Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012622\
Data File : BP008922.D
Acq On : 26 Jan 2022 10:45
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jan 27 05:19:24 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 26 01:52:51 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	98	0.00
2		1,4-Dioxane	40.000	34.243	14.4	99	0.00
3		Pyridine	40.000	36.838	7.9	101	0.00
4		n-Nitrosodimethylamine	40.000	36.884	7.8	101	0.00
5	S	2-Fluorophenol	80.000	72.391	9.5	101	0.00
6		Aniline	40.000	37.582	6.0	102	0.00
7	S	Phenol-d6	80.000	75.241	5.9	102	0.00
8		2-Chlorophenol	40.000	36.954	7.6	101	0.00
9		Benzaldehyde	40.000	36.995	7.5	101	0.00
10	C	Phenol	40.000	37.298	6.8	101	0.00
11		bis(2-Chloroethyl)ether	40.000	37.294	6.8	102	0.00
12		1,3-Dichlorobenzene	40.000	36.026	9.9	100	0.00
13	C	1,4-Dichlorobenzene	40.000	36.030	9.9	100	0.00
14		1,2-Dichlorobenzene	40.000	36.315	9.2	101	0.00
15		Benzyl Alcohol	40.000	38.464	3.8	102	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.467	6.3	101	0.00
17		2-Methylphenol	40.000	37.748	5.6	101	0.00
18		Hexachloroethane	40.000	36.735	8.2	100	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.192	2.0	103	0.00
20		3+4-Methylphenols	40.000	38.646	3.4	103	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	98	0.00
22		Acetophenone	40.000	37.829	5.4	102	0.00
23	S	Nitrobenzene-d5	80.000	75.546	5.6	102	0.00
24		Nitrobenzene	40.000	37.646	5.9	102	0.00
25		Isophorone	40.000	38.858	2.9	103	0.00
26	C	2-Nitrophenol	40.000	39.117	2.2	103	0.00
27		2,4-Dimethylphenol	40.000	38.253	4.4	102	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.950	5.1	102	0.00
29	C	2,4-Dichlorophenol	40.000	37.998	5.0	102	0.00
30		1,2,4-Trichlorobenzene	40.000	36.451	8.9	100	0.00
31		Naphthalene	40.000	36.767	8.1	101	0.00
32		Benzoic acid	40.000	39.622	0.9	99	0.01
33		4-Chloroaniline	40.000	38.050	4.9	102	0.00
34	C	Hexachlorobutadiene	40.000	36.821	7.9	101	0.00
35		Caprolactam	40.000	40.511	-1.3	107	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.193	2.0	104	0.00
37		2-Methylnaphthalene	40.000	37.730	5.7	102	0.00
38		1-Methylnaphthalene	40.000	37.626	5.9	101	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	98	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	36.731	8.2	102	0.00
41	P	Hexachlorocyclopentadiene	40.000	36.390	9.0	93	0.00
42	S	2,4,6-Tribromophenol	80.000	77.338	3.3	106	0.00
43	C	2,4,6-Trichlorophenol	40.000	37.951	5.1	102	0.00
44		2,4,5-Trichlorophenol	40.000	37.818	5.5	102	0.00
45	S	2-Fluorobiphenyl	80.000	73.086	8.6	102	0.00
46		1,1'-Biphenyl	40.000	36.808	8.0	102	0.00
47		2-Chloronaphthalene	40.000	37.096	7.3	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012622\
Data File : BP008922.D
Acq On : 26 Jan 2022 10:45
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jan 27 05:19:24 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 26 01:52:51 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	Amount	Calc.	%Dev	Area%	Dev(min)
48		2-Nitroaniline	40.000	39.993	0.0	107	0.00
49		Acenaphthylene	40.000	37.437	6.4	103	0.00
50		Dimethylphthalate	40.000	37.814	5.5	105	0.00
51		2,6-Dinitrotoluene	40.000	39.188	2.0	106	0.00
52	C	Acenaphthene	40.000	37.757	5.6	104	0.00
53		3-Nitroaniline	40.000	39.310	1.7	106	0.00
54	P	2,4-Dinitrophenol	40.000	38.869	2.8	101	0.00
55		Dibenzofuran	40.000	37.661	5.8	104	0.00
56	P	4-Nitrophenol	40.000	38.176	4.6	101	0.00
57		2,4-Dinitrotoluene	40.000	40.465	-1.2	107	0.00
58		Fluorene	40.000	38.161	4.6	105	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	38.342	4.1	104	0.00
60		Diethylphthalate	40.000	38.341	4.1	106	0.00
61		4-Chlorophenyl-phenylether	40.000	37.938	5.2	104	0.00
62		4-Nitroaniline	40.000	39.410	1.5	106	0.00
63		Azobenzene	40.000	39.168	2.1	105	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	101	0.00
65		4,6-Dinitro-2-methylphenol	40.000	40.025	-0.1	106	0.00
66	c	n-Nitrosodiphenylamine	40.000	37.922	5.2	106	0.00
67		4-Bromophenyl-phenylether	40.000	37.742	5.6	104	0.00
68		Hexachlorobenzene	40.000	37.561	6.1	105	0.00
69		Atrazine	40.000	38.664	3.3	108	0.00
70	C	Pentachlorophenol	40.000	37.590	6.0	103	0.00
71		Phenanthrene	40.000	37.380	6.5	106	0.00
72		Anthracene	40.000	38.232	4.4	107	0.00
73		Carbazole	40.000	37.655	5.9	107	0.00
74		Di-n-butylphthalate	40.000	38.473	3.8	106	0.00
75	C	Fluoranthene	40.000	36.559	8.6	106	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	91	0.00
77		Benzidine	40.000	45.746	-14.4	103	0.00
78		Pyrene	40.000	43.125	-7.8	105	0.00
79	S	Terphenyl-d14	80.000	83.559	-4.4	102	0.00
80		Butylbenzylphthalate	40.000	40.755	-1.9	100	0.00
81		Benzo(a)anthracene	40.000	38.129	4.7	97	-0.01
82		3,3'-Dichlorobenzidine	40.000	37.381	6.5	93	0.00
83		Chrysene	40.000	37.266	6.8	95	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	38.595	3.5	94	-0.01
85	c	Di-n-octyl phthalate	40.000	35.729	10.7	96	0.00
86	I	Perylene-d12	20.000	20.000	0.0	86	-0.01
87		Indeno(1,2,3-cd)pyrene	40.000	39.744	0.6	96	-0.02
88		Benzo(b)fluoranthene	40.000	38.453	3.9	93	-0.01
89		Benzo(k)fluoranthene	40.000	38.087	4.8	95	-0.01
90	C	Benzo(a)pyrene	40.000	38.642	3.4	94	-0.01
91		Dibenzo(a,h)anthracene	40.000	39.713	0.7	95	-0.02
92		Benzo(g,h,i)perylene	40.000	39.330	1.7	94	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012622\
Data File : BP008922.D
Acq On : 26 Jan 2022 10:45
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jan 27 05:19:24 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 26 01:52:51 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	Amount	Calc.	%Dev	Area%	Dev(min)

(#) = Out of Range SPCC's out = 0 CCC's out = 0