

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050219\
 Data File : BF114084.D
 Acq On : 2 May 2019 14:21
 Operator : JU/SJ
 Sample : K2584-03MSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSSI-0419-S-DH-8MSD

Manual Integrations
 APPROVED

Sohil
 5/3/2019 4:32:58 PM

Quant Time: May 03 04:02:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	81868	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	304855	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	168627	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	345659	20.00	ng	0.00
76) Chrysene-d12	14.07	240	240071	20.00	ng	0.00
87) Perylene-d12	15.57	264	250971	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	845114	176.56	ng	0.00
7) Phenol-d6	6.53	99	1143316	190.38	ng	0.01
23) Nitrobenzene-d5	7.45	82	691693	140.09	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	377819	183.92	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1265576	117.38	ng	0.00
79) Terphenyl-d14	13.00	244	1790671	152.92	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.66	88	106392	47.152	ng	96
3) Pyridine	3.44	79	307754	49.969	ng	100
4) n-Nitrosodimethylamine	3.38	42	120608	57.207	ng	97
6) Aniline	6.56	93	172208	20.883	ng	100
8) 2-Chlorophenol	6.68	128	319497	58.464	ng	99
9) Benzaldehyde	6.43	77	56597	12.081	ng	97
10) Phenol	6.55	94	406961	56.671	ng	98
11) bis(2-Chloroethyl)ether	6.63	93	309708	57.313	ng	94
12) 1,3-Dichlorobenzene	6.83	146	340345	54.989	ng	100
13) 1,4-Dichlorobenzene	6.91	146	344513	55.341	ng	100
14) 1,2-Dichlorobenzene	7.06	146	329683	57.625	ng	98
15) Benzyl Alcohol	7.03	79	301481	74.539	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.16	45	347512	55.090	ng	93
17) 2-Methylphenol	7.15	107	276726	57.768	ng	97
18) Hexachloroethane	7.40	117	119966	54.973	ng	97
19) n-Nitroso-di-n-propylamine	7.30	70	224805	57.805	ng	98
20) 3+4-Methylphenols	7.30	107	338128	62.475	ng	# 70
22) Acetophenone	7.30	105	450637	66.426	ng	# 92
24) Nitrobenzene	7.47	77	355685	65.196	ng	100
25) Isophorone	7.71	82	614364	60.812	ng	100
26) 2-Nitrophenol	7.79	139	184094	73.967	ng	94
27) 2,4-Dimethylphenol	7.83	122	288278	71.091	ng	99
28) bis(2-Chloroethoxy)methane	7.92	93	392364	60.948	ng	99
29) 2,4-Dichlorophenol	8.03	162	308523	66.534	ng	97
30) 1,2,4-Trichlorobenzene	8.12	180	332517	64.312	ng	98
31) Naphthalene	8.20	128	959360	66.244	ng	99
32) Benzoic acid	7.93	122	89477	32.674	ng	97
33) 4-Chloroaniline	8.24	127	69100	11.276	ng	97
34) Hexachlorobutadiene	8.32	225	200857	62.317	ng	98
35) Caprolactam	8.61	113	98148m	66.537	ng	
36) 4-Chloro-3-methylphenol	8.73	107	279851	60.916	ng	98
37) 2-Methylnaphthalene	8.89	142	644090	65.952	ng	99
38) 1-Methylnaphthalene	8.99	142	602299	63.899	ng	97
40) 1,2,4,5-Tetrachlorobenzene	9.06	216	339924	62.058	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.05	237	235196	77.000	nq	99
43) 2,4,6-Trichlorophenol	9.17	196	216672	62.294	nq	100
44) 2,4,5-Trichlorophenol	9.21	196	232153	59.506	nq	99
46) 1,1'-Biphenyl	9.35	154	769831	59.761	nq	98
47) 2-Chloronaphthalene	9.37	162	627283	59.880	nq	97
48) 2-Nitroaniline	9.47	65	170041	61.919	nq	88
49) Acenaphthylene	9.79	152	915394	55.817	nq	99
50) Dimethylphthalate	9.65	163	865864	70.999	nq	99
51) 2,6-Dinitrotoluene	9.71	165	167452	64.002	nq	99
52) Acenaphthene	9.96	154	549682	56.714	nq	98
53) 3-Nitroaniline	9.87	138	71001	25.825	nq	96
54) 2,4-Dinitrophenol	9.99	184	137738	117.345	nq	# 1
55) Dibenzofuran	10.13	168	856508	58.996	nq	97
56) 4-Nitrophenol	10.05	139	281044	129.106	nq	98
57) 2,4-Dinitrotoluene	10.12	165	237083	71.791	nq	98
58) Fluorene	10.48	166	658210	60.219	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.26	232	207510	60.526	nq	99
60) Diethylphthalate	10.35	149	772707	66.722	nq	99
61) 4-Chlorophenyl-phenylether	10.47	204	348255	60.162	nq	99
62) 4-Nitroaniline	10.49	138	171365	63.871	nq	99
63) Azobenzene	10.63	77	661271	58.915	nq	97
65) 4,6-Dinitro-2-methylphenol	10.53	198	109373	65.832	nq	98
66) n-Nitrosodiphenylamine	10.59	169	620247	59.803	nq	99
67) 4-Bromophenyl-phenylether	10.96	248	237395	56.851	nq	97
68) Hexachlorobenzene	11.03	284	262014	57.100	nq	97
69) Atrazine	11.11	200	210958	62.526	nq	99
70) Pentachlorophenol	11.23	266	282241	108.636	nq	98
71) Phenanthrene	11.44	178	1115739	64.229	nq	100
72) Anthracene	11.49	178	1129825	64.406	nq	99
73) Carbazole	11.65	167	1048726	67.010	nq	100
74) Di-n-butylphthalate	11.97	149	1364843	74.137	nq	100
75) Fluoranthene	12.63	202	1306108	68.915	nq	99
77) Benzidine	12.74	184	199110	27.836	nq	97
78) Pyrene	12.86	202	1300953	77.491	nq	99
80) Butylbenzylphthalate	13.47	149	582029	88.114	nq	99
81) Benzo(a)anthracene	14.06	228	1068581	75.547	nq	100
82) 3,3'-Dichlorobenzidine	14.01	252	184586	35.604	nq	98
83) Chrysene	14.10	228	1056773	76.714	nq	98
84) Bis(2-ethylhexyl)phthalate	14.04	149	815684	97.057	nq	99
85) Di-n-octyl phthalate	14.68	149	1277222	97.094	nq	98
86) Indeno(1,2,3-cd)pyrene	17.09	276	1251537	95.914	nq	98
88) Benzo(b)fluoranthene	15.14	252	1063734	66.991	nq	100
89) Benzo(k)fluoranthene	15.17	252	1008958	73.870	nq	100
90) Benzo(a)pyrene	15.52	252	1025469	74.664	nq	99
91) Dibenzo(a,h)anthracene	17.11	278	1026096	79.587	nq	100
92) Benzo(g,h,i)perylene	17.55	276	1061772	81.606	nq	99

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

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