

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040919\
 Data File : BG040401.D
 Acq On : 9 Apr 2019 16:24
 Operator : JU/SJ
 Sample : K2317-16MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COMP-3MSD

Manual Integrations
 APPROVED

Sohil
 4/10/2019 11:22:55 AM

Quant Time: Apr 10 08:34:35 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	8.05	152	55267	20.00	ng	-0.02
21) Naphthalene-d8	10.86	136	221382	20.00	ng	-0.02
39) Acenaphthene-d10	14.68	164	134811	20.00	ng	-0.02
64) Phenanthrene-d10	17.42	188	368303	20.00	ng	-0.02
76) Chrysene-d12	21.70	240	396828	20.00	ng	-0.03
87) Perylene-d12	24.98	264	424088	20.00	ng	-0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	377312	113.49	ng	-0.02
7) Phenol-d6	7.21	99	523047	113.84	ng	-0.02
23) Nitrobenzene-d5	9.22	82	303386	72.84	ng	-0.02
42) 2,4,6-Tribromophenol	16.17	330	196124	134.61	ng	-0.02
45) 2-Fluorobiphenyl	13.30	172	652892	71.76	ng	-0.02
79) Terphenyl-d14	20.03	244	1198007	67.92	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	53896	34.477	ng	99
3) Pyridine	3.90	79	126428	30.038	ng	98
4) n-Nitrosodimethylamine	3.82	42	64499	33.129	ng	# 84
6) Aniline	7.38	93	138104	23.136	ng	96
8) 2-Chlorophenol	7.61	128	138315	35.541	ng	97
9) Benzaldehyde	7.20	77	67037	21.271	ng	98
10) Phenol	7.23	94	186065	37.310	ng	98
11) bis(2-Chloroethyl)ether	7.47	93	132273	33.116	ng	97
12) 1,3-Dichlorobenzene	7.94	146	144942	34.261	ng	98
13) 1,4-Dichlorobenzene	8.08	146	149462	35.472	ng	96
14) 1,2-Dichlorobenzene	8.40	146	139644	34.657	ng	96
15) Benzyl Alcohol	8.29	79	120917	32.679	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.56	45	251347	32.788	ng	99
17) 2-Methylphenol	8.49	107	124011	35.936	ng	96
18) Hexachloroethane	9.12	117	52491	32.751	ng	98
19) n-Nitroso-di-n-propylamine	8.85	70	104327	31.670	ng	97
20) 3+4-Methylphenols	8.81	107	170420	36.454	ng	99
22) Acetophenone	8.88	105	207118	37.505	ng	# 98
24) Nitrobenzene	9.26	77	157837	36.596	ng	96
25) Isophorone	9.77	82	300605	35.078	ng	98
26) 2-Nitrophenol	9.97	139	74340	36.376	ng	97
27) 2,4-Dimethylphenol	10.02	122	138271	42.595	ng	98
28) bis(2-Chloroethoxy)methane	10.26	93	182338	35.964	ng	100
29) 2,4-Dichlorophenol	10.51	162	136963	38.871	ng	99
30) 1,2,4-Trichlorobenzene	10.72	180	141342	36.532	ng	95
31) Naphthalene	10.91	128	425550	37.502	ng	100
33) 4-Chloroaniline	11.03	127	122217	25.250	ng	97
34) Hexachlorobutadiene	11.17	225	85515	34.560	ng	99
35) Caprolactam	11.80	113	47930m	34.278	ng	
36) 4-Chloro-3-methylphenol	12.14	107	157521	38.887	ng	97
37) 2-Methylnaphthalene	12.51	142	315732	37.621	ng	99
38) 1-Methylnaphthalene	12.73	142	302917	37.222	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.87	216	162280	37.874	ng	98
41) Hexachlorocyclopentadiene	12.84	237	191097	81.226	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,6-Trichlorophenol	13.12	196	112376	40.679	ng	97
44) 2,4,5-Trichlorophenol	13.19	196	122094	40.548	ng	96
46) 1,1'-Biphenyl	13.51	154	401937	37.734	ng	98
47) 2-Chloronaphthalene	13.56	162	312318	38.225	ng	97
48) 2-Nitroaniline	13.78	65	108489	39.364	ng	95
49) Acenaphthylene	14.40	152	525936	37.965	ng	99
50) Dimethylphthalate	14.13	163	495141	46.929	ng	99
51) 2,6-Dinitrotoluene	14.26	165	94878	40.049	ng	96
52) Acenaphthene	14.75	154	302259	38.077	ng	97
53) 3-Nitroaniline	14.60	138	76762	30.610	ng	97
54) 2,4-Dinitrophenol	14.80	184	79500	63.100	ng	87
55) Dibenzofuran	15.07	168	504079	39.519	ng	99
56) 4-Nitrophenol	14.90	139	167659	82.838	ng	98
57) 2,4-Dinitrotoluene	15.05	165	137982	41.917	ng	97
58) Fluorene	15.73	166	403530	40.239	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.30	232	123207	45.091	ng	97
60) Diethylphthalate	15.48	149	439616	40.718	ng	99
61) 4-Chlorophenyl-phenylether	15.71	204	215543	40.210	ng	100
62) 4-Nitroaniline	15.76	138	111045	41.262	ng	98
63) Azobenzene	16.00	77	406435	39.909	ng	99
65) 4,6-Dinitro-2-methylphenol	15.80	198	73555	35.548	ng	98
66) n-Nitrosodiphenylamine	15.93	169	382294	35.471	ng	97
67) 4-Bromophenyl-phenylether	16.61	248	142617	34.410	ng	95
68) Hexachlorobenzene	16.72	284	149363	35.842	ng	97
69) Atrazine	16.87	200	151717	37.843	ng	97
70) Pentachlorophenol	17.07	266	171324	71.110	ng	99
71) Phenanthrene	17.47	178	721955	37.294	ng	99
72) Anthracene	17.56	178	742416	38.315	ng	99
73) Carbazole	17.84	167	723623	39.461	ng	99
74) Di-n-butylphthalate	18.36	149	856138	39.387	ng	99
75) Fluoranthene	19.48	202	934170	40.752	ng	99
77) Benzidine	19.66	184	312542	21.810	ng	97
78) Pyrene	19.84	202	936404	35.883	ng	99
80) Butylbenzylphthalate	20.70	149	417433	37.382	ng	99
81) Benzo(a)anthracene	21.68	228	875287	35.810	ng	99
82) 3,3'-Dichlorobenzidine	21.60	252	215167	23.141	ng	99
83) Chrysene	21.75	228	861819	36.688	ng	100
84) Bis(2-ethylhexyl)phthalate	21.55	149	610458	38.243	ng	100
85) Di-n-octyl phthalate	22.77	149	1040823	38.271	ng	99
86) Indeno(1,2,3-cd)pyrene	28.74	276	1066837	37.132	ng	# 91
88) Benzo(b)fluoranthene	23.93	252	938425	36.143	ng	100
89) Benzo(k)fluoranthene	24.00	252	919761	38.521	ng	98
90) Benzo(a)pyrene	24.83	252	926343	38.025	ng	98
91) Dibenzo(a,h)anthracene	28.81	278	874385	38.646	ng	98
92) Benzo(g,h,i)perylene	29.93	276	889356	38.248	ng	98

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

