

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010824\
 Data File : BF137000.D
 Acq On : 08 Jan 2024 17:55
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 09 00:41:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 21 01:54:25 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	95	0.00
2	1,4-Dioxane	0.534	0.526	1.5	96	-0.02
3	Pyridine	1.303	1.257	3.5	91	-0.01
4	n-Nitrosodimethylamine	0.696	0.642	7.8	85	0.00
5 S	2-Fluorophenol	1.282	1.290	-0.6	95	0.00
6	Aniline	1.661	1.549	6.7	87	0.00
7 S	Phenol-d6	1.661	1.622	2.3	92	0.00
8	2-Chlorophenol	1.403	1.406	-0.2	94	0.00
9	Benzaldehyde	0.945	0.883	6.6	90	0.00
10 C	Phenol	1.773	1.699	4.2	91	0.00
11	bis(2-Chloroethyl)ether	1.348	1.306	3.1	91	0.00
12	1,3-Dichlorobenzene	1.530	1.555	-1.6	96	0.00
13 C	1,4-Dichlorobenzene	1.564	1.545	1.2	93	0.00
14	1,2-Dichlorobenzene	1.455	1.469	-1.0	95	0.00
15	Benzyl Alcohol	1.121	1.063	5.2	88	0.00
16	2,2'-oxybis(1-Chloropropane	2.205	2.058	6.7	87	0.00
17	2-Methylphenol	1.162	1.137	2.2	93	0.00
18	Hexachloroethane	0.568	0.560	1.4	93	-0.01
19 P	n-Nitroso-di-n-propylamine	1.007	0.946	6.1	90	0.00
20	3+4-Methylphenols	1.452	1.451	0.1	95	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
22	Acetophenone	0.501	0.504	-0.6	92	0.00
23 S	Nitrobenzene-d5	0.391	0.388	0.8	89	0.00
24	Nitrobenzene	0.392	0.391	0.3	89	0.00
25	Isophorone	0.711	0.686	3.5	88	0.00
26 C	2-Nitrophenol	0.187	0.197	-5.3	93	0.00
27	2,4-Dimethylphenol	0.228	0.232	-1.8	92	0.00
28	bis(2-Chloroethoxy)methane	0.432	0.423	2.1	88	0.00
29 C	2,4-Dichlorophenol	0.294	0.296	-0.7	91	0.00
30	1,2,4-Trichlorobenzene	0.327	0.332	-1.5	92	-0.01
31	Naphthalene	1.099	1.097	0.2	91	0.00
32	Benzoic acid	0.221	0.206	6.8	78	0.01
33	4-Chloroaniline	0.406	0.392	3.4	88	0.00
34 C	Hexachlorobutadiene	0.199	0.198	0.5	91	-0.01
35	Caprolactam	0.097	0.099	-2.1	91	0.00
36 C	4-Chloro-3-methylphenol	0.331	0.319	3.6	87	0.00
37	2-Methylnaphthalene	0.707	0.706	0.1	91	-0.01
38	1-Methylnaphthalene	0.694	0.684	1.4	91	-0.01
39 I	Acenaphthene-d10	1.000	1.000	0.0	87	0.00
40	1,2,4,5-Tetrachlorobenzene	0.619	0.649	-4.8	89	0.00
41 P	Hexachlorocyclopentadiene	0.172	0.184	-7.0	88	-0.01
42 S	2,4,6-Tribromophenol	0.236	0.225	4.7	82	0.00
43 C	2,4,6-Trichlorophenol	0.397	0.389	2.0	83	0.00
44	2,4,5-Trichlorophenol	0.443	0.438	1.1	84	0.00
45 S	2-Fluorobiphenyl	1.487	1.524	-2.5	90	-0.01
46	1,1'-Biphenyl	1.678	1.742	-3.8	91	0.00
47	2-Chloronaphthalene	1.276	1.317	-3.2	89	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010824\
 Data File : BF137000.D
 Acq On : 08 Jan 2024 17:55
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 09 00:41:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 21 01:54:25 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	0.400	0.392	2.0	84	0.00
49	Acenaphthylene	1.950	1.939	0.6	85	-0.01
50	Dimethylphthalate	1.461	1.420	2.8	85	-0.01
51	2,6-Dinitrotoluene	0.332	0.334	-0.6	86	0.00
52 C	Acenaphthene	1.209	1.225	-1.3	88	0.00
53	3-Nitroaniline	0.351	0.341	2.8	82	0.00
54 P	2,4-Dinitrophenol	0.165	0.196	-18.8	93	0.00
55	Dibenzofuran	1.832	1.795	2.0	85	0.00
56 P	4-Nitrophenol	0.237	0.296	-24.9	102	0.01
57	2,4-Dinitrotoluene	0.430	0.413	4.0	81	0.00
58	Fluorene	1.454	1.460	-0.4	88	0.00
59	2,3,4,6-Tetrachlorophenol	0.341	0.296	13.2	73	0.00
60	Diethylphthalate	1.516	1.451	4.3	82	-0.01
61	4-Chlorophenyl-phenylether	0.707	0.696	1.6	86	0.00
62	4-Nitroaniline	0.339	0.311	8.3	81	0.00
63	Azobenzene	1.416	1.344	5.1	83	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	84	0.00
65	4,6-Dinitro-2-methylphenol	0.115	0.115	0.0	75	0.00
66 c	n-Nitrosodiphenylamine	0.650	0.675	-3.8	84	0.00
67	4-Bromophenyl-phenylether	0.222	0.228	-2.7	84	0.00
68	Hexachlorobenzene	0.248	0.258	-4.0	85	0.00
69	Atrazine	0.166	0.155	6.6	91	0.00
70 C	Pentachlorophenol	0.115	0.105	8.7	68	0.00
71	Phenanthrene	1.026	1.045	-1.9	83	0.00
72	Anthracene	1.039	1.057	-1.7	84	0.00
73	Carbazole	0.980	0.977	0.3	82	0.00
74	Di-n-butylphthalate	1.196	1.188	0.7	82	0.00
75 C	Fluoranthene	1.098	1.065	3.0	80	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	78	0.00
77	Benzydine	0.590	0.617	-4.6	86	0.00
78	Pyrene	1.963	2.121	-8.0	81	0.00
79 S	Terphenyl-d14	1.431	1.527	-6.7	80	0.00
80	Butylbenzylphthalate	0.714	0.759	-6.3	79	0.00
81	Benzo(a)anthracene	1.389	1.416	-1.9	78	0.00
82	3,3'-Dichlorobenzidine	0.394	0.412	-4.6	82	0.00
83	Chrysene	1.358	1.357	0.1	77	0.00
84	Bis(2-ethylhexyl)phthalate	0.899	0.982	-9.2	81	0.00
85 c	Di-n-octyl phthalate	1.390	1.444	-3.9	78	0.00
86 I	Perylene-d12	1.000	1.000	0.0	90	0.01
87	Indeno(1,2,3-cd)pyrene	1.444	1.588	-10.0	94	0.03
88	Benzo(b)fluoranthene	1.336	1.236	7.5	84	0.00
89	Benzo(k)fluoranthene	1.262	1.203	4.7	83	0.00
90 C	Benzo(a)pyrene	1.128	1.124	0.4	88	0.00
91	Dibenzo(a,h)anthracene	1.185	1.318	-11.2	93	0.02
92	Benzo(g,h,i)perylene	1.213	1.357	-11.9	95	0.03

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010824\
Data File : BF137000.D
Acq On : 08 Jan 2024 17:55
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jan 09 00:41:16 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 21 01:54:25 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