

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110818\
 Data File : BF110581.D
 Acq On : 8 Nov 2018 13:29
 Operator : JU/SJ
 Sample : SSTDIC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleID :
 SSTDIC080

Manual Integrations
 APPROVED

Sohil

11/9/2018 4:26:05 PM

Quant Time: Nov 08 14:15:59 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 08 13:42:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	80266	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	336836	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	158737	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	288917	20.00	ng	0.00
76) Chrysene-d12	14.14	240	175118	20.00	ng	0.00
87) Perylene-d12	15.66	264	137033	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.57	112	771639	156.55	ng	0.00
7) Phenol-d6	6.59	99	968397	153.60	ng	0.01
23) Nitrobenzene-d5	7.53	82	908207	159.93	ng	0.01
42) 2,4,6-Tribromophenol	10.79	330	235484	149.76	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	1558920	154.21	ng	0.00
79) Terphenyl-d14	13.07	244	1478089	175.54	ng	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.79	88	198173	76.739	ng	91
3) Pyridine	3.57	79	442986	77.016	ng	# 85
4) n-Nitrosodimethylamine	3.56	42	198645	82.432	ng	91
6) Aniline	6.63	93	622795	75.833	ng	# 53
8) 2-Chlorophenol	6.74	128	441504	77.693	ng	97
9) Benzaldehyde	6.50	77	207504	54.212	ng	98
10) Phenol	6.60	94	563240	83.578	ng	# 73
11) bis(2-Chloroethyl)ether	6.69	93	429721	77.506	ng	95
12) 1,3-Dichlorobenzene	6.89	146	485540	77.640	ng	99
13) 1,4-Dichlorobenzene	6.97	146	490056	77.481	ng	98
14) 1,2-Dichlorobenzene	7.12	146	446076	75.973	ng	99
15) Benzyl Alcohol	7.10	79	371574	76.630	ng	94
16) 2,2'-oxybis(1-Chloropropan	7.22	45	644023	72.649	ng	70
17) 2-Methylphenol	7.20	107	351709	77.659	ng	# 85
18) Hexachloroethane	7.46	117	184168	79.533	ng	93
19) n-Nitroso-di-n-propylamine	7.37	70	308598	76.298	ng	99
20) 3+4-Methylphenols	7.36	107	428917	73.268	ng	# 87
22) Acetophenone	7.37	105	585331	75.415	ng	96
24) Nitrobenzene	7.55	77	462503	78.884	ng	95
25) Isophorone	7.78	82	756812	78.180	ng	96
26) 2-Nitrophenol	7.85	139	246352	88.297	ng	99
27) 2,4-Dimethylphenol	7.88	122	355455	76.843	ng	98
28) bis(2-Chloroethoxy)methane	7.97	93	499877	76.013	ng	100
29) 2,4-Dichlorophenol	8.09	162	363336	77.746	ng	99
30) 1,2,4-Trichlorobenzene	8.17	180	408030	77.947	ng	95
31) Naphthalene	8.26	128	1187506	76.493	ng	99
32) Benzoic acid	8.05	122	287619	80.449	ng	97
33) 4-Chloroaniline	8.30	127	526003	75.285	ng	100
34) Hexachlorobutadiene	8.36	225	230308	79.238	ng	98
35) Caprolactam	8.72	113	121528m	74.609	ng	
36) 4-Chloro-3-methylphenol	8.79	107	388389	76.854	ng	97
37) 2-Methylnaphthalene	8.95	142	773714	75.327	ng	100
38) 1-Methylnaphthalene	9.05	142	741222	74.675	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.12	216	378079	76.443	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.09	237	158340	62.609	ng	100
43) 2,4,6-Trichlorophenol	9.23	196	240047	75.248	ng	93
44) 2,4,5-Trichlorophenol	9.27	196	256836	76.167	ng	96
46) 1,1'-Biphenyl	9.42	154	1081490	75.342	ng	99
47) 2-Chloronaphthalene	9.45	162	801134	76.085	ng	99
48) 2-Nitroaniline	9.55	65	257457	80.688	ng	92
49) Acenaphthylene	9.86	152	1127533	74.140	ng	99
50) Dimethylphthalate	9.71	163	892986	76.209	ng	99
51) 2,6-Dinitrotoluene	9.79	165	198324	78.575	ng	95
52) Acenaphthene	10.03	154	718612	71.426	ng	98
53) 3-Nitroaniline	9.96	138	231571	77.151	ng	89
54) 2,4-Dinitrophenol	10.07	184	83043	93.647	ng	91
55) Dibenzofuran	10.20	168	1019088	72.346	ng	97
56) 4-Nitrophenol	10.12	139	163442	73.573	ng	96
57) 2,4-Dinitrotoluene	10.19	165	254741	79.459	ng	99
58) Fluorene	10.55	166	790724	75.425	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.32	232	203636	74.090	ng	95
60) Diethylphthalate	10.41	149	862445	75.400	ng	100
61) 4-Chlorophenyl-phenylether	10.53	204	401604	72.617	ng	98
62) 4-Nitroaniline	10.59	138	227719	76.280	ng	91
63) Azobenzene	10.69	77	849186	74.794	ng	97
65) 4,6-Dinitro-2-methylphenol	10.61	198	130107	98.567	ng	88
66) n-Nitrosodiphenylamine	10.66	169	736288	77.696	ng	97
67) 4-Bromophenyl-phenylether	11.02	248	244450	78.002	ng	# 90
68) Hexachlorobenzene	11.10	284	256229	77.057	ng	94
70) Pentachlorophenol	11.29	266	147380	76.011	ng	97
71) Phenanthrene	11.52	178	1134462	75.043	ng	99
72) Anthracene	11.57	178	1137401	75.062	ng	100
73) Carbazole	11.72	167	1104910	74.681	ng	99
74) Di-n-butylphthalate	12.03	149	1306368	78.181	ng	99
75) Fluoranthene	12.70	202	1124073	73.265	ng	99
78) Pyrene	12.94	202	1137999	90.645	ng	99
80) Butylbenzylphthalate	13.53	149	481534	88.368	ng	100
81) Benzo(a)anthracene	14.13	228	820540	79.013	ng	99
82) 3,3'-Dichlorobenzidine	14.08	252	230468	63.732	ng	# 98
83) Chrysene	14.17	228	764825	77.540	ng	100
84) Bis(2-ethylhexyl)phthalate	14.08	149	574624	78.009	ng	97
85) Di-n-octyl phthalate	14.70	149	857343	74.852	ng	98
86) Indeno(1,2,3-cd)pyrene	17.25	276	677594	82.807	ng	98
88) Benzo(b)fluoranthene	15.21	252	621455	78.534	ng	99
89) Benzo(k)fluoranthene	15.24	252	615555m	79.126	ng	
90) Benzo(a)pyrene	15.60	252	585340	80.388	ng	99
91) Dibenzo(a,h)anthracene	17.26	278	555151	88.361	ng	100
92) Benzo(g,h,i)perylene	17.72	276	572857	92.529	ng	98

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

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