Evaluate Continuing Calibration Report

Data Path : V:\HPCHEM1\BNA F\DATA\BF070916\
Data File : BF088864.D
Acq On : 9 Jul 2016 11:40
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 11 04:57:41 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 08 18:51:20 2016
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	75	-0.01
2		1,4-Dioxane	0.662	0.601	9.2	70	0.00
3		Pyridine	1.920	1.659	13.6	67	-0.01
4		n-Nitrosodimethylamine	0.733	0.630	14.1	67	0.00
5	S	2-Fluorophenol	1.334	1.397	-4.7	78	-0.01
6		Aniline	2.513	2.273	9.6	68	-0.01
7	S	Phenol-d6	1.724	1.634	5.2	74	-0.01
8		2-Chlorophenol	1.434	1.508	-5.2	85	0.00
9		Benzaldehyde	1.168	0.935	19.9	65	-0.01
10	C	Phenol	1.968	1.647	16.3	63	-0.01
11		bis(2-Chloroethyl)ether	1.468	1.465	0.2	80	-0.01
12		1,3-Dichlorobenzene	1.578	1.531	3.0	80	-0.01
13	C	1,4-Dichlorobenzene	1.567	1.645	-5.0	84	-0.01
14		1,2-Dichlorobenzene	1.466	1.448	1.2	83	0.00
15		Benzyl Alcohol	1.269	1.208	4.8	70	-0.01
16		2,2'-oxybis(1-Chloropropane	2.125	1.716	19.2	62	0.00
17		2-Methylphenol	1.170	1.205	-3.0	78	0.00
18		Hexachloroethane	0.606	0.596	1.7	76	-0.01
19	P	n-Nitroso-di-n-propylamine	1.158	0.996	14.0	75	-0.01
20		3+4-Methylphenols	1.404	1.478	-5.3	85	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	78	-0.01
22		Acetophenone	0.485	0.457	5.8	75	-0.01
23	S	Nitrobenzene-d5	0.382	0.345	9.7	74	-0.01
24		Nitrobenzene	0.385	0.378	1.8	80	-0.01
25		Isophorone	0.707	0.648	8.3	75	-0.01
26	C	2-Nitrophenol	0.158	0.180	-13.9	96	0.00
27		2,4-Dimethylphenol	0.329	0.334	-1.5	78	0.00
28		bis(2-Chloroethoxy)methane	0.460	0.420	8.7	77	-0.01
29	C	2,4-Dichlorophenol	0.266	0.280	-5.3	87	-0.01
30		1,2,4-Trichlorobenzene	0.291	0.293	-0.7	79	-0.01
31		Naphthalene	0.986	1.026	-4.1	81	-0.01
32		Benzoic acid	0.239	0.226	5.4	75	0.00
33		4-Chloroaniline	0.439	0.418	4.8	75	-0.01
34	C	Hexachlorobutadiene	0.156	0.173	-10.9	87	0.00
35		Caprolactam	0.090	0.083	7.8	69	0.00
36	C	4-Chloro-3-methylphenol	0.314	0.296	5.7	75	-0.01
37		2-Methylnaphthalene	0.635	0.572	9.9	80	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	80	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.665	0.729	-9.6	85	-0.01
40	P	Hexachlorocyclopentadiene	0.327	0.305	6.7	75	-0.01
41	S	2,4,6-Tribromophenol	0.186	0.225	-21.0	93	-0.01
42	C	2,4,6-Trichlorophenol	0.439	0.434	1.1	86	0.00
43		2,4,5-Trichlorophenol	0.377	0.367	2.7	75	-0.01
44	S	2-Fluorobiphenyl	1.266	1.321	-4.3	92	-0.01

