

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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	97	0.00
2	1,4-Dioxane	0.494	0.582	-17.8	94	0.00
3	n-Nitrosodimethylamine	0.517	0.621	-20.1	99	0.00
4 S	2-Fluorophenol	1.154	1.323	-14.6	95	0.00
5 S	Phenol-d6	1.413	1.558	-10.3	93	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
7 S	Nitrobenzene-d5	0.241	0.266	-10.4	93	0.00
8	Naphthalene	1.375	1.565	-13.8	95	0.00
9	2-Methylnaphthalene	0.861	1.000	-16.1	95	0.00
10 I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
11 S	2,4,6-Tribromophenol	0.166	0.162	2.4	84	0.00
12 S	2-Fluorobiphenyl	1.431	1.666	-16.4	96	0.00
13	Acenaphthylene	2.572	2.750	-6.9	93	0.00
14	Acenaphthene	1.702	1.949	-14.5	95	0.00
15	Fluorene	1.908	2.129	-11.6	94	-0.02
16 I	Phenanthrene-d10	1.000	1.000	0.0	94	0.00
17	Hexachlorobenzene	0.308	0.368	-19.5	95	0.00
18	Pentachlorophenol	0.169	0.167	1.2	88	0.00
19	Phenanthrene	2.004	2.397	-19.6	96	0.00
20	Anthracene	1.429	1.595	-11.6	91	0.00
21	Fluoranthene	1.895	2.166	-14.3	93	-0.02
22 I	Chrysene-d12	1.000	1.000	0.0	104	0.00
23	Pyrene	1.835	1.859	-1.3	93	0.00
24 S	Terphenyl-d14	0.632	0.646	-2.2	94	0.00
25	Benzo(a)anthracene	2.000	1.857	7.2	83	0.00
26	Chrysene	1.487	1.746	-17.4	111	0.00
27	Indeno(1,2,3-cd)pyrene	1.674	1.810	-8.1	98	0.00
28 I	Perylene-d12	1.000	1.000	0.0	99	0.00
29	Benzo(b)fluoranthene	1.819	2.027	-11.4	98	0.00
30	Benzo(k)fluoranthene	1.802	2.041	-13.3	98	-0.01
31 C	Benzo(a)pyrene	1.706	1.883	-10.4	96	0.00
32	Dibenzo(a,h)anthracene	1.527	1.712	-12.1	98	0.00
33	Benzo(g,h,i)perylene	1.613	1.817	-12.6	98	0.00

(#) = Out of Range

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