Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP060420\
Data File : BP002367.D
Acq On : 04 Jun 2020 16:56
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP060420

Quant Time: Jun 04 17:54:56 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 04 16:24:14 2020
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	96	0.00
2		1,4-Dioxane	40.000	42.008	-5.0	100	0.00
3		Pyridine	40.000	41.998	-5.0	100	0.00
4		n-Nitrosodimethylamine	40.000	44.391	-11.0	108	0.00
5	S	2-Fluorophenol	80.000	85.014	-6.3	99	0.00
6		Aniline	40.000	43.295	-8.2	101	0.00
7	S	Phenol-d6	80.000	85.729	-7.2	100	0.00
8		2-Chlorophenol	40.000	42.288	-5.7	97	0.00
9		Benzaldehyde	40.000	49.139	-22.8	107	0.00
10	C	Phenol	40.000	43.223	-8.1	101	0.00
11		bis(2-Chloroethyl)ether	40.000	42.860	-7.1	101	0.00
12		1,3-Dichlorobenzene	40.000	41.877	-4.7	97	0.00
13	C	1,4-Dichlorobenzene	40.000	42.070	-5.2	98	0.00
14		1,2-Dichlorobenzene	40.000	42.046	-5.1	96	0.00
15		Benzyl Alcohol	40.000	44.060	-10.2	104	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	43.191	-8.0	103	0.00
17		2-Methylphenol	40.000	43.089	-7.7	100	0.00
18		Hexachloroethane	40.000	42.686	-6.7	100	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	44.218	-10.5	102	0.00
20		3+4-Methylphenols	40.000	43.329	-8.3	100	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	97	0.00
22		Acetophenone	40.000	41.954	-4.9	99	0.00
23	S	Nitrobenzene-d5	80.000	84.294	-5.4	102	0.00
24		Nitrobenzene	40.000	42.352	-5.9	103	0.00
25		Isophorone	40.000	42.141	-5.4	101	0.00
26	C	2-Nitrophenol	40.000	41.575	-3.9	96	0.00
27		2,4-Dimethylphenol	40.000	41.562	-3.9	97	0.00
28		bis(2-Chloroethoxy)methane	40.000	41.972	-4.9	100	0.00
29	C	2,4-Dichlorophenol	40.000	41.006	-2.5	94	0.00
30		1,2,4-Trichlorobenzene	40.000	40.178	-0.4	91	0.00
31		Naphthalene	40.000	40.942	-2.4	96	0.00
32		Benzoic acid	40.000	43.942	-9.9	101	0.01
33		4-Chloroaniline	40.000	41.344	-3.4	97	0.00
34	C	Hexachlorobutadiene	40.000	39.606	1.0	87	0.00
35		Caprolactam	40.000	42.082	-5.2	99	0.01
36	C	4-Chloro-3-methylphenol	40.000	41.939	-4.8	100	0.00
37		2-Methylnaphthalene	40.000	40.790	-2.0	95	0.00
38		1-Methylnaphthalene	40.000	41.149	-2.9	96	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	94	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	39.883	0.3	89	0.00
41	P	Hexachlorocyclopentadiene	40.000	39.061	2.3	85	0.00
42	S	2,4,6-Tribromophenol	80.000	71.052	11.2	72	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.446	-1.1	91	0.00
44		2,4,5-Trichlorophenol	40.000	40.683	-1.7	90	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP060420\
Data File : BP002367.D
Acq On : 04 Jun 2020 16:56
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP060420

Quant Time: Jun 04 17:54:56 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 04 16:24:14 2020
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	80.910	-1.1	91	0.00
46		1,1'-Biphenyl	40.000	40.884	-2.2	93	0.00
47		2-Chloronaphthalene	40.000	40.754	-1.9	93	0.00
48		2-Nitroaniline	40.000	43.766	-9.4	105	0.00
49		Acenaphthylene	40.000	41.007	-2.5	94	0.00
50		Dimethylphthalate	40.000	41.117	-2.8	93	0.00
51		2,6-Dinitrotoluene	40.000	41.386	-3.5	93	0.00
52	C	Acenaphthene	40.000	45.049	-12.6	103	0.00
53		3-Nitroaniline	40.000	42.214	-5.5	96	0.00
54	P	2,4-Dinitrophenol	40.000	42.297	-5.7	93	0.00
55		Dibenzofuran	40.000	40.565	-1.4	92	0.00
56	P	4-Nitrophenol	40.000	41.971	-4.9	95	0.00
57		2,4-Dinitrotoluene	40.000	41.235	-3.1	92	0.00
58		Fluorene	40.000	38.873	2.8	90	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	40.470	-1.2	86	0.00
60		Diethylphthalate	40.000	40.901	-2.3	93	0.00
61		4-Chlorophenyl-phenylether	40.000	37.881	5.3	86	0.00
62		4-Nitroaniline	40.000	41.744	-4.4	95	0.00
63		Azobenzene	40.000	42.394	-6.0	101	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	90	0.01
65		4,6-Dinitro-2-methylphenol	40.000	42.124	-5.3	90	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.471	-3.7	91	0.00
67		4-Bromophenyl-phenylether	40.000	39.410	1.5	82	0.00
68		Hexachlorobenzene	40.000	38.610	3.5	79	0.00
69		Atrazine	40.000	42.866	-7.2	89	0.01
70	C	Pentachlorophenol	40.000	41.500	-3.8	83	0.01
71		Phenanthrene	40.000	40.790	-2.0	90	0.00
72		Anthracene	40.000	40.955	-2.4	90	0.01
73		Carbazole	40.000	41.365	-3.4	92	0.01
74		Di-n-butylphthalate	40.000	42.617	-6.5	95	0.00
75	C	Fluoranthene	40.000	40.589	-1.5	87	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	82	0.01
77		Benzidine	40.000	43.327	-8.3	85	0.00
78		Pyrene	40.000	42.029	-5.1	87	0.00
79	S	Terphenyl-d14	80.000	81.011	-1.3	80	0.01
80		Butylbenzylphthalate	40.000	43.307	-8.3	93	0.00
81		Benzo(a)anthracene	40.000	40.992	-2.5	83	0.00
82		3,3'-Dichlorobenzidine	40.000	42.041	-5.1	81	0.00
83		Chrysene	40.000	41.115	-2.8	82	0.01
84		Bis(2-ethylhexyl)phthalate	40.000	43.577	-8.9	92	0.01
85	c	Di-n-octyl phthalate	40.000	43.864	-9.7	93	0.00
86	I	Perylene-d12	20.000	20.000	0.0	81	0.01
87		Indeno(1,2,3-cd)pyrene	40.000	41.214	-3.0	79	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP060420\
Data File : BP002367.D
Acq On : 04 Jun 2020 16:56
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP060420

Quant Time: Jun 04 17:54:56 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 04 16:24:14 2020
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	43.245	-8.1	84	0.01
89	Benzo(k)fluoranthene	40.000	40.172	-0.4	78	0.00
90 C	Benzo(a)pyrene	40.000	41.376	-3.4	81	0.01
91	Dibenzo(a,h)anthracene	40.000	40.717	-1.8	77	0.02
92	Benzo(g,h,i)perylene	40.000	41.411	-3.5	80	0.02

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