

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012420\
 Data File : BG044209.D
 Acq On : 25 Jan 2020 9:39
 Operator : CG/JU
 Sample : L1247-21MS
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-10-9-COMPMS

Manual Integrations
 APPROVED

mohammad
 1/28/2020 8:11:49 AM

Quant Time: Jan 27 05:37:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 31 13:32:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	84999	20.00	ng	-0.04
21) Naphthalene-d8	10.73	136	383484	20.00	ng	-0.04
39) Acenaphthene-d10	14.57	164	255585	20.00	ng	-0.03
64) Phenanthrene-d10	17.31	188	505394	20.00	ng	-0.02
76) Chrysene-d12	21.56	240	420008	20.00	ng	-0.03
87) Perylene-d12	24.67	264	490603	20.00	ng	-0.05

System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	523453	113.07	ng	-0.03
7) Phenol-d6	7.10	99	708504	109.49	ng	-0.02
23) Nitrobenzene-d5	9.09	82	425620	59.37	ng	-0.04
42) 2,4,6-Tribromophenol	16.05	330	250831	93.78	ng	-0.03
45) 2-Fluorobiphenyl	13.19	172	923402	65.46	ng	-0.04
79) Terphenyl-d14	19.92	244	1233673	66.61	ng	-0.04

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.39	88	78651	38.966	ng	99
3) Pyridine	3.79	79	201360	33.769	ng	95
4) n-Nitrosodimethylamine	3.70	42	105094	39.587	ng	99
6) Aniline	7.25	93	112121	13.192	ng	98
8) 2-Chlorophenol	7.49	128	269255	49.544	ng	97
9) Benzaldehyde	7.06	77	128181	30.950	ng	97
10) Phenol	7.13	94	352393	52.856	ng	98
11) bis(2-Chloroethyl)ether	7.35	93	237960	41.348	ng	99
12) 1,3-Dichlorobenzene	7.82	146	262143	41.220	ng	100
13) 1,4-Dichlorobenzene	7.96	146	266171	42.418	ng	99
14) 1,2-Dichlorobenzene	8.29	146	252213	41.875	ng	100
15) Benzyl Alcohol	8.16	79	221643	43.400	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.46	45	333409	37.446	ng	99
17) 2-Methylphenol	8.38	107	230271	47.728	ng	98
18) Hexachloroethane	9.02	117	101585	41.605	ng	97
19) n-Nitroso-di-n-propylamine	8.73	70	197064	40.893	ng	99
20) 3+4-Methylphenols	8.70	107	316009	46.951	ng	95
22) Acetophenone	8.75	105	364414	38.224	ng	100
24) Nitrobenzene	9.13	77	279450	39.360	ng	97
25) Isophorone	9.65	82	533653	39.680	ng	99
26) 2-Nitrophenol	9.84	139	148907	44.330	ng	98
27) 2,4-Dimethylphenol	9.91	122	258875	51.198	ng	98
28) bis(2-Chloroethoxy)methane	10.14	93	342596	38.210	ng	99
29) 2,4-Dichlorophenol	10.38	162	267831	44.406	ng	97
30) 1,2,4-Trichlorobenzene	10.60	180	263132	38.910	ng	98
31) Naphthalene	10.78	128	784283	42.263	ng	99
32) Benzoic acid	10.06	122	114827	30.510	ng	96
33) 4-Chloroaniline	10.88	127	74721	8.442	ng	96
34) Hexachlorobutadiene	11.08	225	152827	37.013	ng	98
35) Caprolactam	11.66	113	93876m	41.810	ng	
36) 4-Chloro-3-methylphenol	12.02	107	285383	44.447	ng	99
37) 2-Methylnaphthalene	12.39	142	593181	42.458	ng	96
38) 1-Methylnaphthalene	12.61	142	565882	42.736	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.76	216	280726	37.474	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.75	237	326615	84.098	nq	98
43) 2,4,6-Trichlorophenol	13.00	196	210110	40.762	nq	99
44) 2,4,5-Trichlorophenol	13.08	196	223132	41.971	nq	99
46) 1,1'-Biphenyl	13.40	154	751676	41.019	nq	98
47) 2-Chloronaphthalene	13.44	162	588286	39.728	nq	99
48) 2-Nitroaniline	13.64	65	179171	37.615	nq	97
49) Acenaphthylene	14.29	152	927894	43.597	nq	98
50) Dimethylphthalate	14.02	163	874792	48.204	nq	98
51) 2,6-Dinitrotoluene	14.13	165	168417	41.060	nq	96
52) Acenaphthene	14.63	154	572697	38.370	nq	99
53) 3-Nitroaniline	14.46	138	63671	13.669	nq	98
54) 2,4-Dinitrophenol	14.67	184	154660	74.423	nq	97
55) Dibenzofuran	14.97	168	874625	42.567	nq	98
56) 4-Nitrophenol	14.79	139	292204	81.864	nq	98
57) 2,4-Dinitrotoluene	14.93	165	228512	40.943	nq	98
58) Fluorene	15.61	166	687037	42.187	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.20	232	190937	41.767	nq	96
60) Diethylphthalate	15.38	149	738945	40.711	nq	97
61) 4-Chlorophenyl-phenylether	15.61	204	353727	40.001	nq	98
62) 4-Nitroaniline	15.63	138	159275	34.627	nq	95
63) Azobenzene	15.90	77	701778	41.690	nq	99
65) 4,6-Dinitro-2-methylphenol	15.69	198	126828	48.642	nq	97
66) n-Nitrosodiphenylamine	15.82	169	635016	42.667	nq	98
67) 4-Bromophenyl-phenylether	16.50	248	226169	39.789	nq	98
68) Hexachlorobenzene	16.62	284	242952	39.667	nq	98
69) Atrazine	16.77	200	227790	47.085	nq	99
70) Pentachlorophenol	16.97	266	257889	76.382	nq	98
71) Phenanthrene	17.35	178	1060679	43.755	nq	98
72) Anthracene	17.44	178	1088685	45.443	nq	98
73) Carbazole	17.71	167	958163	43.297	nq	97
74) Di-n-butylphthalate	18.28	149	1187237	42.020	nq	97
75) Fluoranthene	19.36	202	1222341	44.091	nq	96
77) Benzidine	19.54	184	293615	27.812	nq	96
78) Pyrene	19.72	202	1221204	44.544	nq	96
80) Butylbenzylphthalate	20.61	149	559278	42.849	nq	96
81) Benzo(a)anthracene	21.54	228	1147511	44.061	nq	96
82) 3,3'-Dichlorobenzidine	21.46	252	173834	16.895	nq	99
83) Chrysene	21.61	228	1097953	44.368	nq	95
84) Bis(2-ethylhexyl)phthalate	21.46	149	794809	44.132	nq	96
85) Di-n-octyl phthalate	22.63	149	1357543	44.837	nq	97
86) Indeno(1,2,3-cd)pyrene	28.19	276	1400154	41.633	nq	# 91
88) Benzo(b)fluoranthene	23.69	252	1199736	41.366	nq	98
89) Benzo(k)fluoranthene	23.75	252	1183712	42.919	nq	99
90) Benzo(a)pyrene	24.52	252	1112406	40.637	nq	97
91) Dibenzo(a,h)anthracene	28.25	278	1147335	41.734	nq	98
92) Benzo(g,h,i)perylene	29.29	276	1179007	43.175	nq	98

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

