Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111915\
Data File : BF083058.D
Acq On : 20 Nov 2015 2:56
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 20 07:39:45 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF111915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 20 01:33:39 2015
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	96	0.00
2		1,4-Dioxane	0.584	0.548	6.2	101	0.00
3		Pyridine	1.583	1.841	-16.3	105	-0.01
4		n-Nitrosodimethylamine	0.801	0.741	7.5	94	0.00
5	S	2-Fluorophenol	1.288	1.302	-1.1	106	0.00
6		Aniline	2.320	2.297	1.0	101	0.00
7	S	Phenol-d6	1.696	1.631	3.8	96	0.00
8		2-Chlorophenol	1.436	1.473	-2.6	105	0.00
9		Benzaldehyde	0.919	0.858	6.6	97	0.00
10	C	Phenol	1.976	1.737	12.1	93	-0.01
11		bis(2-Chloroethyl)ether	1.436	1.422	1.0	104	0.00
12		1,3-Dichlorobenzene	1.608	1.508	6.2	95	-0.01
13	C	1,4-Dichlorobenzene	1.628	1.559	4.2	99	0.00
14		1,2-Dichlorobenzene	1.505	1.619	-7.6	104	0.00
15		Benzyl Alcohol	1.395	1.449	-3.9	107	0.00
16		2,2'-oxybis(1-Chloropropane	2.238	2.164	3.3	94	0.00
17		2-Methylphenol	1.199	1.277	-6.5	108	0.00
18		Hexachloroethane	0.598	0.615	-2.8	105	0.00
19	P	n-Nitroso-di-n-propylamine	1.239	1.179	4.8	102	0.00
20		3+4-Methylphenols	1.563	1.566	-0.2	96	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	96	0.00
22		Acetophenone	0.560	0.531	5.2	96	0.00
23	S	Nitrobenzene-d5	0.437	0.436	0.2	97	0.00
24		Nitrobenzene	0.441	0.419	5.0	105	0.00
25		Isophorone	0.824	0.825	-0.1	96	0.00
26	C	2-Nitrophenol	0.193	0.188	2.6	103	0.00
27		2,4-Dimethylphenol	0.318	0.339	-6.6	97	0.00
28		bis(2-Chloroethoxy)methane	0.462	0.490	-6.1	94	0.00
29	C	2,4-Dichlorophenol	0.309	0.325	-5.2	100	0.00
30		1,2,4-Trichlorobenzene	0.342	0.353	-3.2	102	0.00
31		Naphthalene	1.042	1.043	-0.1	93	0.00
32		Benzoic acid	0.194	0.221	-13.9	116	0.00
33		4-Chloroaniline	0.453	0.423	6.6	101	0.00
34	C	Hexachlorobutadiene	0.214	0.211	1.4	95	0.00
35		Caprolactam	0.098	0.102	-4.1	106	0.00
36	C	4-Chloro-3-methylphenol	0.357	0.328	8.1	86	-0.01
37		2-Methylnaphthalene	0.694	0.677	2.4	103	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
39		1,2,4,5-Tetrachlorobenzene	0.660	0.660	0.0	107	0.00
40	P	Hexachlorocyclopentadiene	0.307	0.309	-0.7	110	0.00
41	S	2,4,6-Tribromophenol	0.212	0.208	1.9	108	0.00
42	C	2,4,6-Trichlorophenol	0.479	0.445	7.1	101	0.00
43		2,4,5-Trichlorophenol	0.477	0.432	9.4	90	0.00
44	S	2-Fluorobiphenyl	1.485	1.241	16.4	91	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111915\
Data File : BF083058.D
Acq On : 20 Nov 2015 2:56
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 20 07:39:45 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF111915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 20 01:33:39 2015
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	1.760	1.775	-0.9	110	0.00
46		2-Chloronaphthalene	1.339	1.350	-0.8	102	0.00
47		2-Nitroaniline	0.505	0.445	11.9	100	0.00
48		Acenaphthylene	2.204	2.184	0.9	106	0.00
49		Dimethylphthalate	1.646	1.594	3.2	102	0.00
50		2,6-Dinitrotoluene	0.355	0.360	-1.4	96	0.00
51	C	Acenaphthene	1.393	1.328	4.7	106	0.00
52		3-Nitroaniline	0.405	0.332	18.0	89	0.00
53	P	2,4-Dinitrophenol	0.181	0.160	11.6	92	-0.01
54		Dibenzofuran	1.866	1.793	3.9	107	0.00
55	P	4-Nitrophenol	0.278	0.232	16.5	77	-0.01
56		2,4-Dinitrotoluene	0.475	0.498	-4.8	105	0.00
57		Fluorene	1.502	1.481	1.4	102	0.00
58		2,3,4,6-Tetrachlorophenol	0.386	0.382	1.0	97	0.00
59		Diethylphthalate	1.605	1.549	3.5	98	0.00
60		4-Chlorophenyl-phenylether	0.703	0.634	9.8	91	0.00
61		4-Nitroaniline	0.413	0.417	-1.0	102	0.00
62		Azobenzene	1.798	1.603	10.8	92	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	96	0.00
64		4,6-Dinitro-2-methylphenol	0.128	0.140	-9.4	89	0.00
65	c	n-Nitrosodiphenylamine	0.701	0.701	0.0	109	0.00
66		4-Bromophenyl-phenylether	0.232	0.220	5.2	101	0.00
67		Hexachlorobenzene	0.253	0.261	-3.2	108	0.00
68		Atrazine	0.220	0.196	10.9	84	-0.01
69	C	Pentachlorophenol	0.129	0.141	-9.3	104	0.00
70		Phenanthrene	1.207	1.104	8.5	88	-0.01
71		Anthracene	1.216	1.195	1.7	108	0.00
72		Carbazole	1.063	1.109	-4.3	101	0.00
73		Di-n-butylphthalate	1.340	1.328	0.9	101	0.00
74	C	Fluoranthene	1.258	1.230	2.2	107	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	89	0.00
76		Benzidine	0.654	0.773	-18.2	104	0.00
77		Pyrene	1.408	1.538	-9.2	105	0.00
78	S	Terphenyl-d14	0.847	0.785	7.3	88	0.00
79		Butylbenzylphthalate	0.622	0.611	1.8	92	0.00
80		Benzo(a)anthracene	1.271	1.320	-3.9	94	0.00
81		3,3'-Dichlorobenzidine	0.428	0.467	-9.1	108	0.00
82		Chrysene	1.165	1.261	-8.2	121	0.00
83		Bis(2-ethylhexyl)phthalate	0.813	0.882	-8.5	102	0.00
84	c	Di-n-octyl phthalate	1.325	1.339	-1.1	98	0.00
85		Indeno(1,2,3-cd)pyrene	0.948	0.976	-3.0	101	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	100	-0.01
87		Benzo(b)fluoranthene	1.399	1.295	7.4	109	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111915\
Data File : BF083058.D
Acq On : 20 Nov 2015 2:56
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 20 07:39:45 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF111915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 20 01:33:39 2015
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	1.065	1.089	-2.3	88	-0.01
89 C	Benzo(a)pyrene	1.122	1.120	0.2	101	-0.01
90	Dibenzo(a,h)anthracene	0.977	0.966	1.1	102	-0.01
91	Benzo(g,h,i)perylene	0.953	0.938	1.6	101	-0.01

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