

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121218\
 Data File : BG038501.D
 Acq On : 12 Dec 2018 13:38
 Operator : JU/SJ
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Sohil
 12/13/2018 3:06:49 PM

Quant Time: Dec 12 15:23:28 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 12:12:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	28078	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	119671	20.00	ng	0.00
39) Acenaphthene-d10	14.70	164	76776	20.00	ng	0.00
64) Phenanthrene-d10	17.45	188	181402	20.00	ng	0.00
76) Chrysene-d12	21.73	240	177865	20.00	ng	0.00
87) Perylene-d12	25.02	264	197435	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	257053	162.01	ng	0.00
7) Phenol-d6	7.22	99	357368	155.84	ng	0.00
23) Nitrobenzene-d5	9.24	82	370670	153.78	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	170376	180.77	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	843748	156.59	ng	0.00
79) Terphenyl-d14	20.05	244	1218264	146.73	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.50	88	57848	79.550	ng	98
3) Pyridine	3.90	79	163534m	75.170	ng	
4) n-Nitrosodimethylamine	3.82	42	84623	88.801	ng	# 93
6) Aniline	7.39	93	223262	76.225	ng	98
8) 2-Chlorophenol	7.62	128	146133	79.303	ng	98
10) Phenol	7.25	94	189548	78.230	ng	98
11) bis(2-Chloroethyl)ether	7.49	93	145767	73.457	ng	99
12) 1,3-Dichlorobenzene	7.95	146	173461	80.542	ng	100
13) 1,4-Dichlorobenzene	8.10	146	176745	79.317	ng	98
14) 1,2-Dichlorobenzene	8.41	146	164738	74.821	ng	97
15) Benzyl Alcohol	8.31	79	146831	77.880	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.59	45	329179	72.930	ng	99
17) 2-Methylphenol	8.50	107	126313	68.682	ng	96
18) Hexachloroethane	9.14	117	65419	81.341	ng	93
19) n-Nitroso-di-n-propylamine	8.87	70	125500	71.098	ng	97
20) 3+4-Methylphenols	8.84	107	181235	74.246	ng	97
22) Acetophenone	8.90	105	233995	74.699	ng	# 99
24) Nitrobenzene	9.29	77	185101	75.585	ng	99
25) Isophorone	9.80	82	403539	95.265	ng	99
26) 2-Nitrophenol	10.00	139	104566	92.153	ng	94
27) 2,4-Dimethylphenol	10.04	122	139642	85.681	ng	99
28) bis(2-Chloroethoxy)methane	10.28	93	191873	79.014	ng	99
29) 2,4-Dichlorophenol	10.52	162	163272	87.760	ng	99
30) 1,2,4-Trichlorobenzene	10.74	180	185358	84.873	ng	99
31) Naphthalene	10.93	128	458510	80.210	ng	98
32) Benzoic acid	10.24	122	101636m	95.749	ng	
33) 4-Chloroaniline	11.05	127	208472	82.381	ng	98
34) Hexachlorobutadiene	11.19	225	128588	94.115	ng	94
35) Caprolactam	11.86	113	60453	78.576	ng	96
36) 4-Chloro-3-methylphenol	12.16	107	175340	84.825	ng	99
37) 2-Methylnaphthalene	12.53	142	333611	81.950	ng	98
38) 1-Methylnaphthalene	12.75	142	318779	81.759	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.89	216	223501	84.600	ng	99
41) Hexachlorocyclopentadiene	12.86	237	133804	92.164	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,6-Trichlorophenol	13.13	196	145417	84.971	nq	98
44) 2,4,5-Trichlorophenol	13.21	196	155190	85.599	nq	98
46) 1,1'-Biphenyl	13.53	154	496462	76.499	nq	98
47) 2-Chloronaphthalene	13.59	162	385827	78.786	nq	97
48) 2-Nitroaniline	13.80	65	129043	78.543	nq	99
49) Acenaphthylene	14.43	152	569801	77.256	nq	97
50) Dimethylphthalate	14.16	163	472790	76.775	nq	100
51) 2,6-Dinitrotoluene	14.28	165	110555	78.592	nq	98
52) Acenaphthene	14.77	154	344115	76.533	nq	99
53) 3-Nitroaniline	14.63	138	115895	76.499	nq	95
54) 2,4-Dinitrophenol	14.82	184	67266	87.232	nq	96
55) Dibenzofuran	15.10	168	575915	77.477	nq	99
56) 4-Nitrophenol	14.92	139	87815	73.377	nq	95
57) 2,4-Dinitrotoluene	15.07	165	152473	78.588	nq	98
58) Fluorene	15.75	166	424534	77.520	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.32	232	131548	83.543	nq	98
60) Diethylphthalate	15.51	149	493438	72.326	nq	99
61) 4-Chlorophenyl-phenylether	15.74	204	267938	81.454	nq	98
62) 4-Nitroaniline	15.79	138	115770	77.693	nq	96
63) Azobenzene	16.03	77	435561	72.948	nq	96
65) 4,6-Dinitro-2-methylphenol	15.83	198	107202	90.195	nq	96
66) n-Nitrosodiphenylamine	15.95	169	422721	82.869	nq	97
67) 4-Bromophenyl-phenylether	16.63	248	176523	90.111	nq	95
68) Hexachlorobenzene	16.74	284	183564	90.847	nq	99
69) Atrazine	16.89	200	135443	71.157	nq	97
70) Pentachlorophenol	17.09	266	118637	85.722	nq	99
71) Phenanthrene	17.49	178	715546	78.765	nq	98
72) Anthracene	17.58	178	707582	80.430	nq	100
73) Carbazole	17.86	167	693081	79.377	nq	99
74) Di-n-butylphthalate	18.39	149	791535	77.597	nq	100
75) Fluoranthene	19.50	202	859457	80.975	nq	99
77) Benzidine	19.68	184	378261	63.023	nq	# 96
78) Pyrene	19.86	202	868409	69.781	nq	98
80) Butylbenzylphthalate	20.73	149	373747	74.669	nq	99
81) Benzo(a)anthracene	21.71	228	846910	77.473	nq	97
82) 3,3'-Dichlorobenzidine	21.63	252	334484	78.655	nq	97
83) Chrysene	21.78	228	812644	78.556	nq	98
84) Bis(2-ethylhexyl)phthalate	21.58	149	527607	74.368	nq	97
85) Di-n-octyl phthalate	22.80	149	886582	72.881	nq	99
86) Indeno(1,2,3-cd)pyrene	28.83	276	1014429	86.285	nq	# 90
88) Benzo(b)fluoranthene	23.97	252	869360	77.929	nq	99
89) Benzo(k)fluoranthene	24.04	252	855474	76.995	nq	100
90) Benzo(a)pyrene	24.87	252	847406	78.312	nq	99
91) Dibenzo(a,h)anthracene	28.89	278	842682	82.108	nq	99
92) Benzo(g,h,i)perylene	30.03	276	831863	81.696	nq	99

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

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