Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020524\
Data File : BF137269.D
Acq On : 05 Feb 2024 10:37
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Feb 05 11:03:14 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 31 04:56:22 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	98	0.00
2		1,4-Dioxane	0.649	0.616	5.1	90	0.00
3		Pyridine	1.129	1.338	-18.5	101	0.00
4		n-Nitrosodimethylamine	0.591	0.633	-7.1	101	0.00
5	S	2-Fluorophenol	1.212	1.337	-10.3	98	0.00
6		Aniline	1.852	1.903	-2.8	91	0.00
7	S	Phenol-d6	1.730	1.776	-2.7	97	0.00
8		2-Chlorophenol	1.355	1.447	-6.8	100	0.00
9		Benzaldehyde	0.659	0.604	8.3	78	0.00
10	C	Phenol	1.931	1.993	-3.2	96	0.00
11		bis(2-Chloroethyl)ether	1.290	1.454	-12.7	107	0.00
12		1,3-Dichlorobenzene	1.589	1.637	-3.0	98	0.00
13	C	1,4-Dichlorobenzene	1.643	1.703	-3.7	99	0.00
14		1,2-Dichlorobenzene	1.536	1.600	-4.2	100	0.00
15		Benzyl Alcohol	1.229	1.377	-12.0	98	0.00
16		2,2'-oxybis(1-Chloropropane	2.324	2.158	7.1	89	0.00
17		2-Methylphenol	1.132	1.218	-7.6	102	0.00
18		Hexachloroethane	0.593	0.609	-2.7	99	0.00
19	P	n-Nitroso-di-n-propylamine	1.186	1.182	0.3	95	0.00
20		3+4-Methylphenols	1.433	1.492	-4.1	96	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
22		Acetophenone	0.538	0.555	-3.2	96	0.00
23	S	Nitrobenzene-d5	0.447	0.463	-3.6	95	0.00
24		Nitrobenzene	0.434	0.451	-3.9	94	0.00
25		Isophorone	0.825	0.823	0.2	95	0.00
26	C	2-Nitrophenol	0.149	0.184	-23.5#	105	0.00
27		2,4-Dimethylphenol	0.236	0.246	-4.2	95	0.00
28		bis(2-Chloroethoxy)methane	0.438	0.455	-3.9	95	0.00
29	C	2,4-Dichlorophenol	0.292	0.317	-8.6	97	0.00
30		1,2,4-Trichlorobenzene	0.363	0.370	-1.9	96	0.00
31		Naphthalene	1.081	1.118	-3.4	97	0.00
32		Benzoic acid	0.140	0.181	-29.3#	117	0.00
33		4-Chloroaniline	0.372	0.394	-5.9	95	0.00
34	C	Hexachlorobutadiene	0.241	0.246	-2.1	96	0.00
35		Caprolactam	0.091	0.100	-9.9	96	0.00
36	C	4-Chloro-3-methylphenol	0.345	0.358	-3.8	96	0.00
37		2-Methylnaphthalene	0.725	0.733	-1.1	95	0.00
38		1-Methylnaphthalene	0.673	0.688	-2.2	95	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
40		1,2,4,5-Tetrachlorobenzene	0.669	0.692	-3.4	94	0.00
41	P	Hexachlorocyclopentadiene	0.319	0.363	-13.8	98	0.00
42	S	2,4,6-Tribromophenol	0.226	0.252	-11.5	99	0.00
43	C	2,4,6-Trichlorophenol	0.380	0.427	-12.4	98	0.00
44		2,4,5-Trichlorophenol	0.415	0.460	-10.8	98	0.00
45	S	2-Fluorobiphenyl	1.455	1.486	-2.1	96	0.00
46		1,1'-Biphenyl	1.591	1.628	-2.3	95	0.00
47		2-Chloronaphthalene	1.178	1.195	-1.4	94	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020524\
Data File : BF137269.D
Acq On : 05 Feb 2024 10:37
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Feb 05 11:03:14 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 31 04:56:22 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	0.381	0.417	-9.4	92	0.00
49		Acenaphthylene	1.885	1.918	-1.8	94	0.00
50		Dimethylphthalate	1.351	1.419	-5.0	96	0.00
51		2,6-Dinitrotoluene	0.288	0.317	-10.1	97	0.00
52	C	Acenaphthene	1.108	1.140	-2.9	95	0.00
53		3-Nitroaniline	0.259	0.301	-16.2	103	0.00
54	P	2,4-Dinitrophenol	0.089	0.139	-56.2#	130	0.00
55		Dibenzofuran	1.763	1.825	-3.5	95	0.00
56	P	4-Nitrophenol	0.169	0.195	-15.4	91	0.00
57		2,4-Dinitrotoluene	0.374	0.428	-14.4	96	0.00
58		Fluorene	1.382	1.428	-3.3	96	0.00
59		2,3,4,6-Tetrachlorophenol	0.386	0.416	-7.8	98	0.00
60		Diethylphthalate	1.330	1.412	-6.2	96	0.00
61		4-Chlorophenyl-phenylether	0.687	0.705	-2.6	95	0.00
62		4-Nitroaniline	0.248	0.263	-6.0	88	0.00
63		Azobenzene	1.473	1.507	-2.3	94	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	96	0.00
65		4,6-Dinitro-2-methylphenol	0.095	0.121	-27.4#	111	0.00
66	c	n-Nitrosodiphenylamine	0.608	0.620	-2.0	95	0.00
67		4-Bromophenyl-phenylether	0.220	0.232	-5.5	97	0.00
68		Hexachlorobenzene	0.246	0.249	-1.2	95	0.00
69		Atrazine	0.201	0.213	-6.0	95	0.00
70	C	Pentachlorophenol	0.141	0.159	-12.8	103	0.00
71		Phenanthrene	1.047	1.074	-2.6	96	0.00
72		Anthracene	1.091	1.108	-1.6	95	0.00
73		Carbazole	0.912	0.957	-4.9	96	0.00
74		Di-n-butylphthalate	0.986	1.105	-12.1	102	0.00
75	C	Fluoranthene	1.096	1.140	-4.0	95	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	98	0.00
77		Benzidine	0.358	0.122	65.9#	28#	0.00
78		Pyrene	1.922	2.003	-4.2	96	0.00
79	S	Terphenyl-d14	1.471	1.573	-6.9	99	0.00
80		Butylbenzylphthalate	0.492	0.656	-33.3#	116	0.00
81		Benzo(a)anthracene	1.416	1.411	0.4	95	0.00
82		3,3'-Dichlorobenzidine	0.386	0.409	-6.0	100	0.00
83		Chrysene	1.393	1.415	-1.6	96	0.00
84		Bis(2-ethylhexyl)phthalate	0.572	0.780	-36.4#	121	0.00
85	c	Di-n-octyl phthalate	0.910	1.166	-28.1#	122	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	97	0.00
87		Indeno(1,2,3-cd)pyrene	1.405	1.608	-14.4	102	0.00
88		Benzo(b)fluoranthene	1.230	1.244	-1.1	101	0.00
89		Benzo(k)fluoranthene	1.301	1.302	-0.1	95	0.00
90	C	Benzo(a)pyrene	1.157	1.195	-3.3	99	0.00
91		Dibenzo(a,h)anthracene	1.126	1.270	-12.8	102	0.00
92		Benzo(g,h,i)perylene	1.105	1.256	-13.7	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020524\
Data File : BF137269.D
Acq On : 05 Feb 2024 10:37
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Feb 05 11:03:14 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 31 04:56:22 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 = 2