

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071720\
 Data File : BF120911.D
 Acq On : 17 Jul 2020 9:13
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 17 10:46:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 16 14:20:32 2020
 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	124	0.00
2	1,4-Dioxane	0.561	0.537	4.3	122	0.04
3	Pyridine	1.494	1.464	2.0	126	0.04
4	n-Nitrosodimethylamine	0.639	0.600	6.1	119	0.05
5 S	2-Fluorophenol	1.220	1.153	5.5	121	0.00
6	Aniline	2.057	1.939	5.7	121	0.00
7 S	Phenol-d6	1.576	1.489	5.5	122	0.00
8	2-Chlorophenol	1.344	1.283	4.5	121	0.00
9	Benzaldehyde	0.811	0.753	7.2	125	0.00
10 C	Phenol	1.734	1.642	5.3	121	0.00
11	bis(2-Chloroethyl)ether	1.329	1.272	4.3	123	0.00
12	1,3-Dichlorobenzene	1.457	1.364	6.4	121	0.00
13 C	1,4-Dichlorobenzene	1.449	1.384	4.5	123	0.00
14	1,2-Dichlorobenzene	1.371	1.293	5.7	120	0.00
15	Benzyl Alcohol	1.114	1.069	4.0	123	0.00
16	2,2'-oxybis(1-Chloropropane	1.915	1.846	3.6	124	0.00
17	2-Methylphenol	1.191	1.138	4.5	124	0.00
18	Hexachloroethane	0.525	0.504	4.0	121	0.00
19 P	n-Nitroso-di-n-propylamine	0.896	0.841	6.1	125	0.00
20	3+4-Methylphenols	1.515	1.311	13.5	121	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	125	0.00
22	Acetophenone	0.449	0.420	6.5	122	0.00
23 S	Nitrobenzene-d5	0.346	0.337	2.6	124	0.00
24	Nitrobenzene	0.373	0.363	2.7	125	0.00
25	Isophorone	0.690	0.665	3.6	125	0.00
26 C	2-Nitrophenol	0.177	0.187	-5.6	129	0.00
27	2,4-Dimethylphenol	0.287	0.280	2.4	126	0.00
28	bis(2-Chloroethoxy)methane	0.421	0.409	2.9	126	0.00
29 C	2,4-Dichlorophenol	0.294	0.288	2.0	124	0.00
30	1,2,4-Trichlorobenzene	0.326	0.310	4.9	120	0.00
31	Naphthalene	1.026	0.965	5.9	121	0.00
32	Benzoic acid	0.216	0.213	1.4	112	0.02
33	4-Chloroaniline	0.458	0.431	5.9	123	0.00
34 C	Hexachlorobutadiene	0.192	0.184	4.2	122	0.00
35	Caprolactam	0.094	0.093	1.1	126	0.01
36 C	4-Chloro-3-methylphenol	0.301	0.291	3.3	123	0.00
37	2-Methylnaphthalene	0.666	0.642	3.6	123	0.00
38	1-Methylnaphthalene	0.631	0.607	3.8	123	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	126	0.00
40	1,2,4,5-Tetrachlorobenzene	0.583	0.562	3.6	123	0.00
41 P	Hexachlorocyclopentadiene	0.322	0.327	-1.6	124	0.00
42 S	2,4,6-Tribromophenol	0.219	0.216	1.4	125	0.00
43 C	2,4,6-Trichlorophenol	0.410	0.407	0.7	126	0.00
44	2,4,5-Trichlorophenol	0.465	0.462	0.6	127	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071720\
 Data File : BF120911.D
 Acq On : 17 Jul 2020 9:13
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 17 10:46:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 16 14:20:32 2020
 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 S	2-Fluorobiphenyl	1.340	1.118	16.6	121	0.00
46	1,1'-Biphenyl	1.526	1.447	5.2	122	0.00
47	2-Chloronaphthalene	1.244	1.187	4.6	122	0.00
48	2-Nitroaniline	0.376	0.379	-0.8	128	0.00
49	Acenaphthylene	1.932	1.855	4.0	124	0.00
50	Dimethylphthalate	1.443	1.370	5.1	124	0.00
51	2,6-Dinitrotoluene	0.310	0.313	-1.0	126	0.00
52 C	Acenaphthene	1.229	1.188	3.3	124	0.00
53	3-Nitroaniline	0.389	0.386	0.8	127	0.00
54 P	2,4-Dinitrophenol	0.149	0.163	-9.4	139	0.00
55	Dibenzofuran	1.777	1.686	5.1	124	0.00
56 P	4-Nitrophenol	0.295	0.296	-0.3	126	0.00
57	2,4-Dinitrotoluene	0.407	0.420	-3.2	127	0.00
58	Fluorene	1.325	1.199	9.5	122	0.00
59	2,3,4,6-Tetrachlorophenol	0.354	0.361	-2.0	130	0.00
60	Diethylphthalate	1.459	1.402	3.9	125	0.00
61	4-Chlorophenyl-phenylether	0.707	0.611	13.6	122	0.00
62	4-Nitroaniline	0.379	0.373	1.6	126	0.00
63	Azobenzene	1.389	1.306	6.0	122	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	125	0.00
65	4,6-Dinitro-2-methylphenol	0.101	0.116	-14.9	137	0.00
66 c	n-Nitrosodiphenylamine	0.629	0.621	1.3	123	0.00
67	4-Bromophenyl-phenylether	0.223	0.225	-0.9	126	0.00
68	Hexachlorobenzene	0.241	0.242	-0.4	126	0.00
69	Atrazine	0.156	0.146	6.4	120	0.00
70 C	Pentachlorophenol	0.138	0.150	-8.7	127	0.00
71	Phenanthrene	1.029	0.980	4.8	122	0.00
72	Anthracene	1.033	0.998	3.4	121	0.00
73	Carbazole	1.025	0.982	4.2	121	0.00
74	Di-n-butylphthalate	1.183	1.143	3.4	122	0.00
75 C	Fluoranthene	1.140	1.085	4.8	120	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	119	0.00
77	Benzidine	0.526	0.566	-7.6	120	0.00
78	Pyrene	1.414	1.420	-0.4	121	0.00
79 S	Terphenyl-d14	0.920	0.823	10.5	115	0.00
80	Butylbenzylphthalate	0.630	0.646	-2.5	121	0.00
81	Benzo(a)anthracene	1.211	1.161	4.1	119	0.00
82	3,3'-Dichlorobenzidine	0.457	0.473	-3.5	124	0.00
83	Chrysene	1.273	1.228	3.5	115	0.00
84	Bis(2-ethylhexyl)phthalate	0.777	0.759	2.3	118	0.00
85 c	Di-n-octyl phthalate	1.546	1.512	2.2	115	0.00
86 I	Perylene-d12	1.000	1.000	0.0	114	0.00
87	Indeno(1,2,3-cd)pyrene	1.391	1.318	5.2	107	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071720\
Data File : BF120911.D
Acq On : 17 Jul 2020 9:13
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jul 17 10:46:09 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 16 14:20:32 2020
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	1.209	1.273	-5.3	122	0.00
89	Benzo(k)fluoranthene	1.103	1.084	1.7	108	0.00
90 C	Benzo(a)pyrene	1.103	1.118	-1.4	114	0.00
91	Dibenzo(a,h)anthracene	1.136	1.077	5.2	107	0.00
92	Benzo(q,h,i)perylene	1.120	1.043	6.9	106	0.00

(#) = Out of Range

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