Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032322\
Data File : BP009510.D
Acq On : 23 Mar 2022 17:31
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Mar 24 06:12:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 24 01:19:11 2022
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	151#	0.00
2		1,4-Dioxane	0.454	0.482	-6.2	169#	0.00
3		Pyridine	1.222	1.350	-10.5	171#	0.00
4		n-Nitrosodimethylamine	0.450	0.499	-10.9	174#	0.00
5	S	2-Fluorophenol	1.146	1.171	-2.2	160#	0.00
6		Aniline	1.766	1.887	-6.9	167#	0.00
7	S	Phenol-d6	1.549	1.631	-5.3	166#	0.00
8		2-Chlorophenol	1.306	1.326	-1.5	160#	0.00
9		Benzaldehyde	0.821	0.786	4.3	171#	0.00
10	C	Phenol	1.554	1.640	-5.5	166#	0.00
11		bis(2-Chloroethyl)ether	1.230	1.311	-6.6	168#	0.00
12		1,3-Dichlorobenzene	1.474	1.454	1.4	155#	0.00
13	C	1,4-Dichlorobenzene	1.507	1.481	1.7	155#	0.00
14		1,2-Dichlorobenzene	1.457	1.439	1.2	156#	0.00
15		Benzyl Alcohol	1.235	1.280	-3.6	162#	0.00
16		2,2'-oxybis(1-Chloropropane	1.333	1.542	-15.7	182#	0.00
17		2-Methylphenol	1.071	1.123	-4.9	165#	0.00
18		Hexachloroethane	0.528	0.520	1.5	155#	0.00
19	P	n-Nitroso-di-n-propylamine	0.980	1.093	-11.5	177#	0.00
20		3+4-Methylphenols	1.473	1.577	-7.1	168#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	163#	0.00
22		Acetophenone	0.495	0.496	-0.2	168#	0.00
23	S	Nitrobenzene-d5	0.376	0.377	-0.3	167#	0.00
24		Nitrobenzene	0.356	0.357	-0.3	168#	0.00
25		Isophorone	0.642	0.689	-7.3	181#	0.00
26	C	2-Nitrophenol	0.185	0.189	-2.2	168#	0.00
27		2,4-Dimethylphenol	0.261	0.273	-4.6	174#	0.00
28		bis(2-Chloroethoxy)methane	0.418	0.445	-6.5	178#	0.00
29	C	2,4-Dichlorophenol	0.319	0.325	-1.9	168#	0.00
30		1,2,4-Trichlorobenzene	0.370	0.355	4.1	160#	0.00
31		Naphthalene	1.030	1.015	1.5	166#	0.00
32		Benzoic acid	0.204	0.210	-2.9	167#	0.00
33		4-Chloroaniline	0.420	0.441	-5.0	176#	0.00
34	C	Hexachlorobutadiene	0.250	0.228	8.8	152#	0.00
35		Caprolactam	0.108	0.120	-11.1	193#	0.00
36	C	4-Chloro-3-methylphenol	0.343	0.356	-3.8	177#	0.00
37		2-Methylnaphthalene	0.723	0.737	-1.9	174#	0.00
38		1-Methylnaphthalene	0.704	0.721	-2.4	174#	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	177#	0.00
40		1,2,4,5-Tetrachlorobenzene	0.664	0.615	7.4	166#	0.00
41	P	Hexachlorocyclopentadiene	0.234	0.266	-13.7	194#	0.00
42	S	2,4,6-Tribromophenol	0.330	0.296	10.3	163#	0.00
43	C	2,4,6-Trichlorophenol	0.440	0.433	1.6	175#	0.00
44		2,4,5-Trichlorophenol	0.478	0.473	1.0	177#	0.00
45	S	2-Fluorobiphenyl	1.443	1.351	6.4	168#	0.00
46		1,1'-Biphenyl	1.492	1.438	3.6	174#	0.00
47		2-Chloronaphthalene	1.175	1.129	3.9	173#	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032322\
Data File : BP009510.D
