Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF032515\
Data File : BF077975.D
Acq On : 26 Mar 2015 3:35
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 26 05:23:28 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 26 00:45:24 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	91	-0.07
2		1,4-Dioxane	0.520	0.505	2.9	88	-0.10
3		Pyridine	1.398	1.416	-1.3	95	-0.13
4		n-Nitrosodimethylamine	0.609	0.534	12.3	74	-0.11
5	S	2-Fluorophenol	1.146	1.177	-2.7	97	-0.06
6		Aniline	2.025	2.034	-0.4	94	-0.07
7	S	Phenol-d6	1.416	1.431	-1.1	93	-0.06
8		2-Chlorophenol	1.364	1.378	-1.0	96	-0.06
9		Benzaldehyde	0.580	0.715	-23.3	113	-0.07
10	C	Phenol	1.673	1.745	-4.3	98	-0.05
11		bis(2-Chloroethyl)ether	1.253	1.231	1.8	91	-0.07
12		1,3-Dichlorobenzene	1.517	1.510	0.5	94	-0.07
13	C	1,4-Dichlorobenzene	1.576	1.588	-0.8	97	-0.07
14		1,2-Dichlorobenzene	1.445	1.453	-0.6	96	-0.07
15		Benzyl Alcohol	1.105	1.160	-5.0	97	-0.06
16		2,2'-oxybis(1-Chloropropane	1.689	1.668	1.2	93	-0.07
17		2-Methylphenol	1.110	1.092	1.6	90	-0.06
18		Hexachloroethane	0.517	0.522	-1.0	98	-0.07
19	P	n-Nitroso-di-n-propylamine	0.979	0.951	2.9	92	-0.07
20		3+4-Methylphenols	1.457	1.500	-3.0	97	-0.06
21	I	Naphthalene-d8	1.000	1.000	0.0	92	-0.07
22		Acetophenone	0.443	0.448	-1.1	95	-0.07
23	S	Nitrobenzene-d5	0.305	0.305	0.0	96	-0.07
24		Nitrobenzene	0.329	0.341	-3.6	101	-0.07
25		Isophorone	0.616	0.604	1.9	91	-0.07
26	C	2-Nitrophenol	0.161	0.163	-1.2	89	-0.06
27		2,4-Dimethylphenol	0.293	0.297	-1.4	93	-0.06
28		bis(2-Chloroethoxy)methane	0.374	0.368	1.6	93	-0.07
29	C	2,4-Dichlorophenol	0.279	0.276	1.1	91	-0.06
30		1,2,4-Trichlorobenzene	0.313	0.305	2.6	91	-0.07
31		Naphthalene	1.027	1.001	2.5	92	-0.07
32		Benzoic acid	0.141	0.155	-9.9	93	-0.05
33		4-Chloroaniline	0.417	0.414	0.7	91	-0.06
34	C	Hexachlorobutadiene	0.177	0.173	2.3	93	-0.07
35		Caprolactam	0.082	0.088	-7.3	99	-0.06
36	C	4-Chloro-3-methylphenol	0.281	0.285	-1.4	96	-0.06
37		2-Methylnaphthalene	0.690	0.679	1.6	90	-0.06
38	I	Acenaphthene-d10	1.000	1.000	0.0	94	-0.07
39		1,2,4,5-Tetrachlorobenzene	0.597	0.537	10.1	85	-0.06
40	P	Hexachlorocyclopentadiene	0.254	0.243	4.3	83	-0.07
41	S	2,4,6-Tribromophenol	0.191	0.182	4.7	89	-0.07
42	C	2,4,6-Trichlorophenol	0.413	0.396	4.1	87	-0.06
43		2,4,5-Trichlorophenol	0.446	0.438	1.8	93	-0.06
44	S	2-Fluorobiphenyl	1.361	1.272	6.5	91	-0.07

