Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010616\
Data File : BF084302.D
Acq On : 6 Jan 2016 21:06
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 07 02:45:57 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	78	-0.03
2		1,4-Dioxane	0.586	0.574	2.0	73	-0.09
3		Pyridine	1.633	1.602	1.9	71	-0.09
4		n-Nitrosodimethylamine	0.838	0.833	0.6	73	-0.09
5	S	2-Fluorophenol	1.236	1.169	5.4	79	-0.02
6		Aniline	2.272	2.008	11.6	67	-0.03
7	S	Phenol-d6	1.553	1.581	-1.8	79	-0.01
8		2-Chlorophenol	1.403	1.464	-4.3	83	-0.02
9		Benzaldehyde	1.011	0.981	3.0	76	-0.03
10	C	Phenol	1.766	1.872	-6.0	77	-0.01
11		bis(2-Chloroethyl)ether	1.368	1.498	-9.5	89	-0.02
12		1,3-Dichlorobenzene	1.447	1.418	2.0	79	-0.03
13	C	1,4-Dichlorobenzene	1.451	1.519	-4.7	78	-0.03
14		1,2-Dichlorobenzene	1.390	1.333	4.1	80	-0.03
15		Benzyl Alcohol	1.306	1.294	0.9	74	-0.02
16		2,2'-oxybis(1-Chloropropane	2.491	2.285	8.3	74	-0.03
17		2-Methylphenol	1.176	1.172	0.3	73	-0.02
18		Hexachloroethane	0.580	0.580	0.0	75	-0.05
19	P	n-Nitroso-di-n-propylamine	1.051	1.046	0.5	80	-0.02
20		3+4-Methylphenols	1.382	1.401	-1.4	85	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	77	-0.03
22		Acetophenone	0.481	0.477	0.8	72	-0.03
23	S	Nitrobenzene-d5	0.359	0.365	-1.7	80	-0.03
24		Nitrobenzene	0.364	0.371	-1.9	71	-0.03
25		Isophorone	0.687	0.666	3.1	75	-0.03
26	C	2-Nitrophenol	0.166	0.195	-17.5	93	-0.02
27		2,4-Dimethylphenol	0.336	0.352	-4.8	76	-0.02
28		bis(2-Chloroethoxy)methane	0.420	0.447	-6.4	89	-0.02
29	C	2,4-Dichlorophenol	0.280	0.268	4.3	76	-0.02
30		1,2,4-Trichlorobenzene	0.289	0.294	-1.7	76	-0.03
31		Naphthalene	0.907	0.881	2.9	76	-0.03
32		Benzoic acid	0.198	0.171	13.6	64	-0.02
33		4-Chloroaniline	0.416	0.453	-8.9	87	-0.02
34	C	Hexachlorobutadiene	0.166	0.168	-1.2	78	-0.03
35		Caprolactam	0.094	0.092	2.1	67	-0.03
36	C	4-Chloro-3-methylphenol	0.313	0.304	2.9	79	-0.02
37		2-Methylnaphthalene	0.651	0.658	-1.1	81	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	77	-0.03
39		1,2,4,5-Tetrachlorobenzene	0.575	0.584	-1.6	78	-0.03
40	P	Hexachlorocyclopentadiene	0.347	0.341	1.7	73	-0.03
41	S	2,4,6-Tribromophenol	0.202	0.206	-2.0	73	-0.03
42	C	2,4,6-Trichlorophenol	0.430	0.389	9.5	71	-0.03
43		2,4,5-Trichlorophenol	0.412	0.422	-2.4	76	-0.02
44	S	2-Fluorobiphenyl	1.317	1.221	7.3	80	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010616\
Data File : BF084302.D
Acq On : 6 Jan 2016 21:06
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 07 02:45:57 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	1.462	8.1	75	-0.03
46		2-Chloronaphthalene	1.249	1.127	9.8	77	-0.03
47		2-Nitroaniline	0.412	0.449	-9.0	87	-0.02
48		Acenaphthylene	1.845	1.905	-3.3	75	-0.03
49		Dimethylphthalate	1.430	1.352	5.5	70	-0.03
50		2,6-Dinitrotoluene	0.314	0.286	8.9	63	-0.03
51	C	Acenaphthene	1.237	1.307	-5.7	75	-0.03
52		3-Nitroaniline	0.389	0.333	14.4	60	-0.03
53	P	2,4-Dinitrophenol	0.093	0.141	-51.6#	81	-0.02
54		Dibenzofuran	1.812	1.743	3.8	76	-0.03
55	P	4-Nitrophenol	0.282	0.295	-4.6	67	-0.01
56		2,4-Dinitrotoluene	0.397	0.382	3.8	63	-0.03
57		Fluorene	1.400	1.269	9.4	74	-0.03
58		2,3,4,6-Tetrachlorophenol	0.352	0.323	8.2	66	-0.03
59		Diethylphthalate	1.410	1.396	1.0	74	-0.03
60		4-Chlorophenyl-phenylether	0.564	0.587	-4.1	80	-0.03
61		4-Nitroaniline	0.346	0.364	-5.2	82	-0.02
62		Azobenzene	1.546	1.564	-1.2	76	-0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	77	-0.03
64		4,6-Dinitro-2-methylphenol	0.109	0.121	-11.0	82	-0.02
65	c	n-Nitrosodiphenylamine	0.639	0.661	-3.4	79	-0.02
66		4-Bromophenyl-phenylether	0.215	0.228	-6.0	77	-0.03
67		Hexachlorobenzene	0.242	0.223	7.9	75	-0.03
68		Atrazine	0.209	0.209	0.0	68	-0.03
69	C	Pentachlorophenol	0.126	0.102	19.0	55	-0.03
70		Phenanthrene	1.046	1.110	-6.1	75	-0.03
71		Anthracene	1.026	0.985	4.0	77	-0.03
72		Carbazole	1.003	0.974	2.9	72	-0.03
73		Di-n-butylphthalate	1.111	1.159	-4.3	78	-0.03
74	C	Fluoranthene	1.093	1.014	7.2	74	-0.03
75	I	Chrysene-d12	1.000	1.000	0.0	81	-0.03
76		Benzidine	0.717	0.569	20.6	63	-0.03
77		Pyrene	1.569	1.274	18.8	74	-0.03
78	S	Terphenyl-d14	0.896	0.707	21.1	74	-0.03
79		Butylbenzylphthalate	0.679	0.671	1.2	78	-0.03
80		Benzo(a)anthracene	1.174	1.024	12.8	75	-0.03
81		3,3'-Dichlorobenzidine	0.410	0.382	6.8	74	-0.03
82		Chrysene	1.128	0.920	18.4	67	-0.05
83		Bis(2-ethylhexyl)phthalate	0.858	0.822	4.2	81	-0.03
84	c	Di-n-octyl phthalate	1.449	1.307	9.8	69	-0.05
85		Indeno(1,2,3-cd)pyrene	0.895	0.959	-7.2	88	-0.06
86	I	Perylene-d12	1.000	1.000	0.0	81	-0.05
87		Benzo(b)fluoranthene	1.308	1.147	12.3	68	-0.05

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010616\
Data File : BF084302.D
Acq On : 6 Jan 2016 21:06
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 07 02:45:57 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	1.201	-12.2	92	-0.03
89 C	Benzo(a)pyrene	1.110	1.118	-0.7	79	-0.05
90	Dibenzo(a,h)anthracene	0.955	0.979	-2.5	87	-0.07
91	Benzo(g,h,i)perylene	0.989	1.022	-3.3	90	-0.07

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