Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM061920\
Data File : BM026462.D
Acq On : 19 Jun 2020 09:05
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jun 19 10:32:25 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 19 10:20:20 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	71	0.00
2		1,4-Dioxane	40.000	42.669	-6.7	70	0.00
3		Pyridine	40.000	41.780	-4.5	70	0.00
4		n-Nitrosodimethylamine	40.000	45.055	-12.6	73	0.00
5	S	2-Fluorophenol	80.000	83.847	-4.8	70	0.00
6		Aniline	40.000	43.101	-7.8	71	0.00
7	S	Phenol-d6	80.000	86.567	-8.2	72	0.00
8		2-Chlorophenol	40.000	43.321	-8.3	73	0.00
9		Benzaldehyde	40.000	33.695	15.8	72	0.00
10	C	Phenol	40.000	42.954	-7.4	72	0.00
11		bis(2-Chloroethyl)ether	40.000	42.760	-6.9	71	0.00
12		1,3-Dichlorobenzene	40.000	42.984	-7.5	71	0.00
13	C	1,4-Dichlorobenzene	40.000	42.619	-6.5	71	0.00
14		1,2-Dichlorobenzene	40.000	43.655	-9.1	73	0.00
15		Benzyl Alcohol	40.000	44.048	-10.1	73	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	44.628	-11.6	73	0.00
17		2-Methylphenol	40.000	43.610	-9.0	73	0.00
18		Hexachloroethane	40.000	44.248	-10.6	73	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	45.489	-13.7	75	0.00
20		3+4-Methylphenols	40.000	44.255	-10.6	74	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	73	0.00
22		Acetophenone	40.000	42.953	-7.4	75	0.00
23	S	Nitrobenzene-d5	80.000	86.177	-7.7	74	0.00
24		Nitrobenzene	40.000	42.405	-6.0	74	0.00
25		Isophorone	40.000	43.670	-9.2	76	0.00
26	C	2-Nitrophenol	40.000	43.397	-8.5	75	0.00
27		2,4-Dimethylphenol	40.000	42.736	-6.8	74	0.00
28		bis(2-Chloroethoxy)methane	40.000	43.159	-7.9	75	0.00
29	C	2,4-Dichlorophenol	40.000	43.052	-7.6	74	0.00
30		1,2,4-Trichlorobenzene	40.000	42.877	-7.2	74	0.00
31		Naphthalene	40.000	42.286	-5.7	74	0.00
32		Benzoic acid	40.000	42.099	-5.2	75	0.00
33		4-Chloroaniline	40.000	41.858	-4.6	73	0.00
34	C	Hexachlorobutadiene	40.000	44.371	-10.9	76	0.00
35		Caprolactam	40.000	43.608	-9.0	76	0.00
36	C	4-Chloro-3-methylphenol	40.000	43.259	-8.1	76	0.00
37		2-Methylnaphthalene	40.000	43.120	-7.8	75	0.00
38		1-Methylnaphthalene	40.000	42.924	-7.3	75	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	76	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	43.002	-7.5	77	0.00
41	P	Hexachlorocyclopentadiene	40.000	37.139	7.2	66	0.00
42	S	2,4,6-Tribromophenol	80.000	87.703	-9.6	81	0.00
43	C	2,4,6-Trichlorophenol	40.000	43.277	-8.2	77	0.00
44		2,4,5-Trichlorophenol	40.000	43.028	-7.6	77	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM061920\
Data File : BM026462.D
Acq On : 19 Jun 2020 09:05
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jun 19 10:32:25 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 19 10:20:20 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	86.859	-8.6	78	0.00
46		1,1'-Biphenyl	40.000	42.296	-5.7	76	0.00
47		2-Chloronaphthalene	40.000	42.529	-6.3	76	0.00
48		2-Nitroaniline	40.000	43.031	-7.6	77	0.00
49		Acenaphthylene	40.000	42.477	-6.2	77	0.00
50		Dimethylphthalate	40.000	43.101	-7.8	78	0.00
51		2,6-Dinitrotoluene	40.000	42.875	-7.2	77	0.00
52	C	Acenaphthene	40.000	42.444	-6.1	76	0.00
53		3-Nitroaniline	40.000	41.079	-2.7	75	0.00
54	P	2,4-Dinitrophenol	40.000	41.004	-2.5	75	0.00
55		Dibenzofuran	40.000	42.192	-5.5	76	0.00
56	P	4-Nitrophenol	40.000	38.203	4.5	70	0.00
57		2,4-Dinitrotoluene	40.000	43.540	-8.8	80	0.00
58		Fluorene	40.000	43.050	-7.6	78	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	42.329	-5.8	77	0.00
60		Diethylphthalate	40.000	43.304	-8.3	80	0.00
61		4-Chlorophenyl-phenylether	40.000	43.554	-8.9	79	0.00
62		4-Nitroaniline	40.000	38.619	3.5	71	0.00
63		Azobenzene	40.000	43.650	-9.1	79	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	78	0.00
65		4,6-Dinitro-2-methylphenol	40.000	43.374	-8.4	80	0.00
66	c	n-Nitrosodiphenylamine	40.000	43.026	-7.6	78	0.00
67		4-Bromophenyl-phenylether	40.000	43.377	-8.4	78	0.00
68		Hexachlorobenzene	40.000	43.555	-8.9	79	0.00
69		Atrazine	40.000	43.599	-9.0	76	0.00
70	C	Pentachlorophenol	40.000	41.185	-3.0	75	0.00
71		Phenanthrene	40.000	42.818	-7.0	79	0.00
72		Anthracene	40.000	42.994	-7.5	79	0.00
73		Carbazole	40.000	41.967	-4.9	79	0.00
74		Di-n-butylphthalate	40.000	45.387	-13.5	85	0.00
75	C	Fluoranthene	40.000	42.480	-6.2	82	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	83	0.00
77		Benzidine	40.000	40.369	-0.9	74	0.00
78		Pyrene	40.000	44.635	-11.6	82	0.00
79	S	Terphenyl-d14	80.000	92.383	-15.5	86	0.00
80		Butylbenzylphthalate	40.000	46.393	-16.0	92	0.00
81		Benzo(a)anthracene	40.000	42.409	-6.0	84	0.00
82		3,3'-Dichlorobenzidine	40.000	43.632	-9.1	85	0.00
83		Chrysene	40.000	42.652	-6.6	83	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	47.121	-17.8	96	0.00
85	c	Di-n-octyl phthalate	40.000	45.655	-14.1	95	0.00
86	I	Perylene-d12	20.000	20.000	0.0	76	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	44.861	-12.2	75	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM061920\
Data File : BM026462.D
Acq On : 19 Jun 2020 09:05
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jun 19 10:32:25 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 19 10:20:20 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	43.359	-8.4	79	0.00
89	Benzo(k)fluoranthene	40.000	43.669	-9.2	79	0.00
90 C	Benzo(a)pyrene	40.000	43.369	-8.4	77	0.00
91	Dibenzo(a,h)anthracene	40.000	44.835	-12.1	75	0.00
92	Benzo(g,h,i)perylene	40.000	45.346	-13.4	74	0.00

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