

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110518\
 Data File : BG037797.D
 Acq On : 5 Nov 2018 11:17
 Operator : JU/SJ
 Sample : SSTDC0040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDC0040

Quant Time: Nov 05 13:48:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 05 13:45:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	56003	20.00	ng	0.00
21) Naphthalene-d8	10.92	136	264992	20.00	ng	0.00
39) Acenaphthene-d10	14.73	164	169187	20.00	ng	0.00
64) Phenanthrene-d10	17.47	188	375322	20.00	ng	0.00
76) Chrysene-d12	21.76	240	339947	20.00	ng	0.00
87) Perylene-d12	25.09	264	341100	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.63	112	260937	77.90	ng	0.00
7) Phenol-d6	7.24	99	395014	77.99	ng	0.00
23) Nitrobenzene-d5	9.27	82	366485	77.47	ng	0.00
42) 2,4,6-Tribromophenol	16.22	330	146556	79.42	ng	0.00
45) 2-Fluorobiphenyl	13.35	172	912347	81.32	ng	0.00
79) Terphenyl-d14	20.08	244	1264881	79.51	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	64180	37.781	ng	93
3) Pyridine	3.93	79	163875	38.274	ng	93
4) n-Nitrosodimethylamine	3.85	42	55018	36.076	ng	94
6) Aniline	7.42	93	242978	39.321	ng	99
8) 2-Chlorophenol	7.66	128	154722	39.483	ng	99
9) Benzaldehyde	7.24	77	108976	34.393	ng	94
10) Phenol	7.27	94	207952	38.910	ng	99
11) bis(2-Chloroethyl)ether	7.51	93	156590	38.578	ng	98
12) 1,3-Dichlorobenzene	7.98	146	179593	41.166	ng	98
13) 1,4-Dichlorobenzene	8.13	146	178537	40.008	ng	98
14) 1,2-Dichlorobenzene	8.45	146	172835	40.263	ng	98
15) Benzyl Alcohol	8.33	79	141679	37.855	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.62	45	278333	38.323	ng	98
17) 2-Methylphenol	8.53	107	138855	39.300	ng	99
18) Hexachloroethane	9.18	117	62905	41.348	ng	96
19) n-Nitroso-di-n-propylamine	8.90	70	132301	37.930	ng	92
20) 3+4-Methylphenols	8.85	107	196442	38.887	ng	99
22) Acetophenone	8.92	105	252478	37.990	ng	# 95
24) Nitrobenzene	9.32	77	191475	38.964	ng	95
25) Isophorone	9.83	82	327035	37.837	ng	95
26) 2-Nitrophenol	10.02	139	92570	41.815	ng	99
27) 2,4-Dimethylphenol	10.06	122	149539	37.832	ng	94
28) bis(2-Chloroethoxy)methane	10.31	93	212878	37.592	ng	98
29) 2,4-Dichlorophenol	10.55	162	174314	39.608	ng	98
30) 1,2,4-Trichlorobenzene	10.77	180	191856	40.088	ng	99
31) Naphthalene	10.97	128	512996	39.414	ng	99
32) Benzoic acid	10.19	122	85002	33.109	ng	94
33) 4-Chloroaniline	11.08	127	238042	38.924	ng	98
34) Hexachlorobutadiene	11.23	225	117259	41.339	ng	96
35) Caprolactam	11.87	113	66120	37.461	ng	95
36) 4-Chloro-3-methylphenol	12.18	107	188631	38.006	ng	99
37) 2-Methylnaphthalene	12.56	142	377548	39.792	ng	99
38) 1-Methylnaphthalene	12.79	142	364904	39.603	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.93	216	231832	42.482	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.89	237	104109	39.938	nq	98
43) 2,4,6-Trichlorophenol	13.16	196	147912	40.570	nq	98
44) 2,4,5-Trichlorophenol	13.23	196	157280	39.622	nq	98
46) 1,1'-Biphenyl	13.57	154	560039	39.970	nq	99
47) 2-Chloronaphthalene	13.61	162	431427	40.699	nq	98
48) 2-Nitroaniline	13.82	65	126809	39.448	nq	93
49) Acenaphthylene	14.45	152	644716	40.432	nq	100
50) Dimethylphthalate	14.18	163	506358	38.687	nq	99
51) 2,6-Dinitrotoluene	14.31	165	116759	40.325	nq	90
52) Acenaphthene	14.79	154	388743	39.585	nq	99
53) 3-Nitroaniline	14.64	138	126602	39.386	nq	# 94
54) 2,4-Dinitrophenol	14.84	184	25946	34.025	nq	96
55) Dibenzofuran	15.13	168	640728	39.379	nq	99
56) 4-Nitrophenol	14.93	139	80892	33.999	nq	96
57) 2,4-Dinitrotoluene	15.09	165	156132	41.079	nq	93
58) Fluorene	15.78	166	474800	38.967	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.35	232	130672	38.448	nq	97
60) Diethylphthalate	15.53	149	522496	38.884	nq	99
61) 4-Chlorophenyl-phenylether	15.76	204	282846	39.827	nq	97
62) 4-Nitroaniline	15.81	138	124212	38.889	nq	88
63) Azobenzene	16.06	77	484699	37.805	nq	96
65) 4,6-Dinitro-2-methylphenol	15.85	198	67939	36.889	nq	96
66) n-Nitrosodiphenylamine	15.98	169	463236	41.032	nq	99
67) 4-Bromophenyl-phenylether	16.66	248	173367	42.063	nq	99
68) Hexachlorobenzene	16.77	284	173233	40.553	nq	99
69) Atrazine	16.92	200	155168	38.014	nq	99
70) Pentachlorophenol	17.12	266	100613	35.163	nq	92
71) Phenanthrene	17.52	178	762932	39.996	nq	100
72) Anthracene	17.61	178	753585	40.257	nq	98
73) Carbazole	17.88	167	747966	39.945	nq	99
74) Di-n-butylphthalate	18.41	149	836843	40.175	nq	99
75) Fluoranthene	19.52	202	859598	39.732	nq	100
77) Benzidine	19.71	184	331405	28.671	nq	99
78) Pyrene	19.89	202	882529	39.713	nq	100
80) Butylbenzylphthalate	20.76	149	370657	40.754	nq	95
81) Benzo(a)anthracene	21.74	228	802484	39.910	nq	99
82) 3,3'-Dichlorobenzidine	21.66	252	305838	39.241	nq	99
83) Chrysene	21.81	228	779591	40.321	nq	99
84) Bis(2-ethylhexyl)phthalate	21.61	149	529492	41.534	nq	99
85) Di-n-octyl phthalate	22.85	149	868724	41.789	nq	97
86) Indeno(1,2,3-cd)pyrene	28.90	276	711131	34.291	nq	# 92
88) Benzo(b)fluoranthene	24.02	252	740756	39.612	nq	99
89) Benzo(k)fluoranthene	24.09	252	799848	43.657	nq	98
90) Benzo(a)pyrene	24.93	252	724056	40.747	nq	97
91) Dibenzo(a,h)anthracene	28.97	278	558906	34.098	nq	98
92) Benzo(g,h,i)perylene	30.10	276	563829	34.000	nq	96

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

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