

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062219\
 Data File : BF115126.D
 Acq On : 22 Jun 2019 17:43
 Operator : HP/JU
 Sample : K3441-04MSD
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAMPLE-3MSD

Quant Time: Jun 24 05:31:55 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 24 04:21:29 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.61	152	311559	20.00	ng	0.00
21) Naphthalene-d8	7.90	136	1215104	20.00	ng	0.00
39) Acenaphthene-d10	9.64	164	598724	20.00	ng	0.00
64) Phenanthrene-d10	11.11	188	1127930	20.00	ng	0.00
76) Chrysene-d12	13.73	240	613735	20.00	ng	-0.02
87) Perylene-d12	15.09	264	669903	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.24	112	2006357	99.38	ng	0.03
7) Phenol-d6	6.26	99	2388002	97.45	ng	0.00
23) Nitrobenzene-d5	7.18	82	1677334	84.92	ng	0.00
42) 2,4,6-Tribromophenol	10.42	330	734630	116.35	ng	0.00
45) 2-Fluorobiphenyl	8.97	172	3026602	85.87	ng	-0.01
79) Terphenyl-d14	12.70	244	2901495	100.52	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.43	88	317769	33.350	ng	# 86
3) Pyridine	3.04	79	683287	25.443	ng	98
4) n-Nitrosodimethylamine	2.97	42	370547	35.196	ng	99
6) Aniline	6.28	93	251635	7.472	ng	# 1
8) 2-Chlorophenol	6.40	128	844220	37.187	ng	99
9) Benzaldehyde	6.16	77	453956	28.395	ng	97
10) Phenol	6.27	94	866444	30.185	ng	90
11) bis(2-Chloroethyl)ether	6.36	93	796407	37.639	ng	99
12) 1,3-Dichlorobenzene	6.55	146	921101	36.559	ng	98
13) 1,4-Dichlorobenzene	6.63	146	931722	36.878	ng	99
14) 1,2-Dichlorobenzene	6.78	146	859335	36.728	ng	98
15) Benzyl Alcohol	6.76	79	676979	37.939	ng	97
16) 2,2'-oxybis(1-Chloropropan	6.90	45	1130523	36.838	ng	99
17) 2-Methylphenol	6.88	107	698612	37.282	ng	99
18) Hexachloroethane	7.13	117	335177	36.798	ng	96
19) n-Nitroso-di-n-propylamine	7.04	70	568798	36.256	ng	98
20) 3+4-Methylphenols	7.03	107	787836	36.411	ng	97
22) Acetophenone	7.02	105	1142296	41.064	ng	# 98
24) Nitrobenzene	7.20	77	873374	40.134	ng	99
25) Isophorone	7.44	82	1600925	39.564	ng	99
26) 2-Nitrophenol	7.51	139	475796	44.274	ng	99
27) 2,4-Dimethylphenol	7.56	122	770536	43.792	ng	99
28) bis(2-Chloroethoxy)methane	7.65	93	1013915	39.644	ng	100
29) 2,4-Dichlorophenol	7.75	162	746066	41.087	ng	99
30) 1,2,4-Trichlorobenzene	7.84	180	809939	40.381	ng	100
31) Naphthalene	7.91	128	2420622	40.583	ng	100
32) Benzoic acid	7.67	122	325577	25.922	ng	99
33) 4-Chloroaniline	7.96	127	136061	4.991	ng	98
34) Hexachlorobutadiene	8.04	225	435543	39.625	ng	98
35) Caprolactam	8.34	113	182134m	29.834	ng	
36) 4-Chloro-3-methylphenol	8.45	107	746726	40.606	ng	97
37) 2-Methylnaphthalene	8.61	142	1662704	41.344	ng	99
38) 1-Methylnaphthalene	8.70	142	1568105	40.826	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.77	216	685607	40.523	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.77	237	698597	69.635	ng	100
43) 2,4,6-Trichlorophenol	8.88	196	544285	41.669	ng	99
44) 2,4,5-Trichlorophenol	8.92	196	563023	41.856	ng	99
46) 1,1'-Biphenyl	9.07	154	1912043	39.912	ng	98
47) 2-Chloronaphthalene	9.09	162	1556356	40.051	ng	98
48) 2-Nitroaniline	9.18	65	458789	40.496	ng	94
49) Acenaphthylene	9.50	152	2439665	40.295	ng	99
50) Dimethylphthalate	9.38	163	1769200	40.164	ng	100
51) 2,6-Dinitrotoluene	9.43	165	425545	41.845	ng	95
52) Acenaphthene	9.67	154	1558905	41.148	ng	99
53) 3-Nitroaniline	9.59	138	184296	14.850	ng	97
54) 2,4-Dinitrophenol	9.70	184	267073	53.695	ng	97
55) Dibenzofuran	9.85	168	2112340	38.847	ng	100
56) 4-Nitrophenol	9.76	139	681951	68.542	ng	97
57) 2,4-Dinitrotoluene	9.83	165	559162	42.583	ng	95
58) Fluorene	10.19	166	1457687	39.452	ng	98
59) 2,3,4,6-Tetrachlorophenol	9.97	232	455102	40.349	ng	99
60) Diethylphthalate	10.08	149	1782969	40.267	ng	99
61) 4-Chlorophenyl-phenylether	10.18	204	705926	40.325	ng	98
62) 4-Nitroaniline	10.21	138	444085	35.670	ng	99
63) Azobenzene	10.34	77	1640038	39.655	ng	98
65) 4,6-Dinitro-2-methylphenol	10.23	198	212624	36.727	ng	79
66) n-Nitrosodiphenylamine	10.30	169	1492891	42.484	ng	100
67) 4-Bromophenyl-phenylether	10.67	248	506969	41.456	ng	92
68) Hexachlorobenzene	10.73	284	549179	41.373	ng	95
69) Atrazine	10.83	200	458798	40.587	ng	98
70) Pentachlorophenol	10.92	266	546067	64.574	ng	98
71) Phenanthrene	11.14	178	2350423	38.703	ng	99
72) Anthracene	11.19	178	2369252	38.649	ng	98
73) Carbazole	11.34	167	2101776	38.395	ng	98
74) Di-n-butylphthalate	11.70	149	2574417	40.699	ng	99
75) Fluoranthene	12.31	202	2248793	37.113	ng	100
77) Benzidine	12.44	184	53470	1.962	ng	# 77
78) Pyrene	12.54	202	2209381	46.843	ng	99
80) Butylbenzylphthalate	13.18	149	1008227	48.438	ng	93
81) Benzo(a)anthracene	13.72	228	1772096	40.241	ng	99
82) 3,3'-Dichlorobenzidine	13.68	252	238736	15.012	ng	99
83) Chrysene	13.75	228	1633108	40.484	ng	100
84) Bis(2-ethylhexyl)phthalate	13.74	149	1283005	47.041	ng	100
85) Di-n-octyl phthalate	14.35	149	2016961	44.014	ng	100
86) Indeno(1,2,3-cd)pyrene	16.36	276	1871533	39.208	ng	100
88) Benzo(b)fluoranthene	14.72	252	1716158	41.056	ng	99
89) Benzo(k)fluoranthene	14.75	252	1581350	42.186	ng	98
90) Benzo(a)pyrene	15.04	252	1629740	43.258	ng	100
91) Dibenzo(a,h)anthracene	16.38	278	1563060	44.440	ng	99
92) Benzo(g,h,i)perylene	16.74	276	1514437	42.543	ng	99

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

