Evaluate Continuing Calibration Report

Data Path : U:\HPCHEM1\BNA F\Data\BF012518\
Data File : BF102431.D
Acq On : 25 Jan 2018 18:44
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 26 00:20:57 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 25 13:44:56 2018
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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	88	0.00
2		1,4-Dioxane	40.000	40.757	-1.9	88	0.00
3		Pyridine	40.000	40.614	-1.5	85	0.00
4		n-Nitrosodimethylamine	40.000	40.885	-2.2	86	0.00
5	S	2-Fluorophenol	80.000	80.252	-0.3	87	0.00
6		Aniline	40.000	39.204	2.0	85	0.00
7	S	Phenol-d6	80.000	78.531	1.8	86	0.00
8		2-Chlorophenol	40.000	39.978	0.1	85	0.00
9		Benzaldehyde	40.000	39.389	1.5	85	0.00
10	C	Phenol	40.000	39.033	2.4	85	0.00
11		bis(2-Chloroethyl)ether	40.000	39.463	1.3	84	0.00
12		1,3-Dichlorobenzene	40.000	38.894	2.8	85	0.00
13	C	1,4-Dichlorobenzene	40.000	39.887	0.3	86	0.00
14		1,2-Dichlorobenzene	40.000	39.894	0.3	86	0.00
15		Benzyl Alcohol	40.000	39.020	2.4	84	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.635	3.4	82	0.00
17		2-Methylphenol	40.000	39.129	2.2	83	0.00
18		Hexachloroethane	40.000	39.996	0.0	86	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.773	3.1	84	0.00
20		3+4-Methylphenols	40.000	40.364	-0.9	85	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	88	0.00
22		Acetophenone	40.000	38.244	4.4	86	0.00
23	S	Nitrobenzene-d5	80.000	77.355	3.3	85	0.00
24		Nitrobenzene	40.000	39.215	2.0	85	0.00
25		Isophorone	40.000	38.001	5.0	83	0.00
26	C	2-Nitrophenol	40.000	39.850	0.4	84	0.00
27		2,4-Dimethylphenol	40.000	38.372	4.1	83	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.364	4.1	83	0.00
29	C	2,4-Dichlorophenol	40.000	38.872	2.8	85	0.00
30		1,2,4-Trichlorobenzene	40.000	38.210	4.5	84	0.00
31		Naphthalene	40.000	38.559	3.6	86	0.00
32		Benzoic acid	40.000	28.967	27.6#	63	-0.02
33		4-Chloroaniline	40.000	38.697	3.3	85	0.00
34	C	Hexachlorobutadiene	40.000	39.128	2.2	85	0.00
35		Caprolactam	40.000	38.053	4.9	82	-0.02
36	C	4-Chloro-3-methylphenol	40.000	38.893	2.8	85	0.00
37		2-Methylnaphthalene	40.000	38.709	3.2	86	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	88	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.803	3.0	85	0.00
40	P	Hexachlorocyclopentadiene	40.000	39.095	2.3	80	0.00
41	S	2,4,6-Tribromophenol	80.000	69.852	12.7	77	0.00
42	C	2,4,6-Trichlorophenol	40.000	37.688	5.8	82	0.00
43		2,4,5-Trichlorophenol	40.000	38.739	3.2	84	0.00
44	S	2-Fluorobiphenyl	80.000	78.586	1.8	86	0.00

Evaluate Continuing Calibration Report

Data Path : U:\HPCHEM1\BNA F\Data\BF012518\
Data File : BF102431.D
Acq On : 25 Jan 2018 18:44
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 26 00:20:57 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 25 13:44:56 2018
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	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	38.665	3.3	86	0.00
46		2-Chloronaphthalene	40.000	38.906	2.7	85	0.00
47		2-Nitroaniline	40.000	39.307	1.7	86	0.00
48		Acenaphthylene	40.000	38.205	4.5	84	0.00
49		Dimethylphthalate	40.000	38.483	3.8	86	0.00
50		2,6-Dinitrotoluene	40.000	38.101	4.7	81	0.00
51	C	Acenaphthene	40.000	38.089	4.8	85	0.00
52		3-Nitroaniline	40.000	38.191	4.5	83	0.00
53	P	2,4-Dinitrophenol	40.000	33.992	15.0	71	0.00
54		Dibenzofuran	40.000	39.754	0.6	85	0.00
55	P	4-Nitrophenol	40.000	34.287	14.3	70	0.00
56		2,4-Dinitrotoluene	40.000	38.318	4.2	82	0.00
57		Fluorene	40.000	40.419	-1.0	89	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	38.867	2.8	86	0.00
59		Diethylphthalate	40.000	37.924	5.2	84	0.00
60		4-Chlorophenyl-phenylether	40.000	40.081	-0.2	86	0.00
61		4-Nitroaniline	40.000	37.605	6.0	81	0.00
62		Azobenzene	40.000	37.908	5.2	83	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	88	0.00
64		4,6-Dinitro-2-methylphenol	40.000	38.308	4.2	79	0.00
65	c	n-Nitrosodiphenylamine	40.000	38.423	3.9	86	0.00
66		4-Bromophenyl-phenylether	40.000	38.600	3.5	84	0.00
67		Hexachlorobenzene	40.000	35.812	10.5	78	0.00
68		Atrazine	40.000	38.259	4.4	85	0.00
69	C	Pentachlorophenol	40.000	40.977	-2.4	82	0.00
70		Phenanthrene	40.000	38.264	4.3	85	0.00
71		Anthracene	40.000	37.739	5.7	84	0.00
72		Carbazole	40.000	38.178	4.6	85	0.00
73		Di-n-butylphthalate	40.000	37.790	5.5	83	0.00
74	C	Fluoranthene	40.000	37.519	6.2	83	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	86	0.00
76		Benzidine	40.000	37.545	6.1	86	0.00
77		Pyrene	40.000	38.948	2.6	83	0.00
78	S	Terphenyl-d14	80.000	75.985	5.0	83	0.00
79		Butylbenzylphthalate	40.000	39.193	2.0	81	0.00
80		Benzo(a)anthracene	40.000	39.457	1.4	86	0.00
81		3,3'-Dichlorobenzidine	40.000	38.681	3.3	82	0.00
82		Chrysene	40.000	38.735	3.2	84	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	38.968	2.6	81	0.00
84	c	Di-n-octyl phthalate	40.000	37.487	6.3	79	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	37.071	7.3	85	0.00
86	I	Perylene-d12	20.000	20.000	0.0	84	0.00
87		Benzo(b)fluoranthene	40.000	38.688	3.3	79	0.00

Evaluate Continuing Calibration Report

Data Path : U:\HPCHEM1\BNA F\Data\BF012518\
Data File : BF102431.D
Acq On : 25 Jan 2018 18:44
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 26 00:20:57 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 25 13:44:56 2018
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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	38.734	3.2	84	0.00
89 C	Benzo(a)pyrene	40.000	39.304	1.7	82	0.00
90	Dibenzo(a,h)anthracene	40.000	38.351	4.1	84	0.00
91	Benzo(g,h,i)perylene	40.000	38.904	2.7	86	0.00

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