Evaluate Continuing Calibration Report

Data Path : V:\HPCHEM1\BNA F\DATA\BF070916\
Data File : BF088864.D
Acq On : 9 Jul 2016 11:40
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 11 04:57:41 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 08 18:51:20 2016
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	1.656	1.697	-2.5	80	-0.01
46		2-Chloronaphthalene	1.300	1.373	-5.6	82	-0.01
47		2-Nitroaniline	0.461	0.394	14.5	61	-0.01
48		Acenaphthylene	2.069	2.106	-1.8	81	-0.01
49		Dimethylphthalate	1.425	1.473	-3.4	90	-0.01
50		2,6-Dinitrotoluene	0.316	0.354	-12.0	96	-0.01
51	C	Acenaphthene	1.268	1.283	-1.2	85	-0.01
52		3-Nitroaniline	0.401	0.423	-5.5	74	0.00
53	P	2,4-Dinitrophenol	0.139	0.165	-18.7	97	0.00
54		Dibenzofuran	1.660	1.676	-1.0	81	-0.01
55	P	4-Nitrophenol	0.305	0.330	-8.2	78	0.00
56		2,4-Dinitrotoluene	0.413	0.455	-10.2	93	0.00
57		Fluorene	1.279	1.486	-16.2	89	-0.01
58		2,3,4,6-Tetrachlorophenol	0.292	0.289	1.0	77	-0.01
59		Diethylphthalate	1.430	1.582	-10.6	89	0.00
60		4-Chlorophenyl-phenylether	0.594	0.582	2.0	85	-0.01
61		4-Nitroaniline	0.386	0.388	-0.5	87	0.00
62		Azobenzene	1.631	1.466	10.1	75	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	83	-0.01
64		4,6-Dinitro-2-methylphenol	0.107	0.137	-28.0#	96	0.00
65	c	n-Nitrosodiphenylamine	0.677	0.686	-1.3	81	-0.01
66		4-Bromophenyl-phenylether	0.202	0.215	-6.4	90	-0.01
67		Hexachlorobenzene	0.223	0.245	-9.9	91	0.00
68		Atrazine	0.211	0.182	13.7	68	-0.01
69	C	Pentachlorophenol	0.132	0.129	2.3	76	-0.01
70		Phenanthrene	1.093	0.981	10.2	78	-0.01
71		Anthracene	1.087	1.173	-7.9	89	0.00
72		Carbazole	1.023	0.963	5.9	87	-0.01
73		Di-n-butylphthalate	1.190	1.146	3.7	84	-0.01
74	C	Fluoranthene	1.108	1.217	-9.8	87	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	88	-0.01
76		Benzidine	1.011	1.045	-3.4	88	0.00
77		Pyrene	1.696	1.664	1.9	82	0.00
78	S	Terphenyl-d14	0.898	0.925	-3.0	92	0.00
79		Butylbenzylphthalate	0.756	0.682	9.8	80	-0.01
80		Benzo(a)anthracene	1.288	1.295	-0.5	88	0.00
81		3,3'-Dichlorobenzidine	0.484	0.422	12.8	79	-0.01
82		Chrysene	1.274	1.009	20.8	70	-0.01
83		Bis(2-ethylhexyl)phthalate	0.864	0.943	-9.1	91	0.00
84	c	Di-n-octyl phthalate	1.467	1.552	-5.8	89	0.00
85		Indeno(1,2,3-cd)pyrene	0.980	0.960	2.0	90	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	84	-0.01
87		Benzo(b)fluoranthene	1.255	1.292	-2.9	91	0.00

Evaluate Continuing Calibration Report

Data Path : V:\HPCHEM1\BNA F\DATA\BF070916\
Data File : BF088864.D
Acq On : 9 Jul 2016 11:40
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 11 04:57:41 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 08 18:51:20 2016
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	1.092	0.968	11.4	75	-0.01
89 C	Benzo(a)pyrene	1.062	1.040	2.1	82	-0.01
90	Dibenzo(a,h)anthracene	0.887	0.929	-4.7	90	-0.01
91	Benzo(g,h,i)perylene	0.918	0.921	-0.3	86	-0.01

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