Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010416\
Data File : BF084232.D
Acq On : 4 Jan 2016 11:52
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 05 02:24:03 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 04 12:22:40 2016
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	76	-0.01
2		1,4-Dioxane	0.586	0.605	-3.2	75	-0.03
3		Pyridine	1.633	1.799	-10.2	77	-0.05
4		n-Nitrosodimethylamine	0.838	0.827	1.3	71	-0.05
5	S	2-Fluorophenol	1.236	1.328	-7.4	87	0.00
6		Aniline	2.272	2.230	1.8	72	-0.01
7	S	Phenol-d6	1.553	1.614	-3.9	78	0.00
8		2-Chlorophenol	1.403	1.435	-2.3	79	0.00
9		Benzaldehyde	1.011	0.928	8.2	70	-0.01
10	C	Phenol	1.766	1.991	-12.7	79	0.00
11		bis(2-Chloroethyl)ether	1.368	1.346	1.6	78	-0.01
12		1,3-Dichlorobenzene	1.447	1.459	-0.8	79	-0.01
13	C	1,4-Dichlorobenzene	1.451	1.527	-5.2	76	-0.01
14		1,2-Dichlorobenzene	1.390	1.295	6.8	75	-0.02
15		Benzyl Alcohol	1.306	1.270	2.8	70	-0.01
16		2,2'-oxybis(1-Chloropropane	2.491	2.318	6.9	72	-0.01
17		2-Methylphenol	1.176	1.101	6.4	66	-0.01
18		Hexachloroethane	0.580	0.584	-0.7	74	-0.02
19	P	n-Nitroso-di-n-propylamine	1.051	0.933	11.2	69	-0.01
20		3+4-Methylphenols	1.382	1.420	-2.7	83	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	76	-0.01
22		Acetophenone	0.481	0.486	-1.0	72	-0.01
23	S	Nitrobenzene-d5	0.359	0.338	5.8	73	-0.01
24		Nitrobenzene	0.364	0.382	-4.9	72	-0.01
25		Isophorone	0.687	0.665	3.2	73	-0.01
26	C	2-Nitrophenol	0.166	0.162	2.4	76	-0.01
27		2,4-Dimethylphenol	0.336	0.327	2.7	69	-0.01
28		bis(2-Chloroethoxy)methane	0.420	0.372	11.4	73	-0.01
29	C	2,4-Dichlorophenol	0.280	0.279	0.4	77	0.00
30		1,2,4-Trichlorobenzene	0.289	0.301	-4.2	76	-0.01
31		Naphthalene	0.907	0.924	-1.9	78	-0.01
32		Benzoic acid	0.198	0.182	8.1	67	-0.01
33		4-Chloroaniline	0.416	0.379	8.9	71	-0.01
34	C	Hexachlorobutadiene	0.166	0.160	3.6	73	-0.02
35		Caprolactam	0.094	0.083	11.7	59	-0.02
36	C	4-Chloro-3-methylphenol	0.313	0.318	-1.6	81	0.00
37		2-Methylnaphthalene	0.651	0.604	7.2	73	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	74	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.575	0.602	-4.7	77	-0.01
40	P	Hexachlorocyclopentadiene	0.347	0.383	-10.4	80	-0.01
41	S	2,4,6-Tribromophenol	0.202	0.212	-5.0	73	-0.01
42	C	2,4,6-Trichlorophenol	0.430	0.420	2.3	74	-0.01
43		2,4,5-Trichlorophenol	0.412	0.400	2.9	69	-0.01
44	S	2-Fluorobiphenyl	1.317	1.139	13.5	72	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010416\
Data File : BF084232.D
Acq On : 4 Jan 2016 11:52
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 05 02:24:03 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 04 12:22:40 2016
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		1,1'-Biphenyl	1.590	1.536	3.4	77	-0.01
46		2-Chloronaphthalene	1.249	1.151	7.8	75	-0.01
47		2-Nitroaniline	0.412	0.419	-1.7	78	0.00
48		Acenaphthylene	1.845	1.856	-0.6	71	-0.01
49		Dimethylphthalate	1.430	1.508	-5.5	75	-0.01
50		2,6-Dinitrotoluene	0.314	0.351	-11.8	75	-0.01
51	C	Acenaphthene	1.237	1.300	-5.1	72	-0.01
52		3-Nitroaniline	0.389	0.414	-6.4	72	-0.01
53	P	2,4-Dinitrophenol	0.093	0.138	-48.4#	77	-0.01
54		Dibenzofuran	1.812	1.695	6.5	72	-0.01
55	P	4-Nitrophenol	0.282	0.310	-9.9	68	0.00
56		2,4-Dinitrotoluene	0.397	0.458	-15.4	73	-0.01
57		Fluorene	1.400	1.264	9.7	71	-0.01
58		2,3,4,6-Tetrachlorophenol	0.352	0.367	-4.3	73	-0.01
59		Diethylphthalate	1.410	1.441	-2.2	74	-0.01
60		4-Chlorophenyl-phenylether	0.564	0.601	-6.6	79	-0.01
61		4-Nitroaniline	0.346	0.321	7.2	70	-0.01
62		Azobenzene	1.546	1.600	-3.5	75	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	74	-0.01
64		4,6-Dinitro-2-methylphenol	0.109	0.106	2.8	69	-0.01
65	c	n-Nitrosodiphenylamine	0.639	0.603	5.6	69	-0.01
66		4-Bromophenyl-phenylether	0.215	0.228	-6.0	73	-0.01
67		Hexachlorobenzene	0.242	0.224	7.4	72	-0.01
68		Atrazine	0.209	0.221	-5.7	69	-0.01
69	C	Pentachlorophenol	0.126	0.114	9.5	59	-0.01
70		Phenanthrene	1.046	1.105	-5.6	71	-0.01
71		Anthracene	1.026	1.023	0.3	76	-0.01
72		Carbazole	1.003	1.018	-1.5	72	-0.01
73		Di-n-butylphthalate	1.111	1.118	-0.6	72	-0.01
74	C	Fluoranthene	1.093	1.000	8.5	70	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	66	-0.01
76		Benzidine	0.717	0.648	9.6	59	-0.01
77		Pyrene	1.569	1.414	9.9	68	-0.01
78	S	Terphenyl-d14	0.896	0.783	12.6	67	-0.01
79		Butylbenzylphthalate	0.679	0.586	13.7	56	-0.01
80		Benzo(a)anthracene	1.174	1.071	8.8	64	-0.01
81		3,3'-Dichlorobenzidine	0.410	0.394	3.9	63	-0.01
82		Chrysene	1.128	0.967	14.3	58	-0.02
83		Bis(2-ethylhexyl)phthalate	0.858	0.675	21.3	54	-0.01
84	c	Di-n-octyl phthalate	1.449	1.165	19.6	51	-0.01
85		Indeno(1,2,3-cd)pyrene	0.895	1.093	-22.1	83	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	71	-0.01
87		Benzo(b)fluoranthene	1.308	1.294	1.1	67	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010416\
Data File : BF084232.D
Acq On : 4 Jan 2016 11:52
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 05 02:24:03 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 04 12:22:40 2016
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(k)fluoranthene	1.070	0.943	11.9	63	-0.01
89 C	Benzo(a)pyrene	1.110	1.115	-0.5	68	-0.01
90	Dibenzo(a,h)anthracene	0.955	1.055	-10.5	82	-0.02
91	Benzo(g,h,i)perylene	0.989	1.110	-12.2	85	-0.02

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