

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN101718\
 Data File : BN003196.D
 Acq On : 18 Oct 2018 00:43
 Operator : MJ/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 18 15:41:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN101718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 18 15:26:23 2018
 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	79	0.00
2	1,4-Dioxane	0.400	0.485	-21.2	105	0.00
3	n-Nitrosodimethylamine	0.400	0.276	31.0#	40	0.02
4 S	2-Fluorophenol	0.400	0.383	4.3	74	0.02
5 S	Phenol-d6	0.400	0.371	7.3	80	0.02
6	bis(2-Chloroethyl)ether	0.400	0.385	3.8	73	0.02
7 I	Naphthalene-d8	0.400	0.400	0.0	86	0.01
8 S	Nitrobenzene-d5	0.400	0.375	6.3	89	0.03
9	Naphthalene	0.400	0.427	-6.7	95	0.01
10	Hexachlorobutadiene	0.400	0.434	-8.5	97	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.428	-7.0	97	0.00
12	2-Methylnaphthalene	0.400	0.434	-8.5	98	0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	92	0.00
14 S	2,4,6-Tribromophenol	0.400	0.393	1.8	104	0.00
15 S	2-Fluorobiphenyl	0.400	0.398	0.5	96	0.00
16	Acenaphthylene	0.400	0.415	-3.7	100	0.01
17	Acenaphthene	0.400	0.424	-6.0	102	0.00
18	Fluorene	0.400	0.422	-5.5	101	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	94	0.00
20	4-Bromophenyl-phenylether	0.400	0.399	0.3	99	0.00
21	Hexachlorobenzene	0.400	0.421	-5.2	103	0.01
22	Pentachlorophenol	0.400	0.388	3.0	103	0.01
23	Phenanthrene	0.400	0.420	-5.0	103	0.00
24	Anthracene	0.400	0.414	-3.5	103	0.01
25 SURR	Fluoranthene-d10	0.400	0.422	-5.5	104	0.00
26	Fluoranthene	0.400	0.423	-5.7	104	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	90	0.00
28	Pvrene	0.400	0.431	-7.7	105	0.00
29 S	Terphenyl-d14	0.400	0.415	-3.7	102	0.00
30	Benzo(a)anthracene	0.400	0.392	2.0	102	0.00
31	Chrysene	0.400	0.437	-9.2	101	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.382	4.5	92	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.396	1.0	97	0.00
34 I	Perylene-d12	0.400	0.400	0.0	90	0.00
35	Benzo(b)fluoranthene	0.400	0.429	-7.2	106	0.00
36	Benzo(k)fluoranthene	0.400	0.429	-7.2	101	0.00
37 C	Benzo(a)pyrene	0.400	0.421	-5.2	101	0.00
38	Dibenzo(a,h)anthracene	0.400	0.401	-0.3	97	0.00
39	Benzo(a,h,i)perylene	0.400	0.403	-0.8	97	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN101718\
Data File : BN003196.D
Acq On : 18 Oct 2018 00:43
Operator : MJ/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_N
LabSampleId :
SSTDCCC0.4

Quant Time: Oct 18 15:41:21 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN101718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 18 15:26:23 2018
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			