

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061719\
 Data File : BE100069.D
 Acq On : 17 Jun 2019 19:25
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jun 18 03:44:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 17 17:23:36 2019
 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	93	0.00
2	1,4-Dioxane	0.400	0.360	10.0	84	0.01
3	n-Nitrosodimethylamine	0.400	0.414	-3.5	69	0.01
4 S	2-Fluorophenol	0.400	0.360	10.0	89	0.00
5 S	Phenol-d6	0.400	0.340	15.0	83	0.01
6	bis(2-Chloroethyl)ether	0.400	0.413	-3.2	93	0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	96	0.00
8 S	Nitrobenzene-d5	0.400	0.355	11.3	94	0.01
9	Naphthalene	0.400	0.406	-1.5	97	0.00
10	Hexachlorobutadiene	0.400	0.425	-6.2	101	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.404	-1.0	100	0.00
12	2-Methylnaphthalene	0.400	0.412	-3.0	102	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	97	0.00
14 S	2,4,6-Tribromophenol	0.400	0.341	14.8	96	0.00
15 S	2-Fluorobiphenyl	0.400	0.395	1.3	94	0.00
16	Acenaphthylene	0.400	0.404	-1.0	100	0.00
17	Acenaphthene	0.400	0.401	-0.3	97	-0.01
18	Fluorene	0.400	0.417	-4.2	102	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	100	0.01
20	4-Bromophenyl-phenylether	0.400	0.390	2.5	101	0.00
21	Hexachlorobenzene	0.400	0.399	0.3	102	0.00
22	Pentachlorophenol	0.400	0.403	-0.8	88	0.01
23	Phenanthrene	0.400	0.383	4.3	99	0.00
24	Anthracene	0.400	0.376	6.0	101	0.00
25 SURR	Fluoranthene-d10	0.400	0.377	5.8	98	0.00
26	Fluoranthene	0.400	0.381	4.8	99	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	94	0.00
28	Pvrene	0.400	0.462	-15.5	98	0.00
29 S	Terphenyl-d14	0.400	0.437	-9.2	94	0.01
30	Benzo(a)anthracene	0.400	0.427	-6.7	94	0.00
31	Chrysene	0.400	0.463	-15.8	97	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.350	12.5	81	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.443	-10.7	93	0.00
34 I	Perylene-d12	0.400	0.400	0.0	92	0.00
35	Benzo(b)fluoranthene	0.400	0.408	-2.0	92	0.00
36	Benzo(k)fluoranthene	0.400	0.408	-2.0	91	0.00
37 C	Benzo(a)pyrene	0.400	0.420	-5.0	96	0.00
38	Dibenzo(a,h)anthracene	0.400	0.406	-1.5	92	0.01
39	Benzo(a,h,i)perylene	0.400	0.411	-2.7	92	0.00

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

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