

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103118\
 Data File : BF110325.D
 Acq On : 31 Oct 2018 11:49
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 11/1/2018 10:45:51 AM

Quant Time: Oct 31 14:10:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 29 12:32:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	88443	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	394560	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	194755	20.00	ng	0.00
64) Phenanthrene-d10	11.52	188	373330	20.00	ng	0.00
76) Chrysene-d12	14.17	240	283091	20.00	ng	0.01
87) Perylene-d12	15.70	264	223047	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.59	112	456426	84.04	ng	0.00
7) Phenol-d6	6.60	99	586658	84.45	ng	0.01
23) Nitrobenzene-d5	7.54	82	562953	84.63	ng	0.00
42) 2,4,6-Tribromophenol	10.82	330	164539	85.29	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	1051689	84.79	ng	0.00
79) Terphenyl-d14	13.10	244	1146125	84.20	ng	0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.79	88	106779	37.526	ng	91
3) Pyridine	3.58	79	257152	40.574	ng	87
4) n-Nitrosodimethylamine	3.55	42	114177	43.000	ng	88
6) Aniline	6.64	93	380187	42.013	ng	# 54
8) 2-Chlorophenol	6.76	128	268334	42.854	ng	97
9) Benzaldehyde	6.53	77	156018	38.168	ng	99
10) Phenol	6.62	94	340434	45.846	ng	# 65
11) bis(2-Chloroethyl)ether	6.71	93	249205	40.792	ng	97
12) 1,3-Dichlorobenzene	6.92	146	280519	40.709	ng	98
13) 1,4-Dichlorobenzene	6.99	146	282990	40.606	ng	98
14) 1,2-Dichlorobenzene	7.15	146	269324	41.629	ng	98
15) Benzyl Alcohol	7.12	79	239004	44.733	ng	91
16) 2,2'-oxybis(1-Chloropropan	7.24	45	398670	40.814	ng	70
17) 2-Methylphenol	7.22	107	213312	42.746	ng	# 82
18) Hexachloroethane	7.49	117	108121	42.375	ng	96
19) n-Nitroso-di-n-propylamine	7.39	70	192185	43.123	ng	# 96
20) 3+4-Methylphenols	7.37	107	280570	43.496	ng	96
22) Acetophenone	7.39	105	376038	41.361	ng	99
24) Nitrobenzene	7.56	77	288683	42.034	ng	92
25) Isophorone	7.79	82	457921	40.384	ng	98
26) 2-Nitrophenol	7.87	139	148228	45.355	ng	94
27) 2,4-Dimethylphenol	7.90	122	219342	40.480	ng	98
28) bis(2-Chloroethoxy)methane	8.00	93	313406	40.685	ng	98
29) 2,4-Dichlorophenol	8.12	162	227796	41.612	ng	100
30) 1,2,4-Trichlorobenzene	8.20	180	247813	40.415	ng	95
31) Naphthalene	8.28	128	739612	40.672	ng	99
32) Benzoic acid	8.02	122	116198m	27.746	ng	
33) 4-Chloroaniline	8.33	127	340349	41.586	ng	99
34) Hexachlorobutadiene	8.39	225	140727	41.334	ng	99
35) Caprolactam	8.71	113	78309	41.042	ng	# 86
36) 4-Chloro-3-methylphenol	8.81	107	249377	42.127	ng	93
37) 2-Methylnaphthalene	8.97	142	499914	41.550	ng	97
38) 1-Methylnaphthalene	9.07	142	479529	41.243	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.14	216	245707	40.491	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.12	237	110734	35.688	nq	100
43) 2,4,6-Trichlorophenol	9.25	196	157038	40.123	nq	94
44) 2,4,5-Trichlorophenol	9.29	196	170207	41.141	nq	96
46) 1,1'-Biphenyl	9.44	154	721724	40.980	nq	99
47) 2-Chloronaphthalene	9.47	162	517584	40.065	nq	99
48) 2-Nitroaniline	9.56	65	166365	42.497	nq	95
49) Acenaphthylene	9.89	152	764901	40.994	nq	99
50) Dimethylphthalate	9.73	163	597003	41.527	nq	100
51) 2,6-Dinitrotoluene	9.80	165	131924	42.601	nq	96
52) Acenaphthene	10.06	154	471330	38.183	nq	98
53) 3-Nitroaniline	9.98	138	157393	42.740	nq	89
54) 2,4-Dinitrophenol	10.09	184	49388	42.760	nq	93
55) Dibenzofuran	10.23	168	702292	40.636	nq	99
56) 4-Nitrophenol	10.13	139	108163	39.685	nq	98
57) 2,4-Dinitrotoluene	10.22	165	177960	45.243	nq	# 85
58) Fluorene	10.57	166	535895	41.664	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.35	232	135115	40.068	nq	94
60) Diethylphthalate	10.43	149	590415	42.071	nq	100
61) 4-Chlorophenyl-phenylether	10.56	204	280196	41.294	nq	99
62) 4-Nitroaniline	10.60	138	158171	43.184	nq	90
63) Azobenzene	10.72	77	563910	40.482	nq	96
65) 4,6-Dinitro-2-methylphenol	10.63	198	70858	41.543	nq	95
66) n-Nitrosodiphenylamine	10.68	169	501485	40.953	nq	96
67) 4-Bromophenyl-phenylether	11.05	248	165438	40.853	nq	93
68) Hexachlorobenzene	11.13	284	177457	41.301	nq	99
69) Atrazine	11.20	200	161957	41.852	nq	99
70) Pentachlorophenol	11.32	266	94167	37.585	nq	99
71) Phenanthrene	11.54	178	797443	40.823	nq	99
72) Anthracene	11.59	178	806699	41.200	nq	100
73) Carbazole	11.75	167	794627	41.565	nq	99
74) Di-n-butylphthalate	12.06	149	925737	42.875	nq	100
75) Fluoranthene	12.74	202	817152	41.218	nq	98
77) Benzidine	12.85	184	357472	44.751	nq	98
78) Pyrene	12.97	202	848423	41.804	nq	98
80) Butylbenzylphthalate	13.56	149	382342	43.404	nq	98
81) Benzo(a)anthracene	14.16	228	689048	41.044	nq	99
82) 3,3'-Dichlorobenzidine	14.12	252	228164	39.030	nq	99
83) Chrysene	14.20	228	652863	40.944	nq	98
84) Bis(2-ethylhexyl)phthalate	14.12	149	481487	40.434	nq	96
85) Di-n-octyl phthalate	14.74	149	790537	42.695	nq	99
86) Indeno(1,2,3-cd)pyrene	17.30	276	517914	39.153	nq	98
88) Benzo(b)fluoranthene	15.25	252	541447	42.037	nq	98
89) Benzo(k)fluoranthene	15.28	252	528254	41.718	nq	98
90) Benzo(a)pyrene	15.64	252	492004	41.513	nq	99
91) Dibenzo(a,h)anthracene	17.32	278	429201	41.970	nq	99
92) Benzo(g,h,i)perylene	17.79	276	424325	42.107	nq	99

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

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