

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032817\
 Data File : BE092836.D
 Acq On : 28 Mar 2017 14:25
 Operator : SJ/MA
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE032817

Quant Time: Mar 28 15:01:45 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE032817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 28 14:58:05 2017
 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	98	0.00
2	1,4-Dioxane	0.400	0.424	-6.0	101	-0.01
3	n-Nitrosodimethylamine	0.400	0.384	4.0	101	0.00
4 S	2-Fluorophenol	0.400	0.385	3.8	96	0.00
5 S	Phenol-d6	0.400	0.379	5.3	100	0.00
6	bis(2-Chloroethyl)ether	0.400	0.400	0.0	100	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	100	0.00
8 S	Nitrobenzene-d5	0.400	0.389	2.8	109	0.00
9	Naphthalene	0.400	0.399	0.3	99	0.00
10	Hexachlorobutadiene	0.400	0.392	2.0	99	0.00
11	2-Methylnaphthalene	0.400	0.387	3.3	99	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	99	0.00
13 S	2,4,6-Tribromophenol	0.400	0.326	18.5	90	0.00
14 S	2-Fluorobiphenyl	0.400	0.405	-1.3	99	-0.01
15	Acenaphthylene	0.400	0.350	12.5	97	0.00
16	Acenaphthene	0.400	0.383	4.3	98	0.00
17	Fluorene	0.400	0.382	4.5	96	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	95	0.00
19	Hexachlorobenzene	0.400	0.414	-3.5	101	0.00
20	Pentachlorophenol	0.400	0.457	-14.2	83	0.00
21	Phenanthrene	0.400	0.438	-9.5	101	0.00
22	Anthracene	0.400	0.390	2.5	100	0.00
23	Fluoranthene	0.400	0.366	8.5	94	0.00
24 I	Chrysene-d12	5.000	5.000	0.0	95	-0.02
25	Pyrene	0.400	0.373	6.8	94	0.00
26 S	Terphenyl-d14	0.400	0.372	7.0	92	0.00
27	Benzo(a)anthracene	0.400	0.373	6.8	97	0.00
28	Chrysene	0.400	0.394	1.5	95	0.00
29	Bis(2-ethylhexyl)phthalate	0.400	0.363	9.3	87	0.00
30	Indeno(1,2,3-cd)pyrene	0.400	0.365	8.8	93	0.00
31 I	Perylene-d12	5.000	5.000	0.0	96	0.00
32	Benzo(b)fluoranthene	0.400	0.342	14.5	85	-0.01
33	Benzo(k)fluoranthene	0.400	0.395	1.3	100	0.00
34 C	Benzo(a)pyrene	0.400	0.348	13.0	90	0.00
35	Dibenzo(a,h)anthracene	0.400	0.364	9.0	92	0.00
36	Benzo(g,h,i)perylene	0.400	0.374	6.5	92	0.00

(#) = Out of Range

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