

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101617\
 Data File : BG029275.D
 Acq On : 16 Oct 2017 15:59
 Operator : SJ/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ICVBG101617

Quant Time: Oct 16 16:43:56 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 16:15:45 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	67605	20.00	ng	0.00
21) Naphthalene-d8	10.99	136	314445	20.00	ng	0.00
38) Acenaphthene-d10	14.79	164	200425	20.00	ng	0.00
63) Phenanthrene-d10	17.53	188	447611	20.00	ng	0.00
75) Chrysene-d12	21.80	240	461189	20.00	ng	0.00
86) Perylene-d12	25.12	264	477511	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.72	112	317221	78.39	ng	0.00
7) Phenol-d6	7.33	99	452409	79.64	ng	0.00
23) Nitrobenzene-d5	9.34	82	400150	80.87	ng	0.00
41) 2,4,6-Tribromophenol	16.27	330	208617	80.62	ng	0.00
44) 2-Fluorobiphenyl	13.42	172	1054391	77.16	ng	0.00
78) Terphenyl-d14	20.12	244	1580670	77.97	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.58	88	62335	37.964	ng	85
3) Pyridine	3.99	79	188259	39.372	ng	92
4) n-Nitrosodimethylamine	3.90	42	88392	39.547	ng	# 93
6) Aniline	7.49	93	275544	39.173	ng	95
8) 2-Chlorophenol	7.73	128	182715	39.390	ng	93
9) Benzaldehyde	7.30	77	111965	39.188	ng	93
10) Phenol	7.35	94	240160	39.216	ng	92
11) bis(2-Chloroethyl)ether	7.59	93	177265	39.729	ng	91
12) 1,3-Dichlorobenzene	8.06	146	204541	39.276	ng	97
13) 1,4-Dichlorobenzene	8.21	146	208176	39.152	ng	95
14) 1,2-Dichlorobenzene	8.53	146	199620	38.924	ng	97
15) Benzyl Alcohol	8.40	79	160378	40.064	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.70	45	298784	39.431	ng	96
17) 2-Methylphenol	8.61	107	159740	39.266	ng	95
18) Hexachloroethane	9.27	117	70484	38.899	ng	96
19) n-Nitroso-di-n-propylamine	8.98	70	151894	39.327	ng	# 96
20) 3+4-Methylphenols	8.94	107	229078	39.964	ng	97
22) Acetophenone	9.00	105	298190	39.350	ng	# 95
24) Nitrobenzene	9.38	77	225358	40.506	ng	93
25) Isophorone	9.91	82	414418	40.118	ng	# 94
26) 2-Nitrophenol	10.09	139	112707	42.386	ng	92
27) 2,4-Dimethylphenol	10.15	122	191187	39.740	ng	90
28) bis(2-Chloroethoxy)methane	10.38	93	250285	39.513	ng	96
29) 2,4-Dichlorophenol	10.63	162	203985	39.760	ng	95
30) 1,2,4-Trichlorobenzene	10.85	180	218292	39.385	ng	98
31) Naphthalene	11.04	128	620599	39.601	ng	99
32) Benzoic acid	10.29	122	171353	42.537	ng	# 85
33) 4-Chloroaniline	11.14	127	264030	40.053	ng	96
34) Hexachlorobutadiene	11.32	225	137051	39.770	ng	98
35) Caprolactam	11.91	113	70077	41.100	ng	# 86
36) 4-Chloro-3-methylphenol	12.25	107	226716	40.180	ng	91
37) 2-Methylnaphthalene	12.63	142	456734	39.989	ng	93
39) 1,2,4,5-Tetrachlorobenzene	13.00	216	282215	38.567	ng	98
40) Hexachlorocyclopentadiene	12.98	237	102465	39.470	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.23	196	183197	39.881	nq	99
43) 2,4,5-Trichlorophenol	13.30	196	186245	39.479	nq	95
45) 1,1'-Biphenyl	13.63	154	671882	39.001	nq	96
46) 2-Chloronaphthalene	13.67	162	488136	38.847	nq	98
47) 2-Nitroaniline	13.87	65	144464	41.856	nq	# 87
48) Acenaphthylene	14.52	152	768087	39.057	nq	98
49) Dimethylphthalate	14.24	163	616503	39.424	nq	99
50) 2,6-Dinitrotoluene	14.36	165	139300	42.494	nq	# 80
51) Acenaphthene	14.86	154	504678	39.442	nq	96
52) 3-Nitroaniline	14.69	138	146340	41.370	nq	# 82
53) 2,4-Dinitrophenol	14.89	184	64791	39.548	nq	# 53
54) Dibenzofuran	15.19	168	704758	38.582	nq	98
55) 4-Nitrophenol	14.99	139	122908	42.145	nq	# 80
56) 2,4-Dinitrotoluene	15.14	165	190795	42.995	nq	# 81
57) Fluorene	15.83	166	589102	39.040	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.41	232	153805	39.078	nq	98
59) Diethylphthalate	15.59	149	613351	39.357	nq	97
60) 4-Chlorophenyl-phenylether	15.82	204	317551	39.470	nq	96
61) 4-Nitroaniline	15.85	138	143858	39.606	nq	86
62) Azobenzene	16.12	77	516696	39.317	nq	96
64) 4,6-Dinitro-2-methylphenol	15.90	198	101829	40.838	nq	93
65) n-Nitrosodiphenylamine	16.03	169	531204	39.617	nq	98
66) 4-Bromophenyl-phenylether	16.72	248	205674	40.044	nq	99
67) Hexachlorobenzene	16.84	284	227293	39.067	nq	98
68) Atrazine	16.97	200	189461	39.600	nq	99
69) Pentachlorophenol	17.17	266	140966	42.178	nq	99
70) Phenanthrene	17.57	178	904053	39.096	nq	98
71) Anthracene	17.66	178	920783	39.656	nq	99
72) Carbazole	17.93	167	875168	39.237	nq	97
73) Di-n-butylphthalate	18.47	149	1035675	40.372	nq	99
74) Fluoranthene	19.57	202	1066979	39.450	nq	97
76) Benzidine	19.74	184	537746	38.617	nq	# 96
77) Pyrene	19.93	202	1104441	39.477	nq	97
79) Butylbenzylphthalate	20.80	149	472663	39.656	nq	94
80) Benzo(a)anthracene	21.79	228	1059603	39.132	nq	97
81) 3,3'-Dichlorobenzidine	21.69	252	435938	39.006	nq	97
82) Chrysene	21.85	228	1011336	39.000	nq	97
83) Bis(2-ethylhexyl)phthalate	21.67	149	671913	39.280	nq	98
84) Di-n-octyl phthalate	22.92	149	1186767	39.808	nq	98
85) Indeno(1,2,3-cd)pyrene	28.91	276	1283792	40.241	nq	# 100
87) Benzo(b)fluoranthene	24.07	252	1101922	40.308	nq	100
88) Benzo(k)fluoranthene	24.14	252	1059699	38.909	nq	98
89) Benzo(a)pyrene	24.96	252	1038060	39.498	nq	98
90) Dibenzo(a,h)anthracene	28.98	278	1040506	39.106	nq	97
91) Benzo(g,h,i)perylene	30.11	276	970518	39.258	nq	97

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

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