Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061215\
Data File : BM001658.D
Acq On : 12 Jun 2015 01:13
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM061215

Quant Time: Jun 12 16:46:12 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 12 16:27:37 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	80	0.00
2		1,4-Dioxane	0.624	0.574	8.0	84	0.00
3		n-Nitrosodimethylamine	0.872	0.882	-1.1	83	0.00
4	S	2-Fluorophenol	1.184	1.139	3.8	81	0.00
5	S	Phenol-d6	1.488	1.418	4.7	81	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	81	0.00
7	S	Nitrobenzene-d5	0.383	0.365	4.7	82	0.00
8		Nitrobenzene	0.414	0.393	5.1	81	0.00
9		Naphthalene	1.212	1.180	2.6	83	0.00
10		Hexachlorobutadiene	0.201	0.196	2.5	85	0.00
11		2-Methylnaphthalene	0.782	0.768	1.8	85	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	82	0.00
13	S	2,4,6-Tribromophenol	0.182	0.187	-2.7	82	0.00
14	S	2-Fluorobiphenyl	1.746	1.727	1.1	86	0.00
15		Acenaphthylene	2.337	2.278	2.5	83	0.00
16		Acenaphthene	1.487	1.419	4.6	84	0.00
17		Fluorene	1.519	1.521	-0.1	87	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	80	0.00
19		4-Bromophenyl-phenylether	0.502	0.516	-2.8	81	0.00
20		Hexachlorobenzene	0.455	0.515	-13.2	91	0.00
21		Pentachlorophenol	0.124	0.113	8.9	81	0.00
22		Phenanthrene	1.640	1.543	5.9	81	0.00
23		Anthracene	1.938	1.908	1.5	79	0.00
24		Fluoranthene	2.429	2.458	-1.2	84	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	79	0.00
26		Pyrene	1.716	1.720	-0.2	83	0.00
27	S	Terphenyl-d14	0.754	0.756	-0.3	82	0.00
28		Benzo(a)anthracene	1.571	1.545	1.7	82	0.00
29		Chrysene	1.503	1.429	4.9	79	0.00
30		Bis(2-ethylhexyl)phthalate	0.979	0.868	11.3	81	0.00
31		Indeno(1,2,3-cd)pyrene	1.181	1.016	14.0	75	0.00
32	I	Perylene-d12	1.000	1.000	0.0	76	0.00
33		Benzo(b)fluoranthene	1.631	1.585	2.8	79	0.00
34		Benzo(k)fluoranthene	1.584	1.560	1.5	78	0.00
35	C	Benzo(a)pyrene	1.359	1.293	4.9	76	0.00
36		Dibenzo(a,h)anthracene	1.142	1.026	10.2	74	0.00
37		Benzo(g,h,i)perylene	1.234	1.143	7.4	77	0.00

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