

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091119\
 Data File : BN007586.D
 Acq On : 11 Sep 2019 13:47
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 11 16:20:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 11 14:14:20 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	104	0.00
2	1,4-Dioxane	0.400	0.464	-16.0	147	0.00
3	n-Nitrosodimethylamine	0.400	0.552	-38.0#	112	0.00
4 S	2-Fluorophenol	0.400	0.429	-7.2	126	0.00
5 S	Phenol-d6	0.400	0.411	-2.7	123	0.00
6	bis(2-Chloroethyl)ether	0.400	0.575	-43.7#	245	-0.03
7 I	Naphthalene-d8	0.400	0.400	0.0	100	0.00
8 S	Nitrobenzene-d5	0.400	0.589	-47.2#	168	0.00
9	Naphthalene	0.400	0.428	-7.0	111	0.00
10	Hexachlorobutadiene	0.400	0.450	-12.5	114	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.424	-6.0	111	0.00
12	2-Methylnaphthalene	0.400	0.423	-5.7	110	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	106	0.00
14 S	2,4,6-Tribromophenol	0.400	0.462	-15.5	152	-0.01
15 S	2-Fluorobiphenyl	0.400	0.401	-0.3	110	-0.01
16	Acenaphthylene	0.400	0.391	2.3	110	-0.01
17	Acenaphthene	0.400	0.389	2.8	108	-0.01
18	Fluorene	0.400	0.389	2.8	108	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	95	0.00
20	4-Bromophenyl-phenylether	0.400	0.442	-10.5	110	0.00
21	Hexachlorobenzene	0.400	0.447	-11.7	110	0.00
22	Pentachlorophenol	0.400	0.553	-38.3#	141	0.00
23	Phenanthrene	0.400	0.429	-7.2	106	-0.01
24	Anthracene	0.400	0.433	-8.2	109	0.00
25 SURR	Fluoranthene-d10	0.400	0.443	-10.7	111	0.00
26	Fluoranthene	0.400	0.438	-9.5	110	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	104	-0.01
28	Pvrene	0.400	0.420	-5.0	111	0.00
29 S	Terphenyl-d14	0.400	0.428	-7.0	113	0.00
30	Benzo(a)anthracene	0.400	0.435	-8.7	117	0.00
31	Chrysene	0.400	0.427	-6.7	114	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.593	-48.2#	182	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.480	-20.0	128	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	119	0.00
35	Benzo(b)fluoranthene	0.400	0.387	3.3	120	0.00
36	Benzo(k)fluoranthene	0.400	0.417	-4.2	131	0.00
37 C	Benzo(a)pyrene	0.400	0.398	0.5	126	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.416	-4.0	130	-0.01
39	Benzo(a,h,i)perylene	0.400	0.399	0.3	123	-0.01

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

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