

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG042919\
 Data File : BG040657.D
 Acq On : 29 Apr 2019 12:17
 Operator : JU/SJ
 Sample : PB119139BSD
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB119139BSD

Manual Integrations
 APPROVED

Sohil
 4/30/2019 6:44:37 PM

Quant Time: Apr 30 05:37:22 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG042319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 30 05:31:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.15	152	66398	20.00	ng	-0.01
21) Naphthalene-d8	10.98	136	302032	20.00	ng	0.00
39) Acenaphthene-d10	14.78	164	216337	20.00	ng	-0.01
64) Phenanthrene-d10	17.53	188	582036	20.00	ng	0.00
76) Chrysene-d12	21.82	240	656861	20.00	ng	0.00
87) Perylene-d12	25.19	264	487731	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.69	112	599080	159.05	ng	0.00
7) Phenol-d6	7.30	99	855752	147.55	ng	0.00
23) Nitrobenzene-d5	9.34	82	531888	81.37	ng	0.00
42) 2,4,6-Tribromophenol	16.27	330	436490	146.79	ng	0.00
45) 2-Fluorobiphenyl	13.41	172	1088807	72.69	ng	-0.01
79) Terphenyl-d14	20.12	244	2135885	72.04	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.55	88	73567	42.946	ng	97
3) Pyridine	3.95	79	129827	26.147	ng	98
4) n-Nitrosodimethylamine	3.88	42	107380	38.990	ng	# 99
6) Aniline	7.47	93	229100	29.803	ng	97
8) 2-Chlorophenol	7.71	128	214687	46.678	ng	94
9) Benzaldehyde	7.29	77	11180	3.109	ng	# 75
10) Phenol	7.33	94	306451	49.155	ng	95
11) bis(2-Chloroethyl)ether	7.57	93	204399	43.722	ng	99
12) 1,3-Dichlorobenzene	8.04	146	206198	41.075	ng	98
13) 1,4-Dichlorobenzene	8.19	146	208891	41.048	ng	97
14) 1,2-Dichlorobenzene	8.51	146	206136	42.002	ng	97
15) Benzyl Alcohol	8.39	79	269047	44.584	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.68	45	392768	42.198	ng	99
17) 2-Methylphenol	8.58	107	209976	47.592	ng	95
18) Hexachloroethane	9.24	117	83218	39.864	ng	99
19) n-Nitroso-di-n-propylamine	8.96	70	210680	40.355	ng	99
20) 3+4-Methylphenols	8.91	107	289539	47.109	ng	98
22) Acetophenone	8.98	105	360489	42.922	ng	98
24) Nitrobenzene	9.38	77	314071	45.033	ng	97
25) Isophorone	9.89	82	578772	39.750	ng	99
26) 2-Nitrophenol	10.09	139	120722	48.290	ng	94
27) 2,4-Dimethylphenol	10.13	122	235713	50.252	ng	99
28) bis(2-Chloroethoxy)methane	10.37	93	295056	40.390	ng	99
29) 2,4-Dichlorophenol	10.62	162	244879	45.921	ng	95
30) 1,2,4-Trichlorobenzene	10.83	180	251631	42.552	ng	98
31) Naphthalene	11.03	128	666951	41.565	ng	99
32) Benzoic acid	10.28	122	173141	53.920	ng	# 86
33) 4-Chloroaniline	11.13	127	179039	25.166	ng	97
34) Hexachlorobutadiene	11.29	225	181632	41.604	ng	99
35) Caprolactam	11.93	113	87862m	41.733	ng	
36) 4-Chloro-3-methylphenol	12.23	107	296898	44.134	ng	96
37) 2-Methylnaphthalene	12.63	142	507892	42.334	ng	98
38) 1-Methylnaphthalene	12.84	142	482403	41.137	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.98	216	336498	41.754	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.95	237	294626	61.955	ng	98
43) 2,4,6-Trichlorophenol	13.21	196	235825	48.422	ng	99
44) 2,4,5-Trichlorophenol	13.29	196	246896	46.537	ng	99
46) 1,1'-Biphenyl	13.62	154	676955	40.303	ng	94
47) 2-Chloronaphthalene	13.67	162	543971	41.383	ng	99
48) 2-Nitroaniline	13.87	65	238477	50.967	ng	96
49) Acenaphthylene	14.51	152	863322	38.711	ng	98
50) Dimethylphthalate	14.24	163	748057	41.328	ng	99
51) 2,6-Dinitrotoluene	14.36	165	166796	47.215	ng	96
52) Acenaphthene	14.85	154	520671	40.089	ng	95
53) 3-Nitroaniline	14.69	138	137001	37.850	ng	98
54) 2,4-Dinitrophenol	14.89	184	213192	104.635	ng	# 77
55) Dibenzofuran	15.18	168	877546	41.252	ng	97
56) 4-Nitrophenol	14.99	139	306585	97.682	ng	96
57) 2,4-Dinitrotoluene	15.15	165	247726	51.606	ng	94
58) Fluorene	15.83	166	709431	41.260	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.40	232	264363	49.507	ng	100
60) Diethylphthalate	15.59	149	797658	41.768	ng	98
61) 4-Chlorophenyl-phenylether	15.82	204	427462	42.745	ng	99
62) 4-Nitroaniline	15.86	138	188234	46.489	ng	94
63) Azobenzene	16.11	77	808375	41.102	ng	99
65) 4,6-Dinitro-2-methylphenol	15.90	198	145136	50.222	ng	99
66) n-Nitrosodiphenylamine	16.03	169	667989	39.009	ng	99
67) 4-Bromophenyl-phenylether	16.71	248	295368	40.733	ng	95
68) Hexachlorobenzene	16.82	284	304033	40.549	ng	98
69) Atrazine	16.97	200	261898	38.602	ng	97
70) Pentachlorophenol	17.17	266	419437	95.597	ng	99
71) Phenanthrene	17.57	178	1258738	40.151	ng	96
72) Anthracene	17.66	178	1287906	40.870	ng	98
73) Carbazole	17.93	167	1188532	41.398	ng	98
74) Di-n-butylphthalate	18.47	149	1426311	41.161	ng	98
75) Fluoranthene	19.57	202	1625242	41.601	ng	98
77) Benzidine	19.75	184	339287	18.865	ng	100
78) Pyrene	19.93	202	1587010	38.548	ng	95
80) Butylbenzylphthalate	20.80	149	673898	41.998	ng	99
81) Benzo(a)anthracene	21.80	228	1622608	39.086	ng	97
82) 3,3'-Dichlorobenzidine	21.71	252	513757	34.721	ng	98
83) Chrysene	21.87	228	1581542	40.058	ng	97
84) Bis(2-ethylhexyl)phthalate	21.67	149	935140	40.373	ng	96
85) Di-n-octyl phthalate	22.94	149	1580411	40.515	ng	99
86) Indeno(1,2,3-cd)pyrene	29.08	276	1979753	41.474	ng	# 90
88) Benzo(b)fluoranthene	24.11	252	1690178	56.709	ng	98
89) Benzo(k)fluoranthene	24.18	252	1585848	56.974	ng	97
90) Benzo(a)pyrene	25.03	252	1614092	57.212	ng	98
91) Dibenzo(a,h)anthracene	29.15	278	1624276	56.892	ng	99
92) Benzo(g,h,i)perylene	30.30	276	1624681	57.179	ng	98

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

