

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG030719\
 Data File : BG039811.D
 Acq On : 7 Mar 2019 20:45
 Operator : JU/SJ
 Sample : K1757-04MSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-6CD-SMSD

Manual Integrations
 APPROVED

Sohil
 3/8/2019 10:48:50 AM

Quant Time: Mar 08 01:43:25 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG030419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 04 14:38:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	45088	20.00	ng	-0.01
21) Naphthalene-d8	10.98	136	199298	20.00	ng	-0.01
39) Acenaphthene-d10	14.78	164	142512	20.00	ng	0.00
64) Phenanthrene-d10	17.53	188	352060	20.00	ng	0.00
76) Chrysene-d12	21.81	240	347572	20.00	ng	-0.01
87) Perylene-d12	25.19	264	357024	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.70	112	362014	136.98	ng	0.00
7) Phenol-d6	7.30	99	500887	127.11	ng	-0.01
23) Nitrobenzene-d5	9.33	82	343315	93.17	ng	-0.01
42) 2,4,6-Tribromophenol	16.27	330	247339	148.90	ng	0.00
45) 2-Fluorobiphenyl	13.40	172	838722	83.80	ng	-0.01
79) Terphenyl-d14	20.12	244	1376034	87.17	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.60	88	37628	30.839	ng	# 36
3) Pyridine	4.00	79	97101	29.726	ng	93
4) n-Nitrosodimethylamine	3.92	42	55946	36.103	ng	92
6) Aniline	7.48	93	39510	7.653	ng	99
8) 2-Chlorophenol	7.72	128	131382	40.976	ng	95
9) Benzaldehyde	7.29	77	79512	31.279	ng	98
10) Phenol	7.33	94	155676	36.466	ng	97
11) bis(2-Chloroethyl)ether	7.57	93	126250	37.931	ng	98
12) 1,3-Dichlorobenzene	8.04	146	126502	34.885	ng	99
13) 1,4-Dichlorobenzene	8.19	146	129439	35.043	ng	98
14) 1,2-Dichlorobenzene	8.51	146	125677	35.216	ng	99
15) Benzyl Alcohol	8.39	79	127846	40.898	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.67	45	247873	37.235	ng	99
17) 2-Methylphenol	8.59	107	116039	42.628	ng	99
18) Hexachloroethane	9.24	117	45026	33.973	ng	96
19) n-Nitroso-di-n-propylamine	8.96	70	108190	38.143	ng	93
20) 3+4-Methylphenols	8.91	107	158411	41.890	ng	94
22) Acetophenone	8.99	105	204972	38.319	ng	# 95
24) Nitrobenzene	9.38	77	157912	40.849	ng	96
25) Isophorone	9.89	82	311360	40.325	ng	96
26) 2-Nitrophenol	10.09	139	78962	42.305	ng	97
27) 2,4-Dimethylphenol	10.12	122	143944	49.587	ng	99
28) bis(2-Chloroethoxy)methane	10.37	93	189723	40.434	ng	99
29) 2,4-Dichlorophenol	10.61	162	150933	46.180	ng	97
30) 1,2,4-Trichlorobenzene	10.83	180	140874	38.305	ng	99
31) Naphthalene	11.03	128	415827	39.407	ng	99
32) Benzoic acid	10.21	122	13908	12.617	ng	91
33) 4-Chloroaniline	11.14	127	7522	1.681	ng	# 83
34) Hexachlorobutadiene	11.29	225	90654	36.072	ng	90
35) Caprolactam	11.92	113	41332m	33.106	ng	
36) 4-Chloro-3-methylphenol	12.24	107	169972	45.368	ng	99
37) 2-Methylnaphthalene	12.62	142	317717	40.274	ng	97
38) 1-Methylnaphthalene	12.84	142	312030	40.700	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.98	216	181539	37.602	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.94	237	192518	74.408	nq	98
43) 2,4,6-Trichlorophenol	13.21	196	132868	47.007	nq	97
44) 2,4,5-Trichlorophenol	13.29	196	144402	45.324	nq	97
46) 1,1'-Biphenyl	13.61	154	453363	38.678	nq	99
47) 2-Chloronaphthalene	13.67	162	348842	39.560	nq	97
48) 2-Nitroaniline	13.87	65	124709	44.008	nq	95
49) Acenaphthylene	14.51	152	570193	39.390	nq	99
50) Dimethylphthalate	14.23	163	488673	41.007	nq	99
51) 2,6-Dinitrotoluene	14.36	165	109574	43.374	nq	93
52) Acenaphthene	14.85	154	345703	39.679	nq	99
53) 3-Nitroaniline	14.70	138	8865	3.491	nq #	93
54) 2,4-Dinitrophenol	14.89	184	34599	25.546	nq	94
55) Dibenzofuran	15.18	168	570356	39.900	nq	97
56) 4-Nitrophenol	14.99	139	159255	70.103	nq #	86
57) 2,4-Dinitrotoluene	15.14	165	156135	43.985	nq	87
58) Fluorene	15.82	166	455166	40.505	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.40	232	142026	45.645	nq	97
60) Diethylphthalate	15.58	149	496188	42.062	nq	98
61) 4-Chlorophenyl-phenylether	15.81	204	242043	40.364	nq	96
62) 4-Nitroaniline	15.86	138	101427	36.842	nq	96
63) Azobenzene	16.10	77	440378	41.182	nq	97
65) 4,6-Dinitro-2-methylphenol	15.90	198	40984	19.689	nq	95
66) n-Nitrosodiphenylamine	16.03	169	425412	41.739	nq	99
67) 4-Bromophenyl-phenylether	16.70	248	160678	40.774	nq	96
68) Hexachlorobenzene	16.82	284	167689	41.052	nq	92
69) Atrazine	16.97	200	170907	42.607	nq	97
70) Pentachlorophenol	17.17	266	126694	49.110	nq	98
71) Phenanthrene	17.57	178	778024	40.121	nq	98
72) Anthracene	17.66	178	785673	41.576	nq	99
73) Carbazole	17.93	167	696478	40.883	nq	98
74) Di-n-butylphthalate	18.46	149	849284	42.371	nq	99
75) Fluoranthene	19.57	202	922832	40.267	nq	99
77) Benzidine	19.75	184	98538	8.934	nq	97
78) Pyrene	19.93	202	919648	40.719	nq	98
80) Butylbenzylphthalate	20.80	149	390675	44.648	nq	95
81) Benzo(a)anthracene	21.79	228	874213	40.132	nq	99
82) 3,3'-Dichlorobenzidine	21.70	252	58291	7.329	nq	98
83) Chrysene	21.86	228	841406	39.842	nq	99
84) Bis(2-ethylhexyl)phthalate	21.66	149	539488	43.875	nq	99
85) Di-n-octyl phthalate	22.91	149	919928	45.184	nq	96
86) Indeno(1,2,3-cd)pyrene	29.06	276	1010434	42.176	nq	100
88) Benzo(b)fluoranthene	24.11	252	878714	41.273	nq	98
89) Benzo(k)fluoranthene	24.18	252	838079	42.289	nq	99
90) Benzo(a)pyrene	25.03	252	859293	43.678	nq	98
91) Dibenzo(a,h)anthracene	29.13	278	815847	42.301	nq	99
92) Benzo(g,h,i)perylene	30.29	276	833965	43.054	nq	98

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

