

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021919\
 Data File : BF112620.D
 Acq On : 19 Feb 2019 14:50
 Operator : JU/SJ
 Sample : K1515-01MSD
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-1MSD

Manual Integrations
 APPROVED

Sohil
 2/20/2019 11:12:45 AM

Quant Time: Feb 19 16:40:15 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	106953	20.00	ng	-0.02
21) Naphthalene-d8	8.10	136	415788	20.00	ng	-0.02
39) Acenaphthene-d10	9.86	164	205996	20.00	ng	-0.02
64) Phenanthrene-d10	11.35	188	384314	20.00	ng	-0.02
76) Chrysene-d12	14.00	240	271299	20.00	ng	-0.01
87) Perylene-d12	15.47	264	262275	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	819858	162.82	ng	-0.02
7) Phenol-d6	6.49	99	1010775	144.46	ng	0.00
23) Nitrobenzene-d5	7.39	82	668130	92.59	ng	-0.02
42) 2,4,6-Tribromophenol	10.66	330	322835	134.64	ng	-0.02
45) 2-Fluorobiphenyl	9.18	172	1302311	89.41	ng	-0.02
79) Terphenyl-d14	12.93	244	1240107	86.09	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.59	88	96524	48.121	ng	95
3) Pyridine	3.36	79	262208	51.389	ng	97
4) n-Nitrosodimethylamine	3.32	42	138990	64.837	ng	84
6) Aniline	6.49	93	386533	46.440	ng	# 54
8) 2-Chlorophenol	6.62	128	336004	49.604	ng	99
9) Benzaldehyde	6.37	77	139403	38.119	ng	99
10) Phenol	6.50	94	405579	52.835	ng	99
11) bis(2-Chloroethyl)ether	6.56	93	293086	48.497	ng	96
12) 1,3-Dichlorobenzene	6.76	146	367362	46.568	ng	97
13) 1,4-Dichlorobenzene	6.84	146	380582	46.876	ng	96
14) 1,2-Dichlorobenzene	6.99	146	352597	46.386	ng	94
15) Benzyl Alcohol	6.97	79	281617	47.747	ng	93
16) 2,2'-oxybis(1-Chloropropan	7.10	45	361588	46.599	ng	# 24
17) 2-Methylphenol	7.09	107	259830	48.388	ng	# 86
18) Hexachloroethane	7.33	117	139831	49.047	ng	98
19) n-Nitroso-di-n-propylamine	7.24	70	234366	48.577	ng	98
20) 3+4-Methylphenols	7.25	107	344533	50.175	ng	# 65
22) Acetophenone	7.24	105	470329	46.455	ng	# 94
24) Nitrobenzene	7.41	77	344317	47.777	ng	99
25) Isophorone	7.65	82	623706	55.095	ng	99
26) 2-Nitrophenol	7.73	139	188073	48.833	ng	# 90
27) 2,4-Dimethylphenol	7.77	122	288366	54.344	ng	92
28) bis(2-Chloroethoxy)methane	7.85	93	387398	50.258	ng	99
29) 2,4-Dichlorophenol	7.98	162	288133	48.522	ng	99
30) 1,2,4-Trichlorobenzene	8.05	180	323089	47.334	ng	99
31) Naphthalene	8.13	128	985261	50.672	ng	99
32) Benzoic acid	7.93	122	118538	39.163	ng	98
33) 4-Chloroaniline	8.18	127	253378	29.927	ng	97
34) Hexachlorobutadiene	8.24	225	188085	46.957	ng	98
35) Caprolactam	8.57	113	87608m	41.821	ng	
36) 4-Chloro-3-methylphenol	8.69	107	297860	47.383	ng	95
37) 2-Methylnaphthalene	8.82	142	703186	52.827	ng	97
38) 1-Methylnaphthalene	8.92	142	653876	51.251	ng	98
40) 1,2,4,5-Tetrachlorobenzene	8.99	216	318619	47.108	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.97	237	178136	70.742	nq	99
43) 2,4,6-Trichlorophenol	9.11	196	194096	46.555	nq	100
44) 2,4,5-Trichlorophenol	9.16	196	195391	44.842	nq	95
46) 1,1'-Biphenyl	9.29	154	849505	47.171	nq	97
47) 2-Chloronaphthalene	9.31	162	650273	46.563	nq	94
48) 2-Nitroaniline	9.42	65	175946	46.224	nq	90
49) Acenaphthylene	9.73	152	1013276	49.503	nq	99
50) Dimethylphthalate	9.58	163	839669	52.463	nq	100
51) 2,6-Dinitrotoluene	9.66	165	169412	47.263	nq #	79
52) Acenaphthene	9.90	154	594862	48.649	nq	98
53) 3-Nitroaniline	9.83	138	115155	29.654	nq	85
54) 2,4-Dinitrophenol	9.96	184	117884	102.476	nq	87
55) Dibenzofuran	10.07	168	878567	47.018	nq #	91
56) 4-Nitrophenol	10.02	139	118333	57.637	nq	86
57) 2,4-Dinitrotoluene	10.07	165	223058	48.881	nq #	80
58) Fluorene	10.42	166	723560	51.745	nq	97
59) 2,3,4,6-Tetrachlorophenol	10.20	232	176050	53.465	nq	97
60) Diethylphthalate	10.28	149	818076	51.506	nq	97
61) 4-Chlorophenyl-phenylether	10.40	204	358111	47.080	nq #	89
62) 4-Nitroaniline	10.45	138	150782	39.586	nq	92
63) Azobenzene	10.56	77	696050	49.829	nq	98
65) 4,6-Dinitro-2-methylphenol	10.49	198	93149	43.230	nq	95
66) n-Nitrosodiphenylamine	10.52	169	620264	47.888	nq	99
67) 4-Bromophenyl-phenylether	10.89	248	217435	48.104	nq	91
68) Hexachlorobenzene	10.97	284	238331	49.145	nq	96
69) Atrazine	11.05	200	174427	41.735	nq	97
70) Pentachlorophenol	11.17	266	140150	74.272	nq	99
71) Phenanthrene	11.38	178	1000196	50.060	nq	99
72) Anthracene	11.43	178	1042082	52.359	nq	99
73) Carbazole	11.59	167	831438	42.351	nq	99
74) Di-n-butylphthalate	11.90	149	1253976	51.459	nq	98
75) Fluoranthene	12.57	202	970574	45.291	nq	99
77) Benzidine	12.69	184	517703	69.331	nq	96
78) Pyrene	12.80	202	956287	50.912	nq	99
80) Butylbenzylphthalate	13.40	149	472074	51.947	nq	95
81) Benzo(a)anthracene	13.99	228	813183	50.988	nq	97
82) 3,3'-Dichlorobenzidine	13.94	252	228403	38.267	nq	99
83) Chrysene	14.03	228	755242	49.610	nq	98
84) Bis(2-ethylhexyl)phthalate	13.95	149	721987	55.969	nq	97
85) Di-n-octyl phthalate	14.56	149	1129585	56.915	nq	98
86) Indeno(1,2,3-cd)pyrene	16.99	276	880201	72.968	nq	96
88) Benzo(b)fluoranthene	15.04	252	750586	47.853	nq	100
89) Benzo(k)fluoranthene	15.07	252	750807	49.973	nq	99
90) Benzo(a)pyrene	15.42	252	737702	52.269	nq	100
91) Dibenzo(a,h)anthracene	16.98	278	735460	64.658	nq	98
92) Benzo(g,h,i)perylene	17.43	276	713753	66.510	nq	97

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

