

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031020\
 Data File : BN010191.D
 Acq On : 11 Mar 2020 14:09
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 11 15:21:27 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN021920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 19 15:57:43 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	90	0.00
2	1,4-Dioxane	0.400	0.449	-12.2	101	0.00
3	n-Nitrosodimethylamine	0.400	0.626	-56.5#	159	0.02
4 S	2-Fluorophenol	0.400	0.441	-10.2	102	0.02
5 S	Phenol-d6	0.400	0.425	-6.2	100	0.03
6	bis(2-Chloroethyl)ether	0.400	0.374	6.5	87	0.02
7 I	Naphthalene-d8	0.400	0.400	0.0	90	0.01
8 S	Nitrobenzene-d5	0.400	0.474	-18.5	113	0.04
9	Naphthalene	0.400	0.418	-4.5	97	0.01
10	Hexachlorobutadiene	0.400	0.450	-12.5	101	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.417	-4.2	97	0.01
12	2-Methylnaphthalene	0.400	0.408	-2.0	95	0.02
13 I	Acenaphthene-d10	0.400	0.400	0.0	96	0.00
14 S	2,4,6-Tribromophenol	0.400	0.436	-9.0	117	0.01
15 S	2-Fluorobiphenyl	0.400	0.369	7.8	92	0.01
16	Acenaphthylene	0.400	0.373	6.8	100	0.00
17	Acenaphthene	0.400	0.407	-1.7	102	0.00
18	Fluorene	0.400	0.403	-0.8	100	0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	100	0.01
20	4-Bromophenyl-phenylether	0.400	0.138	65.5#	54	0.00
21	Hexachlorobenzene	0.400	0.443	-10.7	116	0.01
22	Pentachlorophenol	0.400	0.423	-5.7	107	0.01
23	Phenanthrene	0.400	0.388	3.0	105	0.01
24	Anthracene	0.400	0.354	11.5	98	0.01
25 SURR	Fluoranthene-d10	0.400	0.358	10.5	96	0.00
26	Fluoranthene	0.400	0.353	11.8	95	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	90	0.00
28	Pvrene	0.400	0.408	-2.0	95	0.00
29 S	Terphenyl-d14	0.400	0.421	-5.2	97	0.01
30	Benzo(a)anthracene	0.400	0.382	4.5	97	0.00
31	Chrysene	0.400	0.398	0.5	92	0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.431	-7.7	108	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.408	-2.0	92	0.00
34 I	Perylene-d12	0.400	0.400	0.0	91	0.00
35	Benzo(b)fluoranthene	0.400	0.377	5.8	94	0.00
36	Benzo(k)fluoranthene	0.400	0.405	-1.3	97	0.00
37 C	Benzo(a)pyrene	0.400	0.388	3.0	91	0.00
38	Dibenzo(a,h)anthracene	0.400	0.377	5.8	91	0.01
39	Benzo(a,h,i)perylene	0.400	0.415	-3.7	100	0.00

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

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