Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF062516\
Data File : BF088367.D
Acq On : 25 Jun 2016 10:29
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 26 02:07:13 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 22 14:55:14 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	76	0.01
2		1,4-Dioxane	0.568	0.573	-0.9	79	0.01
3		Pyridine	1.342	1.315	2.0	73	0.03
4		n-Nitrosodimethylamine	0.568	0.466	18.0	63	0.01
5	S	2-Fluorophenol	1.170	1.125	3.8	75	0.01
6		Aniline	2.018	1.660	17.7	63	0.00
7	S	Phenol-d6	1.418	1.314	7.3	74	0.00
8		2-Chlorophenol	1.385	1.395	-0.7	79	0.01
9		Benzaldehyde	0.923	0.851	7.8	75	0.01
10	C	Phenol	1.665	1.756	-5.5	96	0.00
11		bis(2-Chloroethyl)ether	1.243	1.358	-9.3	90	0.00
12		1,3-Dichlorobenzene	1.486	1.439	3.2	74	0.01
13	C	1,4-Dichlorobenzene	1.497	1.538	-2.7	83	0.01
14		1,2-Dichlorobenzene	1.361	1.320	3.0	73	0.01
15		Benzyl Alcohol	1.109	0.990	10.7	66	0.00
16		2,2'-oxybis(1-Chloropropane	1.680	1.261	24.9	57	0.01
17		2-Methylphenol	1.123	1.009	10.2	67	0.00
18		Hexachloroethane	0.531	0.521	1.9	75	0.00
19	P	n-Nitroso-di-n-propylamine	0.907	0.949	-4.6	87	0.01
20		3+4-Methylphenols	1.310	1.313	-0.2	83	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	81	0.01
22		Acetophenone	0.447	0.459	-2.7	85	0.01
23	S	Nitrobenzene-d5	0.343	0.317	7.6	73	0.00
24		Nitrobenzene	0.332	0.338	-1.8	89	0.01
25		Isophorone	0.634	0.614	3.2	76	0.00
26	C	2-Nitrophenol	0.178	0.182	-2.2	72	0.00
27		2,4-Dimethylphenol	0.327	0.281	14.1	67	0.01
28		bis(2-Chloroethoxy)methane	0.393	0.400	-1.8	80	0.01
29	C	2,4-Dichlorophenol	0.273	0.274	-0.4	79	0.01
30		1,2,4-Trichlorobenzene	0.297	0.277	6.7	77	0.01
31		Naphthalene	0.990	0.933	5.8	80	0.01
32		Benzoic acid	0.180	0.153	15.0	59	-0.01
33		4-Chloroaniline	0.442	0.418	5.4	71	0.00
34	C	Hexachlorobutadiene	0.157	0.140	10.8	69	0.01
35		Caprolactam	0.090	0.086	4.4	69	0.00
36	C	4-Chloro-3-methylphenol	0.292	0.296	-1.4	83	0.01
37		2-Methylnaphthalene	0.652	0.589	9.7	69	0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	74	0.01
39		1,2,4,5-Tetrachlorobenzene	0.583	0.568	2.6	83	0.01
40	P	Hexachlorocyclopentadiene	0.323	0.344	-6.5	78	0.01
41	S	2,4,6-Tribromophenol	0.211	0.181	14.2	62	0.01
42	C	2,4,6-Trichlorophenol	0.400	0.351	12.3	64	0.01
43		2,4,5-Trichlorophenol	0.416	0.382	8.2	70	0.00
44	S	2-Fluorobiphenyl	1.502	1.328	11.6	79	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF062516\
Data File : BF088367.D
Acq On : 25 Jun 2016 10:29
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 26 02:07:13 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 22 14:55:14 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	1.613	1.592	1.3	79	0.01
46		2-Chloronaphthalene	1.294	1.267	2.1	78	0.01
47		2-Nitroaniline	0.382	0.376	1.6	67	0.01
48		Acenaphthylene	2.059	1.985	3.6	73	0.01
49		Dimethylphthalate	1.449	1.542	-6.4	87	0.01
50		2,6-Dinitrotoluene	0.332	0.313	5.7	77	0.00
51	C	Acenaphthene	1.284	1.187	7.6	69	0.01
52		3-Nitroaniline	0.383	0.367	4.2	68	0.00
53	P	2,4-Dinitrophenol	0.122	0.179	-46.7#	79	0.01
54		Dibenzofuran	1.769	1.717	2.9	71	0.01
55	P	4-Nitrophenol	0.297	0.337	-13.5	74	0.01
56		2,4-Dinitrotoluene	0.426	0.440	-3.3	83	0.01
57		Fluorene	1.378	1.323	4.0	79	0.01
58		2,3,4,6-Tetrachlorophenol	0.327	0.316	3.4	68	0.01
59		Diethylphthalate	1.466	1.333	9.1	71	0.01
60		4-Chlorophenyl-phenylether	0.588	0.525	10.7	72	0.01
61		4-Nitroaniline	0.363	0.372	-2.5	69	0.00
62		Azobenzene	1.450	1.409	2.8	71	0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	76	0.01
64		4,6-Dinitro-2-methylphenol	0.108	0.115	-6.5	62	0.01
65	c	n-Nitrosodiphenylamine	0.664	0.636	4.2	72	0.01
66		4-Bromophenyl-phenylether	0.215	0.216	-0.5	74	0.01
67		Hexachlorobenzene	0.223	0.202	9.4	75	0.01
68		Atrazine	0.201	0.180	10.4	62	0.01
69	C	Pentachlorophenol	0.118	0.123	-4.2	71	0.01
70		Phenanthrene	1.049	1.025	2.3	83	0.01
71		Anthracene	1.059	1.129	-6.6	77	0.01
72		Carbazole	0.991	1.044	-5.3	88	0.01
73		Di-n-butylphthalate	1.254	1.231	1.8	74	0.01
74	C	Fluoranthene	1.108	1.020	7.9	68	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	72	0.01
76		Benzidine	0.554	0.440	20.6	57	0.01
77		Pyrene	1.484	1.482	0.1	73	0.01
78	S	Terphenyl-d14	0.879	0.817	7.1	70	0.01
79		Butylbenzylphthalate	0.656	0.778	-18.6	82	0.01
80		Benzo(a)anthracene	1.140	1.041	8.7	72	0.02
81		3,3'-Dichlorobenzidine	0.306	0.327	-6.9	72	0.01
82		Chrysene	1.072	1.171	-9.2	84	0.01
83		Bis(2-ethylhexyl)phthalate	0.799	0.856	-7.1	79	0.01
84	c	Di-n-octyl phthalate	1.298	1.542	-18.8	87	0.01
85		Indeno(1,2,3-cd)pyrene	0.996	0.958	3.8	70	0.01
86	I	Perylene-d12	1.000	1.000	0.0	73	0.01
87		Benzo(b)fluoranthene	1.394	1.133	18.7	56	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF062516\
Data File : BF088367.D
Acq On : 25 Jun 2016 10:29
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 26 02:07:13 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 22 14:55:14 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	1.052	1.207	-14.7	103	0.00
89 C	Benzo(a)pyrene	1.128	1.097	2.7	70	0.01
90	Dibenzo(a,h)anthracene	1.047	0.995	5.0	69	0.01
91	Benzo(g,h,i)perylene	1.078	1.068	0.9	71	0.02

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