

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070219\
 Data File : BN006703.D
 Acq On : 02 Jul 2019 21:46
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 05 06:04:37 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	102	0.00
2	1,4-Dioxane	0.625	0.644	-3.0	105	0.00
3	n-Nitrosodimethylamine	0.471	0.391	17.0	91	0.00
4 S	2-Fluorophenol	1.086	1.001	7.8	97	0.00
5 S	Phenol-d6	1.375	1.279	7.0	99	0.00
6	bis(2-Chloroethyl)ether	1.162	1.219	-4.9	120	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
8 S	Nitrobenzene-d5	0.307	0.273	11.1	98	0.00
9	Naphthalene	1.094	1.114	-1.8	103	0.00
10	Hexachlorobutadiene	0.222	0.240	-8.1	104	0.00
11 SURR	2-Methylnaphthalene-d10	0.601	0.611	-1.7	103	0.00
12	2-Methylnaphthalene	0.812	0.709	12.7	97	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
14 S	2,4,6-Tribromophenol	0.197	0.162	17.8	89	0.00
15 S	2-Fluorobiphenyl	1.565	1.625	-3.8	102	0.00
16	Acenaphthylene	1.994	2.002	-0.4	100	0.00
17	Acenaphthene	1.271	1.346	-5.9	101	-0.01
18	Fluorene	1.578	1.612	-2.2	99	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
20	4-Bromophenyl-phenylether	0.236	0.230	2.5	95	0.00
21	Hexachlorobenzene	0.280	0.304	-8.6	100	0.00
22	Pentachlorophenol	0.040	0.029	27.5#	101	-0.02
23	Phenanthrene	1.153	1.168	-1.3	97	0.00
24	Anthracene	1.042	1.021	2.0	95	0.00
25 SURR	Fluoranthene-d10	1.144	1.186	-3.7	96	0.00
26	Fluoranthene	1.354	1.403	-3.6	96	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	97	-0.01
28	Pvrene	1.601	1.650	-3.1	95	0.00
29 S	Terphenyl-d14	0.921	0.962	-4.5	97	0.00
30	Benzo(a)anthracene	1.380	1.378	0.1	90	-0.01
31	Chrysene	1.422	1.542	-8.4	101	-0.01
32	Bis(2-ethylhexyl)phthalate	4.115	1.175	71.4#	44#	-0.01
33	Indeno(1,2,3-cd)pyrene	1.701	1.913	-12.5	100	-0.03
34 I	Perylene-d12	1.000	1.000	0.0	98	-0.03
35	Benzo(b)fluoranthene	1.284	1.384	-7.8	99	-0.02
36	Benzo(k)fluoranthene	1.392	1.511	-8.5	100	-0.02
37 C	Benzo(a)pyrene	1.244	1.360	-9.3	98	-0.02
38	Dibenzo(a,h)anthracene	1.236	1.349	-9.1	100	-0.03
39	Benzo(a,h,i)perylene	1.272	1.392	-9.4	100	-0.03

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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