Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120518\
Data File : BM017902.D
Acq On : 04 Dec 2018 14:20
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Dec 05 10:28:37 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM111218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 12 18:37:25 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	101	-0.04
2		1,4-Dioxane	0.494	0.482	2.4	101	-0.02
3		Pyridine	1.448	1.360	6.1	94	-0.02
4		n-Nitrosodimethylamine	0.638	0.510	20.1	79	-0.02
5	S	2-Fluorophenol	1.185	1.263	-6.6	107	-0.03
6		Aniline	2.053	2.078	-1.2	100	-0.04
7	S	Phenol-d6	1.655	1.702	-2.8	101	-0.04
8		2-Chlorophenol	1.418	1.498	-5.6	106	-0.04
9		Benzaldehyde	1.206	0.964	20.1	79	-0.04
10	C	Phenol	1.737	1.773	-2.1	101	-0.04
11		bis(2-Chloroethyl)ether	1.371	1.379	-0.6	100	-0.04
12		1,3-Dichlorobenzene	1.591	1.674	-5.2	107	-0.04
13	C	1,4-Dichlorobenzene	1.631	1.711	-4.9	106	-0.04
14		1,2-Dichlorobenzene	1.581	1.641	-3.8	104	-0.04
15		Benzyl Alcohol	1.318	1.295	1.7	95	-0.04
16		2,2'-oxybis(1-Chloropropane	2.088	1.554	25.6#	74	-0.03
17		2-Methylphenol	1.170	1.187	-1.5	100	-0.04
18		Hexachloroethane	0.575	0.609	-5.9	104	-0.04
19	P	n-Nitroso-di-n-propylamine	1.188	1.136	4.4	93	-0.04
20		3+4-Methylphenols	1.626	1.639	-0.8	98	-0.04
21	I	Naphthalene-d8	1.000	1.000	0.0	93	-0.04
22		Acetophenone	0.499	0.528	-5.8	96	-0.04
23	S	Nitrobenzene-d5	0.387	0.391	-1.0	91	-0.04
24		Nitrobenzene	0.393	0.384	2.3	90	-0.04
25		Isophorone	0.598	0.622	-4.0	93	-0.04
26	C	2-Nitrophenol	0.181	0.203	-12.2	100	-0.04
27		2,4-Dimethylphenol	0.261	0.283	-8.4	97	-0.04
28		bis(2-Chloroethoxy)methane	0.405	0.417	-3.0	92	-0.04
29	C	2,4-Dichlorophenol	0.303	0.329	-8.6	96	-0.04
30		1,2,4-Trichlorobenzene	0.352	0.367	-4.3	95	-0.04
31		Naphthalene	0.983	1.023	-4.1	95	-0.04
32		Benzoic acid	0.236	0.241	-2.1	92	-0.02
33		4-Chloroaniline	0.431	0.455	-5.6	94	-0.04
34	C	Hexachlorobutadiene	0.219	0.232	-5.9	98	-0.05
35		Caprolactam	0.112	0.101	9.8	83	-0.03
36	C	4-Chloro-3-methylphenol	0.350	0.357	-2.0	91	-0.04
37		2-Methylnaphthalene	0.695	0.714	-2.7	93	-0.04
38		1-Methylnaphthalene	0.681	0.693	-1.8	93	-0.04
39	I	Acenaphthene-d10	1.000	1.000	0.0	86	-0.04
40		1,2,4,5-Tetrachlorobenzene	0.705	0.757	-7.4	91	-0.04
41	P	Hexachlorocyclopentadiene	0.364	0.423	-16.2	98	-0.04
42	S	2,4,6-Tribromophenol	0.287	0.314	-9.4	91	-0.03
43	C	2,4,6-Trichlorophenol	0.442	0.488	-10.4	91	-0.04
44		2,4,5-Trichlorophenol	0.466	0.512	-9.9	92	-0.04

