

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN100719\
 Data File : BN008050.D
 Acq On : 07 Oct 2019 15:11
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Oct 07 23:45:12 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 03 18:39:49 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	36	-0.03
2	1,4-Dioxane	0.400	0.441	-10.2	47	-0.04
3	n-Nitrosodimethylamine	0.400	0.000	100.0#	0	-3.01#
4 S	2-Fluorophenol	0.400	0.362	9.5	38	-0.02
5 S	Phenol-d6	0.400	0.369	7.8	37	0.00
6	bis(2-Chloroethyl)ether	0.400	0.469	-17.2	48	-0.04
7 I	Naphthalene-d8	0.400	0.400	0.0	37	-0.03
8 S	Nitrobenzene-d5	0.400	0.397	0.8	38	-0.03
9	Naphthalene	0.400	0.409	-2.2	38	-0.04
10	Hexachlorobutadiene	0.400	0.449	-12.2	42	-0.04
11 SURR	2-Methylnaphthalene-d10	0.400	0.423	-5.7	41	-0.03
12	2-Methylnaphthalene	0.400	0.423	-5.7	41	-0.03
13 I	Acenaphthene-d10	0.400	0.400	0.0	39	-0.03
14 S	2,4,6-Tribromophenol	0.400	0.398	0.5	46	-0.02
15 S	2-Fluorobiphenyl	0.400	0.364	9.0	38	-0.02
16	Acenaphthylene	0.400	0.375	6.3	40	-0.03
17	Acenaphthene	0.400	0.373	6.8	40	-0.03
18	Fluorene	0.400	0.373	6.8	39	-0.02
19 I	Phenanthrene-d10	0.400	0.400	0.0	38	-0.02
20	4-Bromophenyl-phenylether	0.400	0.386	3.5	37	-0.03
21	Hexachlorobenzene	0.400	0.419	-4.7	40	-0.03
22	Pentachlorophenol	0.400	0.462	-15.5	50	-0.02
23	Phenanthrene	0.400	0.412	-3.0	40	-0.03
24	Anthracene	0.400	0.407	-1.7	40	-0.02
25 SURR	Fluoranthene-d10	0.400	0.424	-6.0	41	-0.03
26	Fluoranthene	0.400	0.444	-11.0	43	-0.03
27 I	Chrysene-d12	0.400	0.400	0.0	53	-0.02
28	Pvrene	0.400	0.321	19.8	44	-0.03
29 S	Terphenyl-d14	0.400	0.360	10.0	50	-0.03
30	Benzo(a)anthracene	0.400	0.320	20.0	44	-0.02
31	Chrysene	0.400	0.356	11.0	49	-0.02
32	Bis(2-ethylhexyl)phthalate	0.400	0.716	-79.0#	106	-0.04
33	Indeno(1,2,3-cd)pyrene	0.400	0.365	8.8	50	-0.05
34 I	Perylene-d12	0.400	0.400	0.0	47	-0.04
35	Benzo(b)fluoranthene	0.400	0.403	-0.8	51	-0.03
36	Benzo(k)fluoranthene	0.400	0.391	2.3	48	-0.03
37 C	Benzo(a)pyrene	0.400	0.385	3.8	48	-0.03
38	Dibenzo(a,h)anthracene	0.400	0.404	-1.0	50	-0.05
39	Benzo(a,h,i)perylene	0.400	0.412	-3.0	51	-0.05

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

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