

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN100719\
 Data File : BNCCC056.D
 Acq On : 07 Oct 2019 19:27
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Oct 07 23:21:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 03 18:39:49 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	40#	-0.03
2	1,4-Dioxane	0.446	0.406	9.0	43#	-0.04
3	n-Nitrosodimethylamine	0.204	0.379	-85.8#	49#	-0.05
4 S	2-Fluorophenol	1.377	1.247	9.4	42#	-0.02
5 S	Phenol-d6	1.962	1.600	18.5	41#	0.00
6	bis(2-Chloroethyl)ether	1.133	1.190	-5.0	48#	-0.04
7 I	Naphthalene-d8	1.000	1.000	0.0	42#	-0.04
8 S	Nitrobenzene-d5	0.350	0.334	4.6	42#	-0.03
9	Naphthalene	1.123	1.156	-2.9	44#	-0.04
10	Hexachlorobutadiene	0.236	0.264	-11.9	47#	-0.05
11 SURR	2-Methylnaphthalene-d10	0.674	0.699	-3.7	45#	-0.03
12	2-Methylnaphthalene	0.759	0.799	-5.3	46#	-0.04
13 I	Acenaphthene-d10	1.000	1.000	0.0	42#	-0.03
14 S	2,4,6-Tribromophenol	0.218	0.206	5.5	47#	-0.03
15 S	2-Fluorobiphenyl	1.647	1.526	7.3	42#	-0.03
16	Acenaphthylene	2.121	2.006	5.4	44#	-0.04
17	Acenaphthene	1.368	1.252	8.5	42#	-0.04
18	Fluorene	1.753	1.623	7.4	42#	-0.03
19 I	Phenanthrene-d10	1.000	1.000	0.0	39#	-0.03
20	4-Bromophenyl-phenylether	0.243	0.232	4.5	38#	-0.04
21	Hexachlorobenzene	0.265	0.283	-6.8	42#	-0.03
22	Pentachlorophenol	0.098	0.111	-13.3	51	-0.03
23	Phenanthrene	1.246	1.246	0.0	40#	-0.03
24	Anthracene	1.125	1.160	-3.1	42#	-0.03
25 SURR	Fluoranthene-d10	1.348	1.437	-6.6	42#	-0.03
26	Fluoranthene	1.549	1.694	-9.4	43#	-0.03
27 I	Chrysene-d12	1.000	1.000	0.0	52	-0.03
28	Pvrene	1.364	1.152	15.5	45#	-0.03
29 S	Terphenyl-d14	0.983	0.857	12.8	47#	-0.03
30	Benzo(a)anthracene	1.473	1.256	14.7	46#	-0.02
31	Chrysene	1.436	1.327	7.6	49#	-0.02
32	Bis(2-ethylhexyl)phthalate	0.758	1.231	-62.4#	94	-0.04
33	Indeno(1,2,3-cd)pyrene	1.797	1.598	11.1	47#	-0.05
34 I	Perylene-d12	1.000	1.000	0.0	45#	-0.04
35	Benzo(b)fluoranthene	1.389	1.448	-4.2	50	-0.03
36	Benzo(k)fluoranthene	1.373	1.371	0.1	47#	-0.04
37 C	Benzo(a)pyrene	1.305	1.268	2.8	46#	-0.04
38	Dibenzo(a,h)anthracene	1.335	1.317	1.3	47#	-0.05
39	Benzo(a,h,i)perylene	1.321	1.337	-1.2	48#	-0.05

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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