

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070919\
 Data File : BN006768.D
 Acq On : 09 Jul 2019 15:58
 Operator : HP/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ICVBN070919

Quant Time: Jul 09 17:36:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN070919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 09 17:31:09 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	98	0.00
2	1,4-Dioxane	0.400	0.371	7.3	92	0.00
3	n-Nitrosodimethylamine	0.400	0.457	-14.2	94	0.00
4 S	2-Fluorophenol	0.400	0.348	13.0	87	0.00
5 S	Phenol-d6	0.400	0.334	16.5	86	0.00
6	bis(2-Chloroethyl)ether	0.400	0.332	17.0	87	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	99	0.01
8 S	Nitrobenzene-d5	0.400	0.343	14.2	87	0.01
9	Naphthalene	0.400	0.384	4.0	88	0.00
10	Hexachlorobutadiene	0.400	0.388	3.0	87	0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.370	7.5	86	0.00
12	2-Methylnaphthalene	0.400	0.381	4.8	89	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	98	0.00
14 S	2,4,6-Tribromophenol	0.400	0.332	17.0	84	0.01
15 S	2-Fluorobiphenyl	0.400	0.380	5.0	88	0.00
16	Acenaphthylene	0.400	0.369	7.8	88	0.00
17	Acenaphthene	0.400	0.384	4.0	87	0.01
18	Fluorene	0.400	0.372	7.0	85	0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	97	0.01
20	4-Bromophenyl-phenylether	0.400	0.367	8.3	84	0.01
21	Hexachlorobenzene	0.400	0.386	3.5	85	0.01
22	Pentachlorophenol	0.400	0.456	-14.0	93	0.00
23	Phenanthrene	0.400	0.374	6.5	85	0.01
24	Anthracene	0.400	0.354	11.5	84	0.00
25 SURR	Fluoranthene-d10	0.400	0.367	8.3	83	0.00
26	Fluoranthene	0.400	0.378	5.5	85	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	93	0.02
28	Pvrene	0.400	0.396	1.0	84	0.00
29 S	Terphenyl-d14	0.400	0.393	1.8	85	0.00
30	Benzo(a)anthracene	0.400	0.362	9.5	81	0.02
31	Chrysene	0.400	0.404	-1.0	84	0.02
32	Bis(2-ethylhexyl)phthalate	0.400	0.332	17.0	81	0.02
33	Indeno(1,2,3-cd)pyrene	0.400	0.404	-1.0	82	0.05
34 I	Perylene-d12	0.400	0.400	0.0	92	0.04
35	Benzo(b)fluoranthene	0.400	0.384	4.0	86	0.03
36	Benzo(k)fluoranthene	0.400	0.406	-1.5	87	0.03
37 C	Benzo(a)pyrene	0.400	0.395	1.3	85	0.04
38	Dibenzo(a,h)anthracene	0.400	0.387	3.3	83	0.05
39	Benzo(a,h,i)perylene	0.400	0.392	2.0	82	0.05

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

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