

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021419\
 Data File : BG039512.D
 Acq On : 14 Feb 2019 18:40
 Operator : JU/SJ
 Sample : K1349-06MS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OR-01-021319-AMS

Manual Integrations
 APPROVED

Sohil
 2/15/2019 9:44:34 AM

Quant Time: Feb 15 04:41:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 29 15:41:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	52627	20.00	ng	-0.01
21) Naphthalene-d8	11.00	136	242305	20.00	ng	-0.01
39) Acenaphthene-d10	14.80	164	167227	20.00	ng	-0.01
64) Phenanthrene-d10	17.54	188	392907	20.00	ng	-0.01
76) Chrysene-d12	21.83	240	397008	20.00	ng	-0.01
87) Perylene-d12	25.21	264	415017	20.00	ng	-0.03

System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	460869	149.88	ng	0.00
7) Phenol-d6	7.32	99	607681	134.26	ng	0.00
23) Nitrobenzene-d5	9.35	82	447579	115.40	ng	-0.01
42) 2,4,6-Tribromophenol	16.28	330	298784	165.92	ng	0.00
45) 2-Fluorobiphenyl	13.42	172	1061096	91.03	ng	-0.01
79) Terphenyl-d14	20.13	244	1633403	87.09	ng	-0.02

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.61	88	47100	35.018	ng	# 3
3) Pyridine	4.02	79	136116	38.223	ng	96
4) n-Nitrosodimethylamine	3.93	42	73022	41.002	ng	86
6) Aniline	7.50	93	43254	7.629	ng	98
8) 2-Chlorophenol	7.74	128	172154	51.219	ng	95
9) Benzaldehyde	7.31	77	137473	85.857	ng	99
10) Phenol	7.35	94	196540	41.656	ng	95
11) bis(2-Chloroethyl)ether	7.59	93	164646	44.162	ng	99
12) 1,3-Dichlorobenzene	8.06	146	173901	42.709	ng	99
13) 1,4-Dichlorobenzene	8.21	146	179204	44.156	ng	98
14) 1,2-Dichlorobenzene	8.53	146	173166	44.076	ng	99
15) Benzyl Alcohol	8.41	79	174193	49.021	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.69	45	315016	37.416	ng	98
17) 2-Methylphenol	8.61	107	152129	49.825	ng	97
18) Hexachloroethane	9.26	117	61690	44.182	ng	94
19) n-Nitroso-di-n-propylamine	8.98	70	137732	42.874	ng	93
20) 3+4-Methylphenols	8.94	107	210901	50.161	ng	98
22) Acetophenone	9.00	105	268874	41.310	ng	# 95
24) Nitrobenzene	9.39	77	207415	49.500	ng	96
25) Isophorone	9.91	82	398534	46.368	ng	98
26) 2-Nitrophenol	10.10	139	103955	58.946	ng	97
27) 2,4-Dimethylphenol	10.15	122	184688	53.118	ng	98
28) bis(2-Chloroethoxy)methane	10.39	93	240303	45.391	ng	97
29) 2,4-Dichlorophenol	10.64	162	205295	54.025	ng	96
30) 1,2,4-Trichlorobenzene	10.85	180	192065	45.019	ng	98
31) Naphthalene	11.04	128	562576	46.801	ng	99
32) Benzoic acid	10.25	122	66366m	32.692	ng	
33) 4-Chloroaniline	11.15	127	10224	1.834	ng	# 90
34) Hexachlorobutadiene	11.31	225	121119	42.689	ng	95
35) Caprolactam	11.94	113	49749	31.637	ng	# 86
36) 4-Chloro-3-methylphenol	12.26	107	221488	50.399	ng	99
37) 2-Methylnaphthalene	12.63	142	429245	48.213	ng	99
38) 1-Methylnaphthalene	12.85	142	407451	47.476	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.99	216	230625	43.345	ng	98

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41) Hexachlorocyclopentadiene	12.96	237	236535	81.583	nq	97
43) 2,4,6-Trichlorophenol	13.23	196	173428	53.013	nq	95
44) 2,4,5-Trichlorophenol	13.30	196	185514	54.106	nq	98
46) 1,1'-Biphenyl	13.63	154	578311	41.564	nq	100
47) 2-Chloronaphthalene	13.68	162	456089	43.574	nq	98
48) 2-Nitroaniline	13.89	65	158237	51.827	nq	92
49) Acenaphthylene	14.53	152	731123	46.292	nq	98
50) Dimethylphthalate	14.24	163	595890	44.121	nq	99
51) 2,6-Dinitrotoluene	14.37	165	135078	52.760	nq	99
52) Acenaphthene	14.86	154	451269	44.017	nq	98
53) 3-Nitroaniline	14.71	138	9525	8.868	nq	# 89
54) 2,4-Dinitrophenol	14.91	184	30420	37.284	nq	89
55) Dibenzofuran	15.19	168	723820	44.035	nq	98
56) 4-Nitrophenol	15.01	139	187234	76.914	nq	90
57) 2,4-Dinitrotoluene	15.16	165	193540	57.924	nq	86
58) Fluorene	15.84	166	570250	46.538	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.41	232	173144	53.933	nq	98
60) Diethylphthalate	15.60	149	602325	43.134	nq	97
61) 4-Chlorophenyl-phenylether	15.82	204	308306	44.435	nq	95
62) 4-Nitroaniline	15.87	138	135683	47.916	nq	89
63) Azobenzene	16.12	77	544029	39.860	nq	97
65) 4,6-Dinitro-2-methylphenol	15.91	198	40606	31.107	nq	94
66) n-Nitrosodiphenylamine	16.04	169	527167	44.126	nq	99
67) 4-Bromophenyl-phenylether	16.72	248	198836	45.616	nq	96
68) Hexachlorobenzene	16.84	284	202488	46.301	nq	93
69) Atrazine	16.98	200	204259	46.785	nq	97
70) Pentachlorophenol	17.18	266	153047	50.353	nq	97
71) Phenanthrene	17.59	178	936683	46.061	nq	98
72) Anthracene	17.68	178	957327	47.793	nq	98
73) Carbazole	17.95	167	839326	41.987	nq	99
74) Di-n-butylphthalate	18.47	149	1026363	44.009	nq	99
75) Fluoranthene	19.58	202	1106144	46.864	nq	99
77) Benzidine	19.76	184	139857	10.745	nq	99
78) Pyrene	19.95	202	1115975	45.154	nq	98
80) Butylbenzylphthalate	20.81	149	487607	46.770	nq	95
81) Benzo(a)anthracene	21.81	228	1109033	47.275	nq	99
82) 3,3'-Dichlorobenzidine	21.72	252	72079	8.286	nq	99
83) Chrysene	21.88	228	1027076	45.992	nq	99
84) Bis(2-ethylhexyl)phthalate	21.68	149	672807	45.235	nq	# 97
85) Di-n-octyl phthalate	22.94	149	1142329	46.488	nq	96
86) Indeno(1,2,3-cd)pyrene	29.11	276	1244115	51.925	nq	98
88) Benzo(b)fluoranthene	24.13	252	1085676	47.968	nq	99
89) Benzo(k)fluoranthene	24.20	252	1054544	46.408	nq	100
90) Benzo(a)pyrene	25.05	252	1057051	48.494	nq	99
91) Dibenzo(a,h)anthracene	29.18	278	1019215	49.929	nq	98
92) Benzo(g,h,i)perylene	30.34	276	1031242	50.684	nq	98

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

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