

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110518\
 Data File : BF110467.D
 Acq On : 5 Nov 2018 16:16
 Operator : JU/SJ
 Sample : J5812-09MS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP2-AMS

Manual Integrations
 APPROVED

Sohil
 11/6/2018 1:19:13 PM

Quant Time: Nov 06 10:23:32 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 16:06:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	89244	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	380315	20.00	ng	0.00
39) Acenaphthene-d10	10.00	164	186447	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	340011	20.00	ng	0.00
76) Chrysene-d12	14.14	240	193434	20.00	ng	0.00
87) Perylene-d12	15.66	264	160267	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.59	112	734143	133.96	ng	0.01
7) Phenol-d6	6.59	99	952624	135.90	ng	0.00
23) Nitrobenzene-d5	7.52	82	604208	94.23	ng	0.00
42) 2,4,6-Tribromophenol	10.80	330	243228	131.70	ng	0.00
45) 2-Fluorobiphenyl	9.32	172	1058076	89.11	ng	0.00
79) Terphenyl-d14	13.07	244	977782	105.13	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.87	88	82144	28.609	ng	88
3) Pyridine	3.64	79	227403	35.558	ng	87
4) n-Nitrosodimethylamine	3.60	42	120726	45.058	ng	88
6) Aniline	6.62	93	306352	33.550	ng	# 54
8) 2-Chlorophenol	6.75	128	274481	43.442	ng	99
9) Benzaldehyde	6.51	77	164368	40.532	ng	98
10) Phenol	6.60	94	390269	52.085	ng	# 63
11) bis(2-Chloroethyl)ether	6.69	93	248362	40.289	ng	96
12) 1,3-Dichlorobenzene	6.90	146	278377	40.036	ng	97
13) 1,4-Dichlorobenzene	6.97	146	278499	39.603	ng	99
14) 1,2-Dichlorobenzene	7.13	146	265116	40.611	ng	99
15) Benzyl Alcohol	7.10	79	240260	44.564	ng	93
16) 2,2'-oxybis(1-Chloropropan	7.22	45	403081	40.896	ng	70
17) 2-Methylphenol	7.20	107	223789	44.443	ng	# 84
18) Hexachloroethane	7.46	117	104262	40.496	ng	93
19) n-Nitroso-di-n-propylamine	7.36	70	190137	42.281	ng	95
20) 3+4-Methylphenols	7.36	107	286150	43.963	ng	91
22) Acetophenone	7.37	105	382449	43.642	ng	98
24) Nitrobenzene	7.54	77	285412	43.114	ng	96
25) Isophorone	7.77	82	525003	48.034	ng	97
26) 2-Nitrophenol	7.86	139	147033	46.674	ng	92
27) 2,4-Dimethylphenol	7.89	122	259034	49.596	ng	99
28) bis(2-Chloroethoxy)methane	7.98	93	332177	44.737	ng	99
29) 2,4-Dichlorophenol	8.10	162	235312	44.596	ng	98
30) 1,2,4-Trichlorobenzene	8.18	180	243839	41.256	ng	98
31) Naphthalene	8.26	128	787633	44.935	ng	100
32) Benzoic acid	7.96	122	4781m	1.184	ng	
33) 4-Chloroaniline	8.31	127	186962	23.700	ng	98
34) Hexachlorobutadiene	8.37	225	130657	39.814	ng	97
35) Caprolactam	8.69	113	71678m	38.974	ng	
36) 4-Chloro-3-methylphenol	8.79	107	261585	45.845	ng	97
37) 2-Methylnaphthalene	8.95	142	552463	47.637	ng	99
38) 1-Methylnaphthalene	9.05	142	517149	46.144	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.12	216	232034	39.942	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.10	237	206925	69.660	nq	100
43) 2,4,6-Trichlorophenol	9.23	196	160464	42.825	nq	95
44) 2,4,5-Trichlorophenol	9.27	196	174233	43.991	nq	96
46) 1,1'-Biphenyl	9.42	154	674852	40.026	nq	100
47) 2-Chloronaphthalene	9.45	162	510797	41.301	nq	100
48) 2-Nitroaniline	9.55	65	167092	44.585	nq	95
49) Acenaphthylene	9.86	152	798858	44.721	nq	99
50) Dimethylphthalate	9.72	163	741710	53.891	nq	100
51) 2,6-Dinitrotoluene	9.79	165	134496	45.367	nq	89
52) Acenaphthene	10.03	154	478756	40.513	nq	98
53) 3-Nitroaniline	9.96	138	116460	33.034	nq	90
54) 2,4-Dinitrophenol	10.07	184	19648	20.794	nq	95
55) Dibenzofuran	10.20	168	688416	41.608	nq	100
56) 4-Nitrophenol	10.13	139	273057	104.649	nq	94
57) 2,4-Dinitrotoluene	10.19	165	176293	46.817	nq	94
58) Fluorene	10.55	166	566279	45.988	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.32	232	136914m	42.411	nq	
60) Diethylphthalate	10.41	149	584076	43.474	nq	99
61) 4-Chlorophenyl-phenylether	10.53	204	265757	40.912	nq	96
62) 4-Nitroaniline	10.58	138	145902	41.610	nq	92
63) Azobenzene	10.70	77	570831	42.805	nq	97
65) 4,6-Dinitro-2-methylphenol	10.60	198	50183	32.305	nq	89
66) n-Nitrosodiphenylamine	10.66	169	493719	44.270	nq	97
67) 4-Bromophenyl-phenylether	11.03	248	157063	42.586	nq	94
68) Hexachlorobenzene	11.10	284	163258	41.720	nq	# 89
69) Atrazine	11.18	200	160804	45.626	nq	99
70) Pentachlorophenol	11.30	266	133243	58.393	nq	96
71) Phenanthrene	11.52	178	806706	45.344	nq	100
72) Anthracene	11.57	178	828064	46.435	nq	99
73) Carbazole	11.72	167	718429	41.262	nq	100
74) Di-n-butylphthalate	12.03	149	889459	45.231	nq	99
75) Fluoranthene	12.71	202	802927	44.469	nq	99
77) Benzidine	12.83	184	347373	Below	Cal	96
78) Pyrene	12.94	202	777529	56.068	nq	98
80) Butylbenzylphthalate	13.54	149	326623	54.264	nq	94
81) Benzo(a)anthracene	14.13	228	535913	46.719	nq	99
82) 3,3'-Dichlorobenzidine	14.08	252	119020	29.797	nq	99
83) Chrysene	14.16	228	498654	45.768	nq	99
84) Bis(2-ethylhexyl)phthalate	14.09	149	429161	52.745	nq	96
85) Di-n-octyl phthalate	14.70	149	606274	47.920	nq	99
86) Indeno(1,2,3-cd)pyrene	17.24	276	489655	54.174	nq	97
88) Benzo(b)fluoranthene	15.21	252	430063	46.469	nq	99
89) Benzo(k)fluoranthene	15.24	252	402542	44.243	nq	99
90) Benzo(a)pyrene	15.60	252	404065	47.448	nq	99
91) Dibenzo(a,h)anthracene	17.24	278	406522	55.324	nq	100
92) Benzo(g,h,i)perylene	17.71	276	412094	56.913	nq	97

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

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