

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091319\
 Data File : BN007660.D
 Acq On : 14 Sep 2019 12:09
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 16 01:39:04 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 12 14:27:51 2019
 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	80	0.00
2	1,4-Dioxane	0.400	0.379	5.3	79	0.00
3	n-Nitrosodimethylamine	0.400	0.560	-40.0#	79	0.00
4 S	2-Fluorophenol	0.400	0.354	11.5	77	0.00
5 S	Phenol-d6	0.400	0.377	5.8	78	0.00
6	bis(2-Chloroethyl)ether	0.400	0.372	7.0	74	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	80	-0.01
8 S	Nitrobenzene-d5	0.400	0.402	-0.5	77	-0.01
9	Naphthalene	0.400	0.425	-6.2	80	-0.01
10	Hexachlorobutadiene	0.400	0.415	-3.7	78	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.415	-3.7	79	0.00
12	2-Methylnaphthalene	0.400	0.422	-5.5	80	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	80	0.00
14 S	2,4,6-Tribromophenol	0.400	0.317	20.8	74	-0.01
15 S	2-Fluorobiphenyl	0.400	0.404	-1.0	81	-0.01
16	Acenaphthylene	0.400	0.366	8.5	75	-0.01
17	Acenaphthene	0.400	0.393	1.8	79	-0.01
18	Fluorene	0.400	0.387	3.3	78	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	78	-0.01
20	4-Bromophenyl-phenylether	0.400	0.422	-5.5	78	-0.01
21	Hexachlorobenzene	0.400	0.424	-6.0	77	-0.01
22	Pentachlorophenol	0.400	0.356	11.0	76	-0.01
23	Phenanthrene	0.400	0.433	-8.2	79	-0.01
24	Anthracene	0.400	0.402	-0.5	76	0.00
25 SURR	Fluoranthene-d10	0.400	0.407	-1.7	76	-0.01
26	Fluoranthene	0.400	0.415	-3.7	77	-0.01
27 I	Chrysene-d12	0.400	0.400	0.0	77	-0.01
28	Pvrene	0.400	0.423	-5.7	77	-0.01
29 S	Terphenyl-d14	0.400	0.419	-4.7	77	-0.01
30	Benzo(a)anthracene	0.400	0.396	1.0	73	-0.01
31	Chrysene	0.400	0.426	-6.5	79	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.341	14.8	70	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.458	-14.5	84	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	77	-0.01
35	Benzo(b)fluoranthene	0.400	0.408	-2.0	78	-0.01
36	Benzo(k)fluoranthene	0.400	0.431	-7.7	84	-0.01
37 C	Benzo(a)pyrene	0.400	0.403	-0.8	79	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.424	-6.0	82	-0.02
39	Benzo(a,h,i)perylene	0.400	0.410	-2.5	80	-0.02

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

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