

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080721\
 Data File : BF124975.D
 Acq On : 7 Aug 2021 15:21
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 07 16:03:39 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 06 11:27:24 2021
 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)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	113	0.00
2	1,4-Dioxane	0.457	0.446	2.4	107	0.04
3	Pyridine	1.061	1.081	-1.9	104	0.03
4	n-Nitrosodimethylamine	0.754	0.753	0.1	101	0.04
5 S	2-Fluorophenol	1.221	1.191	2.5	106	0.00
6	Aniline	1.626	1.585	2.5	103	0.01
7 S	Phenol-d6	1.539	1.523	1.0	104	0.00
8	2-Chlorophenol	1.339	1.304	2.6	106	0.00
9	Benzaldehyde	0.838	0.726	13.4	96	0.00
10 C	Phenol	1.588	1.564	1.5	107	0.01
11	bis(2-Chloroethyl)ether	1.192	1.190	0.2	105	0.01
12	1,3-Dichlorobenzene	1.458	1.405	3.6	103	0.00
13 C	1,4-Dichlorobenzene	1.457	1.452	0.3	105	0.00
14	1,2-Dichlorobenzene	1.366	1.313	3.9	103	0.01
15	Benzyl Alcohol	1.110	1.081	2.6	102	0.00
16	2,2'-oxybis(1-Chloropropane	1.909	1.835	3.9	106	0.00
17	2-Methylphenol	1.002	0.970	3.2	106	0.01
18	Hexachloroethane	0.571	0.572	-0.2	109	0.00
19 P	n-Nitroso-di-n-propylamine	0.913	0.900	1.4	110	0.00
20	3+4-Methylphenols	1.331	1.300	2.3	106	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	105	0.01
22	Acetophenone	0.490	0.491	-0.2	107	0.01
23 S	Nitrobenzene-d5	0.382	0.382	0.0	104	0.00
24	Nitrobenzene	0.370	0.358	3.2	99	0.00
25	Isophorone	0.653	0.635	2.8	102	0.01
26 C	2-Nitrophenol	0.188	0.191	-1.6	103	0.00
27	2,4-Dimethylphenol	0.266	0.270	-1.5	103	0.00
28	bis(2-Chloroethoxy)methane	0.371	0.364	1.9	103	0.00
29 C	2,4-Dichlorophenol	0.297	0.302	-1.7	105	0.00
30	1,2,4-Trichlorobenzene	0.332	0.338	-1.8	105	0.01
31	Naphthalene	1.028	1.049	-2.0	106	0.00
32	Benzoic acid	0.138	0.163	-18.1	122	0.00
33	4-Chloroaniline	0.383	0.387	-1.0	103	0.01
34 C	Hexachlorobutadiene	0.199	0.203	-2.0	104	0.01
35	Caprolactam	0.079	0.080	-1.3	97	0.01
36 C	4-Chloro-3-methylphenol	0.310	0.318	-2.6	107	0.00
37	2-Methylnaphthalene	0.691	0.696	-0.7	106	0.01
38	1-Methylnaphthalene	0.646	0.645	0.2	105	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
40	1,2,4,5-Tetrachlorobenzene	0.574	0.587	-2.3	105	0.01
41 P	Hexachlorocyclopentadiene	0.139	0.194	-39.6#	96	0.00
42 S	2,4,6-Tribromophenol	0.179	0.189	-5.6	105	0.00
43 C	2,4,6-Trichlorophenol	0.382	0.399	-4.5	109	0.00
44	2,4,5-Trichlorophenol	0.387	0.420	-8.5	105	0.00
45 S	2-Fluorobiphenyl	1.367	1.426	-4.3	107	0.00
46	1,1'-Biphenyl	1.490	1.585	-6.4	110	0.01
47	2-Chloronaphthalene	1.186	1.218	-2.7	106	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080721\
 Data File : BF124975.D
 Acq On : 7 Aug 2021 15:21
