Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA F\Data\BF091117\
Data File : BF098438.D
Acq On : 12 Sep 2017 8:23
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 12 11:37:02 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 12 02:14:50 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	99	0.00
2		1,4-Dioxane	0.698	0.647	7.3	90	0.00
3		Pyridine	1.854	1.733	6.5	89	0.00
4		n-Nitrosodimethylamine	0.909	0.810	10.9	83	0.00
5	S	2-Fluorophenol	1.353	1.318	2.6	95	0.00
6		Aniline	2.155	1.923	10.8	91	0.00
7	S	Phenol-d6	1.717	1.581	7.9	93	0.00
8		2-Chlorophenol	1.487	1.439	3.2	95	0.00
9		Benzaldehyde	1.123	1.062	5.4	94	0.00
10	C	Phenol	1.995	1.843	7.6	94	0.00
11		bis(2-Chloroethyl)ether	1.504	1.430	4.9	94	0.00
12		1,3-Dichlorobenzene	1.497	1.462	2.3	98	0.00
13	C	1,4-Dichlorobenzene	1.534	1.473	4.0	96	0.00
14		1,2-Dichlorobenzene	1.397	1.345	3.7	98	0.00
15		Benzyl Alcohol	1.143	1.041	8.9	94	0.00
16		2,2'-oxybis(1-Chloropropane	2.475	2.258	8.8	92	0.00
17		2-Methylphenol	1.213	1.126	7.2	92	0.00
18		Hexachloroethane	0.587	0.440	25.0#	73	0.00
19	P	n-Nitroso-di-n-propylamine	1.080	0.995	7.9	95	0.00
20		3+4-Methylphenols	1.531	1.437	6.1	95	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	91	0.00
22		Acetophenone	0.517	0.511	1.2	94	0.00
23	S	Nitrobenzene-d5	0.359	0.346	3.6	89	0.00
24		Nitrobenzene	0.409	0.394	3.7	89	0.00
25		Isophorone	0.726	0.691	4.8	89	0.00
26	C	2-Nitrophenol	0.165	0.172	-4.2	90	0.00
27		2,4-Dimethylphenol	0.315	0.305	3.2	91	0.00
28		bis(2-Chloroethoxy)methane	0.443	0.434	2.0	91	0.00
29	C	2,4-Dichlorophenol	0.262	0.270	-3.1	93	0.00
30		1,2,4-Trichlorobenzene	0.285	0.289	-1.4	95	0.00
31		Naphthalene	0.989	0.941	4.9	91	0.00
32		Benzoic acid	0.185	0.191	-3.2	88	0.00
33		4-Chloroaniline	0.402	0.389	3.2	89	0.00
34	C	Hexachlorobutadiene	0.146	0.152	-4.1	97	0.00
35		Caprolactam	0.090	0.086	4.4	90	0.00
36	C	4-Chloro-3-methylphenol	0.324	0.302	6.8	86	0.00
37		2-Methylnaphthalene	0.599	0.574	4.2	90	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	82	0.00
39		1,2,4,5-Tetrachlorobenzene	0.623	0.679	-9.0	92	0.00
40	P	Hexachlorocyclopentadiene	0.140	0.014#	90.0#	8#	0.00
41	S	2,4,6-Tribromophenol	0.133	0.127	4.5	76	0.00
42	C	2,4,6-Trichlorophenol	0.366	0.413	-12.8	89	0.00
43		2,4,5-Trichlorophenol	0.382	0.402	-5.2	84	0.00
44	S	2-Fluorobiphenyl	1.180	1.290	-9.3	90	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA F\Data\BF091117\
Data File : BF098438.D
Acq On : 12 Sep 2017 8:23
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 12 11:37:02 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 12 02:14:50 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	1.790	1.876	-4.8	89	0.00
46		2-Chloronaphthalene	1.278	1.317	-3.1	87	0.00
47		2-Nitroaniline	0.491	0.493	-0.4	81	0.00
48		Acenaphthylene	1.901	1.879	1.2	84	0.00
49		Dimethylphthalate	1.548	1.586	-2.5	87	0.00
50		2,6-Dinitrotoluene	0.318	0.331	-4.1	83	0.00
51	C	Acenaphthene	1.208	1.172	3.0	84	0.00
52		3-Nitroaniline	0.384	0.367	4.4	78	0.00
53	P	2,4-Dinitrophenol	0.121	0.016#	86.8#	10#	0.00
54		Dibenzofuran	1.592	1.518	4.6	83	0.00
55	P	4-Nitrophenol	0.221	0.186	15.8	68	0.00
56		2,4-Dinitrotoluene	0.387	0.362	6.5	78	0.00
57		Fluorene	1.318	1.194	9.4	79	0.00
58		2,3,4,6-Tetrachlorophenol	0.257	0.235	8.6	72	0.00
59		Diethylphthalate	1.472	1.458	1.0	85	0.00
60		4-Chlorophenyl-phenylether	0.609	0.590	3.1	84	0.00
61		4-Nitroaniline	0.388	0.335	13.7	71	0.00
62		Azobenzene	1.587	1.365	14.0	73	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	64	0.00
64		4,6-Dinitro-2-methylphenol	0.114	0.031	72.8#	17#	0.00
65	c	n-Nitrosodiphenylamine	0.649	0.773	-19.1	79	0.00
66		4-Bromophenyl-phenylether	0.190	0.226	-18.9	77	0.00
67		Hexachlorobenzene	0.193	0.213	-10.4	73	0.00
68		Atrazine	0.200	0.212	-6.0	70	0.00
69	C	Pentachlorophenol	0.071	0.063	11.3	53	0.00
70		Phenanthrene	1.060	1.062	-0.2	67	0.00
71		Anthracene	1.089	1.072	1.6	65	0.00
72		Carbazole	1.015	0.953	6.1	63	0.00
73		Di-n-butylphthalate	1.216	1.330	-9.4	71	0.00
74	C	Fluoranthene	1.061	0.937	11.7	59	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	65	0.00
76		Benzidine	0.885	0.759	14.2	56	0.00
77		Pyrene	1.578	1.383	12.4	58	0.00
78	S	Terphenyl-d14	0.927	0.861	7.1	63	0.00
79		Butylbenzylphthalate	0.684	0.645	5.7	59	0.00
80		Benzo(a)anthracene	1.173	1.152	1.8	67	0.00
81		3,3'-Dichlorobenzidine	0.456	0.480	-5.3	68	0.00
82		Chrysene	1.185	1.180	0.4	66	0.00
83		Bis(2-ethylhexyl)phthalate	0.859	0.828	3.6	62	0.00
84	c	Di-n-octyl phthalate	1.474	1.545	-4.8	64	0.00
85		Indeno(1,2,3-cd)pyrene	0.832	1.012	-21.6	84	0.00
86	I	Perylene-d12	1.000	1.000	0.0	82	0.00
87		Benzo(b)fluoranthene	1.258	1.112	11.6	75	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA F\Data\BF091117\
Data File : BF098438.D
Acq On : 12 Sep 2017 8:23
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 12 11:37:02 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 12 02:14:50 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	1.174	1.183	-0.8	80	0.00
89 C	Benzo(a)pyrene	1.120	1.098	2.0	80	0.00
90	Dibenzo(a,h)anthracene	0.815	0.808	0.9	85	0.00
91	Benzo(g,h,i)perylene	0.820	0.772	5.9	83	0.00

(#) = Out of Range SPCC's out = 2 CCC's out = 0