Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM081016\
Data File : BM006997.D
Acq On : 10 Aug 2016 09:17
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02041

Quant Time: Aug 10 18:24:42 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080916.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 10 01:30:54 2016
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	53	0.00
2	S	1,4-Dioxane-d8	0.439	0.437	0.5	51	0.00
3	S	Phenol-d5	1.780	1.855	-4.2	54	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	0.995	1.051	-5.6	54	0.00
5	S	2-Chlorophenol-d4	1.350	1.387	-2.7	54	0.00
6	S	4-Methylphenol-d8	1.507	1.575	-4.5	54	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	56	0.00
8	S	Nitrobenzene-d5	0.153	0.142	7.2	51	0.00
9	S	2-Nitrophenol-d4	0.175	0.155	11.4	49	0.00
10	S	2,4-Dichlorophenol-d3	0.342	0.350	-2.3	57	0.00
11		Naphthalene	1.003	1.032	-2.9	57	0.00
12	S	4-Chloroaniline-d4	0.399	0.455	-14.0	58	0.00
13		2-Methylnaphthalene	0.794	0.832	-4.8	57	0.00
14		1-Methylnaphthalene	0.762	0.804	-5.5	58	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	61	0.00
16	S	Dimethylphthalate-d6	1.709	1.793	-4.9	62	0.00
17	S	Acenaphthylene-d8	2.002	2.067	-3.2	60	0.00
18		Acenaphthylene	2.069	2.147	-3.8	61	0.00
19	C	Acenaphthene	1.338	1.399	-4.6	61	0.00
20	S	4-Nitrophenol-d4	0.307	0.334	-8.8	64	0.00
21	S	Fluorene-d10	1.453	1.532	-5.4	62	0.00
22		Fluorene	1.588	1.721	-8.4	63	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	66	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.108	0.095	12.0	59	0.00
25		Phenanthrene	1.072	1.114	-3.9	66	0.00
26	S	Anthracene-d10	0.932	0.972	-4.3	66	0.00
27		Anthracene	1.107	1.164	-5.1	67	0.00
28	C	Fluoranthene	1.350	1.483	-9.9	70	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	79	0.00
30	S	Pyrene-d10	0.909	0.841	7.5	71	0.00
31		Pyrene	1.160	1.088	6.2	71	0.00
32		Benzo(a)anthracene	1.178	1.205	-2.3	79	0.00
33		Chrysene	1.069	1.105	-3.4	79	0.00
34	I	Perylene-d12	1.000	1.000	0.0	98	0.00
35		Benzo(b)fluoranthene	1.242	1.226	1.3	94	0.00
36		Benzo(k)fluoranthene	1.174	1.136	3.2	91	0.00
37	S	Benzo(a)pyrene-d12	0.919	0.943	-2.6	98	0.00
38	C	Benzo(a)pyrene	1.143	1.179	-3.1	98	0.00
39		Indeno(1,2,3-cd)pyrene	1.133	1.390	-22.7#	119	0.00
40		Dibenzo(a,h)anthracene	0.938	1.160	-23.7#	120	0.00
41		Benzo(g,h,i)perylene	0.932	1.169	-25.4#	123	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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