

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN101718\
 Data File : BN003179.D
 Acq On : 17 Oct 2018 13:23
 Operator : MJ/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ICVBN101718

Quant Time: Oct 18 15:30:51 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN101718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 18 15:26:23 2018
 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	95	0.00
2	1,4-Dioxane	0.400	0.410	-2.5	109	0.00
3	n-Nitrosodimethylamine	0.400	0.333	16.8	70	0.02
4 S	2-Fluorophenol	0.400	0.397	0.8	92	0.00
5 S	Phenol-d6	0.400	0.363	9.3	94	0.00
6	bis(2-Chloroethyl)ether	0.400	0.359	10.3	82	0.02
7 I	Naphthalene-d8	0.400	0.400	0.0	97	0.00
8 S	Nitrobenzene-d5	0.400	0.365	8.8	98	0.01
9	Naphthalene	0.400	0.384	4.0	97	0.00
10	Hexachlorobutadiene	0.400	0.384	4.0	97	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.385	3.8	98	0.00
12	2-Methylnaphthalene	0.400	0.389	2.8	99	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	100	0.00
14 S	2,4,6-Tribromophenol	0.400	0.347	13.3	99	0.00
15 S	2-Fluorobiphenyl	0.400	0.376	6.0	99	0.00
16	Acenaphthylene	0.400	0.374	6.5	98	0.01
17	Acenaphthene	0.400	0.387	3.3	101	0.00
18	Fluorene	0.400	0.380	5.0	99	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	100	0.00
20	4-Bromophenyl-phenylether	0.400	0.374	6.5	99	0.00
21	Hexachlorobenzene	0.400	0.386	3.5	100	0.01
22	Pentachlorophenol	0.400	0.335	16.3	90	0.01
23	Phenanthrene	0.400	0.380	5.0	99	0.00
24	Anthracene	0.400	0.374	6.5	98	0.01
25 SURR	Fluoranthene-d10	0.400	0.376	6.0	99	0.00
26	Fluoranthene	0.400	0.378	5.5	99	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	98	0.00
28	Pvrene	0.400	0.375	6.3	99	0.00
29 S	Terphenyl-d14	0.400	0.366	8.5	99	0.00
30	Benzo(a)anthracene	0.400	0.338	15.5	96	0.00
31	Chrysene	0.400	0.391	2.3	99	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.357	10.8	93	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.352	12.0	94	0.00
34 I	Perylene-d12	0.400	0.400	0.0	97	0.00
35	Benzo(b)fluoranthene	0.400	0.379	5.3	100	0.00
36	Benzo(k)fluoranthene	0.400	0.378	5.5	96	0.00
37 C	Benzo(a)pyrene	0.400	0.371	7.3	96	0.00
38	Dibenzo(a,h)anthracene	0.400	0.360	10.0	94	0.00
39	Benzo(a,h,i)perylene	0.400	0.364	9.0	95	0.00

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

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