Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF030915\
Data File : BF077656.D
Acq On : 9 Mar 2015 23:50
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 10 03:56:22 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 10 00:35:08 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	-0.03
2		1,4-Dioxane	0.508	0.596	-17.3	108	-0.05
3		Pyridine	1.454	1.445	0.6	85	-0.06
4		n-Nitrosodimethylamine	0.632	0.708	-12.0	105	-0.06
5	S	2-Fluorophenol	1.198	1.200	-0.2	90	-0.03
6		Aniline	2.129	2.171	-2.0	90	-0.03
7	S	Phenol-d6	1.540	1.536	0.3	88	-0.03
8		2-Chlorophenol	1.413	1.432	-1.3	91	-0.03
9		Benzaldehyde	0.935	0.888	5.0	88	-0.03
10	C	Phenol	1.828	1.704	6.8	79	-0.03
11		bis(2-Chloroethyl)ether	1.313	1.310	0.2	91	-0.03
12		1,3-Dichlorobenzene	1.539	1.549	-0.6	89	-0.03
13	C	1,4-Dichlorobenzene	1.566	1.583	-1.1	90	-0.03
14		1,2-Dichlorobenzene	1.489	1.478	0.7	87	-0.03
15		Benzyl Alcohol	1.171	1.150	1.8	87	-0.03
16		2,2'-oxybis(1-Chloropropane	1.877	1.670	11.0	78	-0.03
17		2-Methylphenol	1.105	1.156	-4.6	88	-0.03
18		Hexachloroethane	0.523	0.487	6.9	83	-0.03
19	P	n-Nitroso-di-n-propylamine	0.978	0.954	2.5	84	-0.03
20		3+4-Methylphenols	1.455	1.510	-3.8	90	-0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	91	-0.03
22		Acetophenone	0.470	0.469	0.2	93	-0.03
23	S	Nitrobenzene-d5	0.322	0.343	-6.5	94	-0.03
24		Nitrobenzene	0.338	0.350	-3.6	92	-0.03
25		Isophorone	0.638	0.626	1.9	84	-0.03
26	C	2-Nitrophenol	0.139	0.136	2.2	80	-0.03
27		2,4-Dimethylphenol	0.310	0.308	0.6	88	-0.03
28		bis(2-Chloroethoxy)methane	0.403	0.360	10.7	79	-0.03
29	C	2,4-Dichlorophenol	0.291	0.293	-0.7	89	-0.03
30		1,2,4-Trichlorobenzene	0.320	0.311	2.8	86	-0.03
31		Naphthalene	1.000	0.992	0.8	90	-0.03
32		Benzoic acid	0.151	0.026	82.8#	14#	0.00
33		4-Chloroaniline	0.445	0.452	-1.6	88	-0.03
34	C	Hexachlorobutadiene	0.183	0.176	3.8	85	-0.03
35		Caprolactam	0.089	0.095	-6.7	92	-0.03
36	C	4-Chloro-3-methylphenol	0.293	0.296	-1.0	86	-0.03
37		2-Methylnaphthalene	0.710	0.663	6.6	81	-0.03
38	I	Acenaphthene-d10	1.000	1.000	0.0	85	-0.03
39		1,2,4,5-Tetrachlorobenzene	0.574	0.549	4.4	78	-0.03
40	P	Hexachlorocyclopentadiene	0.308	0.124	59.7#	31#	-0.03
41	S	2,4,6-Tribromophenol	0.193	0.138	28.5#	57	-0.03
42	C	2,4,6-Trichlorophenol	0.380	0.378	0.5	78	-0.03
43		2,4,5-Trichlorophenol	0.406	0.444	-9.4	87	-0.03
44	S	2-Fluorobiphenyl	1.283	1.309	-2.0	86	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF030915\
Data File : BF077656.D
Acq On : 9 Mar 2015 23:50
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 10 03:56:22 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 10 00:35:08 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	1.515	1.469	3.0	79	-0.03
