

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA M\Data\BM052318\
 Data File : BM015274.D
 Acq On : 23 May 2018 14:00
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: May 24 02:09:23 2018
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM052218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 22 15:47:28 2018
 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	119	0.00
2	1,4-Dioxane	0.505	0.483	4.4	111	0.00
3	Pyridine	1.294	1.330	-2.8	117	-0.01
4	n-Nitrosodimethylamine	0.700	0.749	-7.0	122	0.00
5 S	2-Fluorophenol	1.222	1.243	-1.7	117	0.00
6	Aniline	1.961	2.051	-4.6	121	0.00
7 S	Phenol-d6	1.592	1.659	-4.2	120	0.00
8	2-Chlorophenol	1.380	1.427	-3.4	120	0.00
9	Benzaldehyde	1.007	1.042	-3.5	123	0.00
10 C	Phenol	1.617	1.675	-3.6	120	0.00
11	bis(2-Chloroethyl)ether	1.282	1.339	-4.4	122	0.00
12	1,3-Dichlorobenzene	1.573	1.596	-1.5	119	0.00
13 C	1,4-Dichlorobenzene	1.590	1.627	-2.3	120	0.00
14	1,2-Dichlorobenzene	1.510	1.543	-2.2	120	0.00
15	Benzyl Alcohol	1.151	1.242	-7.9	124	0.00
16	2,2'-oxybis(1-Chloropropane	2.023	2.145	-6.0	127	0.00
17	2-Methylphenol	1.075	1.140	-6.0	123	0.00
18	Hexachloroethane	0.565	0.597	-5.7	122	0.00
19 P	n-Nitroso-di-n-propylamine	1.015	1.097	-8.1	125	0.00
20	3+4-Methylphenols	1.427	1.504	-5.4	122	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	125	0.00
22	Acetophenone	0.480	0.488	-1.7	125	0.00
23 S	Nitrobenzene-d5	0.365	0.378	-3.6	125	0.00
24	Nitrobenzene	0.380	0.390	-2.6	125	0.00
25	Isophorone	0.627	0.660	-5.3	126	0.00
26 C	2-Nitrophenol	0.171	0.175	-2.3	125	0.00
27	2,4-Dimethylphenol	0.309	0.317	-2.6	124	0.00
28	bis(2-Chloroethoxy)methane	0.411	0.426	-3.6	127	0.00
29 C	2,4-Dichlorophenol	0.293	0.303	-3.4	124	0.00
30	1,2,4-Trichlorobenzene	0.345	0.351	-1.7	124	0.00
31	Naphthalene	1.040	1.047	-0.7	125	0.00
32	Benzoic acid	0.179	0.178	0.6	124	0.03
33	4-Chloroaniline	0.415	0.427	-2.9	125	0.00
34 C	Hexachlorobutadiene	0.205	0.211	-2.9	126	0.00
35	Caprolactam	0.083	0.081	2.4	117	0.01
36 C	4-Chloro-3-methylphenol	0.299	0.307	-2.7	123	0.00
37	2-Methylnaphthalene	0.720	0.735	-2.1	125	0.00
38 I	Acenaphthene-d10	1.000	1.000	0.0	124	0.00
39	1,2,4,5-Tetrachlorobenzene	0.704	0.733	-4.1	126	0.00
40 P	Hexachlorocyclopentadiene	0.407	0.455	-11.8	133	0.00
41 S	2,4,6-Tribromophenol	0.251	0.261	-4.0	121	0.00
42 C	2,4,6-Trichlorophenol	0.423	0.453	-7.1	125	0.00
43	2,4,5-Trichlorophenol	0.457	0.480	-5.0	124	0.00
44 S	2-Fluorobiphenyl	1.609	1.665	-3.5	126	0.00

