Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071420\
Data File : BP002741.D
Acq On : 14 Jul 2020 10:09
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Jul 14 12:01:39 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 13 10:54:03 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	108	0.00
2		1,4-Dioxane	40.000	42.557	-6.4	119	0.00
3		Pyridine	40.000	43.777	-9.4	120	0.00
4		n-Nitrosodimethylamine	40.000	43.954	-9.9	120	0.00
5	S	2-Fluorophenol	80.000	80.783	-1.0	112	0.00
6		Aniline	40.000	41.856	-4.6	114	0.00
7	S	Phenol-d6	80.000	82.134	-2.7	113	0.00
8		2-Chlorophenol	40.000	39.750	0.6	110	0.00
9		Benzaldehyde	40.000	37.771	5.6	117	0.00
10	C	Phenol	40.000	41.439	-3.6	114	0.00
11		bis(2-Chloroethyl)ether	40.000	42.068	-5.2	117	0.00
12		1,3-Dichlorobenzene	40.000	38.559	3.6	108	0.00
13	C	1,4-Dichlorobenzene	40.000	38.613	3.5	108	0.00
14		1,2-Dichlorobenzene	40.000	38.576	3.6	107	0.00
15		Benzyl Alcohol	40.000	44.127	-10.3	118	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.573	-3.9	116	0.00
17		2-Methylphenol	40.000	41.454	-3.6	113	0.00
18		Hexachloroethane	40.000	40.238	-0.6	112	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	42.174	-5.4	115	0.00
20		3+4-Methylphenols	40.000	42.022	-5.1	114	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	112	0.00
22		Acetophenone	40.000	39.836	0.4	113	0.00
23	S	Nitrobenzene-d5	80.000	82.423	-3.0	116	0.00
24		Nitrobenzene	40.000	41.256	-3.1	116	0.00
25		Isophorone	40.000	41.846	-4.6	118	0.00
26	C	2-Nitrophenol	40.000	41.603	-4.0	114	0.00
27		2,4-Dimethylphenol	40.000	40.118	-0.3	114	0.00
28		bis(2-Chloroethoxy)methane	40.000	41.457	-3.6	118	0.00
29	C	2,4-Dichlorophenol	40.000	39.904	0.2	111	0.00
30		1,2,4-Trichlorobenzene	40.000	37.870	5.3	108	0.00
31		Naphthalene	40.000	38.745	3.1	112	0.00
32		Benzoic acid	40.000	38.705	3.2	120	0.01
33		4-Chloroaniline	40.000	40.112	-0.3	113	0.00
34	C	Hexachlorobutadiene	40.000	37.786	5.5	108	0.00
35		Caprolactam	40.000	41.146	-2.9	118	0.01
36	C	4-Chloro-3-methylphenol	40.000	41.201	-3.0	115	0.00
37		2-Methylnaphthalene	40.000	38.833	2.9	111	0.00
38		1-Methylnaphthalene	40.000	38.522	3.7	110	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	112	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	38.532	3.7	111	0.00
41	P	Hexachlorocyclopentadiene	40.000	39.114	2.2	119	0.00
42	S	2,4,6-Tribromophenol	80.000	80.731	-0.9	110	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.644	-1.6	111	0.00
44		2,4,5-Trichlorophenol	40.000	39.913	0.2	109	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071420\
Data File : BP002741.D
Acq On : 14 Jul 2020 10:09
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Jul 14 12:01:39 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 13 10:54:03 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	74.101	7.4	105	0.00
46		1,1'-Biphenyl	40.000	38.647	3.4	110	0.00
47		2-Chloronaphthalene	40.000	38.333	4.2	109	0.00
48		2-Nitroaniline	40.000	44.108	-10.3	120	0.00
49		Acenaphthylene	40.000	39.142	2.1	111	0.00
50		Dimethylphthalate	40.000	39.059	2.4	111	0.00
51		2,6-Dinitrotoluene	40.000	40.965	-2.4	113	0.00
52	C	Acenaphthene	40.000	38.816	3.0	111	0.00
53		3-Nitroaniline	40.000	42.117	-5.3	114	0.00
54	P	2,4-Dinitrophenol	40.000	40.264	-0.7	134	0.00
55		Dibenzofuran	40.000	38.377	4.1	110	0.00
56	P	4-Nitrophenol	40.000	39.811	0.5	117	0.00
57		2,4-Dinitrotoluene	40.000	42.470	-6.2	113	0.00
58		Fluorene	40.000	37.861	5.3	106	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	40.945	-2.4	114	0.00
60		Diethylphthalate	40.000	39.374	1.6	111	0.00
61		4-Chlorophenyl-phenylether	40.000	37.066	7.3	104	0.00
62		4-Nitroaniline	40.000	39.847	0.4	114	0.00
63		Azobenzene	40.000	41.759	-4.4	118	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	113	0.00
65		4,6-Dinitro-2-methylphenol	40.000	40.016	-0.0	118	0.00
66	c	n-Nitrosodiphenylamine	40.000	38.521	3.7	108	0.00
67		4-Bromophenyl-phenylether	40.000	39.324	1.7	111	0.00
68		Hexachlorobenzene	40.000	38.583	3.5	109	0.00
69		Atrazine	40.000	40.175	-0.4	109	0.00
70	C	Pentachlorophenol	40.000	40.936	-2.3	125	0.00
71		Phenanthrene	40.000	38.870	2.8	110	0.00
72		Anthracene	40.000	39.513	1.2	110	0.00
73		Carbazole	40.000	39.221	1.9	109	0.00
74		Di-n-butylphthalate	40.000	41.666	-4.2	113	0.00
75	C	Fluoranthene	40.000	38.850	2.9	108	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	108	0.00
77		Benzidine	40.000	42.674	-6.7	110	0.00
78		Pyrene	40.000	39.805	0.5	108	0.00
79	S	Terphenyl-d14	80.000	76.966	3.8	105	0.00
80		Butylbenzylphthalate	40.000	43.719	-9.3	114	0.00
81		Benzo(a)anthracene	40.000	39.631	0.9	106	0.00
82		3,3'-Dichlorobenzidine	40.000	43.070	-7.7	111	0.00
83		Chrysene	40.000	39.201	2.0	107	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	43.599	-9.0	112	0.00
85	c	Di-n-octyl phthalate	40.000	44.096	-10.2	116	0.00
86	I	Perylene-d12	20.000	20.000	0.0	115	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	41.351	-3.4	114	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071420\
Data File : BP002741.D
Acq On : 14 Jul 2020 10:09
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Jul 14 12:01:39 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 13 10:54:03 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	41.149	-2.9	114	0.00
89	Benzo(k)fluoranthene	40.000	39.721	0.7	110	0.00
90 C	Benzo(a)pyrene	40.000	40.749	-1.9	113	0.00
91	Dibenzo(a,h)anthracene	40.000	41.175	-2.9	112	0.01
92	Benzo(g,h,i)perylene	40.000	41.635	-4.1	117	0.01

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