

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022719\
 Data File : BG039639.D
 Acq On : 27 Feb 2019 17:57
 Operator : JU/SJ
 Sample : PB117428BSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB117428BSD

Manual Integrations
 APPROVED

Sohil
 2/28/2019 1:32:06 PM

Quant Time: Feb 28 01:21:33 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	74860	20.00	ng	-0.02
21) Naphthalene-d8	10.99	136	331466	20.00	ng	-0.02
39) Acenaphthene-d10	14.79	164	222585	20.00	ng	-0.02
64) Phenanthrene-d10	17.54	188	529171	20.00	ng	-0.02
76) Chrysene-d12	21.82	240	505909	20.00	ng	-0.02
87) Perylene-d12	25.20	264	498430	20.00	ng	-0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.71	112	607078	138.79	ng	0.00
7) Phenol-d6	7.32	99	909455	141.26	ng	0.00
23) Nitrobenzene-d5	9.35	82	446379	84.13	ng	-0.02
42) 2,4,6-Tribromophenol	16.29	330	392046	163.57	ng	0.00
45) 2-Fluorobiphenyl	13.42	172	1054168	67.94	ng	-0.02
79) Terphenyl-d14	20.13	244	1679328	70.26	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.61	88	75525	39.475	ng	93
3) Pyridine	4.01	79	226663	44.746	ng	93
4) n-Nitrosodimethylamine	3.93	42	112183	44.283	ng	89
6) Aniline	7.50	93	354853	44.002	ng	98
8) 2-Chlorophenol	7.73	128	209049	43.724	ng	99
9) Benzaldehyde	7.31	77	126378	39.033	ng	97
10) Phenol	7.34	94	289572	43.146	ng	98
11) bis(2-Chloroethyl)ether	7.59	93	222284	41.915	ng	98
12) 1,3-Dichlorobenzene	8.06	146	240080	41.451	ng	98
13) 1,4-Dichlorobenzene	8.20	146	243332	42.150	ng	97
14) 1,2-Dichlorobenzene	8.53	146	235877	42.207	ng	97
15) Benzyl Alcohol	8.41	79	219152	43.357	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.68	45	425692	35.545	ng	99
17) 2-Methylphenol	8.60	107	194188	44.711	ng	95
18) Hexachloroethane	9.25	117	87369	43.990	ng	97
19) n-Nitroso-di-n-propylamine	8.97	70	186092	40.724	ng	92
20) 3+4-Methylphenols	8.93	107	268131	44.832	ng	98
22) Acetophenone	9.00	105	354813	39.850	ng	# 93
24) Nitrobenzene	9.39	77	273129	47.649	ng	98
25) Isophorone	9.91	82	535769	45.567	ng	98
26) 2-Nitrophenol	10.10	139	138015	57.671	ng	98
27) 2,4-Dimethylphenol	10.14	122	220496	46.359	ng	99
28) bis(2-Chloroethoxy)methane	10.38	93	309737	42.769	ng	96
29) 2,4-Dichlorophenol	10.63	162	243040	46.754	ng	94
30) 1,2,4-Trichlorobenzene	10.85	180	261135	44.744	ng	99
31) Naphthalene	11.04	128	721865	43.899	ng	98
32) Benzoic acid	10.29	122	155167	49.478	ng	97
33) 4-Chloroaniline	11.15	127	325364	42.674	ng	98
34) Hexachlorobutadiene	11.30	225	164706	42.437	ng	93
35) Caprolactam	11.94	113	87490m	40.672	ng	
36) 4-Chloro-3-methylphenol	12.25	107	262897	43.730	ng	97
37) 2-Methylnaphthalene	12.63	142	547720	44.972	ng	98
38) 1-Methylnaphthalene	12.85	142	524921	44.712	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.99	216	305234	43.100	ng	99

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41) Hexachlorocyclopentadiene	12.96	237	372663	96.567	nq	97
43) 2,4,6-Trichlorophenol	13.22	196	202810	46.576	nq	99
44) 2,4,5-Trichlorophenol	13.30	196	221885	48.619	nq	99
46) 1,1'-Biphenyl	13.63	154	721082	38.936	nq	98
47) 2-Chloronaphthalene	13.68	162	566358	40.652	nq	99
48) 2-Nitroaniline	13.88	65	194962	48.591	nq	94
49) Acenaphthylene	14.52	152	884537	42.076	nq	98
50) Dimethylphthalate	14.24	163	737688	41.035	nq	98
51) 2,6-Dinitrotoluene	14.37	165	169710	50.169	nq	95
52) Acenaphthene	14.86	154	547203	40.100	nq	98
53) 3-Nitroaniline	14.71	138	172049	47.524	nq	# 92
54) 2,4-Dinitrophenol	14.91	184	198599	111.869	nq	99
55) Dibenzofuran	15.19	168	892197	40.779	nq	99
56) 4-Nitrophenol	15.00	139	284522	86.990	nq	# 87
57) 2,4-Dinitrotoluene	15.16	165	243933	55.265	nq	# 84
58) Fluorene	15.84	166	696672	42.715	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.41	232	216543	50.676	nq	97
60) Diethylphthalate	15.59	149	752392	40.480	nq	97
61) 4-Chlorophenyl-phenylether	15.82	204	382141	41.379	nq	98
62) 4-Nitroaniline	15.87	138	185051	48.994	nq	93
63) Azobenzene	16.12	77	669423	36.849	nq	96
65) 4,6-Dinitro-2-methylphenol	15.92	198	144925	62.233	nq	93
66) n-Nitrosodiphenylamine	16.04	169	646614	40.187	nq	99
67) 4-Bromophenyl-phenylether	16.72	248	253351	43.156	nq	94
68) Hexachlorobenzene	16.83	284	261065	44.324	nq	95
69) Atrazine	16.98	200	140590	23.910	nq	96
70) Pentachlorophenol	17.18	266	321686	78.582	nq	99
71) Phenanthrene	17.58	178	1141678	41.685	nq	99
72) Anthracene	17.67	178	1153049	42.741	nq	97
73) Carbazole	17.94	167	1009614	37.500	nq	98
74) Di-n-butylphthalate	18.47	149	1210311	38.533	nq	98
75) Fluoranthene	19.58	202	1324633	41.669	nq	98
77) Benzidine	19.76	184	919479	55.435	nq	98
78) Pyrene	19.94	202	1312280	41.667	nq	98
80) Butylbenzylphthalate	20.81	149	568250	42.772	nq	92
81) Benzo(a)anthracene	21.81	228	1290454	43.168	nq	97
82) 3,3'-Dichlorobenzidine	21.72	252	475451	42.894	nq	98
83) Chrysene	21.88	228	1203952	42.308	nq	99
84) Bis(2-ethylhexyl)phthalate	21.67	149	788917	41.624	nq	98
85) Di-n-octyl phthalate	22.93	149	1336957	42.696	nq	96
86) Indeno(1,2,3-cd)pyrene	29.09	276	1471739	48.203	nq	# 95
88) Benzo(b)fluoranthene	24.12	252	1266799	46.604	nq	99
89) Benzo(k)fluoranthene	24.20	252	1233031	45.182	nq	99
90) Benzo(a)pyrene	25.05	252	1261011	48.170	nq	98
91) Dibenzo(a,h)anthracene	29.16	278	1185905	48.373	nq	99
92) Benzo(g,h,i)perylene	30.33	276	1211871	49.594	nq	99

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

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