Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF031015\
Data File : BF077658.D
Acq On : 10 Mar 2015 13:52
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 11 00:54:39 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 11 00:54:19 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	62	-0.04
2		1,4-Dioxane	0.508	0.571	-12.4	70	-0.06
3		Pyridine	1.454	1.395	4.1	56	-0.06
4		n-Nitrosodimethylamine	0.632	0.676	-7.0	68	-0.07
5	S	2-Fluorophenol	1.198	1.355	-13.1	69	-0.04
6		Aniline	2.129	2.036	4.4	57	-0.04
7	S	Phenol-d6	1.540	1.486	3.5	58	-0.04
8		2-Chlorophenol	1.413	1.352	4.3	58	-0.04
9		Benzaldehyde	0.935	1.074	-14.9	72	-0.04
10	C	Phenol	1.828	1.695	7.3	54	-0.04
11		bis(2-Chloroethyl)ether	1.313	1.309	0.3	62	-0.04
12		1,3-Dichlorobenzene	1.539	1.456	5.4	57	-0.04
13	C	1,4-Dichlorobenzene	1.566	1.512	3.4	58	-0.04
14		1,2-Dichlorobenzene	1.489	1.475	0.9	59	-0.04
15		Benzyl Alcohol	1.171	1.103	5.8	57	-0.04
16		2,2'-oxybis(1-Chloropropane	1.877	1.676	10.7	53	-0.04
17		2-Methylphenol	1.105	1.065	3.6	55	-0.04
18		Hexachloroethane	0.523	0.507	3.1	59	-0.04
19	P	n-Nitroso-di-n-propylamine	0.978	0.908	7.2	55	-0.04
20		3+4-Methylphenols	1.455	1.395	4.1	57	-0.04
21	I	Naphthalene-d8	1.000	1.000	0.0	61	-0.04
22		Acetophenone	0.470	0.464	1.3	62	-0.04
23	S	Nitrobenzene-d5	0.322	0.325	-0.9	60	-0.04
24		Nitrobenzene	0.338	0.336	0.6	60	-0.04
25		Isophorone	0.638	0.595	6.7	53	-0.04
26	C	2-Nitrophenol	0.139	0.162	-16.5	64	-0.04
27		2,4-Dimethylphenol	0.310	0.305	1.6	59	-0.04
28		bis(2-Chloroethoxy)methane	0.403	0.395	2.0	58	-0.04
29	C	2,4-Dichlorophenol	0.291	0.282	3.1	57	-0.04
30		1,2,4-Trichlorobenzene	0.320	0.297	7.2	55	-0.04
31		Naphthalene	1.000	0.952	4.8	58	-0.04
32		Benzoic acid	0.151	0.191	-26.5#	71	-0.04
33		4-Chloroaniline	0.445	0.426	4.3	56	-0.04
34	C	Hexachlorobutadiene	0.183	0.167	8.7	55	-0.04
35		Caprolactam	0.089	0.086	3.4	56	-0.04
36	C	4-Chloro-3-methylphenol	0.293	0.282	3.8	55	-0.04
37		2-Methylnaphthalene	0.710	0.665	6.3	54	-0.04
38	I	Acenaphthene-d10	1.000	1.000	0.0	55	-0.04
39		1,2,4,5-Tetrachlorobenzene	0.574	0.568	1.0	52	-0.04
40	P	Hexachlorocyclopentadiene	0.308	0.293	4.9	47#	-0.04
41	S	2,4,6-Tribromophenol	0.193	0.205	-6.2	54	-0.04
42	C	2,4,6-Trichlorophenol	0.380	0.389	-2.4	52	-0.04
43		2,4,5-Trichlorophenol	0.406	0.439	-8.1	56	-0.04
44	S	2-Fluorobiphenyl	1.283	1.352	-5.4	57	-0.04
45		1,1'-Biphenyl	1.515	1.514	0.1	53	-0.04
46		2-Chloronaphthalene	1.188	1.222	-2.9	56	-0.04
47		2-Nitroaniline	0.293	0.320	-9.2	57	-0.04

Data Path : Z:\HPCHEM1\BNA_F\Data\BF031015\
Data File : BF077658.D
Acq On : 10 Mar 2015 13:52
