

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062618\
 Data File : BF106846.D
 Acq On : 27 Jun 2018 2:29
 Operator : JU/SJ
 Sample : PB110563BSD
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB110563BSD

Manual Integrations
 APPROVED

Sohil
 6/27/2018 2:14:36 PM

Quant Time: Jun 27 05:30:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 25 18:50:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	42212	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	155505	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	70412	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	132474	20.00	ng	0.00
75) Chrysene-d12	14.15	240	122094	20.00	ng	-0.02
86) Perylene-d12	15.68	264	86107	20.00	ng	-0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	201991	81.03	ng	0.00
7) Phenol-d6	6.61	99	247719	79.57	ng	0.00
23) Nitrobenzene-d5	7.52	82	229022	81.85	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	66097	80.99	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	423041	105.20	ng	0.00
78) Terphenyl-d14	13.08	244	472520	93.75	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.87	88	41294	32.220	ng	99
3) Pyridine	3.64	79	118190	34.004	ng	98
4) n-Nitrosodimethylamine	3.61	42	48524	32.870	ng	85
6) Aniline	6.62	93	45268	10.720	ng	# 30
8) 2-Chlorophenol	6.75	128	108586	39.714	ng	95
9) Benzaldehyde	6.51	77	76175	34.801	ng	98
10) Phenol	6.62	94	130489	36.729	ng	93
11) bis(2-Chloroethyl)ether	6.70	93	109234	37.243	ng	96
12) 1,3-Dichlorobenzene	6.90	146	129249	40.360	ng	98
13) 1,4-Dichlorobenzene	6.98	146	130479	40.301	ng	98
14) 1,2-Dichlorobenzene	7.13	146	120273	40.535	ng	98
15) Benzyl Alcohol	7.10	79	92225	36.894	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.23	45	149600	38.875	ng	73
17) 2-Methylphenol	7.22	107	90682	38.755	ng	# 91
18) Hexachloroethane	7.47	117	35825	31.466	ng	89
19) n-Nitroso-di-n-propylamine	7.37	70	74442	36.277	ng	95
20) 3+4-Methylphenols	7.38	107	114977	45.727	ng	96
22) Acetophenone	7.37	105	149698	43.029	ng	# 99
24) Nitrobenzene	7.54	77	112910	39.523	ng	98
25) Isophorone	7.78	82	203799	41.332	ng	97
26) 2-Nitrophenol	7.86	139	44675	33.126	ng	94
27) 2,4-Dimethylphenol	7.90	122	93494	44.852	ng	99
28) bis(2-Chloroethoxy)methane	7.98	93	132521	40.029	ng	98
29) 2,4-Dichlorophenol	8.11	162	88774	40.678	ng	93
30) 1,2,4-Trichlorobenzene	8.18	180	106912	44.471	ng	95
31) Naphthalene	8.27	128	304182	41.839	ng	99
32) Benzoic acid	7.98	122	20388m	17.572	ng	
33) 4-Chloroaniline	8.31	127	20137	6.601	ng	98
34) Hexachlorobutadiene	8.37	225	71325	45.531	ng	99
35) Caprolactam	8.68	113	24260m	34.103	ng	
36) 4-Chloro-3-methylphenol	8.81	107	87491	38.088	ng	95
37) 2-Methylnaphthalene	8.95	142	215820	44.786	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.12	216	112939	47.628	ng	# 96
40) Hexachlorocyclopentadiene	9.10	237	7660	6.279	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.24	196	58369	37.031	nq	95
43) 2,4,5-Trichlorophenol	9.29	196	59328	36.072	nq	# 87
45) 1,1'-Biphenyl	9.42	154	255853	45.594	nq	98
46) 2-Chloronaphthalene	9.45	162	182190	41.247	nq	96
47) 2-Nitroaniline	9.55	65	54514	36.702	nq	99
48) Acenaphthylene	9.87	152	284455	40.790	nq	100
49) Dimethylphthalate	9.71	163	221309	41.907	nq	99
50) 2,6-Dinitrotoluene	9.78	165	46024	41.157	nq	91
51) Acenaphthene	10.04	154	167848	38.222	nq	97
52) 3-Nitroaniline	9.95	138	30548	23.739	nq	98
53) 2,4-Dinitrophenol	10.07	184	4122	11.968	nq	# 21
54) Dibenzofuran	10.21	168	261018	43.408	nq	96
55) 4-Nitrophenol	10.14	139	43752	48.709	nq	93
56) 2,4-Dinitrotoluene	10.19	165	56500	38.708	nq	88
57) Fluorene	10.55	166	195351	40.940	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.33	232	47393	36.249	nq	# 79
59) Diethylphthalate	10.41	149	204511	40.123	nq	# 93
60) 4-Chlorophenyl-phenylether	10.54	204	105097	42.095	nq	88
61) 4-Nitroaniline	10.57	138	48624	38.074	nq	97
62) Azobenzene	10.70	77	177426	35.349	nq	90
64) 4,6-Dinitro-2-methylphenol	10.60	198	5863	7.160	nq	# 74
65) n-Nitrosodiphenylamine	10.66	169	171035	38.385	nq	99
66) 4-Bromophenyl-phenylether	11.03	248	68311	43.207	nq	# 79
67) Hexachlorobenzene	11.10	284	70254	42.456	nq	# 82
68) Atrazine	11.18	200	58133	41.040	nq	95
69) Pentachlorophenol	11.30	266	29513	35.745	nq	97
70) Phenanthrene	11.52	178	277613	43.448	nq	99
71) Anthracene	11.57	178	288289	44.204	nq	99
72) Carbazole	11.73	167	255003	41.763	nq	98
73) Di-n-butylphthalate	12.04	149	293216	40.762	nq	99
74) Fluoranthene	12.71	202	320370	45.491	nq	94
76) Benzidine	12.83	184	122314	35.176	nq	98
77) Pyrene	12.94	202	323801	40.217	nq	99
79) Butylbenzylphthalate	13.54	149	127486	38.512	nq	# 97
80) Benzo(a)anthracene	14.14	228	305388	41.040	nq	100
81) 3,3'-Dichlorobenzidine	14.09	252	76064	29.084	nq	# 96
82) Chrysene	14.17	228	286714	41.881	nq	99
83) Bis(2-ethylhexyl)phthalate	14.09	149	174439	41.218	nq	# 96
84) Di-n-octyl phthalate	14.72	149	303986	44.066	nq	96
85) Indeno(1,2,3-cd)pyrene	17.25	276	202874	34.920	nq	# 90
87) Benzo(b)fluoranthene	15.22	252	231224m	43.228	nq	
88) Benzo(k)fluoranthene	15.25	252	224159m	44.007	nq	
89) Benzo(a)pyrene	15.61	252	209790	44.119	nq	# 97
90) Dibenzo(a,h)anthracene	17.27	278	169621	42.091	nq	# 93
91) Benzo(g,h,i)perylene	17.73	276	166687	42.318	nq	# 91

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

