Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120518\
Data File : BM017919.D
Acq On : 05 Dec 2018 01:46
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Dec 05 04:58:39 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	115	-0.04
2		1,4-Dioxane	0.494	0.425	14.0	102	-0.03
3		Pyridine	1.448	1.285	11.3	101	-0.02
4		n-Nitrosodimethylamine	0.638	0.462	27.6#	82	-0.02
5	S	2-Fluorophenol	1.185	1.221	-3.0	117	-0.03
6		Aniline	2.053	2.122	-3.4	116	-0.03
7	S	Phenol-d6	1.655	1.728	-4.4	117	-0.02
8		2-Chlorophenol	1.418	1.544	-8.9	124	-0.03
9		Benzaldehyde	1.206	0.982	18.6	92	-0.04
10	C	Phenol	1.737	1.783	-2.6	116	-0.02
11		bis(2-Chloroethyl)ether	1.371	1.397	-1.9	115	-0.04
12		1,3-Dichlorobenzene	1.591	1.727	-8.5	126	-0.03
13	C	1,4-Dichlorobenzene	1.631	1.769	-8.5	124	-0.03
14		1,2-Dichlorobenzene	1.581	1.695	-7.2	123	-0.04
15		Benzyl Alcohol	1.318	1.278	3.0	107	-0.03
16		2,2'-oxybis(1-Chloropropane	2.088	1.555	25.5#	84	-0.02
17		2-Methylphenol	1.170	1.167	0.3	112	-0.03
18		Hexachloroethane	0.575	0.398	30.8#	77	-0.04
19	P	n-Nitroso-di-n-propylamine	1.188	1.122	5.6	104	-0.03
20		3+4-Methylphenols	1.626	1.674	-3.0	114	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	105	-0.04
22		Acetophenone	0.499	0.534	-7.0	109	-0.03
23	S	Nitrobenzene-d5	0.387	0.400	-3.4	105	-0.03
24		Nitrobenzene	0.393	0.397	-1.0	105	-0.03
25		Isophorone	0.598	0.627	-4.8	106	-0.04
26	C	2-Nitrophenol	0.181	0.160	11.6	88	-0.03
27		2,4-Dimethylphenol	0.261	0.295	-13.0	115	-0.03
28		bis(2-Chloroethoxy)methane	0.405	0.425	-4.9	106	-0.03
29	C	2,4-Dichlorophenol	0.303	0.344	-13.5	113	-0.03
30		1,2,4-Trichlorobenzene	0.352	0.376	-6.8	110	-0.03
31		Naphthalene	0.983	1.056	-7.4	111	-0.04
32		Benzoic acid	0.236	0.234	0.8	101	0.00
33		4-Chloroaniline	0.431	0.471	-9.3	109	-0.04
34	C	Hexachlorobutadiene	0.219	0.243	-11.0	116	-0.04
35		Caprolactam	0.112	0.118	-5.4	109	-0.01
36	C	4-Chloro-3-methylphenol	0.350	0.362	-3.4	105	-0.02
37		2-Methylnaphthalene	0.695	0.740	-6.5	109	-0.04
38		1-Methylnaphthalene	0.681	0.714	-4.8	108	-0.04
39	I	Acenaphthene-d10	1.000	1.000	0.0	92	-0.03
40		1,2,4,5-Tetrachlorobenzene	0.705	0.819	-16.2	106	-0.04
41	P	Hexachlorocyclopentadiene	0.364	0.009#	97.5#	2#	-0.04
42	S	2,4,6-Tribromophenol	0.287	0.290	-1.0	90	-0.02
43	C	2,4,6-Trichlorophenol	0.442	0.513	-16.1	103	-0.02
44		2,4,5-Trichlorophenol	0.466	0.525	-12.7	101	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.542	1.760	-14.1	104	-0.03
