Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070518\
Data File : BM015784.D
Acq On : 06 Jul 2018 03:58
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jul 06 07:24:54 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 05 17:39:50 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	125	0.00
2		1,4-Dioxane	0.482	0.437	9.3	116	0.00
3		Pyridine	1.273	1.164	8.6	111	0.00
4		n-Nitrosodimethylamine	0.991	0.966	2.5	123	0.00
5	S	2-Fluorophenol	1.166	1.135	2.7	120	0.00
6		Aniline	1.936	1.898	2.0	121	0.00
7	S	Phenol-d6	1.597	1.600	-0.2	122	0.00
8		2-Chlorophenol	1.400	1.386	1.0	122	0.00
9		Benzaldehyde	1.006	1.058	-5.2	131	0.00
10	C	Phenol	1.604	1.581	1.4	121	0.00
11		bis(2-Chloroethyl)ether	1.305	1.273	2.5	121	0.00
12		1,3-Dichlorobenzene	1.552	1.515	2.4	123	0.00
13	C	1,4-Dichlorobenzene	1.605	1.563	2.6	125	0.00
14		1,2-Dichlorobenzene	1.552	1.528	1.5	124	0.00
15		Benzyl Alcohol	1.386	1.398	-0.9	124	0.00
16		2,2'-oxybis(1-Chloropropane	1.425	1.366	4.1	118	0.01
17		2-Methylphenol	1.111	1.106	0.5	123	0.00
18		Hexachloroethane	0.635	0.635	0.0	123	0.00
19	P	n-Nitroso-di-n-propylamine	1.283	1.289	-0.5	122	0.00
20		3+4-Methylphenols	1.544	1.571	-1.7	122	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	122	0.00
22		Acetophenone	0.521	0.517	0.8	123	0.00
23	S	Nitrobenzene-d5	0.410	0.410	0.0	123	0.00
24		Nitrobenzene	0.399	0.393	1.5	122	0.00
25		Isophorone	0.626	0.628	-0.3	122	0.00
26	C	2-Nitrophenol	0.172	0.170	1.2	123	0.00
27		2,4-Dimethylphenol	0.261	0.257	1.5	123	0.00
28		bis(2-Chloroethoxy)methane	0.408	0.402	1.5	121	0.00
29	C	2,4-Dichlorophenol	0.289	0.299	-3.5	125	0.00
30		1,2,4-Trichlorobenzene	0.336	0.338	-0.6	127	0.00
31		Naphthalene	1.037	1.017	1.9	123	0.00
32		Benzoic acid	0.231	0.209	9.5	110	0.01
33		4-Chloroaniline	0.423	0.428	-1.2	123	0.00
34	C	Hexachlorobutadiene	0.224	0.227	-1.3	129	0.00
35		Caprolactam	0.096	0.095	1.0	121	0.01
36	C	4-Chloro-3-methylphenol	0.332	0.341	-2.7	124	0.00
37		2-Methylnaphthalene	0.738	0.738	0.0	125	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	123	0.00
39		1,2,4,5-Tetrachlorobenzene	0.642	0.650	-1.2	127	0.00
40	P	Hexachlorocyclopentadiene	0.236	0.260	-10.2	136	0.00
41	S	2,4,6-Tribromophenol	0.261	0.276	-5.7	127	0.00
42	C	2,4,6-Trichlorophenol	0.406	0.424	-4.4	126	0.00
43		2,4,5-Trichlorophenol	0.438	0.445	-1.6	124	0.00
44	S	2-Fluorobiphenyl	1.611	1.617	-0.4	127	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070518\
Data File : BM015784.D
Acq On : 06 Jul 2018 03:58
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jul 06 07:24:54 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 05 17:39:50 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.729	1.714	0.9	125	0.00
46		2-Chloronaphthalene	1.293	1.293	0.0	125	0.00
47		2-Nitroaniline	0.451	0.452	-0.2	123	0.00
48		Acenaphthylene	2.103	2.119	-0.8	125	0.00
49		Dimethylphthalate	1.764	1.742	1.2	124	0.00
50		2,6-Dinitrotoluene	0.344	0.356	-3.5	126	0.00
51	C	Acenaphthene	1.309	1.304	0.4	125	0.00
52		3-Nitroaniline	0.376	0.373	0.8	123	0.00
53	P	2,4-Dinitrophenol	0.145	0.139	4.1	118	0.00
54		Dibenzofuran	2.017	1.997	1.0	125	0.00
55	P	4-Nitrophenol	0.278	0.271	2.5	119	0.00
56		2,4-Dinitrotoluene	0.497	0.499	-0.4	127	0.00
57		Fluorene	1.663	1.685	-1.3	128	0.00
58		2,3,4,6-Tetrachlorophenol	0.370	0.379	-2.4	126	0.00
59		Diethylphthalate	1.890	1.902	-0.6	126	0.00
60		4-Chlorophenyl-phenylether	0.856	0.865	-1.1	128	0.00
61		4-Nitroaniline	0.391	0.365	6.6	112	0.00
62		Azobenzene	1.772	1.819	-2.7	125	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	122	0.00
64		4,6-Dinitro-2-methylphenol	0.122	0.119	2.5	123	0.00
65	c	n-Nitrosodiphenylamine	0.685	0.687	-0.3	126	0.00
66		4-Bromophenyl-phenylether	0.223	0.230	-3.1	128	0.00
67		Hexachlorobenzene	0.250	0.253	-1.2	127	0.00
68		Atrazine	0.225	0.228	-1.3	123	0.00
69	C	Pentachlorophenol	0.154	0.152	1.3	123	0.00
70		Phenanthrene	1.144	1.131	1.1	125	0.00
71		Anthracene	1.124	1.136	-1.1	125	0.00
72		Carbazole	1.047	1.047	0.0	125	0.00
73		Di-n-butylphthalate	1.411	1.446	-2.5	126	0.00
74	C	Fluoranthene	1.315	1.283	2.4	123	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	106	0.00
76		Benzidine	0.586	0.609	-3.9	138	0.00
77		Pyrene	1.246	1.380	-10.8	120	0.00
78	S	Terphenyl-d14	1.078	1.207	-12.0	124	0.00
79		Butylbenzylphthalate	0.602	0.661	-9.8	117	0.00
80		Benzo(a)anthracene	1.268	1.263	0.4	109	0.00
81		3,3'-Dichlorobenzidine	0.456	0.444	2.6	104	0.00
82		Chrysene	1.181	1.146	3.0	108	0.00
83		Bis(2-ethylhexyl)phthalate	0.890	0.908	-2.0	109	0.00
84	c	Di-n-octyl phthalate	1.451	1.351	6.9	98	0.00
85		Indeno(1,2,3-cd)pyrene	1.097	0.959	12.6	91	0.00
86	I	Perylene-d12	1.000	1.000	0.0	89	0.00
87		Benzo(b)fluoranthene	1.335	1.343	-0.6	93	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070518\
Data File : BM015784.D
Acq On : 06 Jul 2018 03:58
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jul 06 07:24:54 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 05 17:39:50 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.297	1.309	-0.9	92	0.00
89 C	Benzo(a)pyrene	1.175	1.172	0.3	91	0.00
90	Dibenzo(a,h)anthracene	1.058	1.079	-2.0	90	0.00
91	Benzo(g,h,i)perylene	0.929	0.975	-5.0	94	0.00

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