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 11 00:54:39 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 11 00:54:19 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)
48		Acenaphthylene	1.881	1.978	-5.2	56	-0.04
49		Dimethylphthalate	1.406	1.422	-1.1	54	-0.04
50		2,6-Dinitrotoluene	0.282	0.330	-17.0	57	-0.04
51	C	Acenaphthene	1.123	1.114	0.8	52	-0.04
52		3-Nitroaniline	0.325	0.355	-9.2	55	-0.04
53	P	2,4-Dinitrophenol	0.086	0.104	-20.9	63	-0.04
54		Dibenzofuran	1.733	1.710	1.3	52	-0.04
55	P	4-Nitrophenol	0.261	0.276	-5.7	52	-0.04
56		2,4-Dinitrotoluene	0.381	0.435	-14.2	55	-0.04
57		Fluorene	1.386	1.414	-2.0	54	-0.04
58		2,3,4,6-Tetrachlorophenol	0.327	0.340	-4.0	52	-0.04
59		Diethylphthalate	1.386	1.374	0.9	53	-0.04
60		4-Chlorophenyl-phenylether	0.668	0.658	1.5	52	-0.04
61		4-Nitroaniline	0.312	0.354	-13.5	58	-0.04
62		Azobenzene	1.315	1.297	1.4	51	-0.04
63	I	Phenanthrene-d10	1.000	1.000	0.0	59	-0.04
64		4,6-Dinitro-2-methylphenol	0.075	0.080	-6.7	59	-0.04
65	c	n-Nitrosodiphenylamine	0.664	0.604	9.0	52	-0.04
66		4-Bromophenyl-phenylether	0.234	0.198	15.4	50#	-0.04
67		Hexachlorobenzene	0.251	0.217	13.5	52	-0.04
68		Atrazine	0.216	0.172	20.4	50#	-0.04
69	C	Pentachlorophenol	0.132	0.117	11.4	49#	-0.04
70		Phenanthrene	1.159	1.055	9.0	54	-0.04
71		Anthracene	1.150	1.062	7.7	56	-0.04
72		Carbazole	1.094	1.025	6.3	53	-0.04
73		Di-n-butylphthalate	1.284	1.179	8.2	51	-0.04
74	C	Fluoranthene	1.266	1.169	7.7	53	-0.05
75	I	Chrysene-d12	1.000	1.000	0.0	57	-0.03
76		Benzidine	0.676	0.664	1.8	60	-0.04
77		Pyrene	1.223	1.139	6.9	53	-0.05
78	S	Terphenyl-d14	0.856	0.777	9.2	52	-0.04
79		Butylbenzylphthalate	0.503	0.478	5.0	51	-0.04
80		Benzo(a)anthracene	1.172	1.115	4.9	55	-0.03
81		3,3'-Dichlorobenzidine	0.430	0.423	1.6	55	-0.03
82		Chrysene	1.101	1.041	5.4	55	-0.03
83		Bis(2-ethylhexyl)phthalate	0.807	0.711	11.9	49#	-0.02
84	c	Di-n-octyl phthalate	1.348	1.263	6.3	51	0.02
85		Indeno(1,2,3-cd)pyrene	1.255	1.272	-1.4	57	0.04
86	I	Perylene-d12	1.000	1.000	0.0	59	0.05
87		Benzo(b)fluoranthene	1.163	1.096	5.8	53	0.02
88		Benzo(k)fluoranthene	1.140	1.140	0.0	63	0.03
89	C	Benzo(a)pyrene	1.108	1.080	2.5	56	0.04
90		Dibenzo(a,h)anthracene	1.056	1.064	-0.8	58	0.04
91		Benzo(g,h,i)perylene	1.066	1.067	-0.1	58	0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF031015\
Data File : BF077658.D
Acq On : 10 Mar 2015 13:52
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Mar 11 00:54:39 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 11 00:54:19 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)

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