

Data Path : Z:\HPCHEM1\BNA E\Data\BE080717\
 Data File : BE093699.D
 Acq On : 7 Aug 2017 13:38
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE080717

Quant Time: Aug 07 14:33:31 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 07 14:30:21 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	112	0.00
2	1,4-Dioxane	0.400	0.404	-1.0	112	0.00
3	n-Nitrosodimethylamine	0.400	0.402	-0.5	110	0.01
4 S	2-Fluorophenol	0.400	0.401	-0.3	114	0.00
5 S	Phenol-d6	0.400	0.388	3.0	113	0.00
6	bis(2-Chloroethyl)ether	0.400	0.401	-0.3	116	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	114	0.00
8 S	Nitrobenzene-d5	0.400	0.372	7.0	112	0.00
9	Naphthalene	0.400	0.399	0.3	115	0.00
10	Hexachlorobutadiene	0.400	0.397	0.8	115	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.401	-0.3	115	0.00
12	2-Methylnaphthalene	0.400	0.395	1.3	114	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	115	0.00
14 S	2,4,6-Tribromophenol	0.400	0.333	16.8	105	0.00
15 S	2-Fluorobiphenyl	0.400	0.406	-1.5	117	0.00
16	Acenaphthylene	0.400	0.385	3.8	115	0.00
17	Acenaphthene	0.400	0.392	2.0	113	0.00
18	Fluorene	0.400	0.390	2.5	114	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	116	0.00
20	4-Bromophenyl-phenylether	0.400	0.406	-1.5	120	0.00
21	Hexachlorobenzene	0.400	0.393	1.8	110	0.00
22	Pentachlorophenol	0.400	0.330	17.5	102	0.00
23	Phenanthrene	0.400	0.388	3.0	115	0.00
24	Anthracene	0.400	0.381	4.8	113	0.00
25 SURR	Fluoranthene-d10	0.400	0.378	5.5	104	0.00
26	Fluoranthene	0.400	0.375	6.3	104	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	99	0.00
28	Pvrene	0.400	0.403	-0.8	102	0.00
29 S	Terphenyl-d14	0.400	0.409	-2.2	104	0.00
30	Benzo(a)anthracene	0.400	0.387	3.3	99	0.00
31	Chrysene	0.400	0.394	1.5	99	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.336	16.0	82	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.404	-1.0	100	0.00
34 I	Perylene-d12	0.400	0.400	0.0	99	0.00
35	Benzo(b)fluoranthene	0.400	0.400	0.0	98	0.00
36	Benzo(k)fluoranthene	0.400	0.405	-1.3	103	0.00
37 C	Benzo(a)pyrene	0.400	0.399	0.3	99	0.00
38	Dibenzo(a,h)anthracene	0.400	0.398	0.5	99	0.00
39	Benzo(a,h,i)perylene	0.400	0.395	1.3	99	0.00

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

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