Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_F\DATA\BF012315\
Data File : BF077037.D
Acq On : 23 Jan 2015 14:04
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 23 22:17:35 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 23 15:21:24 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	94	0.00
2		1,4-Dioxane	0.521	0.451	13.4	85	0.00
3		Pyridine	1.348	1.411	-4.7	81	0.00
4		n-Nitrosodimethylamine	0.668	0.672	-0.6	93	0.00
5	S	2-Fluorophenol	1.216	1.151	5.3	84	0.00
6		Aniline	2.124	2.038	4.0	84	0.00
7	S	Phenol-d6	1.535	1.469	4.3	85	0.00
8		2-Chlorophenol	1.402	1.366	2.6	86	0.00
9		Benzaldehyde	0.894	0.843	5.7	101	0.00
10	C	Phenol	1.629	1.611	1.1	85	0.00
11		bis(2-Chloroethyl)ether	1.275	1.273	0.2	91	0.00
12		1,3-Dichlorobenzene	1.595	1.554	2.6	89	0.00
13	C	1,4-Dichlorobenzene	1.692	1.613	4.7	87	0.00
14		1,2-Dichlorobenzene	1.559	1.490	4.4	85	0.00
15		Benzyl Alcohol	1.242	1.220	1.8	85	0.00
16		2,2'-oxybis(1-Chloropropane	1.714	1.659	3.2	87	0.00
17		2-Methylphenol	1.146	1.128	1.6	86	0.00
18		Hexachloroethane	0.588	0.569	3.2	86	0.00
19	P	n-Nitroso-di-n-propylamine	1.074	1.046	2.6	85	0.00
20		3+4-Methylphenols	1.517	1.460	3.8	86	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	94	0.00
22		Acetophenone	0.490	0.485	1.0	87	0.00
23	S	Nitrobenzene-d5	0.361	0.366	-1.4	89	0.00
24		Nitrobenzene	0.368	0.360	2.2	84	0.00
25		Isophorone	0.657	0.675	-2.7	90	0.00
26	C	2-Nitrophenol	0.172	0.178	-3.5	85	0.00
27		2,4-Dimethylphenol	0.284	0.294	-3.5	88	0.00
28		bis(2-Chloroethoxy)methane	0.374	0.390	-4.3	89	0.00
29	C	2,4-Dichlorophenol	0.265	0.278	-4.9	90	0.00
30		1,2,4-Trichlorobenzene	0.294	0.296	-0.7	85	0.00
31		Naphthalene	1.030	1.014	1.6	86	0.00
32		Benzoic acid	0.158	0.150	5.1	82	0.00
33		4-Chloroaniline	0.416	0.420	-1.0	89	0.00
34	C	Hexachlorobutadiene	0.175	0.176	-0.6	87	0.00
35		Caprolactam	0.078	0.083	-6.4	88	0.00
36	C	4-Chloro-3-methylphenol	0.303	0.312	-3.0	90	0.00
37		2-Methylnaphthalene	0.703	0.704	-0.1	89	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	92	0.00
39		1,2,4,5-Tetrachlorobenzene	0.494	0.491	0.6	89	0.00
40	P	Hexachlorocyclopentadiene	0.202	0.263	-30.2#	97	0.00
41	S	2,4,6-Tribromophenol	0.162	0.179	-10.5	93	0.00
42	C	2,4,6-Trichlorophenol	0.360	0.365	-1.4	86	0.00
43		2,4,5-Trichlorophenol	0.367	0.394	-7.4	91	0.00
44	S	2-Fluorobiphenyl	1.314	1.312	0.2	88	0.00

Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_F\DATA\BF012315\
Data File : BF077037.D
Acq On : 23 Jan 2015 14:04
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 23 22:17:35 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 23 15:21:24 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	1.483	1.515	-2.2	88	0.00
46		2-Chloronaphthalene	1.220	1.231	-0.9	89	0.00
47		2-Nitroaniline	0.370	0.377	-1.9	86	0.00
48		Acenaphthylene	2.058	2.136	-3.8	90	0.00
49		Dimethylphthalate	1.418	1.436	-1.3	90	0.00
50		2,6-Dinitrotoluene	0.283	0.303	-7.1	88	0.00
51	C	Acenaphthene	1.168	1.172	-0.3	88	0.00
52		3-Nitroaniline	0.339	0.360	-6.2	87	0.00
53	P	2,4-Dinitrophenol	0.095	0.122	-28.4#	105	0.00
54		Dibenzofuran	1.701	1.721	-1.2	89	0.00
55	P	4-Nitrophenol	0.222	0.235	-5.9	90	0.00
56		2,4-Dinitrotoluene	0.384	0.426	-10.9	91	0.00
57		Fluorene	1.441	1.446	-0.3	89	0.00
58		2,3,4,6-Tetrachlorophenol	0.268	0.294	-9.7	93	0.00
59		Diethylphthalate	1.433	1.512	-5.5	92	0.00
60		4-Chlorophenyl-phenylether	0.590	0.605	-2.5	91	0.00
61		4-Nitroaniline	0.328	0.349	-6.4	88	0.00
62		Azobenzene	1.494	1.569	-5.0	92	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
64		4,6-Dinitro-2-methylphenol	0.104	0.132	-26.9#	99	0.00
65	c	n-Nitrosodiphenylamine	0.713	0.706	1.0	90	0.00
66		4-Bromophenyl-phenylether	0.203	0.202	0.5	89	0.00
67		Hexachlorobenzene	0.214	0.216	-0.9	95	0.00
68		Atrazine	0.216	0.226	-4.6	90	0.00
69	C	Pentachlorophenol	0.082	0.107	-30.5#	116	0.00
70		Phenanthrene	1.247	1.194	4.3	90	0.00
71		Anthracene	1.216	1.224	-0.7	95	0.00
72		Carbazole	1.180	1.152	2.4	90	0.00
73		Di-n-butylphthalate	1.365	1.467	-7.5	98	0.00
74	C	Fluoranthene	1.278	1.288	-0.8	92	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
76		Benzidine	0.640	0.599	6.4	96	0.00
77		Pyrene	1.371	1.351	1.5	92	0.00
78	S	Terphenyl-d14	0.869	0.871	-0.2	95	0.00
79		Butylbenzylphthalate	0.621	0.663	-6.8	96	0.00
80		Benzo(a)anthracene	1.255	1.246	0.7	93	0.00
81		3,3'-Dichlorobenzidine	0.399	0.421	-5.5	99	0.00
82		Chrysene	1.144	1.132	1.0	93	0.00
83		Bis(2-ethylhexyl)phthalate	0.859	0.924	-7.6	102	0.00
84	c	Di-n-octyl phthalate	1.491	1.662	-11.5	103	0.00
85		Indeno(1,2,3-cd)pyrene	1.213	1.246	-2.7	95	0.00
86	I	Perylene-d12	1.000	1.000	0.0	99	0.00
87		Benzo(b)fluoranthene	1.347	1.390	-3.2	107	0.00

Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_F\DATA\BF012315\
Data File : BF077037.D
Acq On : 23 Jan 2015 14:04
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 23 22:17:35 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 23 15:21:24 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	1.177	1.161	1.4	86	0.00
89 C	Benzo(a)pyrene	1.188	1.182	0.5	93	0.00
90	Dibenzo(a,h)anthracene	1.090	1.101	-1.0	94	0.00
91	Benzo(g,h,i)perylene	1.084	1.101	-1.6	95	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1