Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010716\
Data File : BF084315.D
Acq On : 7 Jan 2016 19:47
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 08 01:24:38 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	86	-0.03
2		1,4-Dioxane	0.586	0.592	-1.0	83	-0.09
3		Pyridine	1.633	1.591	2.6	77	-0.10
4		n-Nitrosodimethylamine	0.838	0.836	0.2	81	-0.09
5	S	2-Fluorophenol	1.236	1.115	9.8	82	-0.02
6		Aniline	2.272	2.081	8.4	76	-0.03
7	S	Phenol-d6	1.553	1.603	-3.2	88	-0.01
8		2-Chlorophenol	1.403	1.451	-3.4	90	-0.02
9		Benzaldehyde	1.011	1.021	-1.0	87	-0.03
10	C	Phenol	1.766	1.907	-8.0	86	-0.01
11		bis(2-Chloroethyl)ether	1.368	1.493	-9.1	97	-0.02
12		1,3-Dichlorobenzene	1.447	1.395	3.6	85	-0.03
13	C	1,4-Dichlorobenzene	1.451	1.482	-2.1	83	-0.03
14		1,2-Dichlorobenzene	1.390	1.373	1.2	90	-0.03
15		Benzyl Alcohol	1.306	1.275	2.4	79	-0.03
16		2,2'-oxybis(1-Chloropropane	2.491	2.280	8.5	80	-0.03
17		2-Methylphenol	1.176	1.189	-1.1	81	-0.02
18		Hexachloroethane	0.580	0.568	2.1	81	-0.05
19	P	n-Nitroso-di-n-propylamine	1.051	1.066	-1.4	89	-0.02
20		3+4-Methylphenols	1.382	1.359	1.7	90	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	85	-0.03
22		Acetophenone	0.481	0.467	2.9	77	-0.03
23	S	Nitrobenzene-d5	0.359	0.371	-3.3	90	-0.02
24		Nitrobenzene	0.364	0.337	7.4	71	-0.03
25		Isophorone	0.687	0.673	2.0	83	-0.03
26	C	2-Nitrophenol	0.166	0.195	-17.5	101	-0.02
27		2,4-Dimethylphenol	0.336	0.352	-4.8	83	-0.02
28		bis(2-Chloroethoxy)methane	0.420	0.450	-7.1	98	-0.02
29	C	2,4-Dichlorophenol	0.280	0.265	5.4	82	-0.02
30		1,2,4-Trichlorobenzene	0.289	0.288	0.3	81	-0.03
31		Naphthalene	0.907	0.844	6.9	80	-0.03
32		Benzoic acid	0.198	0.153	22.7	62	-0.02
33		4-Chloroaniline	0.416	0.459	-10.3	96	-0.02
34	C	Hexachlorobutadiene	0.166	0.171	-3.0	87	-0.03
35		Caprolactam	0.094	0.090	4.3	71	-0.03
36	C	4-Chloro-3-methylphenol	0.313	0.289	7.7	82	-0.02
37		2-Methylnaphthalene	0.651	0.622	4.5	84	-0.03
38	I	Acenaphthene-d10	1.000	1.000	0.0	85	-0.03
39		1,2,4,5-Tetrachlorobenzene	0.575	0.577	-0.3	85	-0.03
40	P	Hexachlorocyclopentadiene	0.347	0.338	2.6	81	-0.03
41	S	2,4,6-Tribromophenol	0.202	0.206	-2.0	81	-0.03
42	C	2,4,6-Trichlorophenol	0.430	0.401	6.7	81	-0.03
43		2,4,5-Trichlorophenol	0.412	0.425	-3.2	84	-0.02
44	S	2-Fluorobiphenyl	1.317	1.158	12.1	84	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010716\
Data File : BF084315.D
Acq On : 7 Jan 2016 19:47
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 08 01:24:38 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	1.419	10.8	81	-0.03
46		2-Chloronaphthalene	1.249	1.118	10.5	84	-0.03
47		2-Nitroaniline	0.412	0.470	-14.1	100	-0.02
48		Acenaphthylene	1.845	1.865	-1.1	82	-0.03
49		Dimethylphthalate	1.430	1.304	8.8	74	-0.03
50		2,6-Dinitrotoluene	0.314	0.285	9.2	69	-0.03
51	C	Acenaphthene	1.237	1.300	-5.1	83	-0.03
52		3-Nitroaniline	0.389	0.342	12.1	69	-0.03
53	P	2,4-Dinitrophenol	0.093	0.139	-49.5#	88	-0.02
54		Dibenzofuran	1.812	1.672	7.7	81	-0.03
55	P	4-Nitrophenol	0.282	0.316	-12.1	79	-0.01
56		2,4-Dinitrotoluene	0.397	0.371	6.5	68	-0.03
57		Fluorene	1.400	1.286	8.1	83	-0.03
58		2,3,4,6-Tetrachlorophenol	0.352	0.322	8.5	73	-0.03
59		Diethylphthalate	1.410	1.369	2.9	80	-0.03
60		4-Chlorophenyl-phenylether	0.564	0.582	-3.2	88	-0.03
61		4-Nitroaniline	0.346	0.353	-2.0	88	-0.02
62		Azobenzene	1.546	1.536	0.6	83	-0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	83	-0.03
64		4,6-Dinitro-2-methylphenol	0.109	0.123	-12.8	89	-0.02
65	c	n-Nitrosodiphenylamine	0.639	0.656	-2.7	84	-0.02
66		4-Bromophenyl-phenylether	0.215	0.238	-10.7	86	-0.03
67		Hexachlorobenzene	0.242	0.225	7.0	81	-0.03
68		Atrazine	0.209	0.231	-10.5	81	-0.03
69	C	Pentachlorophenol	0.126	0.105	16.7	61	-0.03
70		Phenanthrene	1.046	1.131	-8.1	82	-0.03
71		Anthracene	1.026	0.997	2.8	84	-0.03
72		Carbazole	1.003	1.025	-2.2	81	-0.03
73		Di-n-butylphthalate	1.111	1.203	-8.3	87	-0.03
74	C	Fluoranthene	1.093	1.042	4.7	82	-0.03
75	I	Chrysene-d12	1.000	1.000	0.0	87	-0.03
76		Benzidine	0.717	0.612	14.6	73	-0.03
77		Pyrene	1.569	1.318	16.0	82	-0.03
78	S	Terphenyl-d14	0.896	0.725	19.1	81	-0.03
79		Butylbenzylphthalate	0.679	0.705	-3.8	88	-0.03
80		Benzo(a)anthracene	1.174	1.071	8.8	83	-0.03
81		3,3'-Dichlorobenzidine	0.410	0.388	5.4	81	-0.03
82		Chrysene	1.128	1.045	7.4	81	-0.05
83		Bis(2-ethylhexyl)phthalate	0.858	0.856	0.2	90	-0.03
84	c	Di-n-octyl phthalate	1.449	1.356	6.4	77	-0.05
85		Indeno(1,2,3-cd)pyrene	0.895	0.962	-7.5	95	-0.06
86	I	Perylene-d12	1.000	1.000	0.0	91	-0.05
87		Benzo(b)fluoranthene	1.308	1.132	13.5	76	-0.05

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010716\
Data File : BF084315.D
Acq On : 7 Jan 2016 19:47
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 08 01:24:38 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	1.180	-10.3	101	-0.03
89 C	Benzo(a)pyrene	1.110	1.095	1.4	87	-0.05
90	Dibenzo(a,h)anthracene	0.955	0.949	0.6	95	-0.06
91	Benzo(g,h,i)perylene	0.989	0.979	1.0	97	-0.07

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