

Data Path : Z:\HPCHEM1\BNA M\DATA\BM110515\
 Data File : BM003059.D
 Acq On : 05 Nov 2015 11:21
 Operator : UM/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 ICVBM110415

Quant Time: Nov 16 18:53:40 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM110415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 16 18:52:15 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	134	0.00
2	1,4-Dioxane	1.000	0.987	1.3	129	0.00
3	n-Nitrosodimethylamine	1.000	0.948	5.2	132	0.00
4 S	2-Fluorophenol	1.000	0.970	3.0	136	0.00
5 S	Phenol-d6	1.000	1.005	-0.5	143	0.00
6	bis(2-Chloroethyl)ether	1.000	0.974	2.6	134	-0.02
7 I	Naphthalene-d8	5.000	5.000	0.0	140	0.00
8 S	Nitrobenzene-d5	1.000	0.987	1.3	146	0.00
9	Nitrobenzene	1.000	0.989	1.1	146	0.00
10	Naphthalene	1.000	0.975	2.5	138	0.00
11	Hexachlorobutadiene	1.000	0.961	3.9	135	0.00
12	2-Methylnaphthalene	1.000	0.992	0.8	141	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	148	0.00
14 S	2,4,6-Tribromophenol	1.000	1.004	-0.4	152	0.00
15 S	2-Fluorobiphenyl	1.000	0.946	5.4	142	0.00
16	Acenaphthylene	1.000	0.980	2.0	152	0.00
17	Acenaphthene	1.000	0.970	3.0	156	0.00
18	Fluorene	1.000	0.953	4.7	139	0.02
19 I	Phenanthrene-d10	5.000	5.000	0.0	183	0.00
20	4-Bromophenyl-phenylether	1.000	0.925	7.5	118	0.00
21	Hexachlorobenzene	1.000	0.957	4.3	129	0.00
22	Pentachlorophenol	1.000	0.975	2.5	141	0.02
23	Phenanthrene	1.000	0.875	12.5	112	0.00
24	Anthracene	1.000	0.952	4.8	175	0.00
25	Fluoranthene	1.000	0.925	7.5	147	0.01
26 I	Chrysene-d12	5.000	5.000	0.0	148	0.00
27	Benzydine	1.000	0.988	1.2	129	0.00
28	Pvrene	1.000	0.980	2.0	148	0.00
29 S	Terphenyl-d14	1.000	0.962	3.8	138	0.00
30	Benzo(a)anthracene	1.000	0.974	2.6	145	0.00
31	3,3'-Dichlorobenzidine	1.000	0.981	1.9	126	0.00
32	Chrysene	1.000	0.804	19.6	106	0.01
33	Bis(2-ethylhexyl)phthalate	1.000	1.083	-8.3	180	0.01
34	Indeno(1,2,3-cd)pyrene	1.000	1.074	-7.4	141	0.00
35 I	Perylene-d12	5.000	5.000	0.0	138	0.00
36	Benzo(b)fluoranthene	1.000	0.920	8.0	130	0.00
37	Benzo(k)fluoranthene	1.000	0.975	2.5	136	0.00
38 C	Benzo(a)pyrene	1.000	0.923	7.7	140	0.00
39	Dibenzo(a,h)anthracene	1.000	0.979	2.1	142	0.01
40	Benzo(a,h,i)perylene	1.000	0.938	6.2	139	0.00

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Compound	Amount	Calc.	%Dev Area	% Dev(min)

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