

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061219\
 Data File : BG041208.D
 Acq On : 12 Jun 2019 16:17
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 13 07:41:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 23:08:34 2019
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	71	0.00
2	1,4-Dioxane	0.503	0.476	5.4	67	0.00
3	Pyridine	1.489	1.387	6.9	66	0.00
4	n-Nitrosodimethylamine	0.691	0.718	-3.9	73	0.00
5 S	2-Fluorophenol	1.109	1.087	2.0	69	0.00
6	Aniline	2.286	2.272	0.6	71	0.00
7 S	Phenol-d6	1.713	1.725	-0.7	72	0.00
8	2-Chlorophenol	1.322	1.353	-2.3	73	0.00
9	Benzaldehyde	1.022	1.046	-2.3	72	0.00
10 C	Phenol	1.837	1.856	-1.0	72	0.00
11	bis(2-Chloroethyl)ether	1.332	1.260	5.4	66	0.00
12	1,3-Dichlorobenzene	1.579	1.584	-0.3	71	0.00
13 C	1,4-Dichlorobenzene	1.599	1.615	-1.0	70	0.00
14	1,2-Dichlorobenzene	1.552	1.587	-2.3	72	0.00
15	Benzyl Alcohol	1.585	1.557	1.8	70	0.00
16	2,2'-oxybis(1-Chloropropane	2.177	2.014	7.5	66	-0.01
17	2-Methylphenol	1.330	1.284	3.5	68	0.00
18	Hexachloroethane	0.616	0.628	-1.9	73	0.00
19 P	n-Nitroso-di-n-propylamine	1.318	1.243	5.7	67	0.00
20	3+4-Methylphenols	1.842	1.788	2.9	70	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	72	0.00
22	Acetophenone	0.561	0.554	1.2	70	0.00
23 S	Nitrobenzene-d5	0.451	0.461	-2.2	73	0.00
24	Nitrobenzene	0.459	0.473	-3.1	73	0.00
25	Isophorone	0.857	0.814	5.0	67	0.00
26 C	2-Nitrophenol	0.189	0.207	-9.5	77	0.00
27	2,4-Dimethylphenol	0.313	0.307	1.9	69	0.00
28	bis(2-Chloroethoxy)methane	0.463	0.443	4.3	68	0.00
29 C	2,4-Dichlorophenol	0.397	0.393	1.0	69	0.00
30	1,2,4-Trichlorobenzene	0.452	0.458	-1.3	71	0.00
31	Naphthalene	1.074	1.082	-0.7	71	0.00
32	Benzoic acid	0.223	0.224	-0.4	70	0.00
33	4-Chloroaniline	0.498	0.504	-1.2	71	0.00
34 C	Hexachlorobutadiene	0.346	0.355	-2.6	75	0.00
35	Caprolactam	0.131	0.134	-2.3	71	0.00
36 C	4-Chloro-3-methylphenol	0.453	0.457	-0.9	71	0.00
37	2-Methylnaphthalene	0.827	0.838	-1.3	72	0.00
38	1-Methylnaphthalene	0.779	0.807	-3.6	73	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	74	0.00
40	1,2,4,5-Tetrachlorobenzene	0.806	0.815	-1.1	73	0.00
41 P	Hexachlorocyclopentadiene	0.361	0.365	-1.1	74	0.00
42 S	2,4,6-Tribromophenol	0.297	0.305	-2.7	74	0.00
43 C	2,4,6-Trichlorophenol	0.521	0.529	-1.5	73	0.00
44	2,4,5-Trichlorophenol	0.550	0.541	1.6	72	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061219\
 Data File : BG041208.D
 Acq On : 12 Jun 2019 16:17
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 13 07:41:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 23:08:34 2019
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45 S	2-Fluorobiphenyl	1.389	1.402	-0.9	73	0.00
46	1,1'-Biphenyl	1.480	1.479	0.1	71	0.00
47	2-Chloronaphthalene	1.271	1.304	-2.6	74	0.00
48	2-Nitroaniline	0.456	0.477	-4.6	75	0.00
49	Acenaphthylene	1.913	1.935	-1.2	71	0.00
50	Dimethylphthalate	1.693	1.657	2.1	70	0.00
51	2,6-Dinitrotoluene	0.365	0.365	0.0	72	0.00
52 C	Acenaphthene	1.159	1.146	1.1	71	0.00
53	3-Nitroaniline	0.365	0.370	-1.4	71	0.00
54 P	2,4-Dinitrophenol	0.142	0.142	0.0	77	0.00
55	Dibenzofuran	1.950	1.982	-1.6	72	0.00
56 P	4-Nitrophenol	0.250	0.245	2.0	69	0.00
57	2,4-Dinitrotoluene	0.516	0.525	-1.7	72	0.00
58	Fluorene	1.553	1.562	-0.6	72	0.00
59	2,3,4,6-Tetrachlorophenol	0.540	0.550	-1.9	73	0.00
60	Diethylphthalate	1.728	1.702	1.5	70	0.00
61	4-Chlorophenyl-phenylether	1.009	1.012	-0.3	73	0.00
62	4-Nitroaniline	0.386	0.387	-0.3	72	0.00
63	Azobenzene	1.570	1.554	1.0	70	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	72	0.00
65	4,6-Dinitro-2-methylphenol	0.096	0.112	-16.7	83	0.00
66 c	n-Nitrosodiphenylamine	0.562	0.565	-0.5	71	0.00
67	4-Bromophenyl-phenylether	0.270	0.268	0.7	70	0.00
68	Hexachlorobenzene	0.287	0.285	0.7	71	0.00
69	Atrazine	0.217	0.217	0.0	65	0.00
70 C	Pentachlorophenol	0.139	0.146	-5.0	71	0.00
71	Phenanthrene	1.042	1.076	-3.3	72	0.00
72	Anthracene	1.027	1.067	-3.9	73	0.00
73	Carbazole	0.882	0.928	-5.2	74	0.00
74	Di-n-butylphthalate	1.096	1.131	-3.2	71	0.00
75 C	Fluoranthene	1.287	1.367	-6.2	74	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	72	0.00
77	Benzidine	0.477	0.409	14.3	58	0.00
78	Pyrene	1.300	1.399	-7.6	75	0.00
79 S	Terphenyl-d14	0.942	1.042	-10.6	76	0.00
80	Butylbenzylphthalate	0.506	0.538	-6.3	74	0.00
81	Benzo(a)anthracene	1.331	1.394	-4.7	74	0.00
82	3,3'-Dichlorobenzidine	0.468	0.474	-1.3	73	0.00
83	Chrysene	1.274	1.341	-5.3	73	0.00
84	Bis(2-ethylhexyl)phthalate	0.712	0.739	-3.8	73	0.00
85 c	Di-n-octyl phthalate	1.186	1.240	-4.6	73	0.00
86	Indeno(1,2,3-cd)pyrene	1.561	1.259	19.3	58	0.00
87 I	Perylene-d12	1.000	1.000	0.0	71	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061219\
Data File : BG041208.D
Acq On : 12 Jun 2019 16:17
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Jun 13 07:41:29 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Jun 09 23:08:34 2019
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.213	1.223	-0.8	70	0.00
89	Benzo(k)fluoranthene	1.200	1.297	-8.1	74	0.00
90 C	Benzo(a)pyrene	1.157	1.171	-1.2	70	0.00
91	Dibenzo(a,h)anthracene	1.156	0.987	14.6	59	0.00
92	Benzo(q,h,i)perylene	1.124	0.878	21.9	55	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0