

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012420\
 Data File : BN009433.D
 Acq On : 23 Jan 2020 20:52
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 24 12:17:29 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	112	0.00
2	1,4-Dioxane	0.400	0.399	0.3	107	0.00
3	n-Nitrosodimethylamine	0.400	0.334	16.5	101	-0.02
4 S	2-Fluorophenol	0.400	0.432	-8.0	118	0.00
5 S	Phenol-d6	0.400	0.443	-10.7	122	0.00
6	bis(2-Chloroethyl)ether	0.400	0.394	1.5	121	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	123	-0.01
8 S	Nitrobenzene-d5	0.400	0.403	-0.8	130	-0.01
9	Naphthalene	0.400	0.401	-0.3	122	-0.01
10	Hexachlorobutadiene	0.400	0.400	0.0	119	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.396	1.0	122	-0.01
12	2-Methylnaphthalene	0.400	0.407	-1.7	126	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	133	-0.02
14 S	2,4,6-Tribromophenol	0.400	0.423	-5.7	150	-0.02
15 S	2-Fluorobiphenyl	0.400	0.406	-1.5	134	0.00
16	Acenaphthylene	0.400	0.385	3.8	133	0.00
17	Acenaphthene	0.400	0.385	3.8	127	0.00
18	Fluorene	0.400	0.376	6.0	124	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	130	-0.02
20	4-Bromophenyl-phenylether	0.400	0.393	1.8	130	0.00
21	Hexachlorobenzene	0.400	0.416	-4.0	133	0.00
22	Pentachlorophenol	0.400	0.463	-15.8	158	-0.02
23	Phenanthrene	0.400	0.388	3.0	125	0.00
24	Anthracene	0.400	0.390	2.5	131	-0.02
25 SURR	Fluoranthene-d10	0.400	0.372	7.0	123	-0.01
26	Fluoranthene	0.400	0.379	5.3	125	-0.01
27 I	Chrysene-d12	0.400	0.400	0.0	130	-0.01
28	Pvrene	0.400	0.390	2.5	125	0.00
29 S	Terphenyl-d14	0.400	0.403	-0.8	130	-0.01
30	Benzo(a)anthracene	0.400	0.412	-3.0	139	0.00
31	Chrysene	0.400	0.385	3.8	125	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.551	-37.8#	194	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.369	7.8	117	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	131	-0.01
35	Benzo(b)fluoranthene	0.400	0.369	7.8	126	0.00
36	Benzo(k)fluoranthene	0.400	0.364	9.0	114	0.00
37 C	Benzo(a)pyrene	0.400	0.379	5.3	126	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.358	10.5	119	-0.02
39	Benzo(a,h,i)perylene	0.400	0.346	13.5	113	-0.02

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

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