

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022619\
 Data File : BG039599.D
 Acq On : 26 Feb 2019 14:04
 Operator : JU/SJ
 Sample : PB117390BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB117390BS

Manual Integrations
 APPROVED

Sohil
 2/27/2019 4:21:42 PM

Quant Time: Feb 27 02:55:01 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	52194	20.00	ng	-0.02
21) Naphthalene-d8	10.99	136	219256	20.00	ng	-0.02
39) Acenaphthene-d10	14.79	164	147079	20.00	ng	-0.02
64) Phenanthrene-d10	17.54	188	374143	20.00	ng	-0.02
76) Chrysene-d12	21.83	240	389165	20.00	ng	-0.02
87) Perylene-d12	25.20	264	364219	20.00	ng	-0.04

System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	481383	157.85	ng	0.00
7) Phenol-d6	7.32	99	669019	149.04	ng	0.00
23) Nitrobenzene-d5	9.35	82	329331	93.83	ng	-0.02
42) 2,4,6-Tribromophenol	16.28	330	265830	167.84	ng	-0.01
45) 2-Fluorobiphenyl	13.42	172	790938	77.14	ng	-0.02
79) Terphenyl-d14	20.13	244	1461419	79.49	ng	-0.02

Target Compounds					Qvalue
2) 1,4-Dioxane	3.60	88	52156	39.098 ng	94
3) Pyridine	4.00	79	141124	39.958 ng	94
4) n-Nitrosodimethylamine	3.92	42	71800	40.650 ng	91
6) Aniline	7.49	93	165020	29.349 ng	96
8) 2-Chlorophenol	7.73	128	149677	44.901 ng	97
9) Benzaldehyde	7.31	77	51096	19.361 ng	90
10) Phenol	7.34	94	205280	43.869 ng	97
11) bis(2-Chloroethyl)ether	7.58	93	145587	39.374 ng	99
12) 1,3-Dichlorobenzene	8.05	146	161294	39.942 ng	97
13) 1,4-Dichlorobenzene	8.20	146	162283	40.319 ng	97
14) 1,2-Dichlorobenzene	8.52	146	157193	40.342 ng	99
15) Benzyl Alcohol	8.41	79	143150	40.619 ng	97
16) 2,2'-oxybis(1-Chloropropan	8.68	45	279507	33.474 ng	97
17) 2-Methylphenol	8.60	107	132359	43.710 ng	97
18) Hexachloroethane	9.25	117	57065	41.209 ng	97
19) n-Nitroso-di-n-propylamine	8.97	70	117280	36.810 ng	90
20) 3+4-Methylphenols	8.93	107	184159	44.164 ng	97
22) Acetophenone	8.99	105	223829	38.004 ng	93
24) Nitrobenzene	9.39	77	172531	45.503 ng	96
25) Isophorone	9.90	82	327016	42.047 ng	100
26) 2-Nitrophenol	10.10	139	85223	54.834 ng	97
27) 2,4-Dimethylphenol	10.14	122	156879	49.863 ng	97
28) bis(2-Chloroethoxy)methane	10.38	93	204036	42.592 ng	96
29) 2,4-Dichlorophenol	10.63	162	166148	48.319 ng	93
30) 1,2,4-Trichlorobenzene	10.84	180	163577	42.372 ng	99
31) Naphthalene	11.04	128	461688	42.446 ng	99
32) Benzoic acid	10.27	122	103930	49.987 ng	98
33) 4-Chloroaniline	11.15	127	116350	23.070 ng	96
34) Hexachlorobutadiene	11.30	225	104351	40.646 ng	94
35) Caprolactam	11.93	113	55603m	39.077 ng	
36) 4-Chloro-3-methylphenol	12.25	107	176677	44.428 ng	98
37) 2-Methylnaphthalene	12.63	142	352559	43.762 ng	97
38) 1-Methylnaphthalene	12.85	142	334481	43.071 ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.99	216	192364	41.107 ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.96	237	253579	99.443	nq	99
43) 2,4,6-Trichlorophenol	13.22	196	135149	46.971	nq	99
44) 2,4,5-Trichlorophenol	13.30	196	146317	48.520	nq	98
46) 1,1'-Biphenyl	13.63	154	469980	38.405	nq	98
47) 2-Chloronaphthalene	13.68	162	363774	39.515	nq	98
48) 2-Nitroaniline	13.88	65	125326	47.496	nq	92
49) Acenaphthylene	14.52	152	589822	42.461	nq	99
50) Dimethylphthalate	14.24	163	491954	41.415	nq	99
51) 2,6-Dinitrotoluene	14.37	165	111120	49.772	nq	94
52) Acenaphthene	14.86	154	358169	39.722	nq	99
53) 3-Nitroaniline	14.70	138	75327	33.443	nq	# 96
54) 2,4-Dinitrophenol	14.90	184	132783	112.711	nq	99
55) Dibenzofuran	15.19	168	585378	40.491	nq	99
56) 4-Nitrophenol	15.00	139	205305	94.461	nq	92
57) 2,4-Dinitrotoluene	15.15	165	157315	54.126	nq	88
58) Fluorene	15.84	166	465099	43.156	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.41	232	145856	51.657	nq	96
60) Diethylphthalate	15.59	149	505906	41.192	nq	97
61) 4-Chlorophenyl-phenylether	15.82	204	249901	40.952	nq	96
62) 4-Nitroaniline	15.87	138	120774	48.443	nq	96
63) Azobenzene	16.11	77	444362	37.017	nq	97
65) 4,6-Dinitro-2-methylphenol	15.91	198	96433	59.667	nq	93
66) n-Nitrosodiphenylamine	16.04	169	436449	38.364	nq	99
67) 4-Bromophenyl-phenylether	16.72	248	162999	39.270	nq	92
68) Hexachlorobenzene	16.83	284	169283	40.650	nq	96
69) Atrazine	16.98	200	180801	43.489	nq	95
70) Pentachlorophenol	17.18	266	225522	77.919	nq	98
71) Phenanthrene	17.58	178	803711	41.504	nq	99
72) Anthracene	17.67	178	823919	43.196	nq	98
73) Carbazole	17.94	167	743943	39.082	nq	99
74) Di-n-butylphthalate	18.47	149	908949	40.930	nq	99
75) Fluoranthene	19.58	202	1000052	44.494	nq	98
77) Benzidine	19.76	184	342900	26.875	nq	99
78) Pyrene	19.94	202	1003670	41.428	nq	99
80) Butylbenzylphthalate	20.81	149	427369	41.818	nq	94
81) Benzo(a)anthracene	21.81	228	983423	42.766	nq	99
82) 3,3'-Dichlorobenzidine	21.72	252	247439	29.020	nq	98
83) Chrysene	21.88	228	912513	41.686	nq	99
84) Bis(2-ethylhexyl)phthalate	21.67	149	592555	40.642	nq	# 99
85) Di-n-octyl phthalate	22.93	149	993233	41.235	nq	96
86) Indeno(1,2,3-cd)pyrene	29.09	276	1081540	46.050	nq	100
88) Benzo(b)fluoranthene	24.11	252	952876	47.973	nq	99
89) Benzo(k)fluoranthene	24.19	252	907777	45.521	nq	99
90) Benzo(a)pyrene	25.04	252	926374	48.427	nq	99
91) Dibenzo(a,h)anthracene	29.16	278	882412	49.256	nq	97
92) Benzo(g,h,i)perylene	30.31	276	893097	50.016	nq	100

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