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA M\Data\BM052318\
 Data File : BM015274.D
 Acq On : 23 May 2018 14:00
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: May 24 02:09:23 2018
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM052218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 22 15:47:28 2018
 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)
45	1,1'-Biphenyl	1.750	1.804	-3.1	125	0.00
46	2-Chloronaphthalene	1.347	1.388	-3.0	125	0.00
47	2-Nitroaniline	0.384	0.396	-3.1	124	0.00
48	Acenaphthylene	2.000	2.078	-3.9	124	0.00
49	Dimethylphthalate	1.596	1.611	-0.9	122	0.00
50	2,6-Dinitrotoluene	0.333	0.347	-4.2	121	0.00
51 C	Acenaphthene	1.246	1.271	-2.0	124	0.00
52	3-Nitroaniline	0.351	0.350	0.3	119	0.00
53 P	2,4-Dinitrophenol	0.170	0.171	-0.6	121	0.00
54	Dibenzofuran	1.965	1.969	-0.2	123	0.00
55 P	4-Nitrophenol	0.284	0.276	2.8	116	0.00
56	2,4-Dinitrotoluene	0.456	0.457	-0.2	120	0.00
57	Fluorene	1.553	1.571	-1.2	122	0.00
58	2,3,4,6-Tetrachlorophenol	0.392	0.403	-2.8	121	0.00
59	Diethylphthalate	1.573	1.597	-1.5	122	0.00
60	4-Chlorophenyl-phenylether	0.791	0.805	-1.8	124	0.00
61	4-Nitroaniline	0.367	0.362	1.4	118	0.00
62	Azobenzene	1.495	1.588	-6.2	124	0.00
63 I	Phenanthrene-d10	1.000	1.000	0.0	119	0.00
64	4,6-Dinitro-2-methylphenol	0.110	0.113	-2.7	119	0.00
65 c	n-Nitrosodiphenylamine	0.643	0.677	-5.3	122	0.00
66	4-Bromophenyl-phenylether	0.223	0.237	-6.3	122	0.00
67	Hexachlorobenzene	0.257	0.266	-3.5	121	0.00
68	Atrazine	0.204	0.213	-4.4	117	0.00
69 C	Pentachlorophenol	0.143	0.146	-2.1	118	0.00
70	Phenanthrene	1.139	1.143	-0.4	117	0.00
71	Anthracene	1.136	1.165	-2.6	118	0.00
72	Carbazole	1.005	1.013	-0.8	115	0.00
73	Di-n-butylphthalate	1.184	1.264	-6.8	118	0.00
74 C	Fluoranthene	1.249	1.226	1.8	112	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	107	0.00
76	Benzidine	0.626	0.701	-12.0	109	0.00
77	Pyrene	1.262	1.348	-6.8	113	0.00
78 S	Terphenyl-d14	0.906	0.970	-7.1	113	0.00
79	Butylbenzylphthalate	0.515	0.576	-11.8	111	0.00
80	Benzo(a)anthracene	1.260	1.294	-2.7	108	0.00
81	3,3'-Dichlorobenzidine	0.467	0.488	-4.5	106	0.00
82	Chrysene	1.187	1.200	-1.1	106	0.00
83	Bis(2-ethylhexyl)phthalate	0.750	0.838	-11.7	111	0.00
84 c	Di-n-octyl phthalate	1.311	1.374	-4.8	108	0.00
85	Indeno(1,2,3-cd)pyrene	1.436	1.432	0.3	103	0.00
86 I	Perylene-d12	1.000	1.000	0.0	103	0.00
87	Benzo(b)fluoranthene	1.265	1.297	-2.5	103	0.00

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA M\Data\BM052318\
Data File : BM015274.D
Acq On : 23 May 2018 14:00
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: May 24 02:09:23 2018
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM052218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 22 15:47:28 2018
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(k)fluoranthene	1.197	1.259	-5.2	107	0.00
89 C	Benzo(a)pyrene	1.180	1.226	-3.9	104	0.00
90	Dibenzo(a,h)anthracene	1.204	1.236	-2.7	104	0.00
91	Benzo(a,h,i)perylene	1.173	1.196	-2.0	103	0.00

(#) = Out of Range

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