

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG100719\
 Data File : BG043066.D
 Acq On : 7 Oct 2019 14:03
 Operator : JU
 Sample : PB123594BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB123594BS

Manual Integrations
 APPROVED

mohammad
 10/8/2019 12:22:39 PM

Quant Time: Oct 08 02:02:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG092519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 25 15:35:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	42270	20.00	ng	-0.02
21) Naphthalene-d8	10.86	136	167140	20.00	ng	-0.02
39) Acenaphthene-d10	14.68	164	129592	20.00	ng	-0.01
64) Phenanthrene-d10	17.42	188	333923	20.00	ng	-0.02
76) Chrysene-d12	21.69	240	383667	20.00	ng	-0.02
87) Perylene-d12	24.91	264	434295	20.00	ng	-0.05

System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	278577	117.61	ng	-0.02
7) Phenol-d6	7.23	99	387191	113.52	ng	-0.01
23) Nitrobenzene-d5	9.22	82	199965	58.15	ng	-0.02
42) 2,4,6-Tribromophenol	16.17	330	270068	121.19	ng	-0.01
45) 2-Fluorobiphenyl	13.30	172	538852	60.28	ng	-0.02
79) Terphenyl-d14	20.03	244	1235582	68.62	ng	-0.02

Target Compounds					Qvalue
2) 1,4-Dioxane	3.53	88	33329	33.750 ng	89
3) Pyridine	3.93	79	83829	30.181 ng	90
4) n-Nitrosodimethylamine	3.84	42	45792	34.284 ng	84
6) Aniline	7.38	93	137650	33.432 ng	100
8) 2-Chlorophenol	7.63	128	112893	45.704 ng	93
9) Benzaldehyde	7.19	77	72342	34.708 ng	96
10) Phenol	7.26	94	139943	44.719 ng	93
11) bis(2-Chloroethyl)ether	7.48	93	89603	37.006 ng	93
12) 1,3-Dichlorobenzene	7.95	146	117755	37.370 ng	96
13) 1,4-Dichlorobenzene	8.09	146	122225	38.910 ng	99
14) 1,2-Dichlorobenzene	8.41	146	116020	38.795 ng	94
15) Benzyl Alcohol	8.30	79	102151	39.834 ng	96
16) 2,2'-oxybis(1-Chloropropan	8.59	45	131834	28.352 ng	96
17) 2-Methylphenol	8.51	107	93246	42.585 ng	97
18) Hexachloroethane	9.14	117	41483	35.065 ng	96
19) n-Nitroso-di-n-propylamine	8.86	70	81985	36.606 ng	92
20) 3+4-Methylphenols	8.83	107	131733	43.901 ng	94
22) Acetophenone	8.88	105	157114	36.467 ng	92
24) Nitrobenzene	9.26	77	127916	38.999 ng	95
25) Isophorone	9.78	82	233668	38.685 ng	99
26) 2-Nitrophenol	9.97	139	59649	42.719 ng	97
27) 2,4-Dimethylphenol	10.03	122	106898	49.851 ng	90
28) bis(2-Chloroethoxy)methane	10.26	93	128332	37.447 ng	96
29) 2,4-Dichlorophenol	10.51	162	123980	46.489 ng	97
30) 1,2,4-Trichlorobenzene	10.72	180	131006	38.039 ng	97
31) Naphthalene	10.91	128	333512	40.535 ng	98
32) Benzoic acid	10.17	122	61678	38.636 ng	95
33) 4-Chloroaniline	11.02	127	70606	19.776 ng	97
34) Hexachlorobutadiene	11.20	225	91673	35.550 ng	92
35) Caprolactam	11.79	113	41199m	43.806 ng	
36) 4-Chloro-3-methylphenol	12.14	107	130123	44.875 ng	89
37) 2-Methylnaphthalene	12.51	142	261939	41.732 ng	97
38) 1-Methylnaphthalene	12.73	142	248726	41.879 ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.88	216	171160	35.127 ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.86	237	248871	80.163	ng	94
43) 2,4,6-Trichlorophenol	13.12	196	116922	40.998	ng	98
44) 2,4,5-Trichlorophenol	13.19	196	124833	42.490	ng	96
46) 1,1'-Biphenyl	13.51	154	358050	36.433	ng	99
47) 2-Chloronaphthalene	13.56	162	282062	38.281	ng	99
48) 2-Nitroaniline	13.76	65	91896	37.505	ng	90
49) Acenaphthylene	14.40	152	464667	41.214	ng	98
50) Dimethylphthalate	14.13	163	387278	39.885	ng	99
51) 2,6-Dinitrotoluene	14.25	165	87556	42.343	ng	95
52) Acenaphthene	14.74	154	284872	37.749	ng	100
53) 3-Nitroaniline	14.58	138	57972	27.128	ng #	99
54) 2,4-Dinitrophenol	14.78	184	111471	86.728	ng	97
55) Dibenzofuran	15.08	168	458192	41.044	ng	98
56) 4-Nitrophenol	14.90	139	149785	91.993	ng	96
57) 2,4-Dinitrotoluene	15.03	165	128503	42.437	ng #	97
58) Fluorene	15.72	166	390101	40.930	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.31	232	132210	42.265	ng	100
60) Diethylphthalate	15.49	149	399940	40.473	ng	98
61) 4-Chlorophenyl-phenylether	15.72	204	226741	39.668	ng	98
62) 4-Nitroaniline	15.75	138	92425	39.657	ng	98
63) Azobenzene	16.01	77	342444	38.063	ng	97
65) 4,6-Dinitro-2-methylphenol	15.80	198	82806	43.823	ng	91
66) n-Nitrosodiphenylamine	15.93	169	352905	40.034	ng	99
67) 4-Bromophenyl-phenylether	16.61	248	158697	37.866	ng	99
68) Hexachlorobenzene	16.73	284	183722	39.371	ng	97
69) Atrazine	16.87	200	163993	44.179	ng	95
70) Pentachlorophenol	17.07	266	237546	88.220	ng	98
71) Phenanthrene	17.47	178	661073	40.551	ng	98
72) Anthracene	17.56	178	690865	42.450	ng	100
73) Carbazole	17.83	167	638255	42.291	ng	98
74) Di-n-butylphthalate	18.38	149	743019	41.264	ng	99
75) Fluoranthene	19.47	202	919295	42.634	ng	99
77) Benzidine	19.65	184	639982	66.619	ng	99
78) Pyrene	19.84	202	936577	40.899	ng	100
80) Butylbenzylphthalate	20.72	149	344485	39.632	ng	97
81) Benzo(a)anthracene	21.67	228	962663	40.269	ng	100
82) 3,3'-Dichlorobenzidine	21.59	252	265100	28.274	ng	96
83) Chrysene	21.74	228	917686	39.557	ng	99
84) Bis(2-ethylhexyl)phthalate	21.58	149	496200	40.119	ng	98
85) Di-n-octyl phthalate	22.79	149	838010	41.435	ng	96
86) Indeno(1,2,3-cd)pyrene	28.57	276	1205337	41.735	ng #	98
88) Benzo(b)fluoranthene	23.89	252	1021372	41.564	ng	99
89) Benzo(k)fluoranthene	23.96	252	974419	40.083	ng	99
90) Benzo(a)pyrene	24.76	252	933955	40.284	ng	99
91) Dibenzo(a,h)anthracene	28.64	278	1013019	42.611	ng	98
92) Benzo(g,h,i)perylene	29.72	276	1013471	43.998	ng	98

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