Acq On : 23 Mar 2022 17:31
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Mar 24 06:12:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 24 01:19:11 2022
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	0.320	0.334	-4.4	186#	0.00
49		Acenaphthylene	1.813	1.780	1.8	178#	0.00
50		Dimethylphthalate	1.515	1.496	1.3	180#	0.00
51		2,6-Dinitrotoluene	0.318	0.320	-0.6	180#	0.00
52	C	Acenaphthene	1.083	1.065	1.7	178#	0.00
53		3-Nitroaniline	0.335	0.354	-5.7	190#	0.00
54	P	2,4-Dinitrophenol	0.174	0.190	-9.2	192#	0.00
55		Dibenzofuran	1.820	1.778	2.3	178#	0.00
56	P	4-Nitrophenol	0.250	0.265	-6.0	185#	0.00
57		2,4-Dinitrotoluene	0.437	0.446	-2.1	182#	0.00
58		Fluorene	1.431	1.395	2.5	179#	0.00
59		2,3,4,6-Tetrachlorophenol	0.433	0.426	1.6	177#	0.00
60		Diethylphthalate	1.517	1.504	0.9	181#	0.00
61		4-Chlorophenyl-phenylether	0.788	0.755	4.2	175#	0.00
62		4-Nitroaniline	0.352	0.368	-4.5	186#	0.00
63		Azobenzene	1.319	1.401	-6.2	192#	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	180#	0.00
65		4,6-Dinitro-2-methylphenol	0.129	0.136	-5.4	183#	0.00
66	c	n-Nitrosodiphenylamine	0.592	0.581	1.9	182#	0.00
67		4-Bromophenyl-phenylether	0.245	0.231	5.7	176#	0.00
68		Hexachlorobenzene	0.282	0.259	8.2	170#	0.00
69		Atrazine	0.208	0.216	-3.8	185#	0.00
70	C	Pentachlorophenol	0.151	0.156	-3.3	181#	0.00
71		Phenanthrene	1.071	1.051	1.9	184#	0.00
72		Anthracene	1.057	1.051	0.6	185#	0.00
73		Carbazole	1.034	1.042	-0.8	187#	0.00
74		Di-n-butylphthalate	1.211	1.239	-2.3	188#	0.00
75	C	Fluoranthene	1.296	1.270	2.0	183#	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	176#	0.00
77		Benzidine	0.490	0.331	32.4#	125	0.00
78		Pyrene	1.318	1.363	-3.4	184#	0.00
79	S	Terphenyl-d14	1.114	1.079	3.1	171#	0.00
80		Butylbenzylphthalate	0.560	0.593	-5.9	187#	0.00
81		Benzo(a)anthracene	1.314	1.314	0.0	181#	0.00
82		3,3'-Dichlorobenzidine	0.508	0.473	6.9	169#	0.00
83		Chrysene	1.244	1.212	2.6	176#	0.00
84		Bis(2-ethylhexyl)phthalate	0.781	0.816	-4.5	187#	0.00
85	c	Di-n-octyl phthalate	1.346	1.363	-1.3	184#	0.00
86	I	Perylene-d12	1.000	1.000	0.0	149	0.00
87		Indeno(1,2,3-cd)pyrene	1.517	1.244	18.0	125	0.00
88		Benzo(b)fluoranthene	1.194	1.243	-4.1	161#	0.00
89		Benzo(k)fluoranthene	1.169	1.257	-7.5	164#	0.00
90	C	Benzo(a)pyrene	1.022	1.030	-0.8	155#	0.00
91		Dibenzo(a,h)anthracene	1.265	1.040	17.8	125	0.00
92		Benzo(g,h,i)perylene	1.244	0.989	20.5	121	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032322\
Data File : BP009510.D
Acq On : 23 Mar 2022 17:31
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Mar 24 06:12:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 24 01:19:11 2022
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