

Data Path : Z:\HPCHEM1\BNA E\Data\BE021318\
 Data File : BE095656.D
 Acq On : 12 Feb 2018 11:55
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Feb 13 03:23:15 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE020618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 20:27:34 2018
 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	96	0.00
2	1,4-Dioxane	0.400	0.410	-2.5	98	0.00
3	n-Nitrosodimethylamine	0.400	0.423	-5.7	100	-0.01
4 S	2-Fluorophenol	0.400	0.397	0.8	94	0.00
5 S	Phenol-d6	0.400	0.398	0.5	94	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.408	-2.0	95	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	100	-0.02
8 S	Nitrobenzene-d5	0.400	0.375	6.3	93	-0.02
9	Naphthalene	0.400	0.404	-1.0	98	-0.02
10	Hexachlorobutadiene	0.400	0.424	-6.0	95	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.421	-5.2	103	0.00
12	2-Methylnaphthalene	0.400	0.416	-4.0	100	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	104	0.00
14 S	2,4,6-Tribromophenol	0.400	0.333	16.8	89	0.00
15 S	2-Fluorobiphenyl	0.400	0.405	-1.3	104	-0.02
16	Acenaphthylene	0.400	0.392	2.0	102	0.00
17	Acenaphthene	0.400	0.389	2.8	101	0.00
18	Fluorene	0.400	0.390	2.5	100	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	97	0.00
20	4-Bromophenyl-phenylether	0.400	0.416	-4.0	98	-0.01
21	Hexachlorobenzene	0.400	0.417	-4.2	100	0.00
22	Pentachlorophenol	0.400	0.407	-1.7	106	0.00
23	Phenanthrene	0.400	0.408	-2.0	98	-0.01
24	Anthracene	0.400	0.402	-0.5	99	-0.01
25 SURR	Fluoranthene-d10	0.400	0.391	2.3	96	0.00
26	Fluoranthene	0.400	0.391	2.3	97	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	96	0.00
28	Pvrene	0.400	0.419	-4.7	105	0.00
29 S	Terphenyl-d14	0.400	0.394	1.5	92	0.00
30	Benzo(a)anthracene	0.400	0.416	-4.0	99	0.00
31	Chrysene	0.400	0.410	-2.5	96	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.348	13.0	88	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.432	-8.0	106	0.00
34 I	Perylene-d12	0.400	0.400	0.0	97	0.00
35	Benzo(b)fluoranthene	0.400	0.403	-0.8	97	0.00
36	Benzo(k)fluoranthene	0.400	0.404	-1.0	98	0.00
37 C	Benzo(a)pyrene	0.400	0.407	-1.7	100	0.00
38	Dibenzo(a,h)anthracene	0.400	0.430	-7.5	108	0.00
39	Benzo(a,h,i)perylene	0.400	0.414	-3.5	101	0.00

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

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