

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011520\
 Data File : BN009381.D
 Acq On : 15 Jan 2020 16:51
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 15 17:31:49 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN011420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 14 09:33:55 2020
 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	111	0.00
2	1,4-Dioxane	0.400	0.399	0.3	105	0.00
3	n-Nitrosodimethylamine	0.400	0.368	8.0	109	0.00
4 S	2-Fluorophenol	0.400	0.392	2.0	106	0.00
5 S	Phenol-d6	0.400	0.405	-1.3	110	0.00
6	bis(2-Chloroethyl)ether	0.400	0.375	6.3	113	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	110	0.00
8 S	Nitrobenzene-d5	0.400	0.383	4.3	111	-0.01
9	Naphthalene	0.400	0.415	-3.7	113	-0.01
10	Hexachlorobutadiene	0.400	0.406	-1.5	109	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.402	-0.5	110	0.00
12	2-Methylnaphthalene	0.400	0.408	-2.0	113	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	113	0.00
14 S	2,4,6-Tribromophenol	0.400	0.362	9.5	110	0.00
15 S	2-Fluorobiphenyl	0.400	0.406	-1.5	114	0.00
16	Acenaphthylene	0.400	0.378	5.5	112	0.00
17	Acenaphthene	0.400	0.404	-1.0	114	0.00
18	Fluorene	0.400	0.393	1.8	111	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	116	0.00
20	4-Bromophenyl-phenylether	0.400	0.386	3.5	115	0.00
21	Hexachlorobenzene	0.400	0.383	4.3	110	0.00
22	Pentachlorophenol	0.400	0.396	1.0	112	0.00
23	Phenanthrene	0.400	0.390	2.5	112	0.00
24	Anthracene	0.400	0.372	7.0	112	0.00
25 SURR	Fluoranthene-d10	0.400	0.382	4.5	113	0.00
26	Fluoranthene	0.400	0.382	4.5	113	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	112	0.00
28	Pvrene	0.400	0.409	-2.2	112	0.00
29 S	Terphenyl-d14	0.400	0.403	-0.8	112	0.00
30	Benzo(a)anthracene	0.400	0.377	5.8	109	0.00
31	Chrysene	0.400	0.421	-5.2	118	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.398	0.5	120	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.404	-1.0	110	0.00
34 I	Perylene-d12	0.400	0.400	0.0	114	0.00
35	Benzo(b)fluoranthene	0.400	0.387	3.3	116	0.00
36	Benzo(k)fluoranthene	0.400	0.392	2.0	107	0.00
37 C	Benzo(a)pyrene	0.400	0.404	-1.0	117	0.00
38	Dibenzo(a,h)anthracene	0.400	0.377	5.8	109	0.00
39	Benzo(a,h,i)perylene	0.400	0.384	4.0	109	0.00

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

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