

Data Path : Z:\HPCHEM1\BNA_N\DATA\BN022818\
 Data File : BN000205.D
 Acq On : 28 Feb 2018 17:10
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Feb 28 17:40:46 2018
 Quant Method : Z:\HPCHEM1\BNA_N\METHODS\8270-SIM-BN022718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 15:07:57 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	98	0.00
2	1,4-Dioxane	0.400	0.437	-9.2	113	0.00
3	n-Nitrosodimethylamine	0.400	0.401	-0.3	101	0.00
4 S	2-Fluorophenol	0.400	0.412	-3.0	98	0.00
5 S	Phenol-d6	0.400	0.404	-1.0	100	0.00
6	bis(2-Chloroethyl)ether	0.400	0.432	-8.0	102	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	103	0.00
8 S	Nitrobenzene-d5	0.400	0.384	4.0	98	0.00
9	Naphthalene	0.400	0.403	-0.8	101	0.00
10	Hexachlorobutadiene	0.400	0.409	-2.2	102	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.401	-0.3	101	0.00
12	2-Methylnaphthalene	0.400	0.398	0.5	101	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	105	0.00
14 S	2,4,6-Tribromophenol	0.400	0.358	10.5	96	0.00
15 S	2-Fluorobiphenyl	0.400	0.394	1.5	101	0.00
16	Acenaphthylene	0.400	0.356	11.0	96	0.00
17	Acenaphthene	0.400	0.383	4.3	101	0.00
18	Fluorene	0.400	0.376	6.0	100	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	103	0.00
20	4-Bromophenyl-phenylether	0.400	0.378	5.5	98	0.00
21	Hexachlorobenzene	0.400	0.392	2.0	100	0.00
22	Pentachlorophenol	0.400	0.335	16.3	82	0.00
23	Phenanthrene	0.400	0.379	5.3	97	0.00
24	Anthracene	0.400	0.354	11.5	93	0.00
25 SURR	Fluoranthene-d10	0.400	0.359	10.3	93	0.00
26	Fluoranthene	0.400	0.358	10.5	93	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	91	0.00
28	Pyrene	0.400	0.410	-2.5	92	0.00
29 S	Terphenyl-d14	0.400	0.409	-2.2	91	0.00
30	Benzo(a)anthracene	0.400	0.353	11.8	80	0.01
31	Chrysene	0.400	0.396	1.0	87	0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.365	8.8	90	0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.390	2.5	85	0.02
34 I	Perylene-d12	0.400	0.400	0.0	90	0.02
35	Benzo(b)fluoranthene	0.400	0.382	4.5	87	0.01
36	Benzo(k)fluoranthene	0.400	0.363	9.3	81	0.02
37 C	Benzo(a)pyrene	0.400	0.378	5.5	87	0.01
38	Dibenzo(a,h)anthracene	0.400	0.372	7.0	85	0.02
39	Benzo(g,h,i)perylene	0.400	0.364	9.0	81	0.02

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

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