

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091319\
 Data File : BN007643.D
 Acq On : 14 Sep 2019 00:48
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 14 05:20:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 12 14:27:51 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	72	0.00
2	1,4-Dioxane	0.400	0.382	4.5	72	0.00
3	n-Nitrosodimethylamine	0.400	0.611	-52.7#	78	0.00
4 S	2-Fluorophenol	0.400	0.327	18.3	65	0.00
5 S	Phenol-d6	0.400	0.344	14.0	64	0.00
6	bis(2-Chloroethyl)ether	0.400	0.373	6.8	67	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	72	-0.01
8 S	Nitrobenzene-d5	0.400	0.378	5.5	65	-0.01
9	Naphthalene	0.400	0.427	-6.7	72	-0.01
10	Hexachlorobutadiene	0.400	0.414	-3.5	70	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.417	-4.2	71	0.00
12	2-Methylnaphthalene	0.400	0.421	-5.2	71	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	71	0.00
14 S	2,4,6-Tribromophenol	0.400	0.300	25.0#	62	0.00
15 S	2-Fluorobiphenyl	0.400	0.411	-2.7	73	0.00
16	Acenaphthylene	0.400	0.358	10.5	65	0.00
17	Acenaphthene	0.400	0.393	1.8	70	0.00
18	Fluorene	0.400	0.389	2.8	70	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	70	0.00
20	4-Bromophenyl-phenylether	0.400	0.424	-6.0	70	0.00
21	Hexachlorobenzene	0.400	0.426	-6.5	70	0.00
22	Pentachlorophenol	0.400	0.323	19.3	62	0.00
23	Phenanthrene	0.400	0.426	-6.5	70	0.00
24	Anthracene	0.400	0.395	1.3	66	0.00
25 SURR	Fluoranthene-d10	0.400	0.402	-0.5	67	0.00
26	Fluoranthene	0.400	0.408	-2.0	67	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	67	-0.02
28	Pvrene	0.400	0.427	-6.7	67	0.00
29 S	Terphenyl-d14	0.400	0.428	-7.0	68	0.00
30	Benzo(a)anthracene	0.400	0.380	5.0	61	0.00
31	Chrysene	0.400	0.437	-9.2	70	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.310	22.5	55	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.470	-17.5	74	-0.03
34 I	Perylene-d12	0.400	0.400	0.0	68	-0.02
35	Benzo(b)fluoranthene	0.400	0.401	-0.3	68	-0.02
36	Benzo(k)fluoranthene	0.400	0.424	-6.0	72	-0.02
37 C	Benzo(a)pyrene	0.400	0.385	3.8	66	-0.02
38	Dibenzo(a,h)anthracene	0.400	0.430	-7.5	73	-0.02
39	Benzo(a,h,i)perylene	0.400	0.411	-2.7	70	-0.03

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

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