene	17.18	276	540643	42.345	nq	# 92

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

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