

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061920\
 Data File : BG045807.D
 Acq On : 19 Jun 2020 9:28
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Jun 19 11:12:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 15 18:34:19 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	49507	20.00	ng	0.00
21) Naphthalene-d8	10.72	136	184576	20.00	ng	0.00
39) Acenaphthene-d10	14.57	164	132362	20.00	ng	0.00
64) Phenanthrene-d10	17.32	188	309671	20.00	ng	0.00
76) Chrysene-d12	21.57	240	287141	20.00	ng	0.00
86) Perylene-d12	24.70	264	319900	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.53	112	247310	84.38	ng	0.00
7) Phenol-d6	7.15	99	377120	84.60	ng	-0.02
23) Nitrobenzene-d5	9.08	82	353179	87.41	ng	0.00
42) 2,4,6-Tribromophenol	16.07	330	167450	83.80	ng	0.00
45) 2-Fluorobiphenyl	13.19	172	781129	86.73	ng	0.00
79) Terphenyl-d14	19.94	244	1259781	88.91	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.37	88	54366	40.502	ng	97
3) Pyridine	3.77	79	166051	41.746	ng	99
4) n-Nitrosodimethylamine	3.69	42	56716	42.479	ng	98
6) Aniline	7.25	93	220641	42.035	ng	99
8) 2-Chlorophenol	7.51	128	131543	41.932	ng	95
9) Benzaldehyde	7.05	77	109238	40.798	ng	90
10) Phenol	7.18	94	181613	42.046	ng	98
11) bis(2-Chloroethyl)ether	7.34	93	138984	42.362	ng	98
12) 1,3-Dichlorobenzene	7.80	146	154619	41.360	ng	99
13) 1,4-Dichlorobenzene	7.95	146	159246	42.646	ng	99
14) 1,2-Dichlorobenzene	8.27	146	153115	42.831	ng	98
15) Benzyl Alcohol	8.18	79	146498	42.466	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.45	45	132765	42.937	ng	83
17) 2-Methylphenol	8.41	107	135644	42.197	ng	98
18) Hexachloroethane	9.00	117	60291	42.512	ng	94
19) n-Nitroso-di-n-propylamine	8.73	70	123595	42.171	ng	95
20) 3+4-Methylphenols	8.74	107	179244	42.399	ng	# 72
22) Acetophenone	8.74	105	221681	43.758	ng	# 98
24) Nitrobenzene	9.12	77	187211	43.457	ng	96
25) Isophorone	9.64	82	341255	43.629	ng	99
26) 2-Nitrophenol	9.84	139	79995	44.571	ng	96
27) 2,4-Dimethylphenol	9.93	122	120231	43.878	ng	96
28) bis(2-Chloroethoxy)methane	10.13	93	178700	43.738	ng	99
29) 2,4-Dichlorophenol	10.41	162	138681	43.379	ng	93
30) 1,2,4-Trichlorobenzene	10.59	180	165478	43.723	ng	96
31) Naphthalene	10.77	128	439777	43.581	ng	100
32) Benzoic acid	10.13	122	79233	45.224	ng	94
33) 4-Chloroaniline	10.89	127	184402	43.637	ng	97
34) Hexachlorobutadiene	11.07	225	118163	43.901	ng	97
35) Caprolactam	11.66	113	47830	46.233	ng	89
36) 4-Chloro-3-methylphenol	12.06	107	159074	43.095	ng	98
37) 2-Methylnaphthalene	12.39	142	329795	43.889	ng	99
38) 1-Methylnaphthalene	12.60	142	308262	43.351	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.76	216	186625	42.820	ng	99

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41) Hexachlorocyclopentadiene	12.74	237	84028	39.649	nq	96
43) 2,4,6-Trichlorophenol	13.02	196	124173	42.022	nq	96
44) 2,4,5-Trichlorophenol	13.11	196	143713	42.672	nq	98
46) 1,1'-Biphenyl	13.40	154	409498	43.213	nq	99
47) 2-Chloronaphthalene	13.44	162	328916	42.968	nq	97
48) 2-Nitroaniline	13.66	65	110947	43.572	nq	95
49) Acenaphthylene	14.29	152	532596	43.438	nq	100
50) Dimethylphthalate	14.03	163	445982	43.010	nq	99
51) 2,6-Dinitrotoluene	14.14	165	103030	44.239	nq	96
52) Acenaphthene	14.63	154	315597	42.464	nq	99
53) 3-Nitroaniline	14.49	138	97782	42.034	nq	93
54) 2,4-Dinitrophenol	14.70	184	50191	43.813	nq	94
55) Dibenzofuran	14.97	168	528285	43.756	nq	98
56) 4-Nitrophenol	14.87	139	66183	41.711	nq	96
57) 2,4-Dinitrotoluene	14.94	165	144896	43.423	nq	92
58) Fluorene	15.62	166	431934	43.024	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.21	232	125392	42.898	nq	99
60) Diethylphthalate	15.39	149	463472	43.723	nq	96
61) 4-Chlorophenyl-phenylether	15.61	204	249147	43.058	nq	98
62) 4-Nitroaniline	15.66	138	105671	44.261	nq	98
63) Azobenzene	15.91	77	442774	43.503	nq	98
65) 4,6-Dinitro-2-methylphenol	15.71	198	89269	46.646	nq	95
66) n-Nitrosodiphenylamine	15.83	169	378453	43.147	nq	100
67) 4-Bromophenyl-phenylether	16.50	248	173131	43.027	nq	98
68) Hexachlorobenzene	16.63	284	180148	43.176	nq	97
69) Atrazine	16.78	200	145514	42.753	nq	98
70) Pentachlorophenol	16.99	266	76987	41.476	nq	96
71) Phenanthrene	17.36	178	694043	43.456	nq	100
72) Anthracene	17.45	178	685589	43.108	nq	99
73) Carbazole	17.73	167	659599	43.557	nq	100
74) Di-n-butylphthalate	18.28	149	792712	44.211	nq	100
75) Fluoranthene	19.37	202	846227	43.089	nq	98
77) Benzidine	19.56	184	332546	44.548	nq	98
78) Pyrene	19.74	202	858148	44.630	nq	99
80) Butylbenzylphthalate	20.62	149	356055	44.240	nq	94
81) Benzo(a)anthracene	21.56	228	804180	43.197	nq	99
82) 3,3'-Dichlorobenzidine	21.47	252	307095	43.610	nq	98
83) Chrysene	21.62	228	785615	43.630	nq	99
84) Bis(2-ethylhexyl)phthalate	21.47	149	495852	44.293	nq	97
85) Di-n-octyl phthalate	22.64	149	844851	43.751	nq	100
87) Indeno(1,2,3-cd)pyrene	28.26	276	985494	42.494	nq	# 95
88) Benzo(b)fluoranthene	23.72	252	869065	43.329	nq	98
89) Benzo(k)fluoranthene	23.78	252	818207	42.622	nq	98
90) Benzo(a)pyrene	24.56	252	805900	43.012	nq	98
91) Dibenzo(a,h)anthracene	28.31	278	806917	43.389	nq	100
92) Benzo(g,h,i)perylene	29.36	276	786392	42.247	nq	99

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

