Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090220\
Data File : BP003181.D
Acq On : 02 Sep 2020 11:48
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Sep 02 15:05:54 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 01 14:37:05 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	92	0.00
2		1,4-Dioxane	0.432	0.376	13.0	88	0.00
3		Pyridine	1.138	1.063	6.6	91	0.00
4		n-Nitrosodimethylamine	0.297	0.303	-2.0	99	0.00
5	S	2-Fluorophenol	1.131	1.036	8.4	90	0.00
6		Aniline	1.544	1.497	3.0	95	0.00
7	S	Phenol-d6	1.418	1.363	3.9	95	0.00
8		2-Chlorophenol	1.261	1.188	5.8	94	0.00
9		Benzaldehyde	0.661	0.669	-1.2	98	0.00
10	C	Phenol	1.314	1.255	4.5	94	0.00
11		bis(2-Chloroethyl)ether	1.040	0.984	5.4	94	0.00
12		1,3-Dichlorobenzene	1.533	1.416	7.6	93	0.00
13	C	1,4-Dichlorobenzene	1.548	1.425	7.9	93	0.00
14		1,2-Dichlorobenzene	1.468	1.360	7.4	94	0.00
15		Benzyl Alcohol	0.799	0.828	-3.6	99	0.00
16		2,2'-oxybis(1-Chloropropane	0.881	0.859	2.5	96	0.00
17		2-Methylphenol	0.936	0.918	1.9	95	0.00
18		Hexachloroethane	0.506	0.482	4.7	95	0.00
19	P	n-Nitroso-di-n-propylamine	0.641	0.644	-0.5	98	0.00
20		3+4-Methylphenols	1.261	1.238	1.8	96	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	96	0.00
22		Acetophenone	0.445	0.413	7.2	96	0.00
23	S	Nitrobenzene-d5	0.277	0.265	4.3	97	0.00
24		Nitrobenzene	0.272	0.258	5.1	98	0.00
25		Isophorone	0.489	0.475	2.9	98	0.00
26	C	2-Nitrophenol	0.162	0.162	0.0	97	0.00
27		2,4-Dimethylphenol	0.244	0.229	6.1	95	0.00
28		bis(2-Chloroethoxy)methane	0.365	0.335	8.2	96	0.00
29	C	2,4-Dichlorophenol	0.303	0.289	4.6	97	0.00
30		1,2,4-Trichlorobenzene	0.361	0.328	9.1	96	0.00
31		Naphthalene	1.028	0.930	9.5	96	0.00
32		Benzoic acid	0.159	0.173	-8.8	97	0.01
33		4-Chloroaniline	0.429	0.409	4.7	98	0.00
34	C	Hexachlorobutadiene	0.216	0.202	6.5	99	0.00
35		Caprolactam	0.091	0.094	-3.3	103	0.00
36	C	4-Chloro-3-methylphenol	0.283	0.275	2.8	99	0.00
37		2-Methylnaphthalene	0.735	0.677	7.9	98	0.00
38		1-Methylnaphthalene	0.699	0.642	8.2	97	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
40		1,2,4,5-Tetrachlorobenzene	0.593	0.548	7.6	99	0.00
41	P	Hexachlorocyclopentadiene	0.280	0.246	12.1	87	0.00
42	S	2,4,6-Tribromophenol	0.270	0.268	0.7	106	0.00
43	C	2,4,6-Trichlorophenol	0.390	0.383	1.8	102	0.00
44		2,4,5-Trichlorophenol	0.412	0.398	3.4	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090220\
Data File : BP003181.D
Acq On : 02 Sep 2020 11:48
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Sep 02 15:05:54 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 01 14:37:05 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	1.366	1.243	9.0	99	0.00
46		1,1'-Biphenyl	1.500	1.355	9.7	98	0.00
47		2-Chloronaphthalene	1.228	1.116	9.1	98	0.00
48		2-Nitroaniline	0.230	0.237	-3.0	104	0.00
49		Acenaphthylene	1.857	1.723	7.2	99	0.00
50		Dimethylphthalate	1.525	1.410	7.5	100	0.00
51		2,6-Dinitrotoluene	0.317	0.303	4.4	100	0.00
52	C	Acenaphthene	1.141	1.031	9.6	97	0.00
53		3-Nitroaniline	0.359	0.352	1.9	101	0.00
54	P	2,4-Dinitrophenol	0.142	0.139	2.1	98	0.00
55		Dibenzofuran	1.792	1.622	9.5	99	0.00
56	P	4-Nitrophenol	0.258	0.266	-3.1	102	0.00
57		2,4-Dinitrotoluene	0.426	0.426	0.0	103	0.00
58		Fluorene	1.380	1.269	8.0	101	0.00
59		2,3,4,6-Tetrachlorophenol	0.373	0.365	2.1	103	0.00
60		Diethylphthalate	1.496	1.415	5.4	103	0.00
61		4-Chlorophenyl-phenylether	0.713	0.653	8.4	101	0.00
62		4-Nitroaniline	0.369	0.371	-0.5	105	0.00
63		Azobenzene	0.993	0.938	5.5	102	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
65		4,6-Dinitro-2-methylphenol	0.096	0.094	2.1	102	0.00
66	c	n-Nitrosodiphenylamine	0.585	0.533	8.9	101	0.00
67		4-Bromophenyl-phenylether	0.221	0.205	7.2	103	0.00
68		Hexachlorobenzene	0.263	0.242	8.0	104	0.00
69		Atrazine	0.197	0.189	4.1	102	0.00
70	C	Pentachlorophenol	0.126	0.134	-6.3	107	0.00
71		Phenanthrene	1.084	0.976	10.0	101	0.00
72		Anthracene	1.037	0.953	8.1	102	0.00
73		Carbazole	0.995	0.925	7.0	103	0.00
74		Di-n-butylphthalate	1.170	1.147	2.0	106	0.00
75	C	Fluoranthene	1.253	1.181	5.7	104	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	108	0.00
77		Benzidine	0.457	0.451	1.3	98	0.00
78		Pyrene	1.298	1.146	11.7	104	0.00
79	S	Terphenyl-d14	0.909	0.841	7.5	106	0.00
80		Butylbenzylphthalate	0.534	0.535	-0.2	111	0.00
81		Benzo(a)anthracene	1.253	1.160	7.4	108	0.00
82		3,3'-Dichlorobenzidine	0.461	0.460	0.2	110	0.00
83		Chrysene	1.199	1.086	9.4	107	0.00
84		Bis(2-ethylhexyl)phthalate	0.772	0.762	1.3	110	0.00
85	c	Di-n-octyl phthalate	1.323	1.373	-3.8	116	0.00
86	I	Perylene-d12	1.000	1.000	0.0	116	0.00
87		Indeno(1,2,3-cd)pyrene	1.491	1.301	12.7	114	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090220\
Data File : BP003181.D
Acq On : 02 Sep 2020 11:48
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Sep 02 15:05:54 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 01 14:37:05 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	1.243	1.099	11.6	116	0.00
89	Benzo(k)fluoranthene	1.188	1.027	13.6	111	0.00
90 C	Benzo(a)pyrene	1.153	1.027	10.9	115	0.00
91	Dibenzo(a,h)anthracene	1.225	1.068	12.8	114	0.00
92	Benzo(g,h,i)perylene	1.230	1.061	13.7	113	0.00

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