Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM070918\
Data File : BM015839.D
Acq On : 09 Jul 2018 12:39
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jul 09 15:04:53 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	73	-0.01
2		1,4-Dioxane	0.482	0.456	5.4	71	-0.01
3		Pyridine	1.273	1.165	8.5	65	-0.01
4		n-Nitrosodimethylamine	0.991	0.986	0.5	74	0.00
5	S	2-Fluorophenol	1.166	1.100	5.7	68	-0.01
6		Aniline	1.936	1.970	-1.8	74	-0.01
7	S	Phenol-d6	1.597	1.547	3.1	69	-0.01
8		2-Chlorophenol	1.400	1.380	1.4	71	-0.02
9		Benzaldehyde	1.006	0.976	3.0	71	-0.02
10	C	Phenol	1.604	1.549	3.4	70	-0.01
11		bis(2-Chloroethyl)ether	1.305	1.282	1.8	72	-0.01
12		1,3-Dichlorobenzene	1.552	1.474	5.0	71	-0.01
13	C	1,4-Dichlorobenzene	1.605	1.528	4.8	72	-0.01
14		1,2-Dichlorobenzene	1.552	1.509	2.8	72	-0.01
15		Benzyl Alcohol	1.386	1.408	-1.6	74	-0.01
16		2,2'-oxybis(1-Chloropropane	1.425	1.455	-2.1	74	-0.01
17		2-Methylphenol	1.111	1.136	-2.3	75	-0.02
18		Hexachloroethane	0.635	0.620	2.4	71	-0.01
19	P	n-Nitroso-di-n-propylamine	1.283	1.349	-5.1	75	-0.02
20		3+4-Methylphenols	1.544	1.611	-4.3	74	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	75	-0.02
22		Acetophenone	0.521	0.509	2.3	74	-0.01
23	S	Nitrobenzene-d5	0.410	0.406	1.0	75	-0.01
24		Nitrobenzene	0.399	0.393	1.5	75	-0.01
25		Isophorone	0.626	0.639	-2.1	76	-0.01
26	C	2-Nitrophenol	0.172	0.164	4.7	73	-0.01
27		2,4-Dimethylphenol	0.261	0.268	-2.7	79	-0.02
28		bis(2-Chloroethoxy)methane	0.408	0.402	1.5	75	-0.01
29	C	2,4-Dichlorophenol	0.289	0.290	-0.3	75	-0.01
30		1,2,4-Trichlorobenzene	0.336	0.314	6.5	73	-0.02
31		Naphthalene	1.037	0.998	3.8	74	-0.01
32		Benzoic acid	0.231	0.190	17.7	62	-0.02
33		4-Chloroaniline	0.423	0.431	-1.9	77	-0.02
34	C	Hexachlorobutadiene	0.224	0.208	7.1	73	-0.01
35		Caprolactam	0.096	0.101	-5.2	80	-0.02
36	C	4-Chloro-3-methylphenol	0.332	0.343	-3.3	77	-0.01
37		2-Methylnaphthalene	0.738	0.735	0.4	77	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	79	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.642	0.594	7.5	74	-0.01
40	P	Hexachlorocyclopentadiene	0.236	0.271	-14.8	90	-0.01
41	S	2,4,6-Tribromophenol	0.261	0.264	-1.1	78	-0.01
42	C	2,4,6-Trichlorophenol	0.406	0.391	3.7	74	-0.01
43		2,4,5-Trichlorophenol	0.438	0.408	6.8	72	-0.01
44	S	2-Fluorobiphenyl	1.611	1.487	7.7	74	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.729	1.613	6.7	75	-0.01
46		2-Chloronaphthalene	1.293	1.209	6.5	75	-0.01
47		2-Nitroaniline	0.451	0.444	1.6	77	-0.01
48		Acenaphthylene	2.103	2.087	0.8	79	-0.01
49		Dimethylphthalate	1.764	1.722	2.4	79	-0.01
50		2,6-Dinitrotoluene	0.344	0.346	-0.6	78	-0.01
51	C	Acenaphthene	1.309	1.286	1.8	79	-0.01
52		3-Nitroaniline	0.376	0.381	-1.3	80	-0.01
53	P	2,4-Dinitrophenol	0.145	0.127	12.4	69	-0.01
54		Dibenzofuran	2.017	1.989	1.4	79	-0.01
55	P	4-Nitrophenol	0.278	0.270	2.9	76	-0.01
56		2,4-Dinitrotoluene	0.497	0.504	-1.4	82	-0.01
57		Fluorene	1.663	1.646	1.0	80	-0.01
58		2,3,4,6-Tetrachlorophenol	0.370	0.363	1.9	77	-0.01
59		Diethylphthalate	1.890	1.910	-1.1	81	-0.01
60		4-Chlorophenyl-phenylether	0.856	0.825	3.6	78	-0.01
61		4-Nitroaniline	0.391	0.374	4.3	74	-0.01
62		Azobenzene	1.772	1.873	-5.7	82	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	81	-0.01
64		4,6-Dinitro-2-methylphenol	0.122	0.113	7.4	78	0.00
65	c	n-Nitrosodiphenylamine	0.685	0.655	4.4	80	-0.01
66		4-Bromophenyl-phenylether	0.223	0.213	4.5	79	-0.01
67		Hexachlorobenzene	0.250	0.239	4.4	80	-0.01
68		Atrazine	0.225	0.216	4.0	78	-0.01
69	C	Pentachlorophenol	0.154	0.143	7.1	77	-0.01
70		Phenanthrene	1.144	1.118	2.3	82	-0.01
71		Anthracene	1.124	1.102	2.0	81	-0.01
72		Carbazole	1.047	1.053	-0.6	83	0.00
73		Di-n-butylphthalate	1.411	1.421	-0.7	82	-0.01
74	C	Fluoranthene	1.315	1.316	-0.1	84	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	85	-0.01
76		Benzidine	0.586	0.440	24.9	80	-0.01
77		Pyrene	1.246	1.193	4.3	84	-0.01
78	S	Terphenyl-d14	1.078	1.013	6.0	83	-0.01
79		Butylbenzylphthalate	0.602	0.604	-0.3	86	-0.01
80		Benzo(a)anthracene	1.268	1.229	3.1	85	-0.01
81		3,3'-Dichlorobenzidine	0.456	0.442	3.1	83	-0.01
82		Chrysene	1.181	1.126	4.7	85	0.00
83		Bis(2-ethylhexyl)phthalate	0.890	0.882	0.9	85	-0.01
84	c	Di-n-octyl phthalate	1.451	1.515	-4.4	88	-0.02
85		Indeno(1,2,3-cd)pyrene	1.097	1.253	-14.2	95	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	88	-0.01
87		Benzo(b)fluoranthene	1.335	1.284	3.8	89	-0.01

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88	Benzo(k)fluoranthene	1.297	1.228	5.3	86	-0.01
89 C	Benzo(a)pyrene	1.175	1.150	2.1	89	-0.02
90	Dibenzo(a,h)anthracene	1.058	1.144	-8.1	95	-0.02
91	Benzo(g,h,i)perylene	0.929	1.014	-9.1	97	-0.02

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