

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052919\
 Data File : BF114645.D
 Acq On : 29 May 2019 20:56
 Operator : HP/JU
 Sample : K3068-08MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-4W(3-4)MS

Manual Integrations
 APPROVED

mohammad
 5/30/2019 12:47:55 PM

Quant Time: May 30 04:41:15 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 19:08:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	303991	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	1177785	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	595137	20.00	ng	0.00
64) Phenanthrene-d10	11.19	188	1202050	20.00	ng	0.00
76) Chrysene-d12	13.82	240	759619	20.00	ng	0.00
87) Perylene-d12	15.20	264	825002	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	2154245	124.42	ng	0.01
7) Phenol-d6	6.33	99	2763053	132.43	ng	0.00
23) Nitrobenzene-d5	7.26	82	1702433	90.18	ng	0.00
42) 2,4,6-Tribromophenol	10.51	330	799136	140.92	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	2949479	94.97	ng	0.00
79) Terphenyl-d14	12.78	244	3370511	83.42	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	2.36	88	298115	35.879 ng	98
3) Pyridine	3.06	79	827014	34.200 ng	99
4) n-Nitrosodimethylamine	2.99	42	404448	41.331 ng	99
6) Aniline	6.36	93	698284	22.157 ng	100
8) 2-Chlorophenol	6.48	128	893268	44.673 ng	99
9) Benzaldehyde	6.23	77	387507	30.744 ng	100
10) Phenol	6.35	94	1055264	41.868 ng	99
11) bis(2-Chloroethyl)ether	6.43	93	838800	43.508 ng	100
12) 1,3-Dichlorobenzene	6.63	146	949765	41.691 ng	96
13) 1,4-Dichlorobenzene	6.71	146	946550	41.417 ng	98
14) 1,2-Dichlorobenzene	6.87	146	890591	42.803 ng	99
15) Benzyl Alcohol	6.84	79	754348	45.963 ng	100
16) 2,2'-oxybis(1-Chloropropan	6.98	45	1254864	42.437 ng	100
17) 2-Methylphenol	6.95	107	749435	44.313 ng	99
18) Hexachloroethane	7.21	117	352677	40.758 ng	97
19) n-Nitroso-di-n-propylamine	7.12	70	614040	42.636 ng	98
20) 3+4-Methylphenols	7.10	107	854401	42.954 ng	96
22) Acetophenone	7.10	105	1162799	46.336 ng	97
24) Nitrobenzene	7.27	77	913040	45.530 ng	97
25) Isophorone	7.52	82	1735446	45.807 ng	99
26) 2-Nitrophenol	7.59	139	468449	47.990 ng	97
27) 2,4-Dimethylphenol	7.63	122	828475	51.032 ng	100
28) bis(2-Chloroethoxy)methane	7.73	93	1096553	45.468 ng	100
29) 2,4-Dichlorophenol	7.83	162	777433	47.048 ng	100
30) 1,2,4-Trichlorobenzene	7.92	180	837702	44.497 ng	99
31) Naphthalene	8.00	128	2451710	45.192 ng	99
32) Benzoic acid	7.72	122	168535	15.983 ng	98
33) 4-Chloroaniline	8.04	127	310335	12.007 ng	98
34) Hexachlorobutadiene	8.12	225	449627	42.993 ng	99
35) Caprolactam	8.41	113	267765m	48.017 ng	
36) 4-Chloro-3-methylphenol	8.53	107	817330	49.322 ng	97
37) 2-Methylnaphthalene	8.69	142	1750883	48.574 ng	99
38) 1-Methylnaphthalene	8.79	142	1634938	47.873 ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.86	216	717176	45.251 ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.85	237	774885	73.987	nq	98
43) 2,4,6-Trichlorophenol	8.96	196	567527	45.509	nq	99
44) 2,4,5-Trichlorophenol	9.00	196	611625	49.149	nq	98
46) 1,1'-Biphenyl	9.15	154	2012866	45.563	nq	99
47) 2-Chloronaphthalene	9.17	162	1649830	45.527	nq	98
48) 2-Nitroaniline	9.26	65	500974	44.798	nq	98
49) Acenaphthylene	9.59	152	2534785	45.277	nq	100
50) Dimethylphthalate	9.46	163	2439747	58.832	nq	99
51) 2,6-Dinitrotoluene	9.51	165	466943	48.458	nq	95
52) Acenaphthene	9.76	154	1653935	48.172	nq	100
53) 3-Nitroaniline	9.67	138	261368	22.162	nq	99
54) 2,4-Dinitrophenol	9.77	184	279863	73.631	nq	94
55) Dibenzofuran	9.93	168	2263795	46.320	nq	99
56) 4-Nitrophenol	9.83	139	850770	97.345	nq	90
57) 2,4-Dinitrotoluene	9.91	165	620516	50.175	nq	87
58) Fluorene	10.27	166	1559622	49.546	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.05	232	522421	51.288	nq	100
60) Diethylphthalate	10.15	149	2013222	47.532	nq	100
61) 4-Chlorophenyl-phenylether	10.26	204	778950	48.887	nq	100
62) 4-Nitroaniline	10.28	138	490805	42.708	nq	99
63) Azobenzene	10.42	77	1842115	47.371	nq	99
65) 4,6-Dinitro-2-methylphenol	10.31	198	242843	45.561	nq	96
66) n-Nitrosodiphenylamine	10.38	169	1636312	46.074	nq	99
67) 4-Bromophenyl-phenylether	10.75	248	565572	43.875	nq	98
68) Hexachlorobenzene	10.82	284	598364	42.857	nq	97
69) Atrazine	10.91	200	517788	47.633	nq	99
70) Pentachlorophenol	11.00	266	698416	86.778	nq	99
71) Phenanthrene	11.22	178	2575930	42.009	nq	100
72) Anthracene	11.27	178	2671524	43.047	nq	99
73) Carbazole	11.42	167	2483713	45.537	nq	100
74) Di-n-butylphthalate	11.77	149	3055932	46.880	nq	100
75) Fluoranthene	12.40	202	2729816	42.993	nq	99
77) Benzidine	12.52	184	1083490	34.639	nq	98
78) Pyrene	12.63	202	2761298	42.972	nq	100
80) Butylbenzylphthalate	13.26	149	1467692	45.448	nq	99
81) Benzo(a)anthracene	13.81	228	2461056	42.160	nq	100
82) 3,3'-Dichlorobenzidine	13.77	252	441082	19.566	nq	98
83) Chrysene	13.84	228	2387815	42.814	nq	100
84) Bis(2-ethylhexyl)phthalate	13.82	149	1884408	47.200	nq	100
85) Di-n-octyl phthalate	14.43	149	3205404	46.209	nq	99
86) Indeno(1,2,3-cd)pyrene	16.53	276	2386173	42.718	nq	100
88) Benzo(b)fluoranthene	14.82	252	2304929	44.358	nq	100
89) Benzo(k)fluoranthene	14.84	252	2157600	47.773	nq	100
90) Benzo(a)pyrene	15.14	252	2129946	46.660	nq	97
91) Dibenzo(a,h)anthracene	16.54	278	1967126	48.767	nq	99
92) Benzo(g,h,i)perylene	16.92	276	2011042	51.368	nq	99

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

