Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM060920\
Data File : BM026309.D
Acq On : 09 Jun 2020 19:21
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jun 09 19:52:46 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 09 14:46:47 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	112	0.00
2		1,4-Dioxane	0.510	0.477	6.5	97	0.00
3		Pyridine	1.498	1.405	6.2	99	0.00
4		n-Nitrosodimethylamine	0.742	0.644	13.2	90	0.00
5	S	2-Fluorophenol	1.131	1.097	3.0	103	0.00
6		Aniline	2.149	2.005	6.7	98	0.00
7	S	Phenol-d6	1.638	1.551	5.3	101	0.00
8		2-Chlorophenol	1.305	1.262	3.3	103	0.00
9		Benzaldehyde	1.044	1.086	-4.0	141	0.00
10	C	Phenol	1.744	1.645	5.7	101	0.00
11		bis(2-Chloroethyl)ether	1.405	1.317	6.3	99	0.00
12		1,3-Dichlorobenzene	1.445	1.408	2.6	103	0.00
13	C	1,4-Dichlorobenzene	1.471	1.424	3.2	103	0.00
14		1,2-Dichlorobenzene	1.437	1.376	4.2	101	0.00
15		Benzyl Alcohol	1.391	1.287	7.5	98	0.00
16		2,2'-oxybis(1-Chloropropane	2.624	2.313	11.9	92	0.00
17		2-Methylphenol	1.275	1.198	6.0	100	0.00
18		Hexachloroethane	0.557	0.549	1.4	103	0.00
19	P	n-Nitroso-di-n-propylamine	1.276	1.153	9.6	95	0.00
20		3+4-Methylphenols	1.738	1.629	6.3	99	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
22		Acetophenone	0.514	0.499	2.9	97	0.00
23	S	Nitrobenzene-d5	0.407	0.389	4.4	95	0.00
24		Nitrobenzene	0.426	0.403	5.4	95	0.00
25		Isophorone	0.765	0.744	2.7	97	0.00
26	C	2-Nitrophenol	0.169	0.174	-3.0	103	0.00
27		2,4-Dimethylphenol	0.266	0.262	1.5	99	0.00
28		bis(2-Chloroethoxy)methane	0.434	0.413	4.8	95	0.00
29	C	2,4-Dichlorophenol	0.292	0.290	0.7	99	0.00
30		1,2,4-Trichlorobenzene	0.321	0.320	0.3	100	0.00
31		Naphthalene	1.008	0.976	3.2	98	0.00
32		Benzoic acid	0.198	0.154	22.2	75	0.04
33		4-Chloroaniline	0.440	0.414	5.9	95	0.00
34	C	Hexachlorobutadiene	0.202	0.203	-0.5	100	0.00
35		Caprolactam	0.103	0.095	7.8	94	0.02
36	C	4-Chloro-3-methylphenol	0.356	0.333	6.5	95	0.00
37		2-Methylnaphthalene	0.718	0.682	5.0	96	0.00
38		1-Methylnaphthalene	0.680	0.639	6.0	95	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
40		1,2,4,5-Tetrachlorobenzene	0.569	0.582	-2.3	96	0.00
41	P	Hexachlorocyclopentadiene	0.335	0.340	-1.5	94	0.00
42	S	2,4,6-Tribromophenol	0.242	0.229	5.4	92	0.00
43	C	2,4,6-Trichlorophenol	0.385	0.394	-2.3	96	0.00
44		2,4,5-Trichlorophenol	0.447	0.443	0.9	93	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM060920\
Data File : BM026309.D
Acq On : 09 Jun 2020 19:21
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jun 09 19:52:46 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 09 14:46:47 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.318	1.329	-0.8	95	0.00
46		1,1'-Biphenyl	1.402	1.393	0.6	94	0.00
47		2-Chloronaphthalene	1.139	1.127	1.1	94	0.00
48		2-Nitroaniline	0.451	0.427	5.3	89	0.00
49		Acenaphthylene	1.822	1.777	2.5	93	0.00
50		Dimethylphthalate	1.463	1.397	4.5	91	0.00
51		2,6-Dinitrotoluene	0.321	0.308	4.0	91	0.00
52	C	Acenaphthene	1.094	1.054	3.7	91	0.00
53		3-Nitroaniline	0.357	0.334	6.4	90	0.00
54	P	2,4-Dinitrophenol	0.195	0.174	10.8	86	0.00
55		Dibenzofuran	1.762	1.670	5.2	90	0.00
56	P	4-Nitrophenol	0.283	0.249	12.0	85	0.01
57		2,4-Dinitrotoluene	0.446	0.416	6.7	90	0.00
58		Fluorene	1.431	1.345	6.0	90	0.00
59		2,3,4,6-Tetrachlorophenol	0.383	0.364	5.0	91	0.00
60		Diethylphthalate	1.490	1.385	7.0	90	0.00
61		4-Chlorophenyl-phenylether	0.759	0.722	4.9	91	0.00
62		4-Nitroaniline	0.381	0.337	11.5	86	0.00
63		Azobenzene	1.630	1.500	8.0	87	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
65		4,6-Dinitro-2-methylphenol	0.122	0.119	2.5	88	0.00
66	c	n-Nitrosodiphenylamine	0.579	0.581	-0.3	89	0.00
67		4-Bromophenyl-phenylether	0.220	0.226	-2.7	91	0.00
68		Hexachlorobenzene	0.230	0.231	-0.4	89	0.00
69		Atrazine	0.173	0.182	-5.2	89	0.00
70	C	Pentachlorophenol	0.138	0.130	5.8	83	0.00
71		Phenanthrene	1.042	1.003	3.7	87	0.00
72		Anthracene	1.039	1.009	2.9	87	0.00
73		Carbazole	0.986	0.931	5.6	87	0.00
74		Di-n-butylphthalate	1.162	1.151	0.9	91	0.00
75	C	Fluoranthene	1.194	1.099	8.0	87	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	94	0.00
77		Benzidine	0.415	0.397	4.3	79	0.00
78		Pyrene	1.321	1.357	-2.7	86	0.00
79	S	Terphenyl-d14	1.030	1.057	-2.6	87	0.00
80		Butylbenzylphthalate	0.534	0.575	-7.7	96	0.00
81		Benzo(a)anthracene	1.200	1.170	2.5	87	0.00
82		3,3'-Dichlorobenzidine	0.371	0.407	-9.7	97	0.00
83		Chrysene	1.144	1.124	1.7	87	0.00
84		Bis(2-ethylhexyl)phthalate	0.762	0.803	-5.4	97	0.00
85	c	Di-n-octyl phthalate	1.175	1.286	-9.4	103	0.00
86	I	Perylene-d12	1.000	1.000	0.0	100	0.00
87		Indeno(1,2,3-cd)pyrene	1.157	1.240	-7.2	94	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM060920\
Data File : BM026309.D
Acq On : 09 Jun 2020 19:21
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jun 09 19:52:46 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 09 14:46:47 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.203	1.097	8.8	87	0.00
89	Benzo(k)fluoranthene	1.169	1.115	4.6	91	0.00
90 C	Benzo(a)pyrene	1.069	1.034	3.3	90	0.00
91	Dibenzo(a,h)anthracene	0.966	1.029	-6.5	93	0.00
92	Benzo(g,h,i)perylene	0.919	1.007	-9.6	94	0.01

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