

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062718\
 Data File : BF106864.D
 Acq On : 27 Jun 2018 11:53
 Operator : JU/SJ
 Sample : PB110600BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB110600BS

Manual Integrations
 APPROVED

Sohil
 6/28/2018 2:21:27 PM

Quant Time: Jun 27 14:09:35 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	6.95	152	46164	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	182722	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	99211	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	203403	20.00	ng	-0.01
75) Chrysene-d12	14.15	240	148902	20.00	ng	-0.02
86) Perylene-d12	15.69	264	144294	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	276678	101.49	ng	0.00
7) Phenol-d6	6.60	99	344985	101.33	ng	-0.01
23) Nitrobenzene-d5	7.52	82	220378	67.03	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	129550	112.66	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	440268	73.04	ng	-0.01
78) Terphenyl-d14	13.08	244	533388	86.77	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.87	88	37515	26.766	ng	98
3) Pyridine	3.65	79	111050	29.214	ng	98
4) n-Nitrosodimethylamine	3.61	42	43882	27.180	ng	80
6) Aniline	6.62	93	93143	20.169	ng	95
8) 2-Chlorophenol	6.75	128	103738	34.693	ng	97
9) Benzaldehyde	6.51	77	64839	25.312	ng	93
10) Phenol	6.62	94	124786	32.117	ng	78
11) bis(2-Chloroethyl)ether	6.69	93	101776	31.730	ng	98
12) 1,3-Dichlorobenzene	6.90	146	122200	34.893	ng	98
13) 1,4-Dichlorobenzene	6.97	146	122403	34.570	ng	98
14) 1,2-Dichlorobenzene	7.12	146	114294	35.222	ng	98
15) Benzyl Alcohol	7.10	79	89433	32.715	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.22	45	145040	34.464	ng	75
17) 2-Methylphenol	7.21	107	86870	33.948	ng	# 93
18) Hexachloroethane	7.47	117	41713	33.501	ng	# 83
19) n-Nitroso-di-n-propylamine	7.37	70	70504	31.416	ng	93
20) 3+4-Methylphenols	7.37	107	108777	37.616	ng	# 69
22) Acetophenone	7.36	105	140748	34.430	ng	# 97
24) Nitrobenzene	7.54	77	109179	32.524	ng	98
25) Isophorone	7.78	82	201301	34.744	ng	99
26) 2-Nitrophenol	7.85	139	58415	36.863	ng	97
27) 2,4-Dimethylphenol	7.89	122	94440	38.557	ng	99
28) bis(2-Chloroethoxy)methane	7.98	93	128448	33.019	ng	98
29) 2,4-Dichlorophenol	8.10	162	95552	37.263	ng	94
30) 1,2,4-Trichlorobenzene	8.18	180	106571	37.726	ng	98
31) Naphthalene	8.26	128	304536	35.648	ng	99
32) Benzoic acid	8.02	122	55720	33.272	ng	99
33) 4-Chloroaniline	8.31	127	32979	9.200	ng	97
34) Hexachlorobutadiene	8.37	225	69537	37.778	ng	98
35) Caprolactam	8.69	113	29075m	34.784	ng	
36) 4-Chloro-3-methylphenol	8.80	107	97187	36.007	ng	99
37) 2-Methylnaphthalene	8.95	142	226069	39.925	ng	96
39) 1,2,4,5-Tetrachlorobenzene	9.12	216	116944	35.001	ng	# 95
40) Hexachlorocyclopentadiene	9.10	237	103663	60.304	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.24	196	71947	32.395	nq	92
43) 2,4,5-Trichlorophenol	9.28	196	73661m	31.786	nq	
45) 1,1'-Biphenyl	9.42	154	274950	34.775	nq	96
46) 2-Chloronaphthalene	9.44	162	199651	32.079	nq	96
47) 2-Nitroaniline	9.54	65	59953	28.647	nq	94
48) Acenaphthylene	9.87	152	322132	32.784	nq	99
49) Dimethylphthalate	9.71	163	237021	31.854	nq	98
50) 2,6-Dinitrotoluene	9.78	165	55456	35.196	nq	# 85
51) Acenaphthene	10.04	154	193408	31.258	nq	98
52) 3-Nitroaniline	9.95	138	25623	14.132	nq	96
53) 2,4-Dinitrophenol	10.07	184	56875	70.133	nq	# 13
54) Dibenzofuran	10.21	168	299158	35.309	nq	95
55) 4-Nitrophenol	10.14	139	60422	47.742	nq	96
56) 2,4-Dinitrotoluene	10.19	165	74052	36.006	nq	90
57) Fluorene	10.55	166	237491	35.324	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.32	232	70197	38.106	nq	# 78
59) Diethylphthalate	10.41	149	225699	31.427	nq	# 93
60) 4-Chlorophenyl-phenylether	10.54	204	120790	34.337	nq	# 86
61) 4-Nitroaniline	10.58	138	53152	29.538	nq	94
62) Azobenzene	10.70	77	216662	30.635	nq	95
64) 4,6-Dinitro-2-methylphenol	10.60	198	42484	33.789	nq	79
65) n-Nitrosodiphenylamine	10.66	169	206242	30.146	nq	100
66) 4-Bromophenyl-phenylether	11.03	248	85541	35.238	nq	# 78
67) Hexachlorobenzene	11.10	284	90605	35.661	nq	# 87
68) Atrazine	11.18	200	71080	32.681	nq	97
69) Pentachlorophenol	11.30	266	72613	57.278	nq	99
70) Phenanthrene	11.52	178	349856	35.661	nq	99
71) Anthracene	11.57	178	362147	36.165	nq	100
72) Carbazole	11.72	167	306418	32.684	nq	98
73) Di-n-butylphthalate	12.04	149	355943	32.227	nq	99
74) Fluoranthene	12.71	202	381349	35.267	nq	94
76) Benzidine	12.83	184	67885	16.008	nq	97
77) Pyrene	12.94	202	377303	38.426	nq	99
79) Butylbenzylphthalate	13.54	149	130355	32.289	nq	# 96
80) Benzo(a)anthracene	14.14	228	313926	34.592	nq	99
81) 3,3'-Dichlorobenzidine	14.09	252	58802	18.436	nq	# 97
82) Chrysene	14.18	228	293684	35.176	nq	98
83) Bis(2-ethylhexyl)phthalate	14.10	149	166290	32.219	nq	# 97
84) Di-n-octyl phthalate	14.74	149	252311	29.990	nq	96
85) Indeno(1,2,3-cd)pyrene	17.28	276	336969	47.559	nq	# 89
87) Benzo(b)fluoranthene	15.24	252	272334	30.383	nq	# 95
88) Benzo(k)fluoranthene	15.27	252	275806	32.311	nq	# 95
89) Benzo(a)pyrene	15.63	252	272837	34.240	nq	# 97
90) Dibenzo(a,h)anthracene	17.29	278	279828	41.437	nq	# 93
91) Benzo(g,h,i)perylene	17.76	276	295600	44.784	nq	# 88

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

