Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE092415\
Data File : BE090967.D
Acq On : 24 Sep 2015 5:54
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 24 08:36:38 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	133	0.00
2		1,4-Dioxane	0.795	0.799	-0.5	131	0.00
3		n-Nitrosodimethylamine	1.067	0.823	22.9	105	0.00
4	S	2-Fluorophenol	1.238	0.882	28.8#	96	0.00
5	S	Phenol-d6	1.559	1.031	33.9#	91	0.04
6		bis(2-Chloroethyl)ether	1.546	1.514	2.1	142	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	141	0.00
8	S	Nitrobenzene-d5	0.345	0.280	18.8	115	0.00
9		Nitrobenzene	0.383	0.284	25.8#	109	0.00
10		Naphthalene	1.126	1.099	2.4	139	0.00
11		Hexachlorobutadiene	0.166	0.187	-12.7	157#	0.00
12		2-Methylnaphthalene	0.672	0.651	3.1	140	0.01
13	I	Acenaphthene-d10	1.000	1.000	0.0	151#	0.00
14	S	2,4,6-Tribromophenol	0.196	0.118	39.8#	104	0.00
15	S	2-Fluorobiphenyl	1.418	1.357	4.3	143	0.00
16		Acenaphthylene	8.705	8.316	4.5	145	0.00
17		Acenaphthene	1.341	1.366	-1.9	155#	0.00
18		Fluorene	1.705	1.595	6.5	145	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	137	0.00
20		4-Bromophenyl-phenylether	0.193	0.193	0.0	139	0.00
21		Hexachlorobenzene	0.906	1.024	-13.0	154#	0.00
22		Pentachlorophenol	0.087	0.035	59.8#	60	0.04
23		Phenanthrene	1.292	1.222	5.4	133	0.00
24		Anthracene	1.144	1.032	9.8	128	0.00
25		Fluoranthene	1.525	1.341	12.1	121	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	107	0.00
27		Pyrene	1.141	1.197	-4.9	120	0.00
28	S	Terphenyl-d14	0.638	0.611	4.2	112	0.01
29		Benzo(a)anthracene	1.235	1.048	15.1	95	0.00
30		Chrysene	1.352	1.445	-6.9	115	0.00
31		Bis(2-ethylhexyl)phthalate	0.515	0.365	29.1#	86	0.00
32		Indeno(1,2,3-cd)pyrene	1.455	1.208	17.0	88	0.00
33	I	Perylene-d12	1.000	1.000	0.0	97	0.00
34		Benzo(b)fluoranthene	1.170	1.033	11.7	88	0.00
35		Benzo(k)fluoranthene	1.347	1.468	-9.0	106	0.00
36	C	Benzo(a)pyrene	1.113	1.163	-4.5	105	0.00
37		Dibenzo(a,h)anthracene	1.158	0.996	14.0	84	0.00
38		Benzo(g,h,i)perylene	1.240	1.127	9.1	89	0.00

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