

Data Path : Z:\HPCHEM1\BNA_M\Data\BM060615\
 Data File : BM001578.D
 Acq On : 06 Jun 2015 05:06
 Operator : TP/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM060615

Quant Time: Jun 08 05:41:17 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 08 05:27:25 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	101	0.00
2	1,4-Dioxane	0.523	0.476	9.0	101	0.00
3	n-Nitrosodimethylamine	0.763	0.753	1.3	102	0.00
4 S	2-Fluorophenol	1.104	1.070	3.1	102	0.00
5 S	Phenol-d6	1.357	1.314	3.2	102	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
7 S	Nitrobenzene-d5	0.341	0.324	5.0	100	0.00
8	Nitrobenzene	0.365	0.346	5.2	100	0.00
9	Naphthalene	1.080	1.043	3.4	100	0.00
10	Hexachlorobutadiene	0.188	0.182	3.2	101	0.00
11	2-Methylnaphthalene	0.676	0.644	4.7	99	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
13 S	2,4,6-Tribromophenol	0.207	0.212	-2.4	103	0.00
14 S	2-Fluorobiphenyl	1.541	1.501	2.6	97	0.00
15	Acenaphthylene	2.158	2.077	3.8	96	0.00
16	Acenaphthene	1.359	1.273	6.3	94	0.00
17	Fluorene	0.875	0.859	1.8	102	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	71	0.00
19	4-Bromophenyl-phenylether	0.340	0.443	-30.3#	104	0.00
20	Hexachlorobenzene	0.159	0.211	-32.7#	105	0.00
21	Pentachlorophenol	0.068	0.092	-35.3#	115	0.00
22	Phenanthrene	1.744	1.781	-2.1	77	0.00
23	Anthracene	0.735	0.921	-25.3#	104	0.00
24	Fluoranthene	1.357	1.642	-21.0	95	0.00
25 I	Chrysene-d12	1.000	1.000	0.0	101	0.00
26	Pyrene	1.608	1.504	6.5	96	0.00
27 S	Terphenyl-d14	0.646	0.591	8.5	93	0.00
28	Benzo(a)anthracene	1.408	1.290	8.4	97	0.00
29	Chrysene	1.345	1.257	6.5	98	0.00
30	Bis(2-ethylhexyl)phthalate	0.829	0.734	11.5	98	0.00
31	Indeno(1,2,3-cd)pyrene	1.444	1.275	11.7	96	0.01
32 I	Perylene-d12	1.000	1.000	0.0	99	0.00
33	Benzo(b)fluoranthene	1.494	1.343	10.1	98	0.00
34	Benzo(k)fluoranthene	1.360	1.349	0.8	98	0.01
35 C	Benzo(a)pyrene	1.305	1.221	6.4	98	0.01
36	Dibenzo(a,h)anthracene	1.280	1.139	11.0	96	0.01
37	Benzo(g,h,i)perylene	1.345	1.206	10.3	97	0.00

(#) = Out of Range

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