

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071219\
 Data File : BF115524.D
 Acq On : 12 Jul 2019 12:14
 Operator : HP/JU
 Sample : SSTDICC100
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC100

Manual Integrations
 APPROVED

mohammad
 7/15/2019 2:57:20 PM

Quant Time: Jul 12 14:06:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 13:09:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	192136	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	789519	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	374868	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	699310	20.00	ng	0.00
76) Chrysene-d12	14.06	240	370881	20.00	ng	0.00
87) Perylene-d12	15.52	264	379351	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	2079414	167.21	ng	0.01
7) Phenol-d6	6.54	99	2618438	165.56	ng	0.02
23) Nitrobenzene-d5	7.47	82	2487233	186.13	ng	0.02
42) 2,4,6-Tribromophenol	10.73	330	762095	183.87	ng	0.01
45) 2-Fluorobiphenyl	9.26	172	3546114	165.99	ng	0.00
79) Terphenyl-d14	13.00	244	3574345	197.33	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.71	88	560036	91.813	ng	99
3) Pyridine	3.47	79	1560359	92.511	ng	99
4) n-Nitrosodimethylamine	3.45	42	582493	85.164	ng	97
6) Aniline	6.57	93	1845358	82.683	ng	99
8) 2-Chlorophenol	6.69	128	1181407	84.134	ng	97
10) Phenol	6.55	94	1554331	82.140	ng	83
11) bis(2-Chloroethyl)ether	6.63	93	1213503	86.444	ng	100
12) 1,3-Dichlorobenzene	6.84	146	1342503	87.266	ng	98
13) 1,4-Dichlorobenzene	6.91	146	1315103	85.342	ng	98
14) 1,2-Dichlorobenzene	7.07	146	1135764	79.608	ng	96
15) Benzyl Alcohol	7.05	79	1006187	79.282	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.17	45	1531830	78.041	ng	97
17) 2-Methylphenol	7.14	107	1090954	90.688	ng	98
18) Hexachloroethane	7.41	117	502457	87.528	ng	97
19) n-Nitroso-di-n-propylamine	7.33	70	896424	83.778	ng	99
22) Acetophenone	7.32	105	1531667	80.962	ng	# 92
24) Nitrobenzene	7.49	77	1297557	85.980	ng	99
25) Isophorone	7.73	82	2598581	89.717	ng	100
26) 2-Nitrophenol	7.79	139	724643	116.795	ng	92
27) 2,4-Dimethylphenol	7.83	122	1047888	88.573	ng	98
28) bis(2-Chloroethoxy)methane	7.93	93	1479410	82.831	ng	99
29) 2,4-Dichlorophenol	8.04	162	1043763	88.578	ng	98
30) 1,2,4-Trichlorobenzene	8.12	180	1158650	88.324	ng	99
31) Naphthalene	8.20	128	3125887	79.711	ng	98
32) Benzoic acid	8.02	122	987205	122.119	ng	97
33) 4-Chloroaniline	8.25	127	1543516	88.216	ng	97
34) Hexachlorobutadiene	8.32	225	663868	89.236	ng	100
35) Caprolactam	8.67	113	320322m	83.464	ng	
36) 4-Chloro-3-methylphenol	8.73	107	1111629	89.187	ng	100
37) 2-Methylnaphthalene	8.89	142	2104696	80.296	ng	97
38) 1-Methylnaphthalene	8.99	142	1968694	78.718	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.06	216	956202	88.665	ng	98
41) Hexachlorocyclopentadiene	9.05	237	606980	97.108	ng	99
43) 2,4,6-Trichlorophenol	9.17	196	772684	98.414	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.22	196	759759	93.322	nq	98
46) 1,1'-Biphenyl	9.36	154	2382863	80.856	nq	98
47) 2-Chloronaphthalene	9.39	162	2058918	84.532	nq	98
48) 2-Nitroaniline	9.47	65	722647	100.985	nq	99
49) Acenaphthylene	9.80	152	2958344	80.222	nq	98
50) Dimethylphthalate	9.67	163	2575028	91.519	nq	99
51) 2,6-Dinitrotoluene	9.72	165	601593	98.587	nq	97
52) Acenaphthene	9.97	154	1994866	85.954	nq	99
53) 3-Nitroaniline	9.89	138	691211	97.527	nq	98
54) 2,4-Dinitrophenol	9.99	184	370857	173.684	nq	# 91
56) 4-Nitrophenol	10.05	139	524868	95.720	nq	98
57) 2,4-Dinitrotoluene	10.13	165	751766	96.939	nq	95
58) Fluorene	10.49	166	1901816	73.794	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.26	232	611184	86.903	nq	99
60) Diethylphthalate	10.36	149	2281201	82.370	nq	99
62) 4-Nitroaniline	10.52	138	661122	95.033	nq	98
63) Azobenzene	10.63	77	2172316	76.093	nq	98
65) 4,6-Dinitro-2-methylphenol	10.54	198	451792	148.711	nq	91
66) n-Nitrosodiphenylamine	10.60	169	1868966	84.822	nq	98
67) 4-Bromophenyl-phenylether	10.96	248	703942	89.711	nq	96
68) Hexachlorobenzene	11.04	284	808116	91.779	nq	95
70) Pentachlorophenol	11.22	266	499474	93.080	nq	99
71) Phenanthrene	11.44	178	2968058	80.701	nq	98
73) Carbazole	11.65	167	2735434	81.074	nq	99
78) Pyrene	12.86	202	2999821	104.011	nq	98
80) Butylbenzylphthalate	13.47	149	1289698	102.345	nq	99
81) Benzo(a)anthracene	14.04	228	2153824	90.635	nq	97
82) 3,3'-Dichlorobenzidine	14.00	252	701662	81.373	nq	97
83) Chrysene	14.09	228	2177059	89.163	nq	97
84) Bis(2-ethylhexyl)phthalate	14.03	149	1491220	95.556	nq	# 98
85) Di-n-octyl phthalate	14.64	149	2388190	85.963	nq	99
86) Indeno(1,2,3-cd)pyrene	17.02	276	2347544	111.092	nq	99
88) Benzo(b)fluoranthene	15.10	252	2290874	92.751	nq	99
89) Benzo(k)fluoranthene	15.13	252	1804788	82.163	nq	99
90) Benzo(a)pyrene	15.47	252	1927178	90.464	nq	100
91) Dibenzo(a,h)anthracene	17.04	278	1832665	100.673	nq	98
92) Benzo(g,h,i)perylene	17.48	276	1950784	113.217	nq	98

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

