

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE032415\
 Data File : BE089819.D
 Acq On : 24 Mar 2015 18:12
 Operator : TP/JJ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE032415

Quant Time: Mar 25 16:40:43 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIMPAH-BE032415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 24 18:03:12 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	131	0.00
2	1,4-Dioxane	1.000	0.997	0.3	134	0.03
3	n-Nitrosodimethylamine	1.000	1.042	-4.2	143	-0.03
4 I	Naphthalene-d8	5.000	5.000	0.0	130	0.00
5 S	Nitrobenzene-d5	1.000	1.201	-20.1	191	0.00
6	Naphthalene	1.000	0.983	1.7	133	0.00
7	2-Methylnaphthalene	1.000	1.026	-2.6	129	0.00
8 I	Acenaphthene-d10	5.000	5.000	0.0	116	0.00
9 S	2-Fluorobiphenyl	1.000	1.027	-2.7	123	0.00
10	Acenaphthylene	1.000	1.039	-3.9	125	-0.01
11	Acenaphthene	1.000	1.003	-0.3	122	0.00
12	Fluorene	1.000	1.010	-1.0	122	-0.01
13 I	Phenanthrene-d10	5.000	5.000	0.0	118	0.00
14	Hexachlorobenzene	1.000	0.980	2.0	121	0.00
15	Pentachlorophenol	1.000	0.556	44.4#	97	0.00
16	Phenanthrene	1.000	1.001	-0.1	125	0.00
17	Anthracene	1.000	1.012	-1.2	124	0.00
18	Fluoranthene	1.000	1.082	-8.2	137	0.00
19 I	Chrysene-d12	5.000	5.000	0.0	135	0.00
20	Pyrene	1.000	0.936	6.4	133	0.00
21 S	Terphenyl-d14	1.000	0.930	7.0	131	0.00
22	Benzo(a)anthracene	1.000	0.978	2.2	143	0.00
23	Chrysene	1.000	1.038	-3.8	143	0.00
24	Indeno(1,2,3-cd)pyrene	1.000	0.912	8.8	143	0.00
25 I	Perylene-d12	5.000	5.000	0.0	140	0.00
26	Benzo(b)fluoranthene	1.000	0.879	12.1	132	0.00
27	Benzo(k)fluoranthene	1.000	1.116	-11.6	157	0.00
28 C	Benzo(a)pyrene	1.000	0.898	10.2	136	0.00
29	Dibenzo(a,h)anthracene	1.000	0.859	14.1	123	0.00
30	Benzo(g,h,i)perylene	1.000	0.970	3.0	143	0.00

(#) = Out of Range

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