

Data Path : Z:\HPCHEM1\BNA_N\DATA\BN022718\
 Data File : BN000196.D
 Acq On : 27 Feb 2018 18:13
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Feb 27 18:43:25 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	68	0.00
2	1,4-Dioxane	0.400	0.406	-1.5	73	0.00
3	n-Nitrosodimethylamine	0.400	0.361	9.8	61	0.00
4 S	2-Fluorophenol	0.400	0.387	3.3	64	0.00
5 S	Phenol-d6	0.400	0.386	3.5	66	0.00
6	bis(2-Chloroethyl)ether	0.400	0.408	-2.0	67	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	69	0.00
8 S	Nitrobenzene-d5	0.400	0.372	7.0	64	0.00
9	Naphthalene	0.400	0.396	1.0	67	0.00
10	Hexachlorobutadiene	0.400	0.400	0.0	67	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.398	0.5	68	0.00
12	2-Methylnaphthalene	0.400	0.395	1.3	67	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	68	0.00
14 S	2,4,6-Tribromophenol	0.400	0.379	5.3	66	0.00
15 S	2-Fluorobiphenyl	0.400	0.407	-1.7	68	0.00
16	Acenaphthylene	0.400	0.363	9.3	63	0.00
17	Acenaphthene	0.400	0.398	0.5	67	0.00
18	Fluorene	0.400	0.394	1.5	68	0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	70	0.00
20	4-Bromophenyl-phenylether	0.400	0.386	3.5	68	0.00
21	Hexachlorobenzene	0.400	0.393	1.8	68	0.00
22	Pentachlorophenol	0.400	0.346	13.5	58	0.00
23	Phenanthrene	0.400	0.395	1.3	68	0.00
24	Anthracene	0.400	0.367	8.3	65	0.00
25 SURR	Fluoranthene-d10	0.400	0.383	4.3	68	0.00
26	Fluoranthene	0.400	0.378	5.5	67	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	67	0.00
28	Pyrene	0.400	0.401	-0.3	67	0.00
29 S	Terphenyl-d14	0.400	0.412	-3.0	68	0.00
30	Benzo(a)anthracene	0.400	0.374	6.5	63	0.00
31	Chrysene	0.400	0.404	-1.0	66	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.366	8.5	67	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.426	-6.5	68	0.00
34 I	Perylene-d12	0.400	0.400	0.0	67	0.00
35	Benzo(b)fluoranthene	0.400	0.397	0.8	68	0.00
36	Benzo(k)fluoranthene	0.400	0.402	-0.5	67	0.00
37 C	Benzo(a)pyrene	0.400	0.394	1.5	67	0.00
38	Dibenzo(a,h)anthracene	0.400	0.405	-1.3	69	0.00
39	Benzo(g,h,i)perylene	0.400	0.399	0.3	66	0.00

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

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