

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102919\
 Data File : BG043372.D
 Acq On : 29 Oct 2019 13:49
 Operator : JU
 Sample : PB124300BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB124300BS

Manual Integrations
 APPROVED

mohammad
 10/30/2019 4:54:27 PM

Quant Time: Oct 30 08:09:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 21 16:37:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	35581	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	142823	20.00	ng	0.00
39) Acenaphthene-d10	14.65	164	106696	20.00	ng	0.00
64) Phenanthrene-d10	17.39	188	249915	20.00	ng	0.00
76) Chrysene-d12	21.67	240	287669	20.00	ng	0.00
87) Perylene-d12	24.86	264	329438	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	258571	133.17	ng	0.00
7) Phenol-d6	7.21	99	340824	122.00	ng	0.00
23) Nitrobenzene-d5	9.18	82	175241	63.26	ng	0.00
42) 2,4,6-Tribromophenol	16.14	330	216855	106.80	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	470148	60.89	ng	0.00
79) Terphenyl-d14	20.00	244	946032	61.70	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	30118	39.803	ng	92
3) Pyridine	3.89	79	70299	36.830	ng	98
4) n-Nitrosodimethylamine	3.80	42	43533	45.198	ng	# 96
6) Aniline	7.35	93	108764	33.796	ng	95
8) 2-Chlorophenol	7.60	128	99097	46.233	ng	95
9) Benzaldehyde	7.16	77	55601	40.498	ng	# 80
10) Phenol	7.24	94	119686	46.883	ng	98
11) bis(2-Chloroethyl)ether	7.44	93	80277	40.673	ng	98
12) 1,3-Dichlorobenzene	7.91	146	111777	41.622	ng	95
13) 1,4-Dichlorobenzene	8.06	146	112815	41.520	ng	98
14) 1,2-Dichlorobenzene	8.38	146	108274	40.812	ng	99
15) Benzyl Alcohol	8.27	79	83772	42.697	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.55	45	111514	38.856	ng	97
17) 2-Methylphenol	8.49	107	84422	45.363	ng	99
18) Hexachloroethane	9.11	117	40344	41.155	ng	100
19) n-Nitroso-di-n-propylamine	8.83	70	72423	40.119	ng	95
20) 3+4-Methylphenols	8.82	107	110932	43.470	ng	97
22) Acetophenone	8.85	105	142703	39.016	ng	98
24) Nitrobenzene	9.23	77	111269	42.163	ng	96
25) Isophorone	9.74	82	200233	39.533	ng	100
26) 2-Nitrophenol	9.94	139	56582	44.410	ng	93
27) 2,4-Dimethylphenol	10.01	122	90186	49.387	ng	96
28) bis(2-Chloroethoxy)methane	10.23	93	110355	39.477	ng	99
29) 2,4-Dichlorophenol	10.49	162	105122	44.929	ng	94
30) 1,2,4-Trichlorobenzene	10.69	180	119591	40.554	ng	99
31) Naphthalene	10.88	128	308769	42.591	ng	97
32) Benzoic acid	10.17	122	43132	34.151	ng	96
33) 4-Chloroaniline	11.00	127	62138	20.910	ng	95
34) Hexachlorobutadiene	11.17	225	84535	38.779	ng	92
35) Caprolactam	11.76	113	31590m	39.202	ng	
36) 4-Chloro-3-methylphenol	12.13	107	110034	43.114	ng	97
37) 2-Methylnaphthalene	12.48	142	231578	41.632	ng	97
38) 1-Methylnaphthalene	12.70	142	221169	41.788	ng	96
40) 1,2,4,5-Tetrachlorobenzene	12.85	216	148300	37.557	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.83	237	178617	86.111	nq	97
43) 2,4,6-Trichlorophenol	13.09	196	95076	40.886	nq	99
44) 2,4,5-Trichlorophenol	13.18	196	100284	42.657	nq	98
46) 1,1'-Biphenyl	13.49	154	311097	38.327	nq	100
47) 2-Chloronaphthalene	13.53	162	242867	39.325	nq	99
48) 2-Nitroaniline	13.74	65	70345	38.110	nq	95
49) Acenaphthylene	14.37	152	394799	41.074	nq	97
50) Dimethylphthalate	14.10	163	313343	37.915	nq	99
51) 2,6-Dinitrotoluene	14.22	165	71720	39.946	nq	97
52) Acenaphthene	14.72	154	230531	39.329	nq	97
53) 3-Nitroaniline	14.56	138	45058	25.993	nq	95
54) 2,4-Dinitrophenol	14.77	184	72488	66.052	nq	94
55) Dibenzofuran	15.05	168	379326	39.906	nq	99
56) 4-Nitrophenol	14.90	139	103821	89.375	nq	97
57) 2,4-Dinitrotoluene	15.01	165	102740	40.041	nq	93
58) Fluorene	15.70	166	319152	40.032	nq	96
59) 2,3,4,6-Tetrachlorophenol	15.28	232	103552	41.458	nq	# 99
60) Diethylphthalate	15.47	149	317905	38.284	nq	98
61) 4-Chlorophenyl-phenylether	15.69	204	182247	39.185	nq	97
62) 4-Nitroaniline	15.73	138	66394	37.087	nq	94
63) Azobenzene	15.98	77	265770	39.546	nq	97
65) 4,6-Dinitro-2-methylphenol	15.78	198	60168	40.910	nq	98
66) n-Nitrosodiphenylamine	15.91	169	282687	40.065	nq	98
67) 4-Bromophenyl-phenylether	16.58	248	131490	39.096	nq	97
68) Hexachlorobenzene	16.71	284	146246	37.881	nq	98
69) Atrazine	16.85	200	121294	49.473	nq	94
70) Pentachlorophenol	17.05	266	161246	91.661	nq	97
71) Phenanthrene	17.44	178	516491	40.359	nq	99
72) Anthracene	17.53	178	534541	41.748	nq	99
73) Carbazole	17.81	167	480073	41.403	nq	97
74) Di-n-butylphthalate	18.35	149	567505	40.337	nq	99
75) Fluoranthene	19.45	202	694861	41.648	nq	99
77) Benzidine	19.63	184	299628	51.754	nq	99
78) Pyrene	19.81	202	701561	39.399	nq	98
80) Butylbenzylphthalate	20.69	149	260045	38.547	nq	98
81) Benzo(a)anthracene	21.64	228	721863	40.060	nq	99
82) 3,3'-Dichlorobenzidine	21.56	252	220958	31.703	nq	98
83) Chrysene	21.71	228	711603	39.746	nq	98
84) Bis(2-ethylhexyl)phthalate	21.55	149	385178	40.194	nq	99
85) Di-n-octyl phthalate	22.75	149	663559	40.814	nq	99
86) Indeno(1,2,3-cd)pyrene	28.50	276	931633	41.455	nq	99
88) Benzo(b)fluoranthene	23.85	252	769177	41.224	nq	98
89) Benzo(k)fluoranthene	23.92	252	786110	41.851	nq	100
90) Benzo(a)pyrene	24.72	252	710700	39.888	nq	99
91) Dibenzo(a,h)anthracene	28.56	278	782190	42.920	nq	97
92) Benzo(g,h,i)perylene	29.65	276	771139	42.897	nq	97

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