46		1,1'-Biphenyl	1.795	2.057	-14.6	106	-0.04
47		2-Chloronaphthalene	1.332	1.510	-13.4	104	-0.03
48		2-Nitroaniline	0.412	0.470	-14.1	99	-0.03
49		Acenaphthylene	2.002	2.197	-9.7	100	-0.03
50		Dimethylphthalate	1.626	1.735	-6.7	98	-0.02
51		2,6-Dinitrotoluene	0.362	0.344	5.0	85	-0.02
52	C	Acenaphthene	1.229	1.340	-9.0	100	-0.03
53		3-Nitroaniline	0.394	0.475	-20.6	111	-0.02
54	P	2,4-Dinitrophenol	0.186	0.005#	97.3#	2#	-0.01
55		Dibenzofuran	2.008	2.125	-5.8	98	-0.02
56	P	4-Nitrophenol	0.294	0.298	-1.4	91	-0.01
57		2,4-Dinitrotoluene	0.488	0.435	10.9	78	-0.02
58		Fluorene	1.537	1.577	-2.6	94	-0.02
59		2,3,4,6-Tetrachlorophenol	0.429	0.437	-1.9	91	-0.02
60		Diethylphthalate	1.757	1.788	-1.8	96	-0.02
61		4-Chlorophenyl-phenylether	0.873	0.909	-4.1	95	-0.02
62		4-Nitroaniline	0.372	0.465	-25.0#	113	-0.02
63		Azobenzene	1.624	1.578	2.8	88	-0.02
64	I	Phenanthrene-d10	1.000	1.000	0.0	78	-0.02
65		4,6-Dinitro-2-methylphenol	0.139	0.007	95.0#	4#	-0.02
66	c	n-Nitrosodiphenylamine	0.643	0.801	-24.6#	94	-0.02
67		4-Bromophenyl-phenylether	0.237	0.291	-22.8	93	-0.03
68		Hexachlorobenzene	0.267	0.305	-14.2	88	-0.02
69		Atrazine	0.237	0.206	13.1	65	-0.02
70	C	Pentachlorophenol	0.187	0.191	-2.1	78	-0.02
71		Phenanthrene	1.085	1.177	-8.5	83	-0.02
72		Anthracene	1.056	1.177	-11.5	84	-0.02
73		Carbazole	1.067	1.165	-9.2	82	-0.02
74		Di-n-butylphthalate	1.188	1.473	-24.0	92	-0.02
75	C	Fluoranthene	1.228	1.189	3.2	73	-0.02
76	I	Chrysene-d12	1.000	1.000	0.0	63	-0.02
77		Benzidine	0.651	0.897	-37.8#	82	-0.01
78		Pyrene	1.234	1.385	-12.2	70	-0.02
79	S	Terphenyl-d14	0.882	1.010	-14.5	72	-0.02
80		Butylbenzylphthalate	0.501	0.652	-30.1#	77	-0.02
81		Benzo(a)anthracene	1.152	1.251	-8.6	67	-0.01
82		3,3'-Dichlorobenzidine	0.445	0.563	-26.5#	75	-0.02
83		Chrysene	1.119	1.195	-6.8	66	-0.02
84		Bis(2-ethylhexyl)phthalate	0.741	0.921	-24.3	76	-0.02
85	c	Di-n-octyl phthalate	1.186	1.489	-25.5#	76	-0.02
86		Indeno(1,2,3-cd)pyrene	0.894	1.453	-62.5#	99	-0.02
87	I	Perylene-d12	1.000	1.000	0.0	77	-0.02

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88	Benzo(b)fluoranthene	1.171	1.175	-0.3	76	-0.02
89	Benzo(k)fluoranthene	1.184	1.162	1.9	73	-0.02
90 C	Benzo(a)pyrene	1.068	1.138	-6.6	80	-0.02
91	Dibenzo(a,h)anthracene	0.898	1.156	-28.7#	98	-0.03
92	Benzo(g,h,i)perylene	0.844	1.153	-36.6#	104	-0.02

(#) = Out of Range SPCC's out = 2 CCC's out = 2