

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052519\
 Data File : BF114630.D
 Acq On : 26 May 2019 4:33
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

mohammad
 5/27/2019 3:50:51 AM

Quant Time: May 27 01:37:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 24 07:23:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	149693	20.00	ng	-0.02
21) Naphthalene-d8	8.09	136	584930	20.00	ng	-0.01
39) Acenaphthene-d10	9.84	164	269741	20.00	ng	-0.01
64) Phenanthrene-d10	11.33	188	511338	20.00	ng	-0.01
76) Chrysene-d12	13.97	240	361132	20.00	ng	-0.02
87) Perylene-d12	15.43	264	352385	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	768779	82.65	ng	-0.02
7) Phenol-d6	6.45	99	935342	85.02	ng	-0.01
23) Nitrobenzene-d5	7.37	82	879819	84.41	ng	-0.01
42) 2,4,6-Tribromophenol	10.63	330	211035	75.68	ng	-0.01
45) 2-Fluorobiphenyl	9.16	172	1523112	85.41	ng	-0.02
79) Terphenyl-d14	12.90	244	1464486	83.60	ng	-0.01

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.51	88	178631	34.937	ng	100
3) Pyridine	3.26	79	460706	32.704	ng	97
4) n-Nitrosodimethylamine	3.22	42	239327	35.230	ng	100
6) Aniline	6.47	93	652118	41.034	ng	# 77
8) 2-Chlorophenol	6.59	128	430745	43.377	ng	95
9) Benzaldehyde	6.36	77	269404	34.978	ng	100
10) Phenol	6.46	94	538198	41.585	ng	# 73
11) bis(2-Chloroethyl)ether	6.53	93	409403	41.667	ng	99
12) 1,3-Dichlorobenzene	6.75	146	475533	42.661	ng	97
13) 1,4-Dichlorobenzene	6.82	146	483294	43.026	ng	99
14) 1,2-Dichlorobenzene	6.97	146	443836	43.250	ng	99
15) Benzyl Alcohol	6.95	79	335832	39.138	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.07	45	750842	39.412	ng	60
17) 2-Methylphenol	7.07	107	336098	41.202	ng	# 91
18) Hexachloroethane	7.31	117	176285	41.811	ng	99
19) n-Nitroso-di-n-propylamine	7.22	70	285393	40.926	ng	99
20) 3+4-Methylphenols	7.22	107	429328	45.553	ng	# 87
22) Acetophenone	7.22	105	564158	43.537	ng	# 91
24) Nitrobenzene	7.39	77	450433	42.431	ng	99
25) Isophorone	7.63	82	779809	40.342	ng	99
26) 2-Nitrophenol	7.70	139	223919	43.476	ng	93
27) 2,4-Dimethylphenol	7.75	122	325781	40.274	ng	99
28) bis(2-Chloroethoxy)methane	7.83	93	519202	41.396	ng	99
29) 2,4-Dichlorophenol	7.96	162	347085	43.091	ng	98
30) 1,2,4-Trichlorobenzene	8.03	180	386110	42.155	ng	97
31) Naphthalene	8.10	128	1193962	43.698	ng	98
32) Benzoic acid	7.90	122	168662m	31.986	ng	
33) 4-Chloroaniline	8.16	127	533273	42.830	ng	97
34) Hexachlorobutadiene	8.22	225	222655	42.723	ng	98
35) Caprolactam	8.53	113	112742	42.925	ng	96
36) 4-Chloro-3-methylphenol	8.65	107	358757	44.248	ng	95
37) 2-Methylnaphthalene	8.80	142	774527	43.936	ng	98
38) 1-Methylnaphthalene	8.90	142	727466	43.820	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.97	216	357177	43.713	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.95	237	120412	34.246	nq	97
43) 2,4,6-Trichlorophenol	9.09	196	221149	39.348	nq	98
44) 2,4,5-Trichlorophenol	9.13	196	214785	37.264	nq	94
46) 1,1'-Biphenyl	9.26	154	952377	43.704	nq	97
47) 2-Chloronaphthalene	9.29	162	751561	42.854	nq	99
48) 2-Nitroaniline	9.39	65	262622	42.225	nq	96
49) Acenaphthylene	9.70	152	1134414	42.530	nq	98
50) Dimethylphthalate	9.56	163	817835	41.302	nq	99
51) 2,6-Dinitrotoluene	9.63	165	182051	41.451	nq	86
52) Acenaphthene	9.87	154	673814	41.167	nq	99
53) 3-Nitroaniline	9.80	138	223302	40.958	nq	98
54) 2,4-Dinitrophenol	9.93	184	71726	36.485	nq	97
55) Dibenzofuran	10.05	168	962666	40.109	nq	99
56) 4-Nitrophenol	9.99	139	111779	28.992	nq	94
57) 2,4-Dinitrotoluene	10.05	165	226006	37.913	nq	# 78
58) Fluorene	10.39	166	795421	42.906	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.17	232	187839	37.770	nq	99
60) Diethylphthalate	10.26	149	808613	40.190	nq	99
61) 4-Chlorophenyl-phenylether	10.37	204	387545	42.562	nq	98
62) 4-Nitroaniline	10.42	138	216077	37.717	nq	96
63) Azobenzene	10.54	77	771719	39.627	nq	97
65) 4,6-Dinitro-2-methylphenol	10.46	198	103200	42.002	nq	99
66) n-Nitrosodiphenylamine	10.50	169	682815	43.998	nq	99
67) 4-Bromophenyl-phenylether	10.86	248	241209	42.843	nq	97
68) Hexachlorobenzene	10.94	284	260765	41.867	nq	94
69) Atrazine	11.02	200	198068	41.012	nq	97
70) Pentachlorophenol	11.14	266	87071	29.491	nq	98
71) Phenanthrene	11.35	178	1113574	44.763	nq	98
72) Anthracene	11.40	178	1119749	44.813	nq	99
73) Carbazole	11.56	167	1077910	45.238	nq	98
74) Di-n-butylphthalate	11.88	149	1220285	44.453	nq	99
75) Fluoranthene	12.54	202	1196160	44.650	nq	99
77) Benzidine	12.66	184	537777	33.174	nq	97
78) Pyrene	12.77	202	1211809	41.713	nq	99
80) Butylbenzylphthalate	13.37	149	530973	43.423	nq	98
81) Benzo(a)anthracene	13.96	228	1046641	44.809	nq	99
82) 3,3'-Dichlorobenzidine	13.92	252	401633	44.325	nq	100
83) Chrysene	14.00	228	1001603	43.743	nq	98
84) Bis(2-ethylhexyl)phthalate	13.93	149	709966	44.393	nq	99
85) Di-n-octyl phthalate	14.54	149	1138618	43.920	nq	99
86) Indeno(1,2,3-cd)pyrene	16.90	276	943624	46.631	nq	99
88) Benzo(b)fluoranthene	15.00	252	971654	43.040	nq	99
89) Benzo(k)fluoranthene	15.03	252	870051	41.954	nq	98
90) Benzo(a)pyrene	15.37	252	850328	42.798	nq	98
91) Dibenzo(a,h)anthracene	16.89	278	779612	47.221	nq	97
92) Benzo(g,h,i)perylene	17.33	276	792898	47.923	nq	99

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

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