

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080520\
 Data File : BF121184.D
 Acq On : 5 Aug 2020 9:52
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 05 11:27:11 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 29 14:55:43 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	95	0.00
2	1,4-Dioxane	0.561	0.553	1.4	97	-0.01
3	Pyridine	1.494	1.529	-2.3	101	-0.02
4	n-Nitrosodimethylamine	0.639	0.593	7.2	91	-0.02
5 S	2-Fluorophenol	1.220	1.186	2.8	96	-0.01
6	Aniline	2.057	1.838	10.6	88	-0.01
7 S	Phenol-d6	1.576	1.514	3.9	95	0.00
8	2-Chlorophenol	1.344	1.323	1.6	96	0.00
9	Benzaldehyde	0.811	0.784	3.3	100	0.00
10 C	Phenol	1.734	1.616	6.8	92	0.00
11	bis(2-Chloroethyl)ether	1.329	1.288	3.1	95	-0.01
12	1,3-Dichlorobenzene	1.457	1.431	1.8	97	-0.01
13 C	1,4-Dichlorobenzene	1.449	1.435	1.0	98	0.00
14	1,2-Dichlorobenzene	1.371	1.323	3.5	94	0.00
15	Benzyl Alcohol	1.114	1.012	9.2	89	-0.01
16	2,2'-oxybis(1-Chloropropane	1.915	1.780	7.0	92	-0.01
17	2-Methylphenol	1.191	1.150	3.4	97	-0.01
18	Hexachloroethane	0.525	0.529	-0.8	98	0.00
19 P	n-Nitroso-di-n-propylamine	0.896	0.801	10.6	91	-0.01
20	3+4-Methylphenols	1.515	1.313	13.3	93	-0.01
21 I	Naphthalene-d8	1.000	1.000	0.0	94	0.00
22	Acetophenone	0.449	0.438	2.4	96	0.00
23 S	Nitrobenzene-d5	0.346	0.350	-1.2	97	-0.01
24	Nitrobenzene	0.373	0.374	-0.3	97	-0.01
25	Isophorone	0.690	0.665	3.6	94	-0.01
26 C	2-Nitrophenol	0.177	0.198	-11.9	103	0.00
27	2,4-Dimethylphenol	0.287	0.283	1.4	96	0.00
28	bis(2-Chloroethoxy)methane	0.421	0.414	1.7	96	-0.01
29 C	2,4-Dichlorophenol	0.294	0.298	-1.4	96	-0.01
30	1,2,4-Trichlorobenzene	0.326	0.326	0.0	95	0.00
31	Naphthalene	1.026	1.012	1.4	96	0.00
32	Benzoic acid	0.216	0.206	4.6	81	0.00
33	4-Chloroaniline	0.458	0.444	3.1	95	-0.01
34 C	Hexachlorobutadiene	0.192	0.193	-0.5	96	-0.01
35	Caprolactam	0.094	0.095	-1.1	97	-0.01
36 C	4-Chloro-3-methylphenol	0.301	0.303	-0.7	96	0.00
37	2-Methylnaphthalene	0.666	0.668	-0.3	96	0.00
38	1-Methylnaphthalene	0.631	0.621	1.6	95	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	94	0.00
40	1,2,4,5-Tetrachlorobenzene	0.583	0.607	-4.1	98	0.00
41 P	Hexachlorocyclopentadiene	0.322	0.277	14.0	78	0.00
42 S	2,4,6-Tribromophenol	0.219	0.228	-4.1	97	0.00
43 C	2,4,6-Trichlorophenol	0.410	0.402	2.0	92	0.00
44	2,4,5-Trichlorophenol	0.465	0.465	0.0	95	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080520\
 Data File : BF121184.D
 Acq On : 5 Aug 2020 9:52
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 05 11:27:11 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 29 14:55:43 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	1.206	10.0	96	0.00
46	1,1'-Biphenyl	1.526	1.523	0.2	95	0.00
47	2-Chloronaphthalene	1.244	1.234	0.8	94	0.00
48	2-Nitroaniline	0.376	0.392	-4.3	98	0.00
49	Acenaphthylene	1.932	1.915	0.9	95	0.00
50	Dimethylphthalate	1.443	1.453	-0.7	98	0.00
51	2,6-Dinitrotoluene	0.310	0.325	-4.8	97	0.00
52 C	Acenaphthene	1.229	1.137	7.5	88	0.00
53	3-Nitroaniline	0.389	0.397	-2.1	96	0.00
54 P	2,4-Dinitrophenol	0.149	0.153	-2.7	97	0.00
55	Dibenzofuran	1.777	1.717	3.4	93	0.00
56 P	4-Nitrophenol	0.295	0.249	15.6	78	0.00
57	2,4-Dinitrotoluene	0.407	0.431	-5.9	96	0.00
58	Fluorene	1.325	1.295	2.3	97	0.00
59	2,3,4,6-Tetrachlorophenol	0.354	0.340	4.0	90	0.00
60	Diethylphthalate	1.459	1.409	3.4	93	0.00
61	4-Chlorophenyl-phenylether	0.707	0.658	6.9	97	0.00
62	4-Nitroaniline	0.379	0.399	-5.3	100	0.00
63	Azobenzene	1.389	1.326	4.5	91	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	92	0.00
65	4,6-Dinitro-2-methylphenol	0.101	0.121	-19.8	105	0.00
66 c	n-Nitrosodiphenylamine	0.629	0.641	-1.9	94	0.00
67	4-Bromophenyl-phenylether	0.223	0.232	-4.0	96	0.00
68	Hexachlorobenzene	0.241	0.252	-4.6	96	0.00
69	Atrazine	0.156	0.152	2.6	91	0.00
70 C	Pentachlorophenol	0.138	0.142	-2.9	89	0.00
71	Phenanthrene	1.029	1.023	0.6	93	0.00
72	Anthracene	1.033	1.050	-1.6	94	0.00
73	Carbazole	1.025	1.036	-1.1	94	0.00
74	Di-n-butylphthalate	1.183	1.138	3.8	89	0.00
75 C	Fluoranthene	1.140	1.126	1.2	91	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	86	0.00
77	Benzidine	0.526	0.479	8.9	73	0.00
78	Pyrene	1.414	1.495	-5.7	92	0.00
79 S	Terphenyl-d14	0.920	0.882	4.1	89	0.00
80	Butylbenzylphthalate	0.630	0.638	-1.3	86	0.00
81	Benzo(a)anthracene	1.211	1.274	-5.2	94	0.00
82	3,3'-Dichlorobenzidine	0.457	0.462	-1.1	88	0.00
83	Chrysene	1.273	1.261	0.9	86	0.00
84	Bis(2-ethylhexyl)phthalate	0.777	0.801	-3.1	90	0.00
85 c	Di-n-octyl phthalate	1.546	1.384	10.5	76	0.00
86 I	Perylene-d12	1.000	1.000	0.0	72	0.01
87	Indeno(1,2,3-cd)pyrene	1.391	1.217	12.5	63	0.02

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080520\
Data File : BF121184.D
Acq On : 5 Aug 2020 9:52
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 05 11:27:11 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 29 14:55:43 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	1.316	-8.9	80	0.00
89	Benzo(k)fluoranthene	1.103	1.243	-12.7	79	0.00
90 C	Benzo(a)pyrene	1.103	1.177	-6.7	76	0.01
91	Dibenzo(a,h)anthracene	1.136	0.995	12.4	63	0.01
92	Benzo(q,h,i)perylene	1.120	0.968	13.6	62	0.02

(#) = Out of Range

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