

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022519\
 Data File : BF112686.D
 Acq On : 25 Feb 2019 12:26
 Operator : JU/SJ
 Sample : K1556-01MS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMPMS

Manual Integrations
 APPROVED

Sohil
 2/26/2019 2:38:59 PM

Quant Time: Feb 25 14:18:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 25 14:05:25 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	75782	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	278771	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	117336	20.00	ng	0.00
64) Phenanthrene-d10	11.34	188	185677	20.00	ng	0.00
76) Chrysene-d12	13.99	240	181790	20.00	ng	0.00
87) Perylene-d12	15.46	264	200534	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.47	112	597319	167.42	ng	0.00
7) Phenol-d6	6.49	99	727099	146.66	ng	0.00
23) Nitrobenzene-d5	7.39	82	467475	96.62	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	166498	121.91	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	850812	102.55	ng	0.00
79) Terphenyl-d14	12.92	244	728390	75.47	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.57	88	59225	41.671	ng	94
3) Pyridine	3.35	79	145168	40.154	ng	97
4) n-Nitrosodimethylamine	3.29	42	86667	57.058	ng	83
6) Aniline	6.49	93	163382	27.704	ng	87
8) 2-Chlorophenol	6.62	128	205329	42.781	ng	98
9) Benzaldehyde	6.37	77	83413	32.190	ng	98
10) Phenol	6.50	94	288090	52.967	ng	83
11) bis(2-Chloroethyl)ether	6.55	93	183389	42.827	ng	97
12) 1,3-Dichlorobenzene	6.76	146	224151	40.101	ng	99
13) 1,4-Dichlorobenzene	6.83	146	224494	39.025	ng	98
14) 1,2-Dichlorobenzene	6.99	146	213369	39.615	ng	96
15) Benzyl Alcohol	6.97	79	175583	42.014	ng	92
16) 2,2'-oxybis(1-Chloropropan	7.09	45	210557	38.297	ng	98
17) 2-Methylphenol	7.10	107	154736	40.669	ng	92
18) Hexachloroethane	7.32	117	83614	41.392	ng	93
19) n-Nitroso-di-n-propylamine	7.23	70	137127	40.113	ng	99
20) 3+4-Methylphenols	7.26	107	203931	41.915	ng	# 60
22) Acetophenone	7.23	105	281882	41.526	ng	93
24) Nitrobenzene	7.40	77	203780	42.174	ng	100
25) Isophorone	7.64	82	347712	45.812	ng	98
26) 2-Nitrophenol	7.72	139	106300	41.166	ng	# 90
27) 2,4-Dimethylphenol	7.77	122	167908	47.196	ng	93
28) bis(2-Chloroethoxy)methane	7.85	93	217751	42.134	ng	99
29) 2,4-Dichlorophenol	7.99	162	158992	39.935	ng	97
30) 1,2,4-Trichlorobenzene	8.04	180	179512	39.226	ng	98
31) Naphthalene	8.12	128	584448	44.831	ng	98
32) Benzoic acid	7.92	122	33344	16.431	ng	98
33) 4-Chloroaniline	8.18	127	71012	12.510	ng	# 94
34) Hexachlorobutadiene	8.23	225	106448	39.638	ng	99
35) Caprolactam	8.56	113	39423	28.069	ng	95
36) 4-Chloro-3-methylphenol	8.69	107	158097	37.511	ng	96
37) 2-Methylnaphthalene	8.81	142	384321	43.063	ng	96
38) 1-Methylnaphthalene	8.91	142	360893	42.190	ng	96
40) 1,2,4,5-Tetrachlorobenzene	8.98	216	167453	43.465	ng	97

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41) Hexachlorocyclopentadiene	8.96	237	49847	39.588	nq	99
43) 2,4,6-Trichlorophenol	9.10	196	97358	40.997	nq	99
44) 2,4,5-Trichlorophenol	9.16	196	92026	37.078	nq	# 92
46) 1,1'-Biphenyl	9.27	154	446909	43.567	nq	98
47) 2-Chloronaphthalene	9.30	162	332346	41.780	nq	96
48) 2-Nitroaniline	9.41	65	91707	42.298	nq	88
49) Acenaphthylene	9.72	152	531643	45.598	nq	98
50) Dimethylphthalate	9.57	163	413875	45.399	nq	99
51) 2,6-Dinitrotoluene	9.65	165	80023	39.194	nq	# 77
52) Acenaphthene	9.89	154	305504	43.863	nq	99
53) 3-Nitroaniline	9.83	138	57603	26.042	nq	97
54) 2,4-Dinitrophenol	9.96	184	33411	50.990	nq	# 82
55) Dibenzofuran	10.06	168	444579	41.770	nq	# 91
56) 4-Nitrophenol	10.05	139	40564m	34.687	nq	
57) 2,4-Dinitrotoluene	10.06	165	103172	39.693	nq	# 77
58) Fluorene	10.40	166	351253	44.101	nq	98
59) 2,3,4,6-Tetrachlorophenol	10.19	232	69324	36.961	nq	# 96
60) Diethylphthalate	10.26	149	361380	39.945	nq	97
61) 4-Chlorophenyl-phenylether	10.39	204	163792	37.804	nq	92
62) 4-Nitroaniline	10.45	138	65128	30.018	nq	90
63) Azobenzene	10.55	77	313126	39.354	nq	97
65) 4,6-Dinitro-2-methylphenol	10.48	198	31348	30.113	nq	84
66) n-Nitrosodiphenylamine	10.51	169	283749	45.343	nq	99
67) 4-Bromophenyl-phenylether	10.87	248	89980	41.203	nq	94
68) Hexachlorobenzene	10.96	284	94392	40.287	nq	94
69) Atrazine	11.03	200	78000	38.629	nq	92
70) Pentachlorophenol	11.17	266	54364	59.631	nq	99
71) Phenanthrene	11.36	178	584626	60.564	nq	97
72) Anthracene	11.42	178	499904	51.988	nq	98
73) Carbazole	11.58	167	386011	40.697	nq	99
74) Di-n-butylphthalate	11.89	149	488653	41.505	nq	98
75) Fluoranthene	12.56	202	915460	88.420	nq	99
77) Benzdine	12.68	184	78450	15.679	nq	99
78) Pyrene	12.79	202	890410	70.745	nq	99
80) Butylbenzylphthalate	13.38	149	224203	36.819	nq	95
81) Benzo(a)anthracene	13.97	228	753618	70.520	nq	98
82) 3,3'-Dichlorobenzidine	13.93	252	87478	21.873	nq	# 99
83) Chrysene	14.02	228	665133	65.204	nq	100
84) Bis(2-ethylhexyl)phthalate	13.93	149	310388	35.909	nq	97
85) Di-n-octyl phthalate	14.54	149	578882	43.529	nq	97
86) Indeno(1,2,3-cd)pyrene	16.96	276	674798	83.484	nq	94
88) Benzo(b)fluoranthene	15.03	252	884590	73.760	nq	99
89) Benzo(k)fluoranthene	15.06	252	635657	55.335	nq	100
90) Benzo(a)pyrene	15.40	252	789979	73.207	nq	99
91) Dibenzo(a,h)anthracene	16.95	278	442204	50.845	nq	99
92) Benzo(g,h,i)perylene	17.40	276	540949	65.927	nq	96

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

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