46		2-Chloronaphthalene	1.188	1.247	-5.0	88	-0.03
47		2-Nitroaniline	0.293	0.367	-25.3#	101	-0.03
48		Acenaphthylene	1.881	2.019	-7.3	89	-0.03
49		Dimethylphthalate	1.406	1.330	5.4	79	-0.03
50		2,6-Dinitrotoluene	0.282	0.312	-10.6	83	-0.03
51	C	Acenaphthene	1.123	1.106	1.5	81	-0.03
52		3-Nitroaniline	0.325	0.364	-12.0	87	-0.03
53	P	2,4-Dinitrophenol	0.086	0.006#	93.0#	6#	-0.03
54		Dibenzofuran	1.733	1.683	2.9	80	-0.03
55	P	4-Nitrophenol	0.261	0.203	22.2	59	-0.03
56		2,4-Dinitrotoluene	0.381	0.404	-6.0	79	-0.03
57		Fluorene	1.386	1.411	-1.8	83	-0.03
58		2,3,4,6-Tetrachlorophenol	0.327	0.229	30.0#	54	-0.03
59		Diethylphthalate	1.386	1.330	4.0	79	-0.03
60		4-Chlorophenyl-phenylether	0.668	0.638	4.5	78	-0.03
61		4-Nitroaniline	0.312	0.396	-26.9#	101	-0.03
62		Azobenzene	1.315	1.283	2.4	78	-0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	88	-0.04
64		4,6-Dinitro-2-methylphenol	0.075	0.010	86.7#	11#	-0.03
65	c	n-Nitrosodiphenylamine	0.664	0.630	5.1	81	-0.03
66		4-Bromophenyl-phenylether	0.234	0.207	11.5	78	-0.03
67		Hexachlorobenzene	0.251	0.230	8.4	82	-0.03
68		Atrazine	0.216	0.183	15.3	79	-0.03
69	C	Pentachlorophenol	0.132	0.033	75.0#	21#	-0.03
70		Phenanthrene	1.159	1.094	5.6	84	-0.04
71		Anthracene	1.150	1.150	0.0	89	-0.03
72		Carbazole	1.094	1.105	-1.0	85	-0.03
73		Di-n-butylphthalate	1.284	1.130	12.0	73	-0.04
74	C	Fluoranthene	1.266	1.264	0.2	86	-0.04
75	I	Chrysene-d12	1.000	1.000	0.0	86	-0.04
76		Benzidine	0.676	0.703	-4.0	97	-0.04
77		Pyrene	1.223	1.230	-0.6	88	-0.04
78	S	Terphenyl-d14	0.856	0.737	13.9	74	-0.04
79		Butylbenzylphthalate	0.503	0.480	4.6	78	-0.04
80		Benzo(a)anthracene	1.172	1.176	-0.3	88	-0.04
81		3,3'-Dichlorobenzidine	0.430	0.423	1.6	83	-0.04
82		Chrysene	1.101	1.097	0.4	89	-0.04
83		Bis(2-ethylhexyl)phthalate	0.807	0.646	20.0	68	-0.04
84	c	Di-n-octyl phthalate	1.348	1.154	14.4	70	-0.07
85		Indeno(1,2,3-cd)pyrene	1.255	1.477	-17.7	101	-0.12
86	I	Perylene-d12	1.000	1.000	0.0	90	-0.10
87		Benzo(b)fluoranthene	1.163	1.245	-7.1	91	-0.08

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF030915\
Data File : BF077656.D
Acq On : 9 Mar 2015 23:50
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 10 03:56:22 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 10 00:35:08 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	1.140	1.111	2.5	94	-0.09
89 C	Benzo(a)pyrene	1.108	1.111	-0.3	87	-0.09
90	Dibenzo(a,h)anthracene	1.056	1.215	-15.1	101	-0.12
91	Benzo(g,h,i)perylene	1.066	1.264	-18.6	105	-0.13

(#) = Out of Range SPCC's out = 1 CCC's out = 1