Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031622\
Data File : BM034251.D
Acq On : 16 Mar 2022 10:02
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Mar 16 12:41:20 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 16 02:38:26 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	94	0.00
2		1,4-Dioxane	40.000	39.391	1.5	96	0.00
3		Pyridine	40.000	38.921	2.7	94	0.00
4		n-Nitrosodimethylamine	40.000	38.616	3.5	95	0.00
5	S	2-Fluorophenol	80.000	79.090	1.1	94	0.00
6		Aniline	40.000	38.622	3.4	92	0.00
7	S	Phenol-d6	80.000	77.812	2.7	93	0.00
8		2-Chlorophenol	40.000	38.629	3.4	94	0.00
9		Benzaldehyde	40.000	37.049	7.4	93	0.00
10	C	Phenol	40.000	38.769	3.1	93	0.00
11		bis(2-Chloroethyl)ether	40.000	37.725	5.7	92	0.00
12		1,3-Dichlorobenzene	40.000	38.387	4.0	94	0.00
13	C	1,4-Dichlorobenzene	40.000	38.032	4.9	94	0.00
14		1,2-Dichlorobenzene	40.000	38.049	4.9	94	0.00
15		Benzyl Alcohol	40.000	39.721	0.7	93	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.354	6.6	93	0.00
17		2-Methylphenol	40.000	38.652	3.4	93	0.00
18		Hexachloroethane	40.000	38.744	3.1	95	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	37.930	5.2	92	0.00
20		3+4-Methylphenols	40.000	38.775	3.1	93	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	95	0.00
22		Acetophenone	40.000	38.091	4.8	93	0.00
23	S	Nitrobenzene-d5	80.000	78.089	2.4	94	0.00
24		Nitrobenzene	40.000	38.471	3.8	93	0.00
25		Isophorone	40.000	38.511	3.7	92	0.00
26	C	2-Nitrophenol	40.000	39.538	1.2	92	0.00
27		2,4-Dimethylphenol	40.000	39.099	2.3	94	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.691	3.3	93	0.00
29	C	2,4-Dichlorophenol	40.000	39.666	0.8	94	0.00
30		1,2,4-Trichlorobenzene	40.000	38.514	3.7	94	0.00
31		Naphthalene	40.000	38.139	4.7	94	0.00
32		Benzoic acid	40.000	38.624	3.4	91	0.00
33		4-Chloroaniline	40.000	39.094	2.3	92	0.00
34	C	Hexachlorobutadiene	40.000	38.874	2.8	96	0.00
35		Caprolactam	40.000	39.212	2.0	94	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.294	1.8	93	0.00
37		2-Methylnaphthalene	40.000	38.693	3.3	95	0.00
38		1-Methylnaphthalene	40.000	38.548	3.6	95	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	96	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	38.575	3.6	95	0.00
41	P	Hexachlorocyclopentadiene	40.000	39.624	0.9	94	0.00
42	S	2,4,6-Tribromophenol	80.000	80.435	-0.5	99	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.840	0.4	95	0.00
44		2,4,5-Trichlorophenol	40.000	39.635	0.9	94	0.00
45	S	2-Fluorobiphenyl	80.000	77.908	2.6	95	0.00
46		1,1'-Biphenyl	40.000	38.636	3.4	96	0.00
47		2-Chloronaphthalene	40.000	38.548	3.6	95	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031622\
Data File : BM034251.D
Acq On : 16 Mar 2022 10:02
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Mar 16 12:41:20 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 16 02:38:26 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	40.107	-0.3	95	0.00
49		Acenaphthylene	40.000	38.937	2.7	95	0.00
50		Dimethylphthalate	40.000	38.767	3.1	95	0.00
51		2,6-Dinitrotoluene	40.000	39.601	1.0	95	0.00
52	C	Acenaphthene	40.000	38.983	2.5	95	0.00
53		3-Nitroaniline	40.000	40.866	-2.2	95	0.00
54	P	2,4-Dinitrophenol	40.000	41.143	-2.9	94	0.00
55		Dibenzofuran	40.000	39.018	2.5	97	0.00
56	P	4-Nitrophenol	40.000	40.370	-0.9	96	0.00
57		2,4-Dinitrotoluene	40.000	40.391	-1.0	96	0.00
58		Fluorene	40.000	39.226	1.9	96	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	40.174	-0.4	97	0.00
60		Diethylphthalate	40.000	39.797	0.5	97	0.00
61		4-Chlorophenyl-phenylether	40.000	39.752	0.6	98	0.00
62		4-Nitroaniline	40.000	41.941	-4.9	96	0.00
63		Azobenzene	40.000	39.616	1.0	96	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	99	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.826	0.4	96	0.00
66	c	n-Nitrosodiphenylamine	40.000	38.504	3.7	97	0.00
67		4-Bromophenyl-phenylether	40.000	38.744	3.1	99	0.00
68		Hexachlorobenzene	40.000	38.134	4.7	97	0.00
69		Atrazine	40.000	40.945	-2.4	99	0.00
70	C	Pentachlorophenol	40.000	40.173	-0.4	98	0.00
71		Phenanthrene	40.000	38.387	4.0	98	0.00
72		Anthracene	40.000	39.287	1.8	99	0.00
73		Carbazole	40.000	39.723	0.7	100	0.00
74		Di-n-butylphthalate	40.000	40.776	-1.9	101	0.00
75	C	Fluoranthene	40.000	39.430	1.4	101	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	110	0.00
77		Benzidine	40.000	46.806	-17.0	122	0.00
78		Pyrene	40.000	38.712	3.2	103	0.00
79	S	Terphenyl-d14	80.000	78.764	1.5	103	0.00
80		Butylbenzylphthalate	40.000	40.377	-0.9	107	0.00
81		Benzo(a)anthracene	40.000	39.026	2.4	111	0.00
82		3,3'-Dichlorobenzidine	40.000	39.466	1.3	111	0.00
83		Chrysene	40.000	38.786	3.0	111	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	40.644	-1.6	111	0.00
85	c	Di-n-octyl phthalate	40.000	39.982	0.0	115	0.00
86	I	Perylene-d12	20.000	20.000	0.0	115	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	38.323	4.2	114	0.00
88		Benzo(b)fluoranthene	40.000	40.584	-1.5	118	0.00
89		Benzo(k)fluoranthene	40.000	39.622	0.9	112	0.00
90	C	Benzo(a)pyrene	40.000	39.521	1.2	115	0.00
91		Dibenzo(a,h)anthracene	40.000	38.430	3.9	115	0.00
92		Benzo(g,h,i)perylene	40.000	37.318	6.7	112	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031622\
Data File : BM034251.D
Acq On : 16 Mar 2022 10:02
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Mar 16 12:41:20 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 16 02:38:26 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