Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE092115\
Data File : BE090917.D
Acq On : 21 Sep 2015 14:21
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 22 02:16:22 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE091815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 18 18:17:34 2015
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	128	0.00
2		1,4-Dioxane	0.795	0.719	9.6	113	0.00
3		n-Nitrosodimethylamine	1.067	0.804	24.6	99	0.00
4	S	2-Fluorophenol	1.238	0.873	29.5#	91	0.00
5	S	Phenol-d6	1.559	1.197	23.2	101	0.01
6		bis(2-Chloroethyl)ether	1.546	1.471	4.9	133	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	130	0.00
8	S	Nitrobenzene-d5	0.345	0.295	14.5	112	0.00
9		Nitrobenzene	0.383	0.309	19.3	109	0.00
10		Naphthalene	1.126	1.030	8.5	121	0.00
11		Hexachlorobutadiene	0.166	0.166	0.0	129	0.00
12		2-Methylnaphthalene	0.672	0.595	11.5	118	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	132	0.00
14	S	2,4,6-Tribromophenol	0.196	0.107	45.4#	83	0.00
15	S	2-Fluorobiphenyl	1.418	1.283	9.5	119	0.00
16		Acenaphthylene	8.705	7.655	12.1	117	0.00
17		Acenaphthene	1.341	1.287	4.0	128	0.00
18		Fluorene	1.705	1.542	9.6	123	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	134	0.00
20		4-Bromophenyl-phenylether	0.193	0.169	12.4	119	0.00
21		Hexachlorobenzene	0.906	0.885	2.3	130	0.00
22		Pentachlorophenol	0.087	0.032	63.2#	54	0.02
23		Phenanthrene	1.292	1.185	8.3	126	0.00
24		Anthracene	1.144	1.001	12.5	121	0.00
25		Fluoranthene	1.525	1.254	17.8	111	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	117	0.00
27		Pyrene	1.141	1.006	11.8	111	0.00
28	S	Terphenyl-d14	0.638	0.530	16.9	107	0.00
29		Benzo(a)anthracene	1.235	1.055	14.6	105	0.00
30		Chrysene	1.352	1.254	7.2	109	0.00
31		Bis(2-ethylhexyl)phthalate	0.515	0.242	53.0#	63	0.00
32		Indeno(1,2,3-cd)pyrene	1.455	1.088	25.2#	87	0.00
33	I	Perylene-d12	1.000	1.000	0.0	104	0.00
34		Benzo(b)fluoranthene	1.170	1.012	13.5	93	0.00
35		Benzo(k)fluoranthene	1.347	1.231	8.6	96	0.00
36	C	Benzo(a)pyrene	1.113	0.923	17.1	90	0.00
37		Dibenzo(a,h)anthracene	1.158	0.937	19.1	85	0.00
38		Benzo(g,h,i)perylene	1.240	1.023	17.5	87	0.00

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