

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100223\
 Data File : BF135524.D
 Acq On : 02 Oct 2023 13:23
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Oct 03 02:53:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 30 00:12:20 2023
 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	120	0.00
2	1,4-Dioxane	0.539	0.618	-14.7	139	-0.02
3	Pyridine	1.451	1.653	-13.9	134	-0.02
4	n-Nitrosodimethylamine	0.770	0.805	-4.5	123	-0.03
5 S	2-Fluorophenol	1.230	1.261	-2.5	124	0.00
6	Aniline	2.050	1.973	3.8	116	-0.01
7 S	Phenol-d6	1.564	1.541	1.5	118	-0.01
8	2-Chlorophenol	1.313	1.313	0.0	120	0.00
9	Benzaldehyde	0.891	0.887	0.4	122	0.00
10 C	Phenol	1.715	1.683	1.9	117	0.00
11	bis(2-Chloroethyl)ether	1.286	1.313	-2.1	122	0.00
12	1,3-Dichlorobenzene	1.334	1.287	3.5	116	0.00
13 C	1,4-Dichlorobenzene	1.357	1.298	4.3	116	0.00
14	1,2-Dichlorobenzene	1.260	1.203	4.5	117	0.00
15	Benzyl Alcohol	1.122	0.999	11.0	106	0.00
16	2,2'-oxybis(1-Chloropropane	2.424	2.309	4.7	113	0.00
17	2-Methylphenol	1.158	1.104	4.7	113	0.00
18	Hexachloroethane	0.502	0.476	5.2	114	0.00
19 P	n-Nitroso-di-n-propylamine	0.968	0.886	8.5	112	-0.01
20	3+4-Methylphenols	1.393	1.389	0.3	122	-0.01
21 I	Naphthalene-d8	1.000	1.000	0.0	111	0.00
22	Acetophenone	0.457	0.463	-1.3	116	0.00
23 S	Nitrobenzene-d5	0.368	0.376	-2.2	114	-0.01
24	Nitrobenzene	0.364	0.386	-6.0	118	-0.01
25	Isophorone	0.676	0.688	-1.8	115	0.00
26 C	2-Nitrophenol	0.175	0.185	-5.7	118	0.00
27	2,4-Dimethylphenol	0.294	0.304	-3.4	118	0.00
28	bis(2-Chloroethoxy)methane	0.405	0.419	-3.5	119	0.00
29 C	2,4-Dichlorophenol	0.272	0.273	-0.4	113	0.00
30	1,2,4-Trichlorobenzene	0.284	0.277	2.5	110	0.00
31	Naphthalene	0.972	0.971	0.1	113	0.00
32	Benzoic acid	0.221	0.206	6.8	99	-0.02
33	4-Chloroaniline	0.426	0.434	-1.9	115	0.00
34 C	Hexachlorobutadiene	0.171	0.155	9.4	102	0.00
35	Caprolactam	0.091	0.098	-7.7	122	-0.02
36 C	4-Chloro-3-methylphenol	0.302	0.307	-1.7	114	0.00
37	2-Methylnaphthalene	0.653	0.636	2.6	112	0.00
38	1-Methylnaphthalene	0.612	0.596	2.6	112	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
40	1,2,4,5-Tetrachlorobenzene	0.579	0.559	3.5	106	0.00
41 P	Hexachlorocyclopentadiene	0.210	0.232	-10.5	115	0.00
42 S	2,4,6-Tribromophenol	0.199	0.164	17.6	92	0.00
43 C	2,4,6-Trichlorophenol	0.379	0.370	2.4	107	0.00
44	2,4,5-Trichlorophenol	0.437	0.421	3.7	105	0.00
45 S	2-Fluorobiphenyl	1.326	1.218	8.1	105	0.00
46	1,1'-Biphenyl	1.537	1.524	0.8	109	0.00
47	2-Chloronaphthalene	1.115	1.110	0.4	111	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100223\
 Data File : BF135524.D
 Acq On : 02 Oct 2023 13:23
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Oct 03 02:53:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 30 00:12:20 2023
 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	0.433	0.470	-8.5	118	0.00
49	Acenaphthylene	1.853	1.795	3.1	107	0.00
50	Dimethylphthalate	1.352	1.316	2.7	110	0.00
51	2,6-Dinitrotoluene	0.296	0.300	-1.4	111	-0.01
52 C	Acenaphthene	1.095	1.089	0.5	112	0.00
53	3-Nitroaniline	0.348	0.377	-8.3	119	0.00
54 P	2,4-Dinitrophenol	0.143	0.192	-34.3#	144	0.00
55	Dibenzofuran	1.671	1.567	6.2	105	0.00
56 P	4-Nitrophenol	0.221	0.250	-13.1	119	0.00
57	2,4-Dinitrotoluene	0.371	0.354	4.6	103	0.00
58	Fluorene	1.284	1.264	1.6	111	0.00
59	2,3,4,6-Tetrachlorophenol	0.336	0.304	9.5	100	0.00
60	Diethylphthalate	1.357	1.328	2.1	109	0.00
61	4-Chlorophenyl-phenylether	0.617	0.587	4.9	108	0.00
62	4-Nitroaniline	0.334	0.370	-10.8	120	-0.01
63	Azobenzene	1.328	1.390	-4.7	116	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
65	4,6-Dinitro-2-methylphenol	0.106	0.102	3.8	103	0.00
66 c	n-Nitrosodiphenylamine	0.605	0.610	-0.8	110	0.00
67	4-Bromophenyl-phenylether	0.199	0.189	5.0	105	0.00
68	Hexachlorobenzene	0.203	0.189	6.9	101	0.00
69	Atrazine	0.163	0.154	5.5	104	0.00
70 C	Pentachlorophenol	0.121	0.117	3.3	97	0.00
71	Phenanthrene	0.946	0.969	-2.4	111	0.00
72	Anthracene	0.972	1.000	-2.9	113	0.00
73	Carbazole	0.885	0.944	-6.7	115	0.00
74	Di-n-butylphthalate	1.043	1.086	-4.1	112	0.00
75 C	Fluoranthene	0.986	1.017	-3.1	112	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	106	0.00
77	Benzydine	0.410	0.357	12.9	103	0.00
78	Pyrene	1.904	1.971	-3.5	111	0.00
79 S	Terphenyl-d14	1.290	1.269	1.6	107	0.00
80	Butylbenzylphthalate	0.670	0.762	-13.7	118	0.00
81	Benzo(a)anthracene	1.291	1.304	-1.0	109	0.00
82	3,3'-Dichlorobenzidine	0.403	0.409	-1.5	108	0.00
83	Chrysene	1.284	1.283	0.1	108	0.00
84	Bis(2-ethylhexyl)phthalate	0.688	0.866	-25.9#	132	0.00
85 c	Di-n-octyl phthalate	1.093	1.211	-10.8	115	0.00
86 I	Perylene-d12	1.000	1.000	0.0	110	0.00
87	Indeno(1,2,3-cd)pyrene	1.311	1.235	5.8	99	0.00
88	Benzo(b)fluoranthene	1.083	1.075	0.7	110	0.00
89	Benzo(k)fluoranthene	1.069	0.977	8.6	102	0.00
90 C	Benzo(a)pyrene	1.032	1.026	0.6	108	0.00
91	Dibenzo(a,h)anthracene	1.073	0.998	7.0	99	0.00
92	Benzo(g,h,i)perylene	1.121	1.046	6.7	98	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100223\
Data File : BF135524.D
Acq On : 02 Oct 2023 13:23
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Oct 03 02:53:53 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Sep 30 00:12:20 2023
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