

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012420\
 Data File : BN009427.D
 Acq On : 23 Jan 2020 17:16
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 24 12:16:36 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	93	0.00
2	1,4-Dioxane	0.400	0.394	1.5	87	0.00
3	n-Nitrosodimethylamine	0.400	0.335	16.3	83	0.00
4 S	2-Fluorophenol	0.400	0.408	-2.0	92	0.00
5 S	Phenol-d6	0.400	0.417	-4.2	95	0.00
6	bis(2-Chloroethyl)ether	0.400	0.433	-8.2	109	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	100	-0.01
8 S	Nitrobenzene-d5	0.400	0.394	1.5	104	-0.01
9	Naphthalene	0.400	0.389	2.8	97	-0.01
10	Hexachlorobutadiene	0.400	0.395	1.3	97	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.391	2.3	98	-0.01
12	2-Methylnaphthalene	0.400	0.394	1.5	100	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	107	0.00
14 S	2,4,6-Tribromophenol	0.400	0.391	2.3	112	-0.02
15 S	2-Fluorobiphenyl	0.400	0.410	-2.5	109	0.00
16	Acenaphthylene	0.400	0.383	4.3	107	0.00
17	Acenaphthene	0.400	0.393	1.8	104	0.00
18	Fluorene	0.400	0.380	5.0	101	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	109	-0.02
20	4-Bromophenyl-phenylether	0.400	0.380	5.0	105	0.00
21	Hexachlorobenzene	0.400	0.393	1.8	106	0.00
22	Pentachlorophenol	0.400	0.404	-1.0	108	0.00
23	Phenanthrene	0.400	0.378	5.5	102	0.00
24	Anthracene	0.400	0.377	5.8	106	-0.02
25 SURR	Fluoranthene-d10	0.400	0.364	9.0	101	-0.01
26	Fluoranthene	0.400	0.369	7.8	102	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	106	-0.01
28	Pvrene	0.400	0.397	0.8	103	0.00
29 S	Terphenyl-d14	0.400	0.398	0.5	105	0.00
30	Benzo(a)anthracene	0.400	0.370	7.5	102	0.00
31	Chrysene	0.400	0.397	0.8	105	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.405	-1.3	116	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.385	3.8	99	0.00
34 I	Perylene-d12	0.400	0.400	0.0	108	0.00
35	Benzo(b)fluoranthene	0.400	0.360	10.0	102	0.00
36	Benzo(k)fluoranthene	0.400	0.368	8.0	95	0.00
37 C	Benzo(a)pyrene	0.400	0.379	5.3	104	0.00
38	Dibenzo(a,h)anthracene	0.400	0.365	8.8	100	-0.02
39	Benzo(a,h,i)perylene	0.400	0.357	10.8	96	-0.01

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

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