

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032017\
 Data File : BF093760.D
 Acq On : 20 Mar 2017 12:16
 Operator : SJ/MA
 Sample : SSTDICC025
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC025

Manual Integrations
 APPROVED

mohammad
 3/21/2017 2:44:51 PM

Quant Time: Mar 20 13:21:56 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 20 13:13:09 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	222743	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	890036	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	420832	20.00	ng	0.00
63) Phenanthrene-d10	11.35	188	719361	20.00	ng	0.00
75) Chrysene-d12	13.98	240	557595	20.00	ng	0.00
86) Perylene-d12	15.39	264	517299	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.43	112	692376	51.82	ng	0.00
7) Phenol-d6	6.46	99	784534	47.38	ng	-0.01
23) Nitrobenzene-d5	7.40	82	773980	52.19	ng	-0.01
41) 2,4,6-Tribromophenol	10.65	330	157354	57.13	ng	-0.01
44) 2-Fluorobiphenyl	9.20	172	1437881	64.48	ng	0.00
78) Terphenyl-d14	12.93	244	1253933	48.40	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.44	88	161828	23.46	ng	98
3) Pyridine	3.13	79	420820	22.30	ng	97
4) n-Nitrosodimethylamine	3.06	42	182435	22.09	ng	98
6) Aniline	6.49	93	588042	24.84	ng	96
8) 2-Chlorophenol	6.62	128	385511	25.95	ng	93
9) Benzaldehyde	6.38	77	297497	31.03	ng	92
10) Phenol	6.47	94	423336	21.87	ng	98
11) bis(2-Chloroethyl)ether	6.57	93	375658	24.62	ng	91
12) 1,3-Dichlorobenzene	6.78	146	431613	26.23	ng	98
13) 1,4-Dichlorobenzene	6.86	146	417275	24.76	ng	98
14) 1,2-Dichlorobenzene	7.01	146	406261	25.78	ng	98
15) Benzyl Alcohol	6.97	79	295204	23.81	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.12	45	571112	23.74	ng	99
17) 2-Methylphenol	7.09	107	301047	23.41	ng	98
18) Hexachloroethane	7.35	117	146008	24.39	ng	# 75
19) n-Nitroso-di-n-propylamine	7.25	70	244094	22.20	ng	93
20) 3+4-Methylphenols	7.25	107	392447	25.74	ng	# 87
22) Acetophenone	7.25	105	513665	25.90	ng	# 91
24) Nitrobenzene	7.42	77	418039	26.08	ng	# 89
25) Isophorone	7.66	82	706694	24.76	ng	98
26) 2-Nitrophenol	7.74	139	179489	29.57	ng	# 90
27) 2,4-Dimethylphenol	7.77	122	336399	24.22	ng	98
28) bis(2-Chloroethoxy)methane	7.88	93	499947	27.75	ng	99
29) 2,4-Dichlorophenol	7.98	162	317396	26.60	ng	97
30) 1,2,4-Trichlorobenzene	8.07	180	322851	25.96	ng	97
31) Naphthalene	8.14	128	1064103	24.52	ng	100
32) Benzoic acid	7.88	122	205653	28.30	ng	95
33) 4-Chloroaniline	8.18	127	421137	23.17	ng	98
34) Hexachlorobutadiene	8.28	225	168603	25.88	ng	98
35) Caprolactam	8.54	113	96680	24.14	ng	92
36) 4-Chloro-3-methylphenol	8.66	107	303932	24.19	ng	# 81
37) 2-Methylnaphthalene	8.84	142	760110	27.04	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.01	216	282184	28.64	ng	99
40) Hexachlorocyclopentadiene	9.00	237	151810	30.84	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.11	196	207408	28.55	nq	97
43) 2,4,5-Trichlorophenol	9.14	196	221587m	29.58	nq	
45) 1,1'-Biphenyl	9.29	154	809174	27.42	nq	95
46) 2-Chloronaphthalene	9.32	162	624042	26.36	nq	# 90
47) 2-Nitroaniline	9.41	65	187521	27.37	nq	93
48) Acenaphthylene	9.74	152	1110314	28.76	nq	99
49) Dimethylphthalate	9.60	163	783734	28.51	nq	99
50) 2,6-Dinitrotoluene	9.65	165	163251	27.50	nq	# 64
51) Acenaphthene	9.91	154	687060	30.19	nq	98
52) 3-Nitroaniline	9.82	138	218199	30.97	nq	87
53) 2,4-Dinitrophenol	9.92	184	63306	35.78	nq	# 59
54) Dibenzofuran	10.08	168	885529	28.63	nq	94
55) 4-Nitrophenol	9.97	139	145455	27.37	nq	# 61
56) 2,4-Dinitrotoluene	10.05	165	197543	25.00	nq	# 65
57) Fluorene	10.41	166	655427	26.12	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.20	232	156960	30.47	nq	96
59) Diethylphthalate	10.30	149	743169	28.40	nq	99
60) 4-Chlorophenyl-phenylether	10.41	204	299022	28.29	nq	95
61) 4-Nitroaniline	10.42	138	183672	28.02	nq	94
62) Azobenzene	10.57	77	787507	29.79	nq	94
64) 4,6-Dinitro-2-methylphenol	10.46	198	103775	33.00	nq	# 61
65) n-Nitrosodiphenylamine	10.53	169	607964	25.04	nq	99
66) 4-Bromophenyl-phenylether	10.90	248	189622	27.89	nq	# 86
67) Hexachlorobenzene	10.97	284	182631	25.69	nq	# 85
68) Atrazine	11.05	200	179480	25.67	nq	93
69) Pentachlorophenol	11.16	266	116803	28.68	nq	99
70) Phenanthrene	11.37	178	979474	24.08	nq	99
71) Anthracene	11.42	178	1039281	26.11	nq	99
72) Carbazole	11.58	167	1120213	26.89	nq	98
73) Di-n-butylphthalate	11.92	149	1195940	28.08	nq	99
74) Fluoranthene	12.55	202	1014053	25.73	nq	96
76) Benzidine	12.68	184	654971	30.06	nq	95
77) Pyrene	12.78	202	1002390	21.14	nq	100
79) Butylbenzylphthalate	13.41	149	462502	22.96	nq	# 81
80) Benzo(a)anthracene	13.97	228	838836	23.44	nq	99
81) 3,3'-Dichlorobenzidine	13.93	252	340150	27.10	nq	# 93
82) Chrysene	14.00	228	834552	24.13	nq	100
83) Bis(2-ethylhexyl)phthalate	13.98	149	619406	27.97	nq	# 95
84) Di-n-octyl phthalate	14.59	149	1137867	27.87	nq	99
85) Indeno(1,2,3-cd)pyrene	16.75	276	711217	24.28	nq	99
87) Benzo(b)fluoranthene	14.99	252	811917m	24.76	nq	
88) Benzo(k)fluoranthene	15.01	252	692744m	23.90	nq	
89) Benzo(a)pyrene	15.33	252	735433	25.38	nq	# 97
90) Dibenzo(a,h)anthracene	16.76	278	614488	25.06	nq	96
91) Benzo(g,h,i)perylene	17.15	276	587345	23.56	nq	99

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

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