Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF032717\
Data File : BF093974.D
Acq On : 28 Mar 2017 6:06
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 28 06:43:11 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 27 16:10:49 2017
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	93	-0.01
2		1,4-Dioxane	0.569	0.545	4.2	87	-0.03
3		Pyridine	1.550	1.481	4.5	85	-0.02
4		n-Nitrosodimethylamine	0.638	0.621	2.7	86	-0.02
5	S	2-Fluorophenol	1.184	1.167	1.4	91	-0.01
6		Aniline	2.038	1.948	4.4	86	-0.01
7	S	Phenol-d6	1.465	1.466	-0.1	92	-0.01
8		2-Chlorophenol	1.302	1.315	-1.0	91	-0.01
9		Benzaldehyde	0.989	0.947	4.2	88	-0.01
10	C	Phenol	1.667	1.635	1.9	86	0.00
11		bis(2-Chloroethyl)ether	1.303	1.282	1.6	90	-0.01
12		1,3-Dichlorobenzene	1.460	1.462	-0.1	91	-0.01
13	C	1,4-Dichlorobenzene	1.479	1.474	0.3	91	-0.01
14		1,2-Dichlorobenzene	1.362	1.353	0.7	91	-0.01
15		Benzyl Alcohol	1.072	1.061	1.0	88	-0.01
16		2,2'-oxybis(1-Chloropropane	2.007	1.874	6.6	84	0.00
17		2-Methylphenol	1.115	1.115	0.0	91	0.00
18		Hexachloroethane	0.520	0.489	6.0	84	-0.01
19	P	n-Nitroso-di-n-propylamine	0.941	0.897	4.7	87	-0.01
20		3+4-Methylphenols	1.350	1.368	-1.3	94	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	91	-0.01
22		Acetophenone	0.446	0.456	-2.2	96	-0.01
23	S	Nitrobenzene-d5	0.350	0.350	0.0	90	-0.01
24		Nitrobenzene	0.346	0.342	1.2	88	-0.01
25		Isophorone	0.616	0.605	1.8	89	-0.01
26	C	2-Nitrophenol	0.165	0.160	3.0	82	-0.01
27		2,4-Dimethylphenol	0.284	0.283	0.4	90	-0.01
28		bis(2-Chloroethoxy)methane	0.406	0.396	2.5	88	-0.01
29	C	2,4-Dichlorophenol	0.266	0.270	-1.5	91	0.00
30		1,2,4-Trichlorobenzene	0.274	0.274	0.0	91	-0.01
31		Naphthalene	0.962	0.944	1.9	90	-0.01
32		Benzoic acid	0.189	0.096	49.2#	41#	-0.03
33		4-Chloroaniline	0.390	0.392	-0.5	90	-0.01
34	C	Hexachlorobutadiene	0.140	0.138	1.4	89	-0.01
35		Caprolactam	0.085	0.087	-2.4	90	-0.01
36	C	4-Chloro-3-methylphenol	0.277	0.279	-0.7	90	0.00
37		2-Methylnaphthalene	0.641	0.633	1.2	90	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	89	0.00
39		1,2,4,5-Tetrachlorobenzene	0.547	0.571	-4.4	93	-0.01
40	P	Hexachlorocyclopentadiene	0.256	0.125	51.2#	40#	-0.01
41	S	2,4,6-Tribromophenol	0.158	0.153	3.2	84	-0.01
42	C	2,4,6-Trichlorophenol	0.387	0.394	-1.8	90	-0.01
43		2,4,5-Trichlorophenol	0.419	0.424	-1.2	86	0.00
44	S	2-Fluorobiphenyl	1.311	1.337	-2.0	91	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF032717\
Data File : BF093974.D
Acq On : 28 Mar 2017 6:06
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 28 06:43:11 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 27 16:10:49 2017
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	1.613	1.628	-0.9	90	0.00
46		2-Chloronaphthalene	1.263	1.278	-1.2	91	-0.01
47		2-Nitroaniline	0.375	0.403	-7.5	92	-0.01
48		Acenaphthylene	2.099	2.098	0.0	89	0.00
49		Dimethylphthalate	1.422	1.471	-3.4	92	0.00
50		2,6-Dinitrotoluene	0.326	0.336	-3.1	88	0.00
51	C	Acenaphthene	1.225	1.208	1.4	89	-0.01
52		3-Nitroaniline	0.383	0.407	-6.3	93	0.00
53	P	2,4-Dinitrophenol	0.155	0.021#	86.5#	12#	-0.01
54		Dibenzofuran	1.667	1.647	1.2	87	-0.01
55	P	4-Nitrophenol	0.300	0.270	10.0	76	0.00
56		2,4-Dinitrotoluene	0.409	0.411	-0.5	85	-0.01
57		Fluorene	1.382	1.305	5.6	90	-0.01
58		2,3,4,6-Tetrachlorophenol	0.287	0.260	9.4	78	-0.01
59		Diethylphthalate	1.405	1.424	-1.4	90	-0.01
60		4-Chlorophenyl-phenylether	0.589	0.561	4.8	89	0.00
61		4-Nitroaniline	0.367	0.396	-7.9	92	-0.01
62		Azobenzene	1.329	1.285	3.3	85	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	84	-0.01
64		4,6-Dinitro-2-methylphenol	0.122	0.028	77.0#	18#	-0.02
65	c	n-Nitrosodiphenylamine	0.692	0.734	-6.1	90	0.00
66		4-Bromophenyl-phenylether	0.197	0.206	-4.6	88	0.00
67		Hexachlorobenzene	0.199	0.205	-3.0	87	0.00
68		Atrazine	0.207	0.217	-4.8	87	0.00
69	C	Pentachlorophenol	0.124	0.085	31.5#	56	-0.01
70		Phenanthrene	1.113	1.109	0.4	85	-0.01
71		Anthracene	1.146	1.144	0.2	85	-0.01
72		Carbazole	1.175	1.137	3.2	82	-0.01
73		Di-n-butylphthalate	1.222	1.276	-4.4	87	-0.01
74	C	Fluoranthene	1.139	1.009	11.4	76	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	68	-0.01
76		Benzidine	0.792	0.773	2.4	65	-0.01
77		Pyrene	1.451	1.559	-7.4	73	-0.01
78	S	Terphenyl-d14	0.892	1.003	-12.4	78	-0.01
79		Butylbenzylphthalate	0.618	0.716	-15.9	75	-0.01
80		Benzo(a)anthracene	1.166	1.172	-0.5	68	-0.01
81		3,3'-Dichlorobenzidine	0.417	0.439	-5.3	68	-0.01
82		Chrysene	1.082	1.097	-1.4	69	-0.02
83		Bis(2-ethylhexyl)phthalate	0.844	0.955	-13.2	74	0.00
84	c	Di-n-octyl phthalate	1.410	1.544	-9.5	70	-0.01
85		Indeno(1,2,3-cd)pyrene	0.937	1.030	-9.9	77	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	77	-0.01
87		Benzo(b)fluoranthene	1.226	1.231	-0.4	76	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF032717\
Data File : BF093974.D
Acq On : 28 Mar 2017 6:06
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 28 06:43:11 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 27 16:10:49 2017
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	1.199	1.134	5.4	71	-0.01
89 C	Benzo(a)pyrene	1.129	1.141	-1.1	76	-0.01
90	Dibenzo(a,h)anthracene	0.956	0.971	-1.6	79	-0.02
91	Benzo(g,h,i)perylene	0.964	0.926	3.9	75	-0.02

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