

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101419\
 Data File : BF117399.D
 Acq On : 14 Oct 2019 13:56
 Operator : JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

mohammad
 10/16/2019 8:10:02 AM

Quant Time: Oct 14 14:38:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 14 14:19:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	67906	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	293500	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	148572	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	224489	20.00	ng	0.00
76) Chrysene-d12	13.97	240	136996	20.00	ng	0.00
87) Perylene-d12	15.42	264	111519	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	513440	118.05	ng	0.00
7) Phenol-d6	6.45	99	687156	122.25	ng	0.01
23) Nitrobenzene-d5	7.37	82	678070	121.39	ng	0.01
42) 2,4,6-Tribromophenol	10.63	330	140033	112.77	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1163433	122.44	ng	0.00
79) Terphenyl-d14	12.90	244	1036384	141.41	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.52	88	108413	59.567	ng	99
3) Pyridine	3.26	79	291499	58.515	ng	96
4) n-Nitrosodimethylamine	3.24	42	107470	62.864	ng	93
6) Aniline	6.47	93	431587	59.635	ng	# 78
8) 2-Chlorophenol	6.59	128	289980	61.800	ng	98
9) Benzaldehyde	6.35	77	150046	48.886	ng	97
10) Phenol	6.46	94	378613	61.099	ng	97
11) bis(2-Chloroethyl)ether	6.54	93	283897	62.335	ng	99
12) 1,3-Dichlorobenzene	6.74	146	324423	61.615	ng	98
13) 1,4-Dichlorobenzene	6.81	146	328579	61.150	ng	99
14) 1,2-Dichlorobenzene	6.97	146	306625	61.291	ng	99
15) Benzyl Alcohol	6.95	79	261356	60.632	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.07	45	300838	66.858	ng	# 62
17) 2-Methylphenol	7.07	107	242672	61.709	ng	# 92
18) Hexachloroethane	7.30	117	128752	62.285	ng	95
19) n-Nitroso-di-n-propylamine	7.23	70	225480	65.875	ng	95
20) 3+4-Methylphenols	7.22	107	308923	63.067	ng	93
22) Acetophenone	7.22	105	449492	60.278	ng	# 96
24) Nitrobenzene	7.39	77	332073	60.198	ng	97
25) Isophorone	7.63	82	608764	61.551	ng	99
26) 2-Nitrophenol	7.70	139	154673	65.278	ng	98
27) 2,4-Dimethylphenol	7.75	122	228037	60.689	ng	96
28) bis(2-Chloroethoxy)methane	7.83	93	378481	62.111	ng	98
29) 2,4-Dichlorophenol	7.95	162	239969	60.950	ng	99
30) 1,2,4-Trichlorobenzene	8.02	180	253103	59.503	ng	97
31) Naphthalene	8.10	128	855030	60.574	ng	100
32) Benzoic acid	7.93	122	168223	63.655	ng	89
33) 4-Chloroaniline	8.16	127	355941	56.853	ng	98
34) Hexachlorobutadiene	8.22	225	143483	59.456	ng	97
35) Caprolactam	8.56	113	72594m	54.474	ng	
36) 4-Chloro-3-methylphenol	8.65	107	265418	57.815	ng	96
37) 2-Methylnaphthalene	8.79	142	572474	60.227	ng	98
38) 1-Methylnaphthalene	8.89	142	532722	59.840	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.96	216	233710	60.917	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.94	237	63860	44.254	nq	96
43) 2,4,6-Trichlorophenol	9.08	196	150568	57.831	nq	94
44) 2,4,5-Trichlorophenol	9.13	196	147366	52.977	nq	95
46) 1,1'-Biphenyl	9.26	154	714090	61.274	nq	98
47) 2-Chloronaphthalene	9.29	162	547300	59.505	nq	100
48) 2-Nitroaniline	9.39	65	177773	60.260	nq	98
49) Acenaphthylene	9.70	152	793731	57.607	nq	99
50) Dimethylphthalate	9.56	163	622028	59.130	nq	98
51) 2,6-Dinitrotoluene	9.63	165	131962	61.592	nq	98
52) Acenaphthene	9.87	154	489321	59.910	nq	100
53) 3-Nitroaniline	9.80	138	159808	59.118	nq	90
54) 2,4-Dinitrophenol	9.92	184	53765	70.894	nq	94
55) Dibenzofuran	10.04	168	704173	58.434	nq	98
56) 4-Nitrophenol	9.98	139	74895	42.749	nq	97
57) 2,4-Dinitrotoluene	10.05	165	170921	61.458	nq	89
58) Fluorene	10.39	166	584337	60.196	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.17	232	115070	54.057	nq	99
60) Diethylphthalate	10.26	149	620060	59.912	nq	99
61) 4-Chlorophenyl-phenylether	10.37	204	262171	62.422	nq	99
62) 4-Nitroaniline	10.43	138	142866	50.820	nq	99
63) Azobenzene	10.53	77	632811	59.867	nq	99
65) 4,6-Dinitro-2-methylphenol	10.46	198	75598	80.237	nq	98
66) n-Nitrosodiphenylamine	10.50	169	499081	64.373	nq	98
67) 4-Bromophenyl-phenylether	10.86	248	149513	65.216	nq	95
68) Hexachlorobenzene	10.94	284	159438	63.584	nq	99
69) Atrazine	11.02	200	50393	26.461	nq	98
70) Pentachlorophenol	11.14	266	51330	43.670	nq	90
71) Phenanthrene	11.35	178	779673	61.147	nq	99
72) Anthracene	11.40	178	805200	61.854	nq	99
73) Carbazole	11.56	167	710671	57.514	nq	99
74) Di-n-butylphthalate	11.87	149	979950	66.657	nq	99
75) Fluoranthene	12.54	202	757744	57.836	nq	99
77) Benzidine	12.65	184	112429	25.477	nq	100
78) Pyrene	12.77	202	750713	65.308	nq	99
80) Butylbenzylphthalate	13.37	149	377588	69.517	nq	97
81) Benzo(a)anthracene	13.96	228	552466	60.360	nq	99
82) 3,3'-Dichlorobenzidine	13.91	252	143107	48.520	nq	# 97
83) Chrysene	14.00	228	529218	59.283	nq	98
84) Bis(2-ethylhexyl)phthalate	13.93	149	522679	74.634	nq	97
85) Di-n-octyl phthalate	14.54	149	747445	67.810	nq	99
86) Indeno(1,2,3-cd)pyrene	16.89	276	408453	62.715	nq	99
88) Benzo(b)fluoranthene	15.00	252	433219	64.132	nq	99
89) Benzo(k)fluoranthene	15.03	252	368066	57.520	nq	100
90) Benzo(a)pyrene	15.36	252	357925	60.382	nq	98
91) Dibenzo(a,h)anthracene	16.88	278	339836	71.965	nq	98
92) Benzo(g,h,i)perylene	17.31	276	333447	70.861	nq	98

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

