

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091019\
 Data File : BN007579.D
 Acq On : 10 Sep 2019 16:50
 Operator : HP/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ICVBN091019

Quant Time: Sep 10 17:26:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 10 16:53:18 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	99	0.00
2	1,4-Dioxane	0.400	0.462	-15.5	140	0.00
3	n-Nitrosodimethylamine	0.400	0.509	-27.2#	99	0.00
4 S	2-Fluorophenol	0.400	0.360	10.0	101	0.00
5 S	Phenol-d6	0.400	0.350	12.5	100	0.00
6	bis(2-Chloroethyl)ether	0.400	0.425	-6.2	137	-0.02
7 I	Naphthalene-d8	0.400	0.400	0.0	99	0.00
8 S	Nitrobenzene-d5	0.400	0.409	-2.2	115	0.00
9	Naphthalene	0.400	0.392	2.0	100	0.00
10	Hexachlorobutadiene	0.400	0.394	1.5	99	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.387	3.3	100	0.00
12	2-Methylnaphthalene	0.400	0.387	3.3	100	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	99	0.00
14 S	2,4,6-Tribromophenol	0.400	0.330	17.5	101	-0.01
15 S	2-Fluorobiphenyl	0.400	0.396	1.0	101	0.00
16	Acenaphthylene	0.400	0.376	6.0	98	-0.01
17	Acenaphthene	0.400	0.384	4.0	99	-0.01
18	Fluorene	0.400	0.381	4.8	98	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	99	0.00
20	4-Bromophenyl-phenylether	0.400	0.384	4.0	100	0.00
21	Hexachlorobenzene	0.400	0.388	3.0	99	0.00
22	Pentachlorophenol	0.400	0.451	-12.7	104	0.00
23	Phenanthrene	0.400	0.384	4.0	99	0.00
24	Anthracene	0.400	0.368	8.0	97	0.00
25 SURR	Fluoranthene-d10	0.400	0.373	6.8	97	0.00
26	Fluoranthene	0.400	0.376	6.0	98	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	98	-0.01
28	Pvrene	0.400	0.389	2.8	97	0.00
29 S	Terphenyl-d14	0.400	0.389	2.8	98	0.00
30	Benzo(a)anthracene	0.400	0.366	8.5	94	0.00
31	Chrysene	0.400	0.392	2.0	100	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.322	19.5	94	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.380	5.0	97	0.00
34 I	Perylene-d12	0.400	0.400	0.0	97	0.00
35	Benzo(b)fluoranthene	0.400	0.384	4.0	97	0.00
36	Benzo(k)fluoranthene	0.400	0.388	3.0	99	0.00
37 C	Benzo(a)pyrene	0.400	0.378	5.5	97	0.00
38	Dibenzo(a,h)anthracene	0.400	0.383	4.3	98	0.00
39	Benzo(a,h,i)perylene	0.400	0.387	3.3	97	0.00

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

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