Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP083019\
Data File : BP000012.D
Acq On : 30 Aug 2019 18:35
Operator : HP/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP083019

Quant Time: Aug 30 19:11:19 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 30 17:47:17 2019
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	106	0.00
2		1,4-Dioxane	40.000	39.961	0.1	106	0.00
3		Pyridine	40.000	41.012	-2.5	107	0.00
4		n-Nitrosodimethylamine	40.000	38.528	3.7	104	0.00
5	S	2-Fluorophenol	80.000	82.408	-3.0	108	0.00
6		Aniline	40.000	39.729	0.7	109	0.00
7	S	Phenol-d6	80.000	80.642	-0.8	107	0.00
8		2-Chlorophenol	40.000	41.329	-3.3	109	0.00
9		Benzaldehyde	40.000	35.878	10.3	105	0.00
10	C	Phenol	40.000	39.427	1.4	108	0.00
11		bis(2-Chloroethyl)ether	40.000	39.938	0.2	107	0.00
12		1,3-Dichlorobenzene	40.000	40.940	-2.3	107	0.00
13	C	1,4-Dichlorobenzene	40.000	40.777	-1.9	107	0.00
14		1,2-Dichlorobenzene	40.000	41.186	-3.0	107	0.00
15		Benzyl Alcohol	40.000	40.526	-1.3	108	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.429	1.4	106	0.00
17		2-Methylphenol	40.000	40.463	-1.2	110	0.00
18		Hexachloroethane	40.000	40.734	-1.8	109	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.145	2.1	106	0.00
20		3+4-Methylphenols	40.000	40.990	-2.5	108	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	107	0.00
22		Acetophenone	40.000	40.575	-1.4	106	0.00
23	S	Nitrobenzene-d5	80.000	83.956	-4.9	110	0.00
24		Nitrobenzene	40.000	41.140	-2.9	109	0.00
25		Isophorone	40.000	39.265	1.8	106	0.00
26	C	2-Nitrophenol	40.000	45.331	-13.3	126	0.00
27		2,4-Dimethylphenol	40.000	40.726	-1.8	109	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.075	-0.2	106	0.00
29	C	2,4-Dichlorophenol	40.000	40.771	-1.9	108	0.00
30		1,2,4-Trichlorobenzene	40.000	40.752	-1.9	107	0.00
31		Naphthalene	40.000	41.235	-3.1	104	0.00
32		Benzoic acid	40.000	41.609	-4.0	122	0.00
33		4-Chloroaniline	40.000	39.347	1.6	106	0.00
34	C	Hexachlorobutadiene	40.000	41.047	-2.6	109	0.00
35		Caprolactam	40.000	38.742	3.1	105	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.653	0.9	105	0.00
37		2-Methylnaphthalene	40.000	40.470	-1.2	105	0.00
38		1-Methylnaphthalene	40.000	40.622	-1.6	106	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	106	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	41.449	-3.6	107	0.00
41	P	Hexachlorocyclopentadiene	40.000	42.982	-7.5	112	0.00
42	S	2,4,6-Tribromophenol	80.000	80.501	-0.6	107	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.898	-2.2	107	0.00
44		2,4,5-Trichlorophenol	40.000	41.024	-2.6	108	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP083019\
Data File : BP000012.D
Acq On : 30 Aug 2019 18:35
Operator : HP/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP083019

Quant Time: Aug 30 19:11:19 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 30 17:47:17 2019
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)
45	S	2-Fluorobiphenyl	80.000	78.707	1.6	106	0.00
46		1,1'-Biphenyl	40.000	41.439	-3.6	104	0.00
47		2-Chloronaphthalene	40.000	41.277	-3.2	106	0.00
48		2-Nitroaniline	40.000	42.363	-5.9	110	0.00
49		Acenaphthylene	40.000	41.268	-3.2	104	0.00
50		Dimethylphthalate	40.000	40.457	-1.1	104	0.00
51		2,6-Dinitrotoluene	40.000	41.804	-4.5	107	0.00
52	C	Acenaphthene	40.000	40.703	-1.8	105	0.00
53		3-Nitroaniline	40.000	41.193	-3.0	106	0.00
54	P	2,4-Dinitrophenol	40.000	41.319	-3.3	123	0.00
55		Dibenzofuran	40.000	40.420	-1.1	102	0.00
56	P	4-Nitrophenol	40.000	40.791	-2.0	108	0.00
57		2,4-Dinitrotoluene	40.000	41.217	-3.0	107	0.00
58		Fluorene	40.000	40.259	-0.6	103	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.052	-2.6	108	0.00
60		Diethylphthalate	40.000	39.233	1.9	101	0.00
61		4-Chlorophenyl-phenylether	40.000	40.194	-0.5	103	0.00
62		4-Nitroaniline	40.000	39.531	1.2	106	0.00
63		Azobenzene	40.000	40.115	-0.3	103	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	102	0.00
65		4,6-Dinitro-2-methylphenol	40.000	41.665	-4.2	117	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.984	-5.0	103	0.00
67		4-Bromophenyl-phenylether	40.000	40.769	-1.9	102	0.00
68		Hexachlorobenzene	40.000	40.513	-1.3	102	0.00
69		Atrazine	40.000	37.490	6.3	100	0.00
70	C	Pentachlorophenol	40.000	41.326	-3.3	103	0.00
71		Phenanthrene	40.000	41.208	-3.0	99	0.00
72		Anthracene	40.000	41.586	-4.0	101	0.00
73		Carbazole	40.000	40.461	-1.2	98	0.00
74		Di-n-butylphthalate	40.000	40.473	-1.2	97	0.00
75	C	Fluoranthene	40.000	40.333	-0.8	97	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	96	0.00
77		Benzidine	40.000	37.006	7.5	94	0.00
78		Pyrene	40.000	42.365	-5.9	99	0.00
79	S	Terphenyl-d14	80.000	82.920	-3.7	96	0.00
80		Butylbenzylphthalate	40.000	41.150	-2.9	94	0.00
81		Benzo(a)anthracene	40.000	41.560	-3.9	97	0.00
82		3,3'-Dichlorobenzidine	40.000	40.026	-0.1	92	0.00
83		Chrysene	40.000	41.425	-3.6	96	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	39.524	1.2	91	0.00
85	c	Di-n-octyl phthalate	40.000	40.620	-1.5	90	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	40.144	-0.4	98	0.00
87	I	Perylene-d12	20.000	20.000	0.0	95	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP083019\
Data File : BP000012.D
Acq On : 30 Aug 2019 18:35
Operator : HP/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP083019

Quant Time: Aug 30 19:11:19 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 30 17:47:17 2019
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)
88	Benzo(b)fluoranthene	40.000	41.098	-2.7	98	0.00
89	Benzo(k)fluoranthene	40.000	40.045	-0.1	92	0.00
90 C	Benzo(a)pyrene	40.000	40.977	-2.4	97	0.00
91	Dibenzo(a,h)anthracene	40.000	40.663	-1.7	97	0.00
92	Benzo(g,h,i)perylene	40.000	39.839	0.4	97	0.01

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