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF032515\
Data File : BF077975.D
Acq On : 26 Mar 2015 3:35
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 26 05:23:28 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 26 00:45:24 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	1.599	1.515	5.3	89	-0.06
46		2-Chloronaphthalene	1.299	1.226	5.6	91	-0.07
47		2-Nitroaniline	0.323	0.336	-4.0	92	-0.07
48		Acenaphthylene	2.161	2.113	2.2	95	-0.07
49		Dimethylphthalate	1.483	1.446	2.5	94	-0.07
50		2,6-Dinitrotoluene	0.307	0.305	0.7	93	-0.07
51	C	Acenaphthene	1.239	1.172	5.4	89	-0.07
52		3-Nitroaniline	0.357	0.378	-5.9	97	-0.06
53	P	2,4-Dinitrophenol	0.093	0.103	-10.8	104	-0.06
54		Dibenzofuran	1.837	1.717	6.5	88	-0.07
55	P	4-Nitrophenol	0.265	0.268	-1.1	89	-0.06
56		2,4-Dinitrotoluene	0.401	0.403	-0.5	95	-0.07
57		Fluorene	1.526	1.447	5.2	89	-0.07
58		2,3,4,6-Tetrachlorophenol	0.344	0.340	1.2	92	-0.07
59		Diethylphthalate	1.445	1.436	0.6	94	-0.07
60		4-Chlorophenyl-phenylether	0.696	0.649	6.8	89	-0.07
61		4-Nitroaniline	0.356	0.412	-15.7	108	-0.06
62		Azobenzene	1.381	1.338	3.1	84	-0.07
63	I	Phenanthrene-d10	1.000	1.000	0.0	97	-0.08
64		4,6-Dinitro-2-methylphenol	0.078	0.097	-24.4	105	-0.07
65	c	n-Nitrosodiphenylamine	0.666	0.631	5.3	90	-0.07
66		4-Bromophenyl-phenylether	0.213	0.205	3.8	91	-0.07
67		Hexachlorobenzene	0.235	0.223	5.1	91	-0.07
68		Atrazine	0.187	0.173	7.5	85	-0.07
69	C	Pentachlorophenol	0.104	0.102	1.9	86	-0.07
70		Phenanthrene	1.148	1.144	0.3	96	-0.08
71		Anthracene	1.134	1.127	0.6	99	-0.07
72		Carbazole	1.079	1.102	-2.1	102	-0.07
73		Di-n-butylphthalate	1.210	1.226	-1.3	98	-0.07
74	C	Fluoranthene	1.266	1.302	-2.8	94	-0.08
75	I	Chrysene-d12	1.000	1.000	0.0	104	-0.18
76		Benzidine	0.546	0.555	-1.6	113	-0.08
77		Pyrene	1.258	1.175	6.6	90	-0.08
78	S	Terphenyl-d14	0.819	0.736	10.1	89	-0.09
79		Butylbenzylphthalate	0.496	0.497	-0.2	103	-0.13
80		Benzo(a)anthracene	1.186	1.154	2.7	105	-0.18
81		3,3'-Dichlorobenzidine	0.391	0.417	-6.6	117	-0.17
82		Chrysene	1.066	1.094	-2.6	108	-0.18
83		Bis(2-ethylhexyl)phthalate	0.751	0.746	0.7	106	-0.19
84	c	Di-n-octyl phthalate	1.284	1.340	-4.4	112	-0.32
85		Indeno(1,2,3-cd)pyrene	1.331	1.405	-5.6	119	-0.58#
86	I	Perylene-d12	1.000	1.000	0.0	112	-0.46
87		Benzo(b)fluoranthene	1.306	1.302	0.3	112	-0.38

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF032515\
Data File : BF077975.D
Acq On : 26 Mar 2015 3:35
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 26 05:23:28 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 26 00:45:24 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	1.020	0.983	3.6	113	-0.38
89 C	Benzo(a)pyrene	1.089	1.111	-2.0	115	-0.43
90	Dibenzo(a,h)anthracene	1.081	1.108	-2.5	117	-0.59#
91	Benzo(g,h,i)perylene	1.100	1.116	-1.5	114	-0.61#

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