

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061217\
 Data File : BE093102.D
 Acq On : 12 Jun 2017 13:27
 Operator : SJ/MA
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE061217

Quant Time: Jun 12 14:07:47 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE061217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 12 14:07:54 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	0.400	0.400	0.0	104	0.00
2	1,4-Dioxane	0.400	0.390	2.5	120	0.00
3	n-Nitrosodimethylamine	0.400	0.398	0.5	114	0.00
4 S	2-Fluorophenol	0.400	0.396	1.0	120	0.00
5 S	Phenol-d6	0.400	0.386	3.5	114	0.00
6	bis(2-Chloroethyl)ether	0.400	0.403	-0.8	117	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	105	-0.02
8 S	Nitrobenzene-d5	0.400	0.400	0.0	121	0.00
9	Naphthalene	0.400	0.394	1.5	117	0.00
10	Hexachlorobutadiene	0.400	0.388	3.0	118	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.391	2.3	119	0.00
12	2-Methylnaphthalene	0.400	0.393	1.8	118	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	112	0.00
14 S	2,4,6-Tribromophenol	0.400	0.379	5.3	128	0.00
15 S	2-Fluorobiphenyl	0.400	0.389	2.8	121	0.00
16	Acenaphthylene	0.400	0.415	-3.7	139	0.00
17	Acenaphthene	0.400	0.404	-1.0	130	0.00
18	Fluorene	0.400	0.401	-0.3	129	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	113	0.00
20	4-Bromophenyl-phenylether	0.400	0.397	0.8	130	0.00
21	Hexachlorobenzene	0.400	0.341	14.8	125	0.00
22	Pentachlorophenol	0.400	0.339	15.3	121	0.00
23	Phenanthrene	0.400	0.376	6.0	125	0.00
24	Anthracene	0.400	0.378	5.5	138	0.00
25 SURR	Fluoranthene-d10	0.400	0.375	6.3	126	0.00
26	Fluoranthene	0.400	0.373	6.8	129	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	99	0.00
28	Pyrene	0.400	0.414	-3.5	112	0.00
29 S	Terphenyl-d14	0.400	0.429	-7.2	116	0.00
30	Benzo(a)anthracene	0.400	0.397	0.8	110	0.00
31	Chrysene	0.400	0.399	0.3	110	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.441	-10.2	115	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.383	4.3	107	0.00
34 I	Perylene-d12	0.400	0.400	0.0	94	0.00
35	Benzo(b)fluoranthene	0.400	0.393	1.8	108	0.00
36	Benzo(k)fluoranthene	0.400	0.397	0.8	108	0.00
37 C	Benzo(a)pyrene	0.400	0.392	2.0	108	0.00
38	Dibenzo(a,h)anthracene	0.400	0.393	1.8	108	0.00
39	Benzo(g,h,i)perylene	0.400	0.384	4.0	106	0.02

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

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