

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011819\
 Data File : BF112134.D
 Acq On : 18 Jan 2019 19:41
 Operator : JU/SJ
 Sample : K1011-06MSD
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-03-011719-AMSD

Manual Integrations
 APPROVED

Sohil
 1/21/2019 12:00:57 PM

Quant Time: Jan 19 00:28:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 18 23:33:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	273806	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1049829	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	527132	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	944295	20.00	ng	0.00
76) Chrysene-d12	14.03	240	595242	20.00	ng	0.00
87) Perylene-d12	15.50	264	593879	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.51	112	1684922	111.69	ng	0.02
7) Phenol-d6	6.51	99	1990539	105.46	ng	0.00
23) Nitrobenzene-d5	7.42	82	1489927	98.05	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	740256	130.18	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	2935147	81.64	ng	0.00
79) Terphenyl-d14	12.97	244	2615121	93.46	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.79	88	223250	31.554	ng	# 84
3) Pyridine	3.52	79	528076	29.801	ng	99
4) n-Nitrosodimethylamine	3.44	42	283136	39.353	ng	98
6) Aniline	6.52	93	260375	11.342	ng	# 1
8) 2-Chlorophenol	6.65	128	706213	40.308	ng	95
9) Benzaldehyde	6.41	77	168550	12.720	ng	96
10) Phenol	6.52	94	677736	32.856	ng	81
11) bis(2-Chloroethyl)ether	6.59	93	647623	39.486	ng	98
12) 1,3-Dichlorobenzene	6.80	146	773684	38.695	ng	98
13) 1,4-Dichlorobenzene	6.87	146	784068	38.337	ng	99
14) 1,2-Dichlorobenzene	7.03	146	731126	38.809	ng	99
15) Benzyl Alcohol	7.00	79	586833	41.732	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.13	45	853738	36.096	ng	83
17) 2-Methylphenol	7.12	107	544061	41.350	ng	95
18) Hexachloroethane	7.37	117	271531	39.778	ng	97
19) n-Nitroso-di-n-propylamine	7.27	70	476863	41.473	ng	96
20) 3+4-Methylphenols	7.27	107	679866	42.616	ng	# 75
22) Acetophenone	7.26	105	944272	38.824	ng	96
24) Nitrobenzene	7.44	77	724572	44.993	ng	98
25) Isophorone	7.68	82	1308646	44.587	ng	98
26) 2-Nitrophenol	7.76	139	385974	46.821	ng	98
27) 2,4-Dimethylphenol	7.80	122	621253	45.600	ng	96
28) bis(2-Chloroethoxy)methane	7.89	93	842635	41.854	ng	100
29) 2,4-Dichlorophenol	8.00	162	627105	42.136	ng	99
30) 1,2,4-Trichlorobenzene	8.08	180	685603	39.695	ng	100
31) Naphthalene	8.16	128	2048103	42.893	ng	99
32) Benzoic acid	7.93	122	304088	48.274	ng	99
33) 4-Chloroaniline	8.21	127	156648	7.300	ng	98
34) Hexachlorobutadiene	8.28	225	371402	38.641	ng	100
35) Caprolactam	8.58	113	162799m	31.254	ng	
36) 4-Chloro-3-methylphenol	8.70	107	623511	42.894	ng	100
37) 2-Methylnaphthalene	8.86	142	1468126	45.279	ng	98
38) 1-Methylnaphthalene	8.95	142	1391559	44.656	ng	97
40) 1,2,4,5-Tetrachlorobenzene	9.02	216	684920	40.463	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.01	237	267769	43.562	ng	97
43) 2,4,6-Trichlorophenol	9.14	196	456382	45.070	ng	100
44) 2,4,5-Trichlorophenol	9.19	196	481938	44.687	ng	99
46) 1,1'-Biphenyl	9.32	154	1788508	40.577	ng	100
47) 2-Chloronaphthalene	9.35	162	1371064	39.738	ng	98
48) 2-Nitroaniline	9.44	65	382165	49.176	ng	95
49) Acenaphthylene	9.76	152	2160484	42.911	ng	99
50) Dimethylphthalate	9.62	163	1693691	44.058	ng	100
51) 2,6-Dinitrotoluene	9.68	165	378731	49.332	ng	92
52) Acenaphthene	9.93	154	1266697	41.119	ng	100
53) 3-Nitroaniline	9.85	138	124897	14.548	ng	89
54) 2,4-Dinitrophenol	9.96	184	175126	73.530	ng	94
55) Dibenzofuran	10.10	168	1908427	40.966	ng	97
56) 4-Nitrophenol	10.03	139	458054	75.029	ng	97
57) 2,4-Dinitrotoluene	10.09	165	491029	49.651	ng	91
58) Fluorene	10.45	166	1500744	43.116	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.23	232	395677	48.619	ng	99
60) Diethylphthalate	10.32	149	1642206	43.763	ng	99
61) 4-Chlorophenyl-phenylether	10.43	204	751160	39.993	ng	95
62) 4-Nitroaniline	10.46	138	322390	37.818	ng	97
63) Azobenzene	10.60	77	1422806	39.593	ng	93
65) 4,6-Dinitro-2-methylphenol	10.50	198	136355	37.428	ng	91
66) n-Nitrosodiphenylamine	10.56	169	1368463	43.836	ng	99
67) 4-Bromophenyl-phenylether	10.92	248	475977	43.395	ng	97
68) Hexachlorobenzene	11.00	284	484122	40.974	ng	96
69) Atrazine	11.08	200	468662	46.524	ng	98
70) Pentachlorophenol	11.20	266	441112	77.828	ng	100
71) Phenanthrene	11.41	178	2077139	43.709	ng	100
72) Anthracene	11.46	178	2088330	43.380	ng	99
73) Carbazole	11.62	167	1865710	38.688	ng	99
74) Di-n-butylphthalate	11.94	149	2457537	45.304	ng	100
75) Fluoranthene	12.60	202	1912486	38.116	ng	99
77) Benzidine	12.71	184	447129	22.514	ng	95
78) Pyrene	12.83	202	1866475	47.527	ng	100
80) Butylbenzylphthalate	13.44	149	895688	51.989	ng	99
81) Benzo(a)anthracene	14.02	228	1524091	44.656	ng	99
82) 3,3'-Dichlorobenzidine	13.97	252	380568	29.775	ng	98
83) Chrysene	14.05	228	1435435	44.449	ng	100
84) Bis(2-ethylhexyl)phthalate	14.00	149	1273291	51.568	ng	99
85) Di-n-octyl phthalate	14.60	149	1942211	51.052	ng	99
86) Indeno(1,2,3-cd)pyrene	16.99	276	1326156	44.511	ng	99
88) Benzo(b)fluoranthene	15.07	252	1489435	44.183	ng	99
89) Benzo(k)fluoranthene	15.10	252	1439556	44.120	ng	99
90) Benzo(a)pyrene	15.44	252	1410185	45.614	ng	99
91) Dibenzo(a,h)anthracene	16.99	278	1115971	40.642	ng	100
92) Benzo(g,h,i)perylene	17.43	276	1033923	38.221	ng	99

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

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