

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031816\
 Data File : BM004687.D
 Acq On : 18 Mar 2016 17:44
 Operator : UM/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM031816

Quant Time: Mar 19 07:55:07 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM031816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 19 07:50:46 2016
 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	97	0.00
2	1,4-Dioxane	0.400	0.461	-15.3	94	0.00
3	n-Nitrosodimethylamine	0.400	0.481	-20.2	99	0.00
4 S	2-Fluorophenol	0.400	0.459	-14.8	95	0.00
5 S	Phenol-d6	0.400	0.441	-10.2	93	0.00
6 I	Naphthalene-d8	5.000	5.000	0.0	97	0.00
7 S	Nitrobenzene-d5	0.400	0.442	-10.5	93	0.00
8	Naphthalene	0.400	0.455	-13.7	95	0.00
9	2-Methylnaphthalene	0.400	0.464	-16.0	95	0.00
10 I	Acenaphthene-d10	5.000	5.000	0.0	98	0.00
11 S	2,4,6-Tribromophenol	0.400	0.390	2.5	84	0.00
12 S	2-Fluorobiphenyl	0.400	0.466	-16.5	96	0.00
13	Acenaphthylene	0.400	0.428	-7.0	93	0.00
14	Acenaphthene	0.400	0.458	-14.5	95	0.00
15	Fluorene	0.400	0.446	-11.5	94	-0.02
16 I	Phenanthrene-d10	5.000	5.000	0.0	94	0.00
17	Hexachlorobenzene	0.400	0.478	-19.5	95	0.00
18	Pentachlorophenol	0.400	0.462	-15.5	88	0.00
19	Phenanthrene	0.400	0.478	-19.5	96	0.00
20	Anthracene	0.400	0.446	-11.5	91	0.00
21	Fluoranthene	0.400	0.457	-14.2	93	-0.02
22 I	Chrysene-d12	5.000	5.000	0.0	104	0.00
23	Pyrene	0.400	0.405	-1.3	93	0.00
24 S	Terphenyl-d14	0.400	0.409	-2.2	94	0.00
25	Benzo(a)anthracene	0.400	0.371	7.3	83	0.00
26	Chrysene	0.400	0.470	-17.5	111	0.00
27	Indeno(1,2,3-cd)pyrene	0.400	0.433	-8.2	98	0.00
28 I	Perylene-d12	5.000	5.000	0.0	99	0.00
29	Benzo(b)fluoranthene	0.400	0.446	-11.5	98	0.00
30	Benzo(k)fluoranthene	0.400	0.453	-13.2	98	-0.01
31 C	Benzo(a)pyrene	0.400	0.442	-10.5	96	0.00
32	Dibenzo(a,h)anthracene	0.400	0.449	-12.2	98	0.00
33	Benzo(g,h,i)perylene	0.400	0.451	-12.7	98	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0