Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF020515\
Data File : BF077197.D
Acq On : 5 Feb 2015 23:23
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 06 11:22:39 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Feb 06 03:00:41 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	73	0.00
2		1,4-Dioxane	0.521	0.529	-1.5	79	0.00
3		Pyridine	1.348	1.360	-0.9	61	0.00
4		n-Nitrosodimethylamine	0.668	0.635	4.9	69	0.00
5	S	2-Fluorophenol	1.216	1.184	2.6	68	0.00
6		Aniline	2.124	1.963	7.6	64	0.00
7	S	Phenol-d6	1.535	1.429	6.9	65	0.00
8		2-Chlorophenol	1.402	1.335	4.8	65	0.00
9		Benzaldehyde	0.894	0.693	22.5	65	0.00
10	C	Phenol	1.629	1.427	12.4	59	-0.02
11		bis(2-Chloroethyl)ether	1.275	1.264	0.9	70	0.00
12		1,3-Dichlorobenzene	1.595	1.570	1.6	70	0.00
13	C	1,4-Dichlorobenzene	1.692	1.587	6.2	67	0.00
14		1,2-Dichlorobenzene	1.559	1.506	3.4	67	0.00
15		Benzyl Alcohol	1.242	1.143	8.0	63	0.00
16		2,2'-oxybis(1-Chloropropane	1.714	1.738	-1.4	72	0.00
17		2-Methylphenol	1.146	1.053	8.1	63	0.00
18		Hexachloroethane	0.588	0.599	-1.9	71	0.00
19	P	n-Nitroso-di-n-propylamine	1.074	0.992	7.6	63	0.00
20		3+4-Methylphenols	1.517	1.305	14.0	60	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	68	0.00
22		Acetophenone	0.490	0.498	-1.6	65	0.00
23	S	Nitrobenzene-d5	0.361	0.372	-3.0	65	0.00
24		Nitrobenzene	0.368	0.359	2.4	61	0.00
25		Isophorone	0.657	0.660	-0.5	63	0.00
26	C	2-Nitrophenol	0.172	0.164	4.7	57	0.00
27		2,4-Dimethylphenol	0.284	0.277	2.5	60	0.00
28		bis(2-Chloroethoxy)methane	0.374	0.382	-2.1	63	0.00
29	C	2,4-Dichlorophenol	0.265	0.236	10.9	55	0.00
30		1,2,4-Trichlorobenzene	0.294	0.313	-6.5	65	0.00
31		Naphthalene	1.030	1.052	-2.1	65	0.00
32		Benzoic acid	0.158	0.104	34.2#	41#	-0.02
33		4-Chloroaniline	0.416	0.380	8.7	58	0.00
34	C	Hexachlorobutadiene	0.175	0.182	-4.0	65	0.00
35		Caprolactam	0.078	0.067	14.1	52	-0.03
36	C	4-Chloro-3-methylphenol	0.303	0.285	5.9	59	0.00
37		2-Methylnaphthalene	0.703	0.727	-3.4	66	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	63	0.00
39		1,2,4,5-Tetrachlorobenzene	0.494	0.518	-4.9	64	0.00
40	P	Hexachlorocyclopentadiene	0.202	0.250	-23.8	63	0.00
41	S	2,4,6-Tribromophenol	0.162	0.159	1.9	57	0.00
42	C	2,4,6-Trichlorophenol	0.360	0.363	-0.8	59	0.00
43		2,4,5-Trichlorophenol	0.367	0.338	7.9	54	0.00
44	S	2-Fluorobiphenyl	1.314	1.427	-8.6	65	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF020515\
Data File : BF077197.D
Acq On : 5 Feb 2015 23:23
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 06 11:22:39 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Feb 06 03:00:41 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	1.587	-7.0	63	0.00
46		2-Chloronaphthalene	1.220	1.302	-6.7	65	0.00
47		2-Nitroaniline	0.370	0.381	-3.0	59	0.00
48		Acenaphthylene	2.058	2.207	-7.2	64	0.00
49		Dimethylphthalate	1.418	1.549	-9.2	67	0.00
50		2,6-Dinitrotoluene	0.283	0.303	-7.1	61	0.00
51	C	Acenaphthene	1.168	1.209	-3.5	62	0.00
52		3-Nitroaniline	0.339	0.345	-1.8	58	0.00
53	P	2,4-Dinitrophenol	0.095	0.067	29.5#	40#	0.02
54		Dibenzofuran	1.701	1.795	-5.5	64	0.00
55	P	4-Nitrophenol	0.222	0.228	-2.7	60	0.02
56		2,4-Dinitrotoluene	0.384	0.412	-7.3	61	0.00
57		Fluorene	1.441	1.521	-5.6	64	0.00
58		2,3,4,6-Tetrachlorophenol	0.268	0.190	29.1#	41#	0.00
59		Diethylphthalate	1.433	1.596	-11.4	67	0.00
60		4-Chlorophenyl-phenylether	0.590	0.646	-9.5	67	0.00
61		4-Nitroaniline	0.328	0.333	-1.5	58	0.00
62		Azobenzene	1.494	1.690	-13.1	68	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	67	0.00
64		4,6-Dinitro-2-methylphenol	0.104	0.102	1.9	52	0.00
65	c	n-Nitrosodiphenylamine	0.713	0.727	-2.0	64	0.00
66		4-Bromophenyl-phenylether	0.203	0.207	-2.0	63	0.00
67		Hexachlorobenzene	0.214	0.227	-6.1	68	0.00
68		Atrazine	0.216	0.212	1.9	58	0.00
69	C	Pentachlorophenol	0.082	0.075	8.5	56	0.00
70		Phenanthrene	1.247	1.261	-1.1	65	0.00
71		Anthracene	1.216	1.221	-0.4	65	0.00
72		Carbazole	1.180	1.227	-4.0	66	0.00
73		Di-n-butylphthalate	1.365	1.519	-11.3	69	0.00
74	C	Fluoranthene	1.278	1.343	-5.1	66	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	73	0.00
76		Benzidine	0.640	0.584	8.8	66	0.00
77		Pyrene	1.371	1.428	-4.2	69	0.00
78	S	Terphenyl-d14	0.869	0.905	-4.1	70	0.00
79		Butylbenzylphthalate	0.621	0.672	-8.2	69	0.00
80		Benzo(a)anthracene	1.255	1.304	-3.9	70	0.00
81		3,3'-Dichlorobenzidine	0.399	0.409	-2.5	69	0.00
82		Chrysene	1.144	1.178	-3.0	69	0.00
83		Bis(2-ethylhexyl)phthalate	0.859	0.978	-13.9	77	0.00
84	c	Di-n-octyl phthalate	1.491	1.681	-12.7	74	-0.02
85		Indeno(1,2,3-cd)pyrene	1.213	1.303	-7.4	71	-0.09
86	I	Perylene-d12	1.000	1.000	0.0	70	-0.03
87		Benzo(b)fluoranthene	1.347	1.354	-0.5	74	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF020515\
Data File : BF077197.D
Acq On : 5 Feb 2015 23:23
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 06 11:22:39 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Feb 06 03:00:41 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	1.374	-16.7	72	0.00
89 C	Benzo(a)pyrene	1.188	1.272	-7.1	71	-0.02
90	Dibenzo(a,h)anthracene	1.090	1.162	-6.6	70	-0.09
91	Benzo(g,h,i)perylene	1.084	1.196	-10.3	73	-0.07

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