

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100824\
 Data File : BF139701.D
 Acq On : 08 Oct 2024 09:39
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Oct 08 11:37:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 2024
 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	82	0.00
2	1,4-Dioxane	0.499	0.483	3.2	81	-0.02
3	Pyridine	1.213	1.233	-1.6	87	-0.01
4	n-Nitrosodimethylamine	0.794	0.764	3.8	80	-0.02
5 S	2-Fluorophenol	1.154	1.159	-0.4	85	0.00
6	Aniline	1.377	1.240	9.9	79	0.00
7 S	Phenol-d6	1.552	1.556	-0.3	85	0.00
8	2-Chlorophenol	1.236	1.238	-0.2	85	0.00
9	Benzaldehyde	0.822	0.781	5.0	81	0.00
10 C	Phenol	1.632	1.649	-1.0	84	0.00
11	bis(2-Chloroethyl)ether	1.329	1.224	7.9	80	0.00
12	1,3-Dichlorobenzene	1.391	1.370	1.5	83	0.00
13 C	1,4-Dichlorobenzene	1.409	1.378	2.2	84	0.00
14	1,2-Dichlorobenzene	1.320	1.289	2.3	83	0.00
15	Benzyl Alcohol	1.186	1.162	2.0	81	0.00
16	2,2'-oxybis(1-Chloropropane	2.300	2.156	6.3	79	0.00
17	2-Methylphenol	1.032	1.016	1.6	82	0.00
18	Hexachloroethane	0.525	0.503	4.2	81	0.00
19 P	n-Nitroso-di-n-propylamine	0.998	0.936	6.2	80	0.00
20	3+4-Methylphenols	1.296	1.284	0.9	84	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	82	0.00
22	Acetophenone	0.475	0.455	4.2	83	0.00
23 S	Nitrobenzene-d5	0.395	0.382	3.3	82	0.00
24	Nitrobenzene	0.409	0.396	3.2	82	0.00
25	Isophorone	0.702	0.659	6.1	80	0.00
26 C	2-Nitrophenol	0.176	0.175	0.6	80	0.00
27	2,4-Dimethylphenol	0.215	0.208	3.3	83	0.00
28	bis(2-Chloroethoxy)methane	0.419	0.393	6.2	80	0.00
29 C	2,4-Dichlorophenol	0.281	0.278	1.1	83	0.00
30	1,2,4-Trichlorobenzene	0.318	0.306	3.8	82	0.00
31	Naphthalene	1.023	0.997	2.5	83	0.00
32	Benzoic acid	0.214	0.225	-5.1	82	0.00
33	4-Chloroaniline	0.346	0.322	6.9	77	0.00
34 C	Hexachlorobutadiene	0.205	0.194	5.4	81	0.00
35	Caprolactam	0.092	0.094	-2.2	87	0.00
36 C	4-Chloro-3-methylphenol	0.318	0.310	2.5	84	0.00
37	2-Methylnaphthalene	0.666	0.630	5.4	81	0.00
38	1-Methylnaphthalene	0.651	0.623	4.3	82	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	83	0.00
40	1,2,4,5-Tetrachlorobenzene	0.555	0.527	5.0	82	0.00
41 P	Hexachlorocyclopentadiene	0.170	0.141	17.1	64	0.00
42 S	2,4,6-Tribromophenol	0.193	0.183	5.2	81	0.00
43 C	2,4,6-Trichlorophenol	0.374	0.350	6.4	79	0.00
44	2,4,5-Trichlorophenol	0.404	0.389	3.7	83	0.00
45 S	2-Fluorobiphenyl	1.303	1.211	7.1	82	0.00
46	1,1'-Biphenyl	1.488	1.403	5.7	82	0.00
47	2-Chloronaphthalene	1.115	1.061	4.8	82	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100824\
 Data File : BF139701.D
 Acq On : 08 Oct 2024 09:39
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Oct 08 11:37:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 2024
 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	0.396	1.0	84	0.00
49	Acenaphthylene	1.696	1.622	4.4	83	0.00
50	Dimethylphthalate	1.309	1.237	5.5	82	0.00
51	2,6-Dinitrotoluene	0.296	0.293	1.0	84	0.00
52 C	Acenaphthene	1.164	1.106	5.0	82	0.00
53	3-Nitroaniline	0.302	0.296	2.0	84	0.00
54 P	2,4-Dinitrophenol	0.159	0.159	0.0	81	0.00
55	Dibenzofuran	1.630	1.554	4.7	83	0.00
56 P	4-Nitrophenol	0.230	0.217	5.7	79	0.01
57	2,4-Dinitrotoluene	0.387	0.385	0.5	84	0.00
58	Fluorene	1.298	1.253	3.5	85	0.00
59	2,3,4,6-Tetrachlorophenol	0.326	0.317	2.8	84	0.00
60	Diethylphthalate	1.352	1.275	5.7	82	0.00
61	4-Chlorophenyl-phenylether	0.650	0.611	6.0	84	0.00
62	4-Nitroaniline	0.309	0.304	1.6	85	0.00
63	Azobenzene	1.360	1.269	6.7	80	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
65	4,6-Dinitro-2-methylphenol	0.113	0.115	-1.8	83	0.00
66 c	n-Nitrosodiphenylamine	0.590	0.555	5.9	84	0.00
67	4-Bromophenyl-phenylether	0.205	0.189	7.8	82	0.00
68	Hexachlorobenzene	0.228	0.212	7.0	83	0.00
69	Atrazine	0.158	0.114	27.8#	74	0.00
70 C	Pentachlorophenol	0.135	0.127	5.9	78	0.00
71	Phenanthrene	0.943	0.898	4.8	85	0.00
72	Anthracene	0.933	0.897	3.9	86	0.00
73	Carbazole	0.896	0.890	0.7	88	0.00
74	Di-n-butylphthalate	1.090	1.048	3.9	84	0.00
75 C	Fluoranthene	1.031	1.016	1.5	88	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	90	0.00
77	Benzydine	0.353	0.281	20.4	73	0.00
78	Pyrene	1.789	1.775	0.8	88	0.00
79 S	Terphenyl-d14	1.219	1.190	2.4	89	0.00
80	Butylbenzylphthalate	0.640	0.662	-3.4	92	0.00
81	Benzo(a)anthracene	1.315	1.265	3.8	91	0.00
82	3,3'-Dichlorobenzidine	0.402	0.364	9.5	83	0.00
83	Chrysene	1.202	1.177	2.1	89	0.00
84	Bis(2-ethylhexyl)phthalate	0.799	0.821	-2.8	91	0.00
85 c	Di-n-octyl phthalate	1.174	1.104	6.0	84	0.00
86 I	Perylene-d12	1.000	1.000	0.0	76	0.00
87	Indeno(1,2,3-cd)pyrene	1.177	1.190	-1.1	75	0.01
88	Benzo(b)fluoranthene	1.293	1.195	7.6	79	0.00
89	Benzo(k)fluoranthene	1.097	1.144	-4.3	77	0.00
90 C	Benzo(a)pyrene	1.020	0.998	2.2	77	0.00
91	Dibenzo(a,h)anthracene	0.976	0.984	-0.8	74	0.01
92	Benzo(g,h,i)perylene	0.979	1.024	-4.6	77	0.01

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100824\
Data File : BF139701.D
Acq On : 08 Oct 2024 09:39
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Oct 08 11:37:13 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 02 04:58:47 2024
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