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.542	1.634	-6.0	90	-0.04
46		1,1'-Biphenyl	1.795	1.913	-6.6	91	-0.04
47		2-Chloronaphthalene	1.332	1.409	-5.8	90	-0.04
48		2-Nitroaniline	0.412	0.410	0.5	81	-0.04
49		Acenaphthylene	2.002	2.125	-6.1	90	-0.04
50		Dimethylphthalate	1.626	1.699	-4.5	89	-0.03
51		2,6-Dinitrotoluene	0.362	0.389	-7.5	90	-0.04
52	C	Acenaphthene	1.229	1.291	-5.0	90	-0.04
53		3-Nitroaniline	0.394	0.420	-6.6	91	-0.03
54	P	2,4-Dinitrophenol	0.186	0.231	-24.2	102	-0.03
55		Dibenzofuran	2.008	2.081	-3.6	89	-0.03
56	P	4-Nitrophenol	0.294	0.328	-11.6	94	-0.03
57		2,4-Dinitrotoluene	0.488	0.541	-10.9	90	-0.03
58		Fluorene	1.537	1.587	-3.3	88	-0.03
59		2,3,4,6-Tetrachlorophenol	0.429	0.464	-8.2	90	-0.04
60		Diethylphthalate	1.757	1.804	-2.7	90	-0.03
61		4-Chlorophenyl-phenylether	0.873	0.898	-2.9	88	-0.03
62		4-Nitroaniline	0.372	0.407	-9.4	92	-0.02
63		Azobenzene	1.624	1.615	0.6	84	-0.03
64	I	Phenanthrene-d10	1.000	1.000	0.0	84	-0.03
65		4,6-Dinitro-2-methylphenol	0.139	0.160	-15.1	94	-0.03
66	c	n-Nitrosodiphenylamine	0.643	0.697	-8.4	89	-0.04
67		4-Bromophenyl-phenylether	0.237	0.255	-7.6	88	-0.04
68		Hexachlorobenzene	0.267	0.280	-4.9	88	-0.03
69		Atrazine	0.237	0.247	-4.2	84	-0.03
70	C	Pentachlorophenol	0.187	0.209	-11.8	92	-0.03
71		Phenanthrene	1.085	1.140	-5.1	88	-0.03
72		Anthracene	1.056	1.129	-6.9	87	-0.04
73		Carbazole	1.067	1.162	-8.9	88	-0.03
74		Di-n-butylphthalate	1.188	1.384	-16.5	93	-0.03
75	C	Fluoranthene	1.228	1.299	-5.8	87	-0.03
76	I	Chrysene-d12	1.000	1.000	0.0	83	-0.02
77		Benzidine	0.651	0.604	7.2	73	-0.03
78		Pyrene	1.234	1.292	-4.7	86	-0.03
79	S	Terphenyl-d14	0.882	0.929	-5.3	87	-0.02
80		Butylbenzylphthalate	0.501	0.583	-16.4	91	-0.03
81		Benzo(a)anthracene	1.152	1.197	-3.9	85	-0.02
82		3,3'-Dichlorobenzidine	0.445	0.488	-9.7	85	-0.03
83		Chrysene	1.119	1.149	-2.7	85	-0.03
84		Bis(2-ethylhexyl)phthalate	0.741	0.849	-14.6	93	-0.02
85	c	Di-n-octyl phthalate	1.186	1.344	-13.3	91	-0.03
86		Indeno(1,2,3-cd)pyrene	0.894	1.042	-16.6	93	-0.05
87	I	Perylene-d12	1.000	1.000	0.0	86	-0.04

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88	Benzo(b)fluoranthene	1.171	1.183	-1.0	86	-0.03
89	Benzo(k)fluoranthene	1.184	1.190	-0.5	84	-0.03
90 C	Benzo(a)pyrene	1.068	1.117	-4.6	88	-0.04
91	Dibenzo(a,h)anthracene	0.898	1.007	-12.1	95	-0.05
92	Benzo(g,h,i)perylene	0.844	0.951	-12.7	96	-0.06

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