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 07 16:03:39 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 06 11:27:24 2021
 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)
48	2-Nitroaniline	0.348	0.346	0.6	102	0.00
49	Acenaphthylene	1.801	1.890	-4.9	111	0.01
50	Dimethylphthalate	1.355	1.403	-3.5	107	0.00
51	2,6-Dinitrotoluene	0.302	0.305	-1.0	102	0.01
52 C	Acenaphthene	1.092	1.107	-1.4	106	0.01
53	3-Nitroaniline	0.318	0.326	-2.5	108	0.01
54 P	2,4-Dinitrophenol	0.132	0.139	-5.3	106	0.00
55	Dibenzofuran	1.706	1.778	-4.2	109	0.01
56 P	4-Nitrophenol	0.210	0.218	-3.8	102	0.00
57	2,4-Dinitrotoluene	0.410	0.448	-9.3	111	0.00
58	Fluorene	1.308	1.342	-2.6	106	0.00
59	2,3,4,6-Tetrachlorophenol	0.290	0.314	-8.3	108	0.00
60	Diethylphthalate	1.391	1.419	-2.0	108	0.00
61	4-Chlorophenyl-phenylether	0.618	0.652	-5.5	107	0.01
62	4-Nitroaniline	0.309	0.308	0.3	100	0.00
63	Azobenzene	1.360	1.383	-1.7	106	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
65	4,6-Dinitro-2-methylphenol	0.125	0.131	-4.8	110	0.00
66 c	n-Nitrosodiphenylamine	0.648	0.650	-0.3	102	0.00
67	4-Bromophenyl-phenylether	0.219	0.231	-5.5	108	0.01
68	Hexachlorobenzene	0.238	0.241	-1.3	105	0.00
69	Atrazine	0.195	0.197	-1.0	104	0.00
70 C	Pentachlorophenol	0.102	0.111	-8.8	108	0.00
71	Phenanthrene	1.085	1.099	-1.3	104	0.01
72	Anthracene	1.088	1.108	-1.8	106	0.00
73	Carbazole	0.998	1.004	-0.6	105	0.00
74	Di-n-butylphthalate	1.268	1.339	-5.6	111	0.00
75 C	Fluoranthene	1.122	1.144	-2.0	106	0.01
76 I	Chrysene-d12	1.000	1.000	0.0	101	0.00
77	Benzydine	0.535	0.483	9.7	96	0.00
78	Pyrene	1.575	1.631	-3.6	103	0.00
79 S	Terphenyl-d14	1.186	1.240	-4.6	107	0.01
80	Butylbenzylphthalate	0.701	0.725	-3.4	105	0.01
81	Benzo(a)anthracene	1.303	1.276	2.1	100	0.00
82	3,3'-Dichlorobenzidine	0.344	0.341	0.9	97	0.01
83	Chrysene	1.260	1.264	-0.3	102	0.01
84	Bis(2-ethylhexyl)phthalate	0.973	1.037	-6.6	104	0.00
85 c	Di-n-octyl phthalate	1.534	1.641	-7.0	108	0.00
86 I	Perylene-d12	1.000	1.000	0.0	96	0.00
87	Indeno(1,2,3-cd)pyrene	1.191	1.207	-1.3	99	0.01
88	Benzo(b)fluoranthene	1.404	1.519	-8.2	99	0.00
89	Benzo(k)fluoranthene	1.277	1.353	-6.0	99	0.00
90 C	Benzo(a)pyrene	1.196	1.231	-2.9	97	0.01
91	Dibenzo(a,h)anthracene	1.047	1.040	0.7	95	0.01
92	Benzo(g,h,i)perylene	1.004	0.973	3.1	96	0.02

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080721\
Data File : BF124975.D
Acq On : 7 Aug 2021 15:21
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 07 16:03:39 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 06 11:27:24 2021
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)
----------	-------	------	-----------	------------

(#) = Out of Range

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