

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN032420\
 Data File : BN010337.D
 Acq On : 24 Mar 2020 21:40
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 25 04:13:25 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN031920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 19 15:26:45 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	100	0.00
2	1,4-Dioxane	0.400	0.386	3.5	94	0.00
3	n-Nitrosodimethylamine	0.400	0.374	6.5	99	0.02
4 S	2-Fluorophenol	0.400	0.402	-0.5	101	0.00
5 S	Phenol-d6	0.400	0.405	-1.3	104	0.00
6	bis(2-Chloroethyl)ether	0.400	0.406	-1.5	100	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	100	0.00
8 S	Nitrobenzene-d5	0.400	0.387	3.3	101	0.00
9	Naphthalene	0.400	0.407	-1.7	101	-0.01
10	Hexachlorobutadiene	0.400	0.415	-3.7	104	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.447	-11.7	114	0.00
12	2-Methylnaphthalene	0.400	0.401	-0.3	102	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	97	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.354	11.5	96	0.00
15 S	2-Fluorobiphenyl	0.400	0.409	-2.2	98	-0.01
16	Acenaphthylene	0.400	0.376	6.0	98	-0.01
17	Acenaphthene	0.400	0.411	-2.7	101	0.00
18	Fluorene	0.400	0.399	0.3	100	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	97	0.00
20	4-Bromophenyl-phenylether	0.400	0.383	4.3	95	0.00
21	Hexachlorobenzene	0.400	0.429	-7.2	105	0.00
22	Pentachlorophenol	0.400	0.329	17.8	93	0.00
23	Phenanthrene	0.400	0.402	-0.5	99	-0.01
24	Anthracene	0.400	0.373	6.8	96	0.00
25 SURR	Fluoranthene-d10	0.400	0.435	-8.7	111	0.00
26	Fluoranthene	0.400	0.394	1.5	100	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	94	0.00
28	Pvrene	0.400	0.415	-3.7	98	0.00
29 S	Terphenyl-d14	0.400	0.411	-2.7	96	0.00
30	Benzo(a)anthracene	0.400	0.377	5.8	92	-0.01
31	Chrysene	0.400	0.406	-1.5	95	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.299	25.3#	78	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.410	-2.5	89	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	86	-0.01
35	Benzo(b)fluoranthene	0.400	0.404	-1.0	96	-0.01
36	Benzo(k)fluoranthene	0.400	0.422	-5.5	93	-0.01
37 C	Benzo(a)pyrene	0.400	0.391	2.3	93	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.410	-2.5	88	-0.02
39	Benzo(a,h,i)perylene	0.400	0.416	-4.0	95	-0.01

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN032420\
Data File : BN010337.D
Acq On : 24 Mar 2020 21:40
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_N
LabSampleId :
SSTDCCC0.4

Quant Time: Mar 25 04:13:25 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN031920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 19 15:26:45 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)

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