

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121618\
 Data File : BG038637.D
 Acq On : 16 Dec 2018 12:46
 Operator : JU/SJ
 Sample : PB115713BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115713BS

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:14:01 PM

Quant Time: Dec 17 05:58:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 15:26:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	22007	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	86550	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	54214	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	136892	20.00	ng	0.00
76) Chrysene-d12	21.73	240	147204	20.00	ng	0.00
87) Perylene-d12	25.02	264	158295	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	175955	141.81	ng	0.00
7) Phenol-d6	7.21	99	230814	135.51	ng	0.00
23) Nitrobenzene-d5	9.24	82	121723	71.85	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	102975	137.82	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	274045	69.35	ng	0.00
79) Terphenyl-d14	20.05	244	559669	82.74	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	19984	34.606	ng	92
3) Pyridine	3.89	79	55620	34.983	ng	91
4) n-Nitrosodimethylamine	3.82	42	37865	48.606	ng	# 86
6) Aniline	7.38	93	69303	31.872	ng	97
8) 2-Chlorophenol	7.62	128	62230	43.340	ng	98
9) Benzaldehyde	7.20	77	21253	19.234	ng	95
10) Phenol	7.24	94	77674	42.923	ng	95
11) bis(2-Chloroethyl)ether	7.48	93	55169	38.292	ng	96
12) 1,3-Dichlorobenzene	7.94	146	66362	39.089	ng	99
13) 1,4-Dichlorobenzene	8.10	146	66396	38.186	ng	98
14) 1,2-Dichlorobenzene	8.41	146	63185	38.546	ng	96
15) Benzyl Alcohol	8.30	79	56014	42.131	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.58	45	123760	37.498	ng	98
17) 2-Methylphenol	8.50	107	53326	43.983	ng	97
18) Hexachloroethane	9.14	117	23890	37.600	ng	91
19) n-Nitroso-di-n-propylamine	8.86	70	44977	37.596	ng	98
20) 3+4-Methylphenols	8.82	107	72201	43.120	ng	97
22) Acetophenone	8.89	105	89550	41.423	ng	98
24) Nitrobenzene	9.28	77	70476	41.121	ng	99
25) Isophorone	9.79	82	126269	41.482	ng	# 98
26) 2-Nitrophenol	9.99	139	35394	40.349	ng	96
27) 2,4-Dimethylphenol	10.03	122	60672	48.261	ng	98
28) bis(2-Chloroethoxy)methane	10.28	93	71065	41.370	ng	99
29) 2,4-Dichlorophenol	10.52	162	63389	45.324	ng	99
30) 1,2,4-Trichlorobenzene	10.73	180	68029	40.270	ng	100
31) Naphthalene	10.93	128	185880	43.589	ng	98
32) Benzoic acid	10.19	122	24431m	34.281	ng	
33) 4-Chloroaniline	11.05	127	41397	22.360	ng	98
34) Hexachlorobutadiene	11.19	225	45718	39.995	ng	95
35) Caprolactam	11.83	113	20500m	35.419	ng	
36) 4-Chloro-3-methylphenol	12.15	107	66276	41.526	ng	95
37) 2-Methylnaphthalene	12.53	142	136151	44.753	ng	95
38) 1-Methylnaphthalene	12.75	142	128967	43.686	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.89	216	80087	39.520	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.85	237	109752	98.578	nq	93
43) 2,4,6-Trichlorophenol	13.13	196	53410	42.864	nq	99
44) 2,4,5-Trichlorophenol	13.20	196	62073	47.052	nq	97
46) 1,1'-Biphenyl	13.53	154	177406	39.034	nq	97
47) 2-Chloronaphthalene	13.58	162	145542	41.457	nq	97
48) 2-Nitroaniline	13.79	65	48390	43.273	nq	98
49) Acenaphthylene	14.42	152	230157	43.853	nq	98
50) Dimethylphthalate	14.15	163	186179	42.052	nq	99
51) 2,6-Dinitrotoluene	14.28	165	41704	41.567	nq	97
52) Acenaphthene	14.76	154	128461	40.933	nq	98
53) 3-Nitroaniline	14.61	138	30433	29.756	nq	97
54) 2,4-Dinitrophenol	14.82	184	47254	79.211	nq	98
55) Dibenzofuran	15.09	168	219051	40.719	nq	99
56) 4-Nitrophenol	14.91	139	67484	86.977	nq	98
57) 2,4-Dinitrotoluene	15.06	165	59932	42.411	nq	98
58) Fluorene	15.74	166	174993	43.906	nq	97
59) 2,3,4,6-Tetrachlorophenol	15.32	232	54270	45.724	nq	99
60) Diethylphthalate	15.50	149	192322	39.571	nq	99
61) 4-Chlorophenyl-phenylether	15.73	204	99190	39.887	nq	99
62) 4-Nitroaniline	15.77	138	44831	42.577	nq	96
63) Azobenzene	16.02	77	167481	41.250	nq	99
65) 4,6-Dinitro-2-methylphenol	15.82	198	37285	38.181	nq	93
66) n-Nitrosodiphenylamine	15.95	169	161632	40.185	nq	99
67) 4-Bromophenyl-phenylether	16.63	248	65122	38.952	nq	98
68) Hexachlorobenzene	16.74	284	67454	38.922	nq	96
69) Atrazine	16.89	200	68762	46.066	nq	99
70) Pentachlorophenol	17.09	266	77163	75.275	nq	94
71) Phenanthrene	17.49	178	313386	44.147	nq	99
72) Anthracene	17.58	178	321366	45.857	nq	99
73) Carbazole	17.85	167	290258	41.880	nq	98
74) Di-n-butylphthalate	18.38	149	353716	44.478	nq	99
75) Fluoranthene	19.49	202	405208	46.135	nq	97
77) Benzidine	19.68	184	194892	44.768	nq	98
78) Pyrene	19.86	202	409232	42.082	nq	98
80) Butylbenzylphthalate	20.72	149	160388	42.215	nq	97
81) Benzo(a)anthracene	21.70	228	399862	44.578	nq	99
82) 3,3'-Dichlorobenzidine	21.62	252	105091	29.541	nq	95
83) Chrysene	21.77	228	369514	42.769	nq	99
84) Bis(2-ethylhexyl)phthalate	21.57	149	227107	42.147	nq	100
85) Di-n-octyl phthalate	22.80	149	390309	43.251	nq	99
86) Indeno(1,2,3-cd)pyrene	28.80	276	456178	44.911	nq	# 97
88) Benzo(b)fluoranthene	23.97	252	397612	45.750	nq	99
89) Benzo(k)fluoranthene	24.04	252	386397	44.647	nq	97
90) Benzo(a)pyrene	24.86	252	386948	46.150	nq	98
91) Dibenzo(a,h)anthracene	28.87	278	375759	45.010	nq	98
92) Benzo(g,h,i)perylene	29.99	276	373618	45.412	nq	97

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

