Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091815\
Data File : BE090890.D
Acq On : 18 Sep 2015 17:49
Operator : UM/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE091815

Quant Time: Sep 18 18:33:10 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	102	0.00
2		1,4-Dioxane	0.795	0.738	7.2	93	0.00
3		n-Nitrosodimethylamine	1.067	0.938	12.1	92	0.00
4	S	2-Fluorophenol	1.238	1.074	13.2	90	0.00
5	S	Phenol-d6	1.559	1.382	11.4	93	0.00
6		bis(2-Chloroethyl)ether	1.546	1.459	5.6	105	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
8	S	Nitrobenzene-d5	0.345	0.317	8.1	94	0.00
9		Nitrobenzene	0.383	0.347	9.4	95	0.00
10		Naphthalene	1.126	1.039	7.7	94	0.00
11		Hexachlorobutadiene	0.166	0.157	5.4	95	0.00
12		2-Methylnaphthalene	0.672	0.617	8.2	95	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
14	S	2,4,6-Tribromophenol	0.196	0.150	23.5	90	-0.02
15	S	2-Fluorobiphenyl	1.418	1.331	6.1	96	0.00
16		Acenaphthylene	8.705	7.970	8.4	95	0.00
17		Acenaphthene	1.341	1.247	7.0	97	0.00
18		Fluorene	1.705	1.567	8.1	97	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
20		4-Bromophenyl-phenylether	0.193	0.180	6.7	96	-0.02
21		Hexachlorobenzene	0.906	0.853	5.8	96	0.00
22		Pentachlorophenol	0.087	0.059	32.2#	77	0.00
23		Phenanthrene	1.292	1.227	5.0	99	0.00
24		Anthracene	1.144	1.031	9.9	95	0.00
25		Fluoranthene	1.525	1.370	10.2	93	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	95	0.00
27		Pyrene	1.141	1.030	9.7	92	0.00
28	S	Terphenyl-d14	0.638	0.552	13.5	91	0.00
29		Benzo(a)anthracene	1.235	1.080	12.6	88	0.00
30		Chrysene	1.352	1.245	7.9	88	0.00
31		Bis(2-ethylhexyl)phthalate	0.515	0.353	31.5#	75	0.00
32		Indeno(1,2,3-cd)pyrene	1.455	1.267	12.9	83	0.00
33	I	Perylene-d12	1.000	1.000	0.0	92	0.00
34		Benzo(b)fluoranthene	1.170	1.016	13.2	83	0.00
35		Benzo(k)fluoranthene	1.347	1.244	7.6	86	0.00
36	C	Benzo(a)pyrene	1.113	0.964	13.4	83	0.00
37		Dibenzo(a,h)anthracene	1.158	1.009	12.9	81	0.00
38		Benzo(g,h,i)perylene	1.240	1.124	9.4	85	0.00

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