

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012020\
 Data File : BN009405.D
 Acq On : 20 Jan 2020 16:38
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 20 17:51:16 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	95	0.00
2	1,4-Dioxane	0.400	0.404	-1.0	91	0.00
3	n-Nitrosodimethylamine	0.400	0.360	10.0	92	0.00
4 S	2-Fluorophenol	0.400	0.357	10.8	82	0.00
5 S	Phenol-d6	0.400	0.356	11.0	83	0.00
6	bis(2-Chloroethyl)ether	0.400	0.339	15.3	88	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	96	0.00
8 S	Nitrobenzene-d5	0.400	0.348	13.0	88	-0.01
9	Naphthalene	0.400	0.387	3.3	92	-0.01
10	Hexachlorobutadiene	0.400	0.398	0.5	93	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.383	4.3	92	0.00
12	2-Methylnaphthalene	0.400	0.386	3.5	94	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	102	0.00
14 S	2,4,6-Tribromophenol	0.400	0.338	15.5	93	0.00
15 S	2-Fluorobiphenyl	0.400	0.379	5.3	96	0.00
16	Acenaphthylene	0.400	0.349	12.8	93	0.00
17	Acenaphthene	0.400	0.377	5.8	96	0.00
18	Fluorene	0.400	0.388	3.0	99	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	105	0.00
20	4-Bromophenyl-phenylether	0.400	0.369	7.8	98	0.00
21	Hexachlorobenzene	0.400	0.368	8.0	95	0.00
22	Pentachlorophenol	0.400	0.378	5.5	94	0.00
23	Phenanthrene	0.400	0.369	7.8	96	0.00
24	Anthracene	0.400	0.350	12.5	95	0.00
25 SURR	Fluoranthene-d10	0.400	0.360	10.0	96	0.00
26	Fluoranthene	0.400	0.357	10.8	95	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	95	0.00
28	Pvrene	0.400	0.404	-1.0	95	0.00
29 S	Terphenyl-d14	0.400	0.389	2.8	92	0.00
30	Benzo(a)anthracene	0.400	0.347	13.3	86	0.00
31	Chrysene	0.400	0.405	-1.3	97	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.371	7.3	96	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.378	5.5	88	0.00
34 I	Perylene-d12	0.400	0.400	0.0	96	0.00
35	Benzo(b)fluoranthene	0.400	0.359	10.3	90	0.00
36	Benzo(k)fluoranthene	0.400	0.390	2.5	90	0.00
37 C	Benzo(a)pyrene	0.400	0.373	6.8	91	0.00
38	Dibenzo(a,h)anthracene	0.400	0.356	11.0	87	0.00
39	Benzo(a,h,i)perylene	0.400	0.374	6.5	90	0.00

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

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