

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG032919\
 Data File : BG040230.D
 Acq On : 29 Mar 2019 17:55
 Operator : JU/SJ
 Sample : PB118394BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB118394BS

Manual Integrations
 APPROVED

Sohil
 4/1/2019 10:06:01 AM

Quant Time: Mar 30 04:38:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 28 05:47:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	54119	20.00	ng	-0.01
21) Naphthalene-d8	10.87	136	240814	20.00	ng	-0.01
39) Acenaphthene-d10	14.69	164	159004	20.00	ng	-0.01
64) Phenanthrene-d10	17.44	188	408455	20.00	ng	-0.01
76) Chrysene-d12	21.71	240	392861	20.00	ng	-0.02
87) Perylene-d12	25.00	264	359438	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	460013	141.31	ng	-0.01
7) Phenol-d6	7.21	99	637753	141.75	ng	-0.01
23) Nitrobenzene-d5	9.23	82	345856	76.34	ng	-0.01
42) 2,4,6-Tribromophenol	16.18	330	252120	146.72	ng	-0.01
45) 2-Fluorobiphenyl	13.31	172	811696	75.64	ng	-0.01
79) Terphenyl-d14	20.03	244	1314671	75.28	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.51	88	57304	37.435	ng	98
3) Pyridine	3.90	79	141826	34.411	ng	99
4) n-Nitrosodimethylamine	3.83	42	67804	35.565	ng	# 88
6) Aniline	7.38	93	178503	30.538	ng	96
8) 2-Chlorophenol	7.62	128	162329	42.597	ng	97
9) Benzaldehyde	7.20	77	46096	14.937	ng	96
10) Phenol	7.24	94	213504	43.720	ng	98
11) bis(2-Chloroethyl)ether	7.48	93	153391	39.217	ng	99
12) 1,3-Dichlorobenzene	7.95	146	155072	37.434	ng	97
13) 1,4-Dichlorobenzene	8.09	146	162628	39.416	ng	97
14) 1,2-Dichlorobenzene	8.41	146	151961	38.514	ng	97
15) Benzyl Alcohol	8.30	79	147339	40.664	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.58	45	282448	37.627	ng	99
17) 2-Methylphenol	8.49	107	150791	44.623	ng	98
18) Hexachloroethane	9.13	117	58249	37.115	ng	98
19) n-Nitroso-di-n-propylamine	8.86	70	122963	38.119	ng	98
20) 3+4-Methylphenols	8.82	107	206653	45.143	ng	100
22) Acetophenone	8.89	105	239815	39.922	ng	# 98
24) Nitrobenzene	9.28	77	180433	38.459	ng	97
25) Isophorone	9.79	82	363951	39.043	ng	99
26) 2-Nitrophenol	9.99	139	90255	40.600	ng	96
27) 2,4-Dimethylphenol	10.03	122	170430	48.265	ng	95
28) bis(2-Chloroethoxy)methane	10.27	93	219485	39.797	ng	99
29) 2,4-Dichlorophenol	10.51	162	168901	44.067	ng	99
30) 1,2,4-Trichlorobenzene	10.73	180	166644	39.596	ng	99
31) Naphthalene	10.93	128	498909	40.419	ng	98
32) Benzoic acid	10.14	122	35591m	16.682	ng	
33) 4-Chloroaniline	11.04	127	149081	28.315	ng	97
34) Hexachlorobutadiene	11.18	225	101023	37.533	ng	96
35) Caprolactam	11.85	113	69704	45.827	ng	# 84
36) 4-Chloro-3-methylphenol	12.15	107	191320	43.420	ng	98
37) 2-Methylnaphthalene	12.52	142	383999	42.063	ng	99
38) 1-Methylnaphthalene	12.74	142	364896	41.220	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.88	216	190610	37.717	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.85	237	223085	80.395	ng	96
43) 2,4,6-Trichlorophenol	13.12	196	142558	43.752	ng	95
44) 2,4,5-Trichlorophenol	13.20	196	149246	42.024	ng	97
46) 1,1'-Biphenyl	13.52	154	495280	39.422	ng	98
47) 2-Chloronaphthalene	13.57	162	383295	39.774	ng	100
48) 2-Nitroaniline	13.78	65	131966	40.597	ng	97
49) Acenaphthylene	14.42	152	644413	39.440	ng	100
50) Dimethylphthalate	14.14	163	511849	41.131	ng	99
51) 2,6-Dinitrotoluene	14.28	165	118064	42.253	ng	95
52) Acenaphthene	14.76	154	380842	40.677	ng	96
53) 3-Nitroaniline	14.60	138	89965	30.417	ng	99
54) 2,4-Dinitrophenol	14.81	184	125162	84.227	ng	96
55) Dibenzofuran	15.09	168	624362	41.501	ng	99
56) 4-Nitrophenol	14.91	139	211043	88.408	ng	99
57) 2,4-Dinitrotoluene	15.06	165	167514	43.145	ng	# 95
58) Fluorene	15.74	166	491145	41.523	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.31	232	148731	46.150	ng	98
60) Diethylphthalate	15.49	149	548883	43.103	ng	99
61) 4-Chlorophenyl-phenylether	15.72	204	267004	42.231	ng	97
62) 4-Nitroaniline	15.77	138	134612	42.409	ng	98
63) Azobenzene	16.01	77	499733	41.604	ng	98
65) 4,6-Dinitro-2-methylphenol	15.81	198	97716	42.582	ng	97
66) n-Nitrosodiphenylamine	15.94	169	472730	39.551	ng	98
67) 4-Bromophenyl-phenylether	16.62	248	177310	38.575	ng	99
68) Hexachlorobenzene	16.73	284	182982	39.593	ng	96
69) Atrazine	16.88	200	169152	38.044	ng	98
70) Pentachlorophenol	17.08	266	213423	79.876	ng	97
71) Phenanthrene	17.48	178	857633	39.947	ng	100
72) Anthracene	17.57	178	862473	40.136	ng	99
73) Carbazole	17.84	167	814462	40.049	ng	99
74) Di-n-butylphthalate	18.38	149	945605	39.227	ng	100
75) Fluoranthene	19.49	202	1031428	40.572	ng	100
77) Benzidine	19.67	184	375338	26.457	ng	97
78) Pyrene	19.85	202	1019810	39.474	ng	99
80) Butylbenzylphthalate	20.71	149	443884	40.152	ng	98
81) Benzo(a)anthracene	21.69	228	946143	39.100	ng	99
82) 3,3'-Dichlorobenzidine	21.61	252	248312	26.976	ng	98
83) Chrysene	21.76	228	917113	39.436	ng	98
84) Bis(2-ethylhexyl)phthalate	21.56	149	627171	39.687	ng	100
85) Di-n-octyl phthalate	22.78	149	1070698	39.767	ng	99
86) Indeno(1,2,3-cd)pyrene	28.78	276	1125408	39.567	ng	99
88) Benzo(b)fluoranthene	23.95	252	985968	44.804	ng	97
89) Benzo(k)fluoranthene	24.02	252	952230	47.054	ng	98
90) Benzo(a)pyrene	24.85	252	966046	46.787	ng	99
91) Dibenzo(a,h)anthracene	28.83	278	921622	48.060	ng	99
92) Benzo(g,h,i)perylene	29.97	276	944488	47.925	ng	100

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

