Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM080117\
Data File : BM011076.D
Acq On : 01 Aug 2017 11:04
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 01 16:21:16 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM072617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 26 17:29:06 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	67	-0.01
2		1,4-Dioxane	0.531	0.511	3.8	67	0.00
3		Pyridine	1.451	1.587	-9.4	73	0.00
4		n-Nitrosodimethylamine	0.739	0.831	-12.4	76	0.00
5	S	2-Fluorophenol	1.183	1.210	-2.3	70	0.00
6		Aniline	2.119	2.352	-11.0	76	-0.01
7	S	Phenol-d6	1.681	1.871	-11.3	75	-0.01
8		2-Chlorophenol	1.201	1.235	-2.8	70	-0.01
9		Benzaldehyde	1.088	1.221	-12.2	79	0.00
10	C	Phenol	1.826	1.993	-9.1	74	-0.01
11		bis(2-Chloroethyl)ether	1.423	1.507	-5.9	72	0.00
12		1,3-Dichlorobenzene	1.464	1.500	-2.5	71	-0.01
13	C	1,4-Dichlorobenzene	1.501	1.566	-4.3	71	-0.01
14		1,2-Dichlorobenzene	1.435	1.481	-3.2	71	-0.01
15		Benzyl Alcohol	1.613	1.892	-17.3	80	-0.01
16		2,2'-oxybis(1-Chloropropane	1.280	1.471	-14.9	79	-0.02
17		2-Methylphenol	1.167	1.313	-12.5	76	-0.01
18		Hexachloroethane	0.654	0.701	-7.2	73	-0.01
19	P	n-Nitroso-di-n-propylamine	1.428	1.744	-22.1	83	-0.01
20		3+4-Methylphenols	1.622	1.830	-12.8	77	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	74	-0.01
22		Acetophenone	0.600	0.607	-1.2	78	-0.01
23	S	Nitrobenzene-d5	0.533	0.583	-9.4	82	0.00
24		Nitrobenzene	0.552	0.594	-7.6	82	0.00
25		Isophorone	0.888	0.996	-12.2	86	-0.01
26	C	2-Nitrophenol	0.176	0.182	-3.4	77	-0.01
27		2,4-Dimethylphenol	0.309	0.313	-1.3	77	-0.01
28		bis(2-Chloroethoxy)methane	0.496	0.512	-3.2	78	-0.01
29	C	2,4-Dichlorophenol	0.348	0.361	-3.7	78	-0.01
30		1,2,4-Trichlorobenzene	0.430	0.428	0.5	77	-0.01
31		Naphthalene	1.006	1.028	-2.2	79	-0.01
32		Benzoic acid	0.249	0.271	-8.8	83	-0.01
33		4-Chloroaniline	0.401	0.431	-7.5	82	-0.01
34	C	Hexachlorobutadiene	0.338	0.344	-1.8	77	-0.01
35		Caprolactam	0.099	0.120	-21.2	97	0.00
36	C	4-Chloro-3-methylphenol	0.449	0.516	-14.9	88	-0.01
37		2-Methylnaphthalene	0.739	0.797	-7.8	82	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	88	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.859	0.820	4.5	86	-0.01
40	P	Hexachlorocyclopentadiene	0.501	0.461	8.0	81	-0.01
41	S	2,4,6-Tribromophenol	0.321	0.354	-10.3	101	-0.01
42	C	2,4,6-Trichlorophenol	0.520	0.519	0.2	87	-0.01
43		2,4,5-Trichlorophenol	0.536	0.533	0.6	88	-0.01
44	S	2-Fluorobiphenyl	1.595	1.540	3.4	86	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM080117\
Data File : BM011076.D
Acq On : 01 Aug 2017 11:04
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 01 16:21:16 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM072617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 26 17:29:06 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	1.729	1.657	4.2	87	-0.01
46		2-Chloronaphthalene	1.274	1.226	3.8	87	-0.01
47		2-Nitroaniline	0.451	0.518	-14.9	101	0.00
48		Acenaphthylene	1.902	1.919	-0.9	91	-0.01
49		Dimethylphthalate	1.711	1.733	-1.3	94	-0.01
50		2,6-Dinitrotoluene	0.346	0.371	-7.2	95	-0.01
51	C	Acenaphthene	1.177	1.191	-1.2	92	-0.01
52		3-Nitroaniline	0.300	0.323	-7.7	97	-0.01
53	P	2,4-Dinitrophenol	0.246	0.247	-0.4	92	-0.01
54		Dibenzofuran	1.908	1.948	-2.1	93	-0.01
55	P	4-Nitrophenol	0.264	0.289	-9.5	99	0.00
56		2,4-Dinitrotoluene	0.486	0.532	-9.5	98	-0.01
57		Fluorene	1.661	1.771	-6.6	98	-0.01
58		2,3,4,6-Tetrachlorophenol	0.502	0.532	-6.0	97	0.00
59		Diethylphthalate	1.747	1.854	-6.1	97	0.00
60		4-Chlorophenyl-phenylether	1.003	1.049	-4.6	97	-0.01
61		4-Nitroaniline	0.319	0.363	-13.8	105	0.00
62		Azobenzene	1.870	2.116	-13.2	103	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	100	-0.01
64		4,6-Dinitro-2-methylphenol	0.135	0.136	-0.7	99	-0.01
65	c	n-Nitrosodiphenylamine	0.572	0.562	1.7	100	0.00
66		4-Bromophenyl-phenylether	0.257	0.254	1.2	101	0.00
67		Hexachlorobenzene	0.291	0.287	1.4	99	-0.01
68		Atrazine	0.229	0.232	-1.3	100	-0.01
69	C	Pentachlorophenol	0.184	0.194	-5.4	103	0.00
70		Phenanthrene	1.047	1.087	-3.8	105	-0.01
71		Anthracene	1.044	1.090	-4.4	105	0.00
72		Carbazole	0.913	0.980	-7.3	110	-0.01
73		Di-n-butylphthalate	1.112	1.196	-7.6	107	0.00
74	C	Fluoranthene	1.298	1.451	-11.8	114	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	113	0.00
76		Benzidine	0.478	0.450	5.9	104	0.00
77		Pyrene	1.188	1.171	1.4	113	0.00
78	S	Terphenyl-d14	1.010	1.036	-2.6	115	0.00
79		Butylbenzylphthalate	0.423	0.461	-9.0	119	0.00
80		Benzo(a)anthracene	1.142	1.179	-3.2	119	0.00
81		3,3'-Dichlorobenzidine	0.448	0.462	-3.1	115	0.00
82		Chrysene	1.096	1.123	-2.5	119	0.00
83		Bis(2-ethylhexyl)phthalate	0.622	0.647	-4.0	116	0.00
84	c	Di-n-octyl phthalate	1.009	1.094	-8.4	120	0.00
85		Indeno(1,2,3-cd)pyrene	1.135	1.020	10.1	106	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	114	0.00
87		Benzo(b)fluoranthene	1.284	1.348	-5.0	120	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM080117\
Data File : BM011076.D
Acq On : 01 Aug 2017 11:04
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 01 16:21:16 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM072617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 26 17:29:06 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	1.230	1.243	-1.1	118	-0.01
89 C	Benzo(a)pyrene	1.162	1.177	-1.3	117	-0.01
90	Dibenzo(a,h)anthracene	1.077	0.957	11.1	105	-0.01
91	Benzo(g,h,i)perylene	1.057	0.919	13.1	102	-0.01

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