

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG040715\
 Data File : BG016274.D
 Acq On : 8 Apr 2015 1:20
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Apr 08 03:49:29 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 08 03:25:31 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	76512	20.00	ng	0.00
21) Naphthalene-d8	10.48	136	392249	20.00	ng	0.00
38) Acenaphthene-d10	14.34	164	235885	20.00	ng	0.00
63) Phenanthrene-d10	17.08	188	462536	20.00	ng	0.00
75) Chrysene-d12	21.25	240	387927	20.00	ng	0.00
86) Perylene-d12	23.50	264	368961	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.28	112	373072	76.95	ng	0.00
7) Phenol-d6	6.88	99	574766	78.98	ng	0.00
23) Nitrobenzene-d5	8.86	82	513012	75.76	ng	0.00
41) 2,4,6-Tribromophenol	15.83	330	132278	82.92	ng	0.00
44) 2-Fluorobiphenyl	12.96	172	1041794	75.20	ng	0.00
78) Terphenyl-d14	19.70	244	1257682	75.88	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.18	88	75425	34.22	ng	98
3) Pyridine	3.57	79	248273	37.69	ng	98
4) n-Nitrosodimethylamine	3.49	42	72260	36.90	ng	99
6) Aniline	7.03	93	399252	38.41	ng	99
8) 2-Chlorophenol	7.26	128	233310	40.07	ng	98
9) Benzaldehyde	6.84	77	153246	35.95	ng	97
10) Phenol	6.90	94	309063	39.69	ng	100
11) bis(2-Chloroethyl)ether	7.13	93	232955	37.52	ng	99
12) 1,3-Dichlorobenzene	7.59	146	227719	38.73	ng	98
13) 1,4-Dichlorobenzene	7.73	146	233412	38.27	ng	98
14) 1,2-Dichlorobenzene	8.05	146	225567	38.67	ng	98
15) Benzyl Alcohol	7.94	79	203761	40.16	ng	100
16) 2,2'-oxybis(1-Chloropropan	8.23	45	258109	36.91	ng	99
17) 2-Methylphenol	8.15	107	209695	40.16	ng	97
18) Hexachloroethane	8.77	117	87201	37.37	ng	100
19) n-Nitroso-di-n-propylamine	8.50	70	202262	38.79	ng	98
20) 3+4-Methylphenols	8.48	107	294239	40.68	ng	97
22) Acetophenone	8.52	105	339256	37.30	ng	# 99
24) Nitrobenzene	8.90	77	270683	37.38	ng	93
25) Isophorone	9.42	82	562169	37.39	ng	99
26) 2-Nitrophenol	9.60	139	149730	44.35	ng	97
27) 2,4-Dimethylphenol	9.67	122	248484	39.51	ng	98
28) bis(2-Chloroethoxy)methane	9.90	93	323937	37.74	ng	100
29) 2,4-Dichlorophenol	10.14	162	210650	42.33	ng	99
30) 1,2,4-Trichlorobenzene	10.35	180	201405	38.76	ng	98
31) Naphthalene	10.54	128	774478	37.89	ng	100
32) Benzoic acid	9.81	122	180532	44.73	ng	99
33) 4-Chloroaniline	10.65	127	378667	39.48	ng	100
34) Hexachlorobutadiene	10.82	225	89809	39.35	ng	100
35) Caprolactam	11.42	113	126793	40.47	ng	98
36) 4-Chloro-3-methylphenol	11.78	107	267010	40.62	ng	98
37) 2-Methylnaphthalene	12.15	142	563722	38.34	ng	99
39) 1,2,4,5-Tetrachlorobenzene	12.52	216	202444	38.98	ng	99
40) Hexachlorocyclopentadiene	12.50	237	91084	38.32	ng	98

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG040715\
 Data File : BG016274.D
 Acq On : 8 Apr 2015 1:20
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Apr 08 03:49:29 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 08 03:25:31 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.77	196	163266	42.31	ng	95
43) 2,4,5-Trichlorophenol	12.84	196	174506	42.32	ng	98
45) 1,1'-Biphenyl	13.17	154	680475	37.61	ng	99
46) 2-Chloronaphthalene	13.21	162	519574	37.69	ng	98
47) 2-Nitroaniline	13.42	65	172810	38.90	ng	94
48) Acenaphthylene	14.06	152	919577	37.52	ng	98
49) Dimethylphthalate	13.80	163	641851	37.13	ng	99
50) 2,6-Dinitrotoluene	13.92	165	163365	39.07	ng	95
51) Acenaphthene	14.40	154	547171	37.36	ng	98
52) 3-Nitroaniline	14.25	138	208261	39.00	ng	93
53) 2,4-Dinitrophenol	14.46	184	81855	38.60	ng	97
54) Dibenzofuran	14.74	168	757311	37.26	ng	98
55) 4-Nitrophenol	14.56	139	168467	38.18	ng	98
56) 2,4-Dinitrotoluene	14.71	165	223975	39.33	ng	95
57) Fluorene	15.38	166	667037	37.20	ng	99
58) 2,3,4,6-Tetrachlorophenol	14.97	232	131881	41.44	ng	94
59) Diethylphthalate	15.17	149	673426	36.59	ng	99
60) 4-Chlorophenyl-phenylether	15.38	204	276654	37.72	ng	99
61) 4-Nitroaniline	15.41	138	233579	38.54	ng	96
62) Azobenzene	15.67	77	690331	35.25	ng	99
64) 4,6-Dinitro-2-methylphenol	15.47	198	124243	38.96	ng	96
65) n-Nitrosodiphenylamine	15.60	169	614350	39.29	ng	99
66) 4-Bromophenyl-phenylether	16.27	248	147563	39.96	ng	97
67) Hexachlorobenzene	16.39	284	153489	39.99	ng	96
68) Atrazine	16.55	200	171543	39.36	ng	100
69) Pentachlorophenol	16.73	266	100736	42.98	ng	99
70) Phenanthrene	17.12	178	980563	37.69	ng	100
71) Anthracene	17.21	178	991331	38.45	ng	100
72) Carbazole	17.49	167	977688	37.79	ng	99
73) Di-n-butylphthalate	18.05	149	1191683	37.64	ng	99
74) Fluoranthene	19.14	202	1000295	37.30	ng	98
76) Benzidine	19.33	184	588763	35.79	ng	99
77) Pyrene	19.50	202	1030748	38.16	ng	99
79) Butylbenzylphthalate	20.40	149	540991	40.76	ng	98
80) Benzo(a)anthracene	21.24	228	872195	38.04	ng	100
81) 3,3'-Dichlorobenzidine	21.18	252	303511	40.26	ng	98
82) Chrysene	21.30	228	829885	37.49	ng	99
83) Bis(2-ethylhexyl)phthalate	21.17	149	729011	40.58	ng	98
84) Di-n-octyl phthalate	22.04	149	1196466	40.87	ng	99
85) Indeno(1,2,3-cd)pyrene	25.80	276	919719	40.00	ng	98
87) Benzo(b)fluoranthene	22.82	252	796207	36.84	ng	98
88) Benzo(k)fluoranthene	22.86	252	817000	38.63	ng	98
89) Benzo(a)pyrene	23.40	252	771815	38.11	ng	99
90) Dibenzo(a,h)anthracene	25.80	278	758856	38.51	ng	99
91) Benzo(g,h,i)perylene	26.49	276	758792	38.55	ng	99

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

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG040715\
Data File : BG016274.D
Acq On : 8 Apr 2015 1:20
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SSTDCCC040EC

Quant Time: Apr 08 03:49:29 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 08 03:25:31 2015
Response via : Initial Calibration

