Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041023\
Data File : BM039371.D
Acq On : 10 Apr 2023 10:53
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Apr 11 02:07:19 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 11 02:05:37 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	89	0.00
2		1,4-Dioxane	0.543	0.640	-17.9	106	0.00
3		Pyridine	1.520	1.717	-13.0	99	0.00
4		n-Nitrosodimethylamine	0.597	0.669	-12.1	100	0.00
5	S	2-Fluorophenol	1.327	1.441	-8.6	97	0.00
6		Aniline	2.125	2.267	-6.7	94	0.00
7	S	Phenol-d6	1.717	1.848	-7.6	95	0.01
8		2-Chlorophenol	1.359	1.445	-6.3	95	0.00
9		Benzaldehyde	0.978	0.861	12.0	78	0.00
10	C	Phenol	1.721	1.844	-7.1	96	0.00
11		bis(2-Chloroethyl)ether	1.384	1.538	-11.1	99	0.01
12		1,3-Dichlorobenzene	1.441	1.527	-6.0	95	0.00
13	C	1,4-Dichlorobenzene	1.457	1.523	-4.5	93	0.00
14		1,2-Dichlorobenzene	1.393	1.486	-6.7	96	0.00
15		Benzyl Alcohol	1.181	1.166	1.3	87	0.01
16		2,2'-oxybis(1-Chloropropane	1.771	2.012	-13.6	102	0.00
17		2-Methylphenol	1.204	1.262	-4.8	93	0.00
18		Hexachloroethane	0.564	0.587	-4.1	94	0.00
19	P	n-Nitroso-di-n-propylamine	1.098	1.174	-6.9	96	0.00
20		3+4-Methylphenols	1.618	1.689	-4.4	92	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	87	0.00
22		Acetophenone	0.535	0.570	-6.5	93	0.00
23	S	Nitrobenzene-d5	0.412	0.450	-9.2	95	0.00
24		Nitrobenzene	0.409	0.448	-9.5	95	0.00
25		Isophorone	0.768	0.823	-7.2	93	0.00
26	C	2-Nitrophenol	0.189	0.189	0.0	85	0.00
27		2,4-Dimethylphenol	0.314	0.323	-2.9	89	0.00
28		bis(2-Chloroethoxy)methane	0.465	0.496	-6.7	93	0.00
29	C	2,4-Dichlorophenol	0.307	0.323	-5.2	90	0.00
30		1,2,4-Trichlorobenzene	0.328	0.348	-6.1	93	0.00
31		Naphthalene	1.078	1.143	-6.0	93	0.00
32		Benzoic acid	0.245	0.207	15.5	71	0.00
33		4-Chloroaniline	0.472	0.463	1.9	85	0.00
34	C	Hexachlorobutadiene	0.193	0.206	-6.7	94	0.00
35		Caprolactam	0.106	0.102	3.8	81	0.00
36	C	4-Chloro-3-methylphenol	0.338	0.343	-1.5	87	0.00
37		2-Methylnaphthalene	0.750	0.783	-4.4	91	0.00
38		1-Methylnaphthalene	0.702	0.731	-4.1	91	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	86	0.00
40		1,2,4,5-Tetrachlorobenzene	0.609	0.656	-7.7	92	0.00
41	P	Hexachlorocyclopentadiene	0.333	0.342	-2.7	83	0.00
42	S	2,4,6-Tribromophenol	0.254	0.266	-4.7	88	0.00
43	C	2,4,6-Trichlorophenol	0.415	0.435	-4.8	87	0.00
44		2,4,5-Trichlorophenol	0.487	0.513	-5.3	87	0.00
45	S	2-Fluorobiphenyl	1.476	1.604	-8.7	93	0.00
46		1,1'-Biphenyl	1.579	1.689	-7.0	91	0.00
47		2-Chloronaphthalene	1.196	1.259	-5.3	90	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041023\
Data File : BM039371.D
Acq On : 10 Apr 2023 10:53
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Apr 11 02:07:19 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 11 02:05:37 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	0.402	0.406	-1.0	84	0.00
49		Acenaphthylene	1.965	2.080	-5.9	90	0.00
50		Dimethylphthalate	1.541	1.615	-4.8	90	0.00
51		2,6-Dinitrotoluene	0.331	0.347	-4.8	87	0.00
52	C	Acenaphthene	1.161	1.236	-6.5	91	0.00
53		3-Nitroaniline	0.383	0.352	8.1	76	0.00
54	P	2,4-Dinitrophenol	0.196	0.178	9.2	74	0.00
55		Dibenzofuran	1.868	1.960	-4.9	90	0.00
56	P	4-Nitrophenol	0.295	0.278	5.8	74	0.00
57		2,4-Dinitrotoluene	0.448	0.457	-2.0	84	0.00
58		Fluorene	1.480	1.571	-6.1	91	0.00
59		2,3,4,6-Tetrachlorophenol	0.387	0.408	-5.4	89	0.00
60		Diethylphthalate	1.551	1.631	-5.2	90	0.00
61		4-Chlorophenyl-phenylether	0.733	0.781	-6.5	92	0.00
62		4-Nitroaniline	0.393	0.327	16.8	67	0.00
63		Azobenzene	1.554	1.712	-10.2	93	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	84	0.00
65		4,6-Dinitro-2-methylphenol	0.132	0.136	-3.0	84	0.00
66	c	n-Nitrosodiphenylamine	0.630	0.670	-6.3	88	0.00
67		4-Bromophenyl-phenylether	0.214	0.225	-5.1	88	0.00
68		Hexachlorobenzene	0.232	0.247	-6.5	89	0.00
69		Atrazine	0.200	0.208	-4.0	87	0.00
70	C	Pentachlorophenol	0.166	0.181	-9.0	86	0.00
71		Phenanthrene	1.095	1.158	-5.8	88	0.00
72		Anthracene	1.117	1.206	-8.0	89	0.00
73		Carbazole	1.044	1.063	-1.8	84	0.00
74		Di-n-butylphthalate	1.291	1.417	-9.8	90	0.00
75	C	Fluoranthene	1.249	1.337	-7.0	89	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	82	0.00
77		Benzidine	0.528	0.475	10.0	69	0.00
78		Pyrene	1.424	1.540	-8.1	89	0.00
79	S	Terphenyl-d14	1.083	1.190	-9.9	90	0.00
80		Butylbenzylphthalate	0.652	0.700	-7.4	87	0.00
81		Benzo(a)anthracene	1.411	1.466	-3.9	86	0.00
82		3,3'-Dichlorobenzidine	0.470	0.468	0.4	81	0.00
83		Chrysene	1.362	1.463	-7.4	88	0.00
84		Bis(2-ethylhexyl)phthalate	0.963	1.053	-9.3	90	0.00
85	c	Di-n-octyl phthalate	1.736	1.814	-4.5	86	0.00
86	I	Perylene-d12	1.000	1.000	0.0	84	0.00
87		Indeno(1,2,3-cd)pyrene	1.474	1.636	-11.0	91	-0.01
88		Benzo(b)fluoranthene	1.224	1.232	-0.7	85	0.00
89		Benzo(k)fluoranthene	1.247	1.344	-7.8	90	0.00
90	C	Benzo(a)pyrene	1.202	1.289	-7.2	89	0.00
91		Dibenzo(a,h)anthracene	1.237	1.364	-10.3	91	0.00
92		Benzo(g,h,i)perylene	1.217	1.345	-10.5	91	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041023\
Data File : BM039371.D
Acq On : 10 Apr 2023 10:53
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Apr 11 02:07:19 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 11 02:05:37 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