Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111516\
Data File : BF091185.D
Acq On : 15 Nov 2016 18:10
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 16 06:48:58 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 08 12:59:29 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	98	-0.02
2		1,4-Dioxane	0.693	0.680	1.9	97	-0.06
3		Pyridine	1.621	1.857	-14.6	106	-0.06
4		n-Nitrosodimethylamine	0.641	0.608	5.1	87	-0.06
5	S	2-Fluorophenol	1.211	1.322	-9.2	100	-0.02
6		Aniline	1.939	1.959	-1.0	89	-0.02
7	S	Phenol-d6	1.644	1.559	5.2	87	-0.02
8		2-Chlorophenol	1.445	1.487	-2.9	97	-0.03
9		Benzaldehyde	0.912	0.884	3.1	91	-0.02
10	C	Phenol	1.819	1.787	1.8	95	-0.02
11		bis(2-Chloroethyl)ether	1.533	1.523	0.7	95	-0.02
12		1,3-Dichlorobenzene	1.461	1.501	-2.7	95	-0.03
13	C	1,4-Dichlorobenzene	1.568	1.573	-0.3	98	-0.03
14		1,2-Dichlorobenzene	1.439	1.541	-7.1	97	-0.02
15		Benzyl Alcohol	1.159	1.208	-4.2	97	-0.02
16		2,2'-oxybis(1-Chloropropane	2.195	1.958	10.8	87	-0.03
17		2-Methylphenol	1.144	1.144	0.0	90	-0.02
18		Hexachloroethane	0.607	0.601	1.0	92	-0.03
19	P	n-Nitroso-di-n-propylamine	1.139	1.128	1.0	98	-0.02
20		3+4-Methylphenols	1.373	1.516	-10.4	97	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	91	-0.02
22		Acetophenone	0.460	0.493	-7.2	97	-0.02
23	S	Nitrobenzene-d5	0.367	0.369	-0.5	94	-0.03
24		Nitrobenzene	0.405	0.430	-6.2	90	-0.02
25		Isophorone	0.706	0.701	0.7	93	-0.02
26	C	2-Nitrophenol	0.171	0.193	-12.9	102	-0.02
27		2,4-Dimethylphenol	0.283	0.286	-1.1	86	-0.02
28		bis(2-Chloroethoxy)methane	0.386	0.421	-9.1	90	-0.02
29	C	2,4-Dichlorophenol	0.247	0.257	-4.0	92	-0.03
30		1,2,4-Trichlorobenzene	0.278	0.313	-12.6	99	-0.02
31		Naphthalene	0.956	1.032	-7.9	95	-0.02
32		Benzoic acid	0.199	0.179	10.1	76	-0.02
33		4-Chloroaniline	0.398	0.445	-11.8	100	-0.02
34	C	Hexachlorobutadiene	0.150	0.165	-10.0	102	-0.03
35		Caprolactam	0.078	0.084	-7.7	98	-0.02
36	C	4-Chloro-3-methylphenol	0.299	0.273	8.7	79	-0.02
37		2-Methylnaphthalene	0.615	0.581	5.5	88	-0.03
38	I	Acenaphthene-d10	1.000	1.000	0.0	94	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.510	0.584	-14.5	95	-0.02
40	P	Hexachlorocyclopentadiene	0.164	0.124	24.4	67	-0.02
41	S	2,4,6-Tribromophenol	0.181	0.217	-19.9	117	-0.02
42	C	2,4,6-Trichlorophenol	0.329	0.346	-5.2	90	-0.02
43		2,4,5-Trichlorophenol	0.346	0.380	-9.8	97	-0.02
44	S	2-Fluorobiphenyl	1.162	1.069	8.0	87	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111516\
Data File : BF091185.D
Acq On : 15 Nov 2016 18:10
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 16 06:48:58 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 08 12:59:29 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	1.460	1.538	-5.3	94	-0.02
46		2-Chloronaphthalene	1.109	1.183	-6.7	94	-0.02
47		2-Nitroaniline	0.411	0.364	11.4	74	-0.02
48		Acenaphthylene	1.728	1.799	-4.1	96	-0.02
49		Dimethylphthalate	1.311	1.341	-2.3	95	-0.02
50		2,6-Dinitrotoluene	0.281	0.296	-5.3	103	-0.03
51	C	Acenaphthene	1.085	1.127	-3.9	101	-0.02
52		3-Nitroaniline	0.333	0.317	4.8	86	-0.02
53	P	2,4-Dinitrophenol	0.132	0.095	28.0#	68	-0.02
54		Dibenzofuran	1.460	1.545	-5.8	101	-0.02
55	P	4-Nitrophenol	0.186	0.155	16.7	71	-0.02
56		2,4-Dinitrotoluene	0.358	0.397	-10.9	92	-0.02
57		Fluorene	1.194	1.305	-9.3	100	-0.02
58		2,3,4,6-Tetrachlorophenol	0.271	0.324	-19.6	98	-0.02
59		Diethylphthalate	1.345	1.426	-6.0	95	-0.02
60		4-Chlorophenyl-phenylether	0.543	0.559	-2.9	87	-0.03
61		4-Nitroaniline	0.337	0.345	-2.4	89	-0.02
62		Azobenzene	1.583	1.192	24.7	71	-0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	94	-0.02
64		4,6-Dinitro-2-methylphenol	0.122	0.113	7.4	88	-0.02
65	c	n-Nitrosodiphenylamine	0.636	0.651	-2.4	102	-0.02
66		4-Bromophenyl-phenylether	0.208	0.251	-20.7	123	-0.02
67		Hexachlorobenzene	0.229	0.264	-15.3	103	-0.02
68		Atrazine	0.183	0.198	-8.2	91	-0.02
69	C	Pentachlorophenol	0.102	0.092	9.8	82	-0.02
70		Phenanthrene	1.063	1.036	2.5	86	-0.03
71		Anthracene	1.130	1.229	-8.8	109	-0.02
72		Carbazole	1.019	1.061	-4.1	105	-0.02
73		Di-n-butylphthalate	1.281	1.211	5.5	83	-0.03
74	C	Fluoranthene	1.103	1.240	-12.4	112	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	93	-0.03
76		Benzidine	0.447	0.483	-8.1	92	-0.02
77		Pyrene	1.310	1.348	-2.9	105	-0.03
78	S	Terphenyl-d14	0.801	0.922	-15.1	102	-0.02
79		Butylbenzylphthalate	0.627	0.714	-13.9	106	-0.03
80		Benzo(a)anthracene	1.136	1.202	-5.8	99	-0.03
81		3,3'-Dichlorobenzidine	0.399	0.466	-16.8	109	-0.02
82		Chrysene	1.120	1.310	-17.0	105	-0.02
83		Bis(2-ethylhexyl)phthalate	0.848	0.969	-14.3	103	-0.02
84	c	Di-n-octyl phthalate	1.394	1.606	-15.2	101	-0.03
85		Indeno(1,2,3-cd)pyrene	0.929	0.948	-2.0	99	-0.06
86	I	Perylene-d12	1.000	1.000	0.0	96	-0.03
87		Benzo(b)fluoranthene	1.356	1.330	1.9	85	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111516\
Data File : BF091185.D
Acq On : 15 Nov 2016 18:10
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 16 06:48:58 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 08 12:59:29 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	1.190	1.433	-20.4	116	-0.03
89 C	Benzo(a)pyrene	1.181	1.255	-6.3	94	-0.03
90	Dibenzo(a,h)anthracene	1.021	1.083	-6.1	96	-0.06
91	Benzo(g,h,i)perylene	0.923	1.001	-8.5	99	-0.06

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