Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM080618\
Data File : BM016193.D
Acq On : 06 Aug 2018 16:17
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 07 13:13:12 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 07 13:02:43 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	79	0.04
2		1,4-Dioxane	0.564	0.589	-4.4	78	0.03
3		Pyridine	1.359	1.352	0.5	74	0.03
4		n-Nitrosodimethylamine	1.060	1.159	-9.3	80	0.03
5	S	2-Fluorophenol	1.204	1.234	-2.5	76	0.03
6		Aniline	1.810	1.787	1.3	74	0.03
7	S	Phenol-d6	1.572	1.564	0.5	74	0.03
8		2-Chlorophenol	1.431	1.427	0.3	75	0.04
9		Benzaldehyde	1.105	1.068	3.3	72	0.03
10	C	Phenol	1.572	1.557	1.0	74	0.02
11		bis(2-Chloroethyl)ether	1.242	1.175	5.4	73	0.03
12		1,3-Dichlorobenzene	1.675	1.616	3.5	75	0.04
13	C	1,4-Dichlorobenzene	1.699	1.620	4.6	73	0.04
14		1,2-Dichlorobenzene	1.662	1.580	4.9	75	0.04
15		Benzyl Alcohol	1.252	1.193	4.7	72	0.03
16		2,2'-oxybis(1-Chloropropane	1.901	1.696	10.8	69	0.03
17		2-Methylphenol	1.075	1.080	-0.5	74	0.03
18		Hexachloroethane	0.734	0.738	-0.5	75	0.03
19	P	n-Nitroso-di-n-propylamine	1.149	1.124	2.2	71	0.02
20		3+4-Methylphenols	1.546	1.467	5.1	71	0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	74	0.03
22		Acetophenone	0.504	0.508	-0.8	73	0.03
23	S	Nitrobenzene-d5	0.430	0.464	-7.9	77	0.04
24		Nitrobenzene	0.433	0.444	-2.5	72	0.03
25		Isophorone	0.588	0.609	-3.6	73	0.02
26	C	2-Nitrophenol	0.181	0.184	-1.7	74	0.04
27		2,4-Dimethylphenol	0.252	0.249	1.2	70	0.04
28		bis(2-Chloroethoxy)methane	0.382	0.380	0.5	70	0.04
29	C	2,4-Dichlorophenol	0.299	0.306	-2.3	71	0.03
30		1,2,4-Trichlorobenzene	0.363	0.350	3.6	70	0.04
31		Naphthalene	1.012	0.979	3.3	71	0.03
32		Benzoic acid	0.192	0.194	-1.0	75	0.04
33		4-Chloroaniline	0.413	0.417	-1.0	70	0.03
34	C	Hexachlorobutadiene	0.252	0.253	-0.4	73	0.04
35		Caprolactam	0.083	0.085	-2.4	73	0.00
36	C	4-Chloro-3-methylphenol	0.329	0.346	-5.2	73	0.02
37		2-Methylnaphthalene	0.678	0.664	2.1	70	0.03
38	I	Acenaphthene-d10	1.000	1.000	0.0	73	0.02
39		1,2,4,5-Tetrachlorobenzene	0.681	0.673	1.2	72	0.03
40	P	Hexachlorocyclopentadiene	0.278	0.285	-2.5	74	0.03
41	S	2,4,6-Tribromophenol	0.254	0.236	7.1	68	0.02
42	C	2,4,6-Trichlorophenol	0.426	0.436	-2.3	71	0.02
43		2,4,5-Trichlorophenol	0.439	0.450	-2.5	72	0.03
44	S	2-Fluorobiphenyl	1.644	1.703	-3.6	76	0.03

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.805	1.765	2.2	71	0.03
46		2-Chloronaphthalene	1.404	1.379	1.8	71	0.02
47		2-Nitroaniline	0.464	0.485	-4.5	74	0.03
48		Acenaphthylene	2.055	2.060	-0.2	71	0.02
49		Dimethylphthalate	1.750	1.753	-0.2	72	0.02
50		2,6-Dinitrotoluene	0.352	0.354	-0.6	70	0.02
51	C	Acenaphthene	1.264	1.239	2.0	72	0.02
52		3-Nitroaniline	0.380	0.370	2.6	69	0.02
53	P	2,4-Dinitrophenol	0.131	0.153	-16.8	84	0.03
54		Dibenzofuran	2.083	2.032	2.4	71	0.02
55	P	4-Nitrophenol	0.273	0.264	3.3	69	0.02
56		2,4-Dinitrotoluene	0.504	0.505	-0.2	72	0.02
57		Fluorene	1.548	1.544	0.3	72	0.02
58		2,3,4,6-Tetrachlorophenol	0.378	0.385	-1.9	71	0.02
59		Diethylphthalate	1.887	1.946	-3.1	74	0.02
60		4-Chlorophenyl-phenylether	0.862	0.848	1.6	72	0.03
61		4-Nitroaniline	0.365	0.353	3.3	70	0.02
62		Azobenzene	1.987	2.183	-9.9	77	0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	74	0.03
64		4,6-Dinitro-2-methylphenol	0.121	0.128	-5.8	77	0.02
65	c	n-Nitrosodiphenylamine	0.659	0.654	0.8	72	0.02
66		4-Bromophenyl-phenylether	0.226	0.225	0.4	72	0.02
67		Hexachlorobenzene	0.249	0.239	4.0	69	0.02
68		Atrazine	0.236	0.236	0.0	71	0.02
69	C	Pentachlorophenol	0.154	0.145	5.8	69	0.03
70		Phenanthrene	1.120	1.085	3.1	72	0.02
71		Anthracene	1.088	1.086	0.2	73	0.02
72		Carbazole	1.127	1.122	0.4	71	0.02
73		Di-n-butylphthalate	1.405	1.540	-9.6	79	0.02
74	C	Fluoranthene	1.249	1.265	-1.3	74	0.02
75	I	Chrysene-d12	1.000	1.000	0.0	80	0.02
76		Benzidine	0.559	0.565	-1.1	77	0.02
77		Pyrene	1.199	1.136	5.3	74	0.02
78	S	Terphenyl-d14	0.955	0.981	-2.7	81	0.02
79		Butylbenzylphthalate	0.609	0.625	-2.6	79	0.02
80		Benzo(a)anthracene	1.232	1.182	4.1	77	0.02
81		3,3'-Dichlorobenzidine	0.496	0.480	3.2	75	0.02
82		Chrysene	1.182	1.121	5.2	76	0.02
83		Bis(2-ethylhexyl)phthalate	0.882	0.933	-5.8	82	0.02
84	c	Di-n-octyl phthalate	1.483	1.558	-5.1	81	0.02
85		Indeno(1,2,3-cd)pyrene	1.402	1.339	4.5	77	0.05
86	I	Perylene-d12	1.000	1.000	0.0	80	0.04
87		Benzo(b)fluoranthene	1.178	1.153	2.1	78	0.03

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88	Benzo(k)fluoranthene	1.193	1.131	5.2	74	0.03
89 C	Benzo(a)pyrene	1.132	1.103	2.6	77	0.04
90	Dibenzo(a,h)anthracene	1.173	1.148	2.1	77	0.04
91	Benzo(g,h,i)perylene	1.109	1.054	5.0	76	0.04

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