

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070219\
 Data File : BN006696.D
 Acq On : 02 Jul 2019 17:43
 Operator : HP/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ICVBN070219

Quant Time: Jul 05 05:48:55 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN070219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 05 05:30:28 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	92	0.00
2	1,4-Dioxane	0.400	0.409	-2.2	94	0.00
3	n-Nitrosodimethylamine	0.400	0.373	6.8	93	0.00
4 S	2-Fluorophenol	0.400	0.381	4.8	91	0.00
5 S	Phenol-d6	0.400	0.373	6.8	90	0.00
6	bis(2-Chloroethyl)ether	0.400	0.415	-3.7	107	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	95	0.01
8 S	Nitrobenzene-d5	0.400	0.364	9.0	91	0.00
9	Naphthalene	0.400	0.407	-1.7	93	0.00
10	Hexachlorobutadiene	0.400	0.432	-8.0	95	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.416	-4.0	96	0.00
12	2-Methylnaphthalene	0.400	0.374	6.5	90	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	95	0.00
14 S	2,4,6-Tribromophenol	0.400	0.342	14.5	87	0.01
15 S	2-Fluorobiphenyl	0.400	0.411	-2.7	95	0.00
16	Acenaphthylene	0.400	0.398	0.5	93	0.00
17	Acenaphthene	0.400	0.427	-6.7	96	0.00
18	Fluorene	0.400	0.416	-4.0	95	0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	95	0.01
20	4-Bromophenyl-phenylether	0.400	0.402	-0.5	95	0.00
21	Hexachlorobenzene	0.400	0.439	-9.7	98	0.01
22	Pentachlorophenol	0.400	0.391	2.3	91	0.00
23	Phenanthrene	0.400	0.408	-2.0	94	0.01
24	Anthracene	0.400	0.389	2.8	92	0.01
25 SURR	Fluoranthene-d10	0.400	0.418	-4.5	94	0.00
26	Fluoranthene	0.400	0.422	-5.5	94	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	95	0.01
28	Pvrene	0.400	0.412	-3.0	94	0.00
29 S	Terphenyl-d14	0.400	0.409	-2.2	94	0.00
30	Benzo(a)anthracene	0.400	0.390	2.5	86	0.00
31	Chrysene	0.400	0.430	-7.5	98	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.197	50.7#	74	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.434	-8.5	95	0.01
34 I	Perylene-d12	0.400	0.400	0.0	95	0.00
35	Benzo(b)fluoranthene	0.400	0.415	-3.7	93	0.00
36	Benzo(k)fluoranthene	0.400	0.442	-10.5	99	0.01
37 C	Benzo(a)pyrene	0.400	0.434	-8.5	94	0.00
38	Dibenzo(a,h)anthracene	0.400	0.428	-7.0	94	0.01
39	Benzo(a,h,i)perylene	0.400	0.433	-8.2	96	0.01

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

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