Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF061816\
Data File : BF088127.D
Acq On : 18 Jun 2016 11:52
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 19 07:58:25 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 17 16:09:13 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	72	-0.01
2		1,4-Dioxane	0.568	0.577	-1.6	75	0.00
3		Pyridine	1.342	1.349	-0.5	71	0.02
4		n-Nitrosodimethylamine	0.568	0.516	9.2	66	0.00
5	S	2-Fluorophenol	1.170	1.120	4.3	70	0.00
6		Aniline	2.018	1.714	15.1	61	-0.01
7	S	Phenol-d6	1.418	1.261	11.1	67	-0.01
8		2-Chlorophenol	1.385	1.374	0.8	73	-0.01
9		Benzaldehyde	0.923	0.711	23.0	59	-0.01
10	C	Phenol	1.665	1.368	17.8	70	-0.01
11		bis(2-Chloroethyl)ether	1.243	1.321	-6.3	82	-0.01
12		1,3-Dichlorobenzene	1.486	1.469	1.1	71	-0.01
13	C	1,4-Dichlorobenzene	1.497	1.452	3.0	73	-0.01
14		1,2-Dichlorobenzene	1.361	1.353	0.6	70	-0.01
15		Benzyl Alcohol	1.109	0.951	14.2	60	-0.01
16		2,2'-oxybis(1-Chloropropane	1.680	1.449	13.7	62	-0.01
17		2-Methylphenol	1.123	1.029	8.4	64	0.00
18		Hexachloroethane	0.531	0.534	-0.6	72	-0.01
19	P	n-Nitroso-di-n-propylamine	0.907	0.861	5.1	74	-0.01
20		3+4-Methylphenols	1.310	1.269	3.1	75	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	74	-0.01
22		Acetophenone	0.447	0.438	2.0	74	-0.01
23	S	Nitrobenzene-d5	0.343	0.301	12.2	63	-0.02
24		Nitrobenzene	0.332	0.347	-4.5	84	-0.01
25		Isophorone	0.634	0.580	8.5	66	-0.02
26	C	2-Nitrophenol	0.178	0.184	-3.4	67	-0.01
27		2,4-Dimethylphenol	0.327	0.283	13.5	62	-0.01
28		bis(2-Chloroethoxy)methane	0.393	0.407	-3.6	74	-0.01
29	C	2,4-Dichlorophenol	0.273	0.261	4.4	69	-0.01
30		1,2,4-Trichlorobenzene	0.297	0.291	2.0	74	-0.01
31		Naphthalene	0.990	0.935	5.6	73	-0.01
32		Benzoic acid	0.180	0.134	25.6#	47#	-0.02
33		4-Chloroaniline	0.442	0.425	3.8	66	-0.01
34	C	Hexachlorobutadiene	0.157	0.149	5.1	67	-0.01
35		Caprolactam	0.090	0.080	11.1	59	-0.02
36	C	4-Chloro-3-methylphenol	0.292	0.255	12.7	66	-0.01
37		2-Methylnaphthalene	0.652	0.603	7.5	64	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	67	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.583	0.602	-3.3	79	-0.01
40	P	Hexachlorocyclopentadiene	0.323	0.290	10.2	59	-0.01
41	S	2,4,6-Tribromophenol	0.211	0.178	15.6	55	-0.01
42	C	2,4,6-Trichlorophenol	0.400	0.369	7.8	60	-0.01
43		2,4,5-Trichlorophenol	0.416	0.394	5.3	65	-0.01
44	S	2-Fluorobiphenyl	1.502	1.180	21.4	63	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF061816\
Data File : BF088127.D
Acq On : 18 Jun 2016 11:52
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 19 07:58:25 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 17 16:09:13 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	1.706	-5.8	77	-0.01
46		2-Chloronaphthalene	1.294	1.300	-0.5	72	-0.01
47		2-Nitroaniline	0.382	0.368	3.7	59	-0.01
48		Acenaphthylene	2.059	1.990	3.4	66	-0.01
49		Dimethylphthalate	1.449	1.467	-1.2	75	-0.01
50		2,6-Dinitrotoluene	0.332	0.305	8.1	68	-0.02
51	C	Acenaphthene	1.284	1.178	8.3	62	-0.01
52		3-Nitroaniline	0.383	0.361	5.7	61	-0.01
53	P	2,4-Dinitrophenol	0.122	0.141	-15.6	56	-0.01
54		Dibenzofuran	1.769	1.704	3.7	63	-0.01
55	P	4-Nitrophenol	0.297	0.314	-5.7	62	0.00
56		2,4-Dinitrotoluene	0.426	0.365	14.3	62	-0.02
57		Fluorene	1.378	1.309	5.0	71	-0.01
58		2,3,4,6-Tetrachlorophenol	0.327	0.310	5.2	60	-0.01
59		Diethylphthalate	1.466	1.471	-0.3	70	-0.01
60		4-Chlorophenyl-phenylether	0.588	0.522	11.2	65	-0.02
61		4-Nitroaniline	0.363	0.382	-5.2	64	-0.01
62		Azobenzene	1.450	1.381	4.8	63	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	65	-0.01
64		4,6-Dinitro-2-methylphenol	0.108	0.105	2.8	48#	-0.02
65	c	n-Nitrosodiphenylamine	0.664	0.683	-2.9	66	-0.01
66		4-Bromophenyl-phenylether	0.215	0.215	0.0	63	-0.01
67		Hexachlorobenzene	0.223	0.200	10.3	64	-0.02
68		Atrazine	0.201	0.208	-3.5	62	-0.01
69	C	Pentachlorophenol	0.118	0.132	-11.9	65	-0.01
70		Phenanthrene	1.049	0.969	7.6	67	-0.02
71		Anthracene	1.059	1.125	-6.2	66	-0.01
72		Carbazole	0.991	1.054	-6.4	76	-0.01
73		Di-n-butylphthalate	1.254	1.202	4.1	62	-0.01
74	C	Fluoranthene	1.108	1.055	4.8	61	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	60	-0.02
76		Benzidine	0.554	0.503	9.2	54	-0.01
77		Pyrene	1.484	1.453	2.1	59	-0.02
78	S	Terphenyl-d14	0.879	0.859	2.3	61	-0.02
79		Butylbenzylphthalate	0.656	0.710	-8.2	62	-0.02
80		Benzo(a)anthracene	1.140	1.062	6.8	61	-0.02
81		3,3'-Dichlorobenzidine	0.306	0.313	-2.3	57	-0.02
82		Chrysene	1.072	1.084	-1.1	65	-0.02
83		Bis(2-ethylhexyl)phthalate	0.799	0.848	-6.1	65	-0.02
84	c	Di-n-octyl phthalate	1.298	1.589	-22.4#	74	-0.02
85		Indeno(1,2,3-cd)pyrene	0.996	0.911	8.5	55	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	59	-0.02
87		Benzo(b)fluoranthene	1.394	1.180	15.4	47#	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF061816\
Data File : BF088127.D
Acq On : 18 Jun 2016 11:52
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 19 07:58:25 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 17 16:09:13 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	1.205	-14.5	82	-0.02
89 C	Benzo(a)pyrene	1.128	1.139	-1.0	58	-0.02
90	Dibenzo(a,h)anthracene	1.047	0.993	5.2	55	-0.03
91	Benzo(g,h,i)perylene	1.078	1.037	3.8	55	-0.05

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