Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072823\
Data File : BP016411.D
Acq On : 28 Jul 2023 08:03
Operator : MA/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jul 28 08:40:25 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 28 01:48:25 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	112	0.00
2		1,4-Dioxane	0.478	0.507	-6.1	124	0.00
3		Pyridine	1.429	1.533	-7.3	122	0.00
4		n-Nitrosodimethylamine	0.539	0.581	-7.8	125	0.00
5	S	2-Fluorophenol	1.226	1.299	-6.0	119	0.00
6		Aniline	2.076	2.215	-6.7	119	0.00
7	S	Phenol-d6	1.680	1.786	-6.3	119	0.00
8		2-Chlorophenol	1.285	1.350	-5.1	116	0.00
9		Benzaldehyde	0.866	0.889	-2.7	125	0.00
10	C	Phenol	1.632	1.737	-6.4	119	0.00
11		bis(2-Chloroethyl)ether	1.326	1.399	-5.5	119	0.00
12		1,3-Dichlorobenzene	1.380	1.405	-1.8	115	0.00
13	C	1,4-Dichlorobenzene	1.400	1.426	-1.9	115	0.00
14		1,2-Dichlorobenzene	1.348	1.370	-1.6	114	0.00
15		Benzyl Alcohol	1.148	1.238	-7.8	119	0.00
16		2,2'-oxybis(1-Chloropropane	1.504	1.643	-9.2	125	0.00
17		2-Methylphenol	1.176	1.244	-5.8	117	0.00
18		Hexachloroethane	0.499	0.519	-4.0	118	0.00
19	P	n-Nitroso-di-n-propylamine	1.053	1.118	-6.2	119	0.00
20		3+4-Methylphenols	1.597	1.691	-5.9	116	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	114	0.00
22		Acetophenone	0.505	0.518	-2.6	117	0.00
23	S	Nitrobenzene-d5	0.357	0.376	-5.3	119	0.00
24		Nitrobenzene	0.367	0.388	-5.7	120	0.00
25		Isophorone	0.721	0.750	-4.0	117	0.00
26	C	2-Nitrophenol	0.141	0.153	-8.5	115	0.00
27		2,4-Dimethylphenol	0.323	0.330	-2.2	115	0.00
28		bis(2-Chloroethoxy)methane	0.442	0.460	-4.1	118	0.00
29	C	2,4-Dichlorophenol	0.302	0.308	-2.0	111	0.00
30		1,2,4-Trichlorobenzene	0.327	0.320	2.1	108	0.00
31		Naphthalene	1.046	1.055	-0.9	114	0.00
32		Benzoic acid	0.191	0.199	-4.2	116	0.01
33		4-Chloroaniline	0.472	0.480	-1.7	114	0.00
34	C	Hexachlorobutadiene	0.194	0.182	6.2	102	0.00
35		Caprolactam	0.110	0.112	-1.8	112	0.00
36	C	4-Chloro-3-methylphenol	0.336	0.346	-3.0	114	0.00
37		2-Methylnaphthalene	0.760	0.762	-0.3	112	0.00
38		1-Methylnaphthalene	0.713	0.715	-0.3	112	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	108	0.00
40		1,2,4,5-Tetrachlorobenzene	0.608	0.595	2.1	102	0.00
41	P	Hexachlorocyclopentadiene	0.305	0.309	-1.3	101	0.00
42	S	2,4,6-Tribromophenol	0.294	0.259	11.9	88	0.00
43	C	2,4,6-Trichlorophenol	0.390	0.402	-3.1	104	0.00
44		2,4,5-Trichlorophenol	0.473	0.484	-2.3	105	0.00
45	S	2-Fluorobiphenyl	1.395	1.387	0.6	106	0.00
46		1,1'-Biphenyl	1.525	1.551	-1.7	109	0.00
47		2-Chloronaphthalene	1.150	1.171	-1.8	109	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072823\
Data File : BP016411.D
Acq On : 28 Jul 2023 08:03
Operator : MA/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jul 28 08:40:25 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 28 01:48:25 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	0.324	0.366	-13.0	118	0.00
49		Acenaphthylene	1.908	1.958	-2.6	110	0.00
50		Dimethylphthalate	1.559	1.562	-0.2	108	0.00
51		2,6-Dinitrotoluene	0.319	0.335	-5.0	108	0.00
52	C	Acenaphthene	1.135	1.159	-2.1	109	0.00
53		3-Nitroaniline	0.361	0.387	-7.2	112	0.00
54	P	2,4-Dinitrophenol	0.121	0.124	-2.5	110	0.00
55		Dibenzofuran	1.861	1.866	-0.3	108	0.00
56	P	4-Nitrophenol	0.283	0.301	-6.4	110	0.00
57		2,4-Dinitrotoluene	0.416	0.449	-7.9	108	0.00
58		Fluorene	1.462	1.453	0.6	105	0.00
59		2,3,4,6-Tetrachlorophenol	0.389	0.380	2.3	99	0.00
60		Diethylphthalate	1.549	1.566	-1.1	109	0.00
61		4-Chlorophenyl-phenylether	0.738	0.714	3.3	101	0.00
62		4-Nitroaniline	0.358	0.379	-5.9	109	0.00
63		Azobenzene	1.462	1.540	-5.3	116	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
65		4,6-Dinitro-2-methylphenol	0.084	0.092	-9.5	110	0.00
66	c	n-Nitrosodiphenylamine	0.589	0.616	-4.6	107	0.00
67		4-Bromophenyl-phenylether	0.212	0.211	0.5	97	0.00
68		Hexachlorobenzene	0.236	0.226	4.2	93	0.00
69		Atrazine	0.175	0.179	-2.3	102	0.00
70	C	Pentachlorophenol	0.163	0.159	2.5	96	0.00
71		Phenanthrene	1.062	1.076	-1.3	104	0.00
72		Anthracene	1.068	1.096	-2.6	104	0.00
73		Carbazole	1.001	1.026	-2.5	105	0.00
74		Di-n-butylphthalate	1.158	1.238	-6.9	107	0.00
75	C	Fluoranthene	1.245	1.231	1.1	99	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	91	0.00
77		Benzidine	0.347	0.465	-34.0#	111	0.00
78		Pyrene	1.341	1.487	-10.9	99	0.00
79	S	Terphenyl-d14	0.947	1.013	-7.0	93	0.00
80		Butylbenzylphthalate	0.527	0.619	-17.5	104	0.00
81		Benzo(a)anthracene	1.342	1.394	-3.9	93	0.00
82		3,3'-Dichlorobenzidine	0.441	0.458	-3.9	89	0.00
83		Chrysene	1.281	1.312	-2.4	92	0.00
84		Bis(2-ethylhexyl)phthalate	0.791	0.901	-13.9	101	0.00
85	c	Di-n-octyl phthalate	1.330	1.443	-8.5	97	0.00
86	I	Perylene-d12	1.000	1.000	0.0	79	0.00
87		Indeno(1,2,3-cd)pyrene	1.348	1.237	8.2	73	0.00
88		Benzo(b)fluoranthene	1.159	1.283	-10.7	84	0.00
89		Benzo(k)fluoranthene	1.179	1.279	-8.5	83	0.00
90	C	Benzo(a)pyrene	1.123	1.187	-5.7	81	0.00
91		Dibenzo(a,h)anthracene	1.124	1.048	6.8	73	0.00
92		Benzo(g,h,i)perylene	1.084	0.980	9.6	72	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072823\
Data File : BP016411.D
Acq On : 28 Jul 2023 08:03
Operator : MA/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jul 28 08:40:25 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 28 01:48:25 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