

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062619\
 Data File : BG041463.D
 Acq On : 26 Jun 2019 10:19
 Operator : HP/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

mohammad
 6/28/2019 8:26:44 AM

Quant Time: Jun 26 13:49:04 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG062619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 26 13:18:22 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	24526	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	97260	20.00	ng	0.00
39) Acenaphthene-d10	14.68	164	67427	20.00	ng	0.00
64) Phenanthrene-d10	17.43	188	183826	20.00	ng	0.00
76) Chrysene-d12	21.71	240	200261	20.00	ng	0.00
87) Perylene-d12	25.02	264	212997	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.57	112	24414	18.29	ng	0.00
7) Phenol-d6	7.19	99	35948	18.58	ng	0.00
23) Nitrobenzene-d5	9.22	82	42923	18.56	ng	0.00
42) 2,4,6-Tribromophenol	16.17	330	20555	19.86	ng	0.00
45) 2-Fluorobiphenyl	13.30	172	104369	21.04	ng	0.00
79) Terphenyl-d14	20.03	244	214962	21.27	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.41	88	5511	8.234	ng	# 90
3) Pyridine	3.84	79	13289	7.484	ng	88
4) n-Nitrosodimethylamine	3.76	42	7874	8.315	ng	# 69
6) Aniline	7.35	93	22198	8.254	ng	95
8) 2-Chlorophenol	7.59	128	15125	9.644	ng	92
9) Benzaldehyde	7.18	77	13852	11.193	ng	87
10) Phenol	7.22	94	18763	9.018	ng	96
11) bis(2-Chloroethyl)ether	7.45	93	17602	11.150	ng	95
12) 1,3-Dichlorobenzene	7.92	146	18803	9.575	ng	91
13) 1,4-Dichlorobenzene	8.07	146	19675	9.808	ng	92
14) 1,2-Dichlorobenzene	8.39	146	19293m	10.069	ng	
15) Benzyl Alcohol	8.29	79	10203	5.504	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.55	45	22042	8.529	ng	98
17) 2-Methylphenol	8.48	107	12861	8.646	ng	93
18) Hexachloroethane	9.11	117	8026	10.097	ng	94
19) n-Nitroso-di-n-propylamine	8.84	70	14831	9.477	ng	98
20) 3+4-Methylphenols	8.81	107	17344	8.759	ng	91
22) Acetophenone	8.87	105	27652	9.770	ng	# 93
24) Nitrobenzene	9.26	77	22356	9.612	ng	93
25) Isophorone	9.77	82	40987	9.660	ng	98
26) 2-Nitrophenol	9.98	139	8176	8.671	ng	97
27) 2,4-Dimethylphenol	10.02	122	12985	9.187	ng	99
28) bis(2-Chloroethoxy)methane	10.26	93	20493	9.270	ng	# 97
29) 2,4-Dichlorophenol	10.52	162	16715	9.537	ng	96
30) 1,2,4-Trichlorobenzene	10.72	180	21373	9.522	ng	96
31) Naphthalene	10.91	128	50852	9.620	ng	98
32) Benzoic acid	10.14	122	6712m	7.295	ng	
33) 4-Chloroaniline	11.03	127	20933	9.325	ng	94
34) Hexachlorobutadiene	11.16	225	16890	9.480	ng	91
35) Caprolactam	11.79	113	5233	8.673	ng	# 81
36) 4-Chloro-3-methylphenol	12.15	107	18303	9.041	ng	90
37) 2-Methylnaphthalene	12.51	142	38003	9.761	ng	97
38) 1-Methylnaphthalene	12.73	142	35704	9.585	ng	# 99
40) 1,2,4,5-Tetrachlorobenzene	12.87	216	29098	10.451	ng	99

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41) Hexachlorocyclopentadiene	12.84	237	12254	6.534	nq	99
43) 2,4,6-Trichlorophenol	13.12	196	16948	9.840	nq	94
44) 2,4,5-Trichlorophenol	13.21	196	15138	8.542	nq	94
46) 1,1'-Biphenyl	13.51	154	51290	10.062	nq	92
47) 2-Chloronaphthalene	13.57	162	45991	10.480	nq	94
48) 2-Nitroaniline	13.78	65	13826	8.602	nq	89
49) Acenaphthylene	14.41	152	67598	10.151	nq	97
50) Dimethylphthalate	14.12	163	61697	10.284	nq	99
51) 2,6-Dinitrotoluene	14.26	165	12975	10.257	nq	92
52) Acenaphthene	14.75	154	38242	9.779	nq	96
53) 3-Nitroaniline	14.60	138	11557	9.273	nq	83
54) 2,4-Dinitrophenol	14.83	184	5439	6.712	nq	87
55) Dibenzofuran	15.08	168	69255	10.186	nq	97
56) 4-Nitrophenol	14.95	139	6444m	7.442	nq	
57) 2,4-Dinitrotoluene	15.05	165	17488	9.457	nq	# 83
58) Fluorene	15.73	166	56422	10.549	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.31	232	18894	10.310	nq	# 97
60) Diethylphthalate	15.48	149	62877	10.311	nq	98
61) 4-Chlorophenyl-phenylether	15.71	204	34841	10.012	nq	95
62) 4-Nitroaniline	15.77	138	11484	8.615	nq	97
63) Azobenzene	16.01	77	53487	9.832	nq	97
65) 4,6-Dinitro-2-methylphenol	15.82	198	9219	7.726	nq	89
66) n-Nitrosodiphenylamine	15.93	169	48724	9.982	nq	97
67) 4-Bromophenyl-phenylether	16.61	248	23205	9.763	nq	98
68) Hexachlorobenzene	16.73	284	25798	10.136	nq	99
69) Atrazine	16.87	200	22294	14.004	nq	97
70) Pentachlorophenol	17.08	266	9637	7.394	nq	89
71) Phenanthrene	17.47	178	98171	10.013	nq	99
72) Anthracene	17.57	178	97953	10.134	nq	98
73) Carbazole	17.84	167	82026	9.701	nq	99
74) Di-n-butylphthalate	18.37	149	106235	10.070	nq	99
75) Fluoranthene	19.48	202	133470	10.247	nq	99
77) Benzidine	19.68	184	45973	9.491	nq	98
78) Pyrene	19.85	202	133387	9.876	nq	99
80) Butylbenzylphthalate	20.71	149	51004	9.984	nq	95
81) Benzo(a)anthracene	21.69	228	134056	9.887	nq	99
82) 3,3'-Dichlorobenzidine	21.61	252	43606	9.161	nq	99
83) Chrysene	21.76	228	132697	10.176	nq	98
84) Bis(2-ethylhexyl)phthalate	21.56	149	71143	10.219	nq	97
85) Di-n-octyl phthalate	22.78	149	119647	10.158	nq	99
86) Indeno(1,2,3-cd)pyrene	28.82	276	152639	9.866	nq	98
88) Benzo(b)fluoranthene	23.95	252	130196	9.863	nq	# 96
89) Benzo(k)fluoranthene	24.02	252	131137	10.028	nq	100
90) Benzo(a)pyrene	24.85	252	123767	9.928	nq	97
91) Dibenzo(a,h)anthracene	28.88	278	124215	9.897	nq	97
92) Benzo(g,h,i)perylene	30.01	276	123895	9.989	nq	# 94

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

