

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112219\
 Data File : BN008721.D
 Acq On : 23 Nov 2019 00:35
 Operator : JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ICVBN112219

Quant Time: Nov 25 10:58:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN112219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 25 10:49:40 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	0.394	1.5	103	0.00
3	n-Nitrosodimethylamine	0.400	0.389	2.8	109	0.00
4 S	2-Fluorophenol	0.400	0.396	1.0	103	0.00
5 S	Phenol-d6	0.400	0.393	1.8	102	0.00
6	bis(2-Chloroethyl)ether	0.400	0.402	-0.5	103	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	104	0.00
8 S	Nitrobenzene-d5	0.400	0.395	1.3	103	0.00
9	Naphthalene	0.400	0.406	-1.5	103	0.00
10	Hexachlorobutadiene	0.400	0.403	-0.8	102	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.404	-1.0	105	0.00
12	2-Methylnaphthalene	0.400	0.402	-0.5	103	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	103	0.00
14 S	2,4,6-Tribromophenol	0.400	0.371	7.3	100	0.00
15 S	2-Fluorobiphenyl	0.400	0.402	-0.5	103	0.00
16	Acenaphthylene	0.400	0.384	4.0	101	0.00
17	Acenaphthene	0.400	0.393	1.8	104	0.00
18	Fluorene	0.400	0.397	0.8	103	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	103	0.00
20	4-Bromophenyl-phenylether	0.400	0.396	1.0	103	0.00
21	Hexachlorobenzene	0.400	0.402	-0.5	104	0.00
22	Pentachlorophenol	0.400	0.322	19.5	97	0.00
23	Phenanthrene	0.400	0.396	1.0	103	-0.01
24	Anthracene	0.400	0.387	3.3	102	0.00
25 SURR	Fluoranthene-d10	0.400	0.397	0.8	103	0.00
26	Fluoranthene	0.400	0.395	1.3	103	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	108	0.00
28	Pvrene	0.400	0.383	4.3	102	0.00
29 S	Terphenyl-d14	0.400	0.387	3.3	103	0.00
30	Benzo(a)anthracene	0.400	0.381	4.8	104	-0.01
31	Chrysene	0.400	0.396	1.0	106	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.353	11.8	103	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.395	1.3	105	0.00
34 I	Perylene-d12	0.400	0.400	0.0	105	-0.01
35	Benzo(b)fluoranthene	0.400	0.395	1.3	104	0.00
36	Benzo(k)fluoranthene	0.400	0.404	-1.0	104	0.00
37 C	Benzo(a)pyrene	0.400	0.395	1.3	104	0.00
38	Dibenzo(a,h)anthracene	0.400	0.405	-1.3	106	-0.01
39	Benzo(a,h,i)perylene	0.400	0.402	-0.5	105	0.00

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

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