

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082319\
 Data File : BF116477.D
 Acq On : 23 Aug 2019 14:45
 Operator : HP/JU
 Sample : K4086-01RE
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NB-08-082019RE

Quant Time: Aug 23 16:40:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 23 13:38:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	6392	20.00	ng	0.00
21) Naphthalene-d8	8.43	136	143	20.00	ng	0.27
39) Acenaphthene-d10	10.23	164	670	20.00	ng	0.32
64) Phenanthrene-d10	11.76	188	370	20.00	ng	0.36
76) Chrysene-d12	14.45	240	2465	20.00	ng	0.42
87) Perylene-d12	15.99	264	908	20.00	ng	0.48

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.50	112	701024	1602.47	ng	0.00
7) Phenol-d6	6.50	99	1095942	2043.16	ng	0.00
23) Nitrobenzene-d5	7.44	82	710983	272012.91	ng	0.00
42) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng	
45) 2-Fluorobiphenyl	0.00	172	0	0.00	ng	
79) Terphenyl-d14	13.34	244	762	5.77	ng	0.36

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
9) Benzaldehyde	6.43	77	1202	3.354	ng	# 78
10) Phenol	6.52	94	16075	27.597	ng	# 92
15) Benzyl Alcohol	7.04	79	13732	31.634	ng	# 46
18) Hexachloroethane	7.42	117	436	2.338	ng	# 1
19) n-Nitroso-di-n-propylamine	7.44	70	99249	271.161	ng	# 77
20) 3+4-Methylphenols	7.27	107	1872	4.018	ng	# 72
22) Acetophenone	7.53	105	600	167.795	ng	# 55
24) Nitrobenzene	7.72	77	120	44.332	ng	# 38
25) Isophorone	7.97	82	932	180.940	ng	# 1
27) 2,4-Dimethylphenol	7.94	122	268	129.907	ng	# 64
28) bis(2-Chloroethoxy)methane	8.18	93	829	260.613	ng	# 1
29) 2,4-Dichlorophenol	8.33	162	804	409.798	ng	# 25
31) Naphthalene	8.18	128	10811	1562.801	ng	# 98
32) Benzoic acid	7.94	122	268	203.203	ng	# 46
33) 4-Chloroaniline	8.55	127	281	92.530	ng	# 1
35) Caprolactam	8.87	113	128	187.300	ng	# 1
36) 4-Chloro-3-methylphenol	9.00	107	570	257.489	ng	# 22
37) 2-Methylnaphthalene	9.23	142	675	142.112	ng	# 1
38) 1-Methylnaphthalene	9.23	142	675	152.132	ng	# 66
43) 2,4,6-Trichlorophenol	9.35	196	134	10.622	ng	# 11
44) 2,4,5-Trichlorophenol	9.35	196	134	10.439	ng	# 1
46) 1,1'-Biphenyl	9.63	154	3113	58.114	ng	# 41
47) 2-Chloronaphthalene	9.63	162	146	3.516	ng	# 1
48) 2-Nitroaniline	9.75	65	137	10.774	ng	# 53
49) Acenaphthylene	10.06	152	839	13.809	ng	# 1
51) 2,6-Dinitrotoluene	10.02	165	491	51.316	ng	# 87
52) Acenaphthene	10.26	154	728	18.673	ng	# 1
53) 3-Nitroaniline	10.19	138	749	65.671	ng	# 56
54) 2,4-Dinitrophenol	10.32	184	605	152.908	ng	# 38
55) Dibenzofuran	10.47	168	603	11.095	ng	# 1
56) 4-Nitrophenol	10.32	139	742	83.701	ng	# 1
57) 2,4-Dinitrotoluene	10.41	165	3113	255.684	ng	# 49
58) Fluorene	10.78	166	413	9.781	ng	# 1
59) 2,3,4,6-Tetrachlorophenol	10.58	232	172	16.318	ng	# 1

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
60) Diethylphthalate	10.62	149	490	10.665	nq	# 38
61) 4-Chlorophenyl-phenylether	10.79	204	524	25.438	nq	# 1
62) 4-Nitroaniline	10.66	138	601	53.878	nq	# 91
63) Azobenzene	10.95	77	1542	32.641	nq	# 40
65) 4,6-Dinitro-2-methylphenol	10.82	198	1608	912.192	nq	# 1
66) n-Nitrosodiphenylamine	10.90	169	6305	526.272	nq	# 45
67) 4-Bromophenyl-phenylether	11.31	248	116	29.904	nq	# 1
69) Atrazine	11.38	200	4071	1134.120	nq	# 64
70) Pentachlorophenol	11.76	266	112	44.744	nq	# 1
71) Phenanthrene	11.81	178	1377	68.317	nq	# 1
72) Anthracene	11.81	178	1377	67.986	nq	# 1
73) Carbazole	12.00	167	787	43.288	nq	# 1
74) Di-n-butylphthalate	12.49	149	57873	2568.892	nq	# 1
75) Fluoranthene	12.84	202	412670	19803.120	nq	# 97
77) Benzidine	13.07	184	556	5.606	nq	# 1
78) Pyrene	13.22	202	1273	6.458	nq	# 1
80) Butylbenzylphthalate	13.90	149	10055	118.108	nq	# 83
81) Benzo(a)anthracene	14.45	228	1084	6.790	nq	# 1
82) 3,3'-Dichlorobenzidine	14.39	252	1478	25.978	nq	# 9
83) Chrysene	14.49	228	1327	8.201	nq	# 1
84) Bis(2-ethylhexyl)phthalate	14.43	149	407	3.585	nq	# 59
85) Di-n-octyl phthalate	15.10	149	11118	57.688	nq	# 76
86) Indeno(1,2,3-cd)pyrene	17.43	276	120689	870.689	nq	# 93
88) Benzo(b)fluoranthene	15.44	252	212186	3782.820	nq	# 95
89) Benzo(k)fluoranthene	15.53	252	53857	1059.835	nq	# 68
90) Benzo(a)pyrene	15.94	252	874	17.630	nq	# 1
91) Dibenzo(a,h)anthracene	17.54	278	2424	59.505	nq	# 1
92) Benzo(a,h,i)perylene	17.93	276	634	16.325	nq	# 1

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

