

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN032020\
 Data File : BN010313.D
 Acq On : 20 Mar 2020 21:31
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 20 23:01:07 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN031920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 19 15:26:45 2020
 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	99	0.00
2	1,4-Dioxane	0.376	0.369	1.9	94	0.00
3	n-Nitrosodimethylamine	0.317	0.283	10.7	93	0.02
4 S	2-Fluorophenol	0.792	0.772	2.5	97	0.00
5 S	Phenol-d6	0.985	0.977	0.8	100	0.00
6	bis(2-Chloroethyl)ether	0.946	0.949	-0.3	98	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
8 S	Nitrobenzene-d5	0.256	0.245	4.3	98	0.00
9	Naphthalene	0.997	1.000	-0.3	97	0.00
10	Hexachlorobutadiene	0.242	0.251	-3.7	102	0.00
11 SURR	2-Methylnaphthalene-d10	0.639	0.619	3.1	96	0.00
12	2-Methylnaphthalene	0.692	0.681	1.6	98	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	96	-0.01
14 S	2,4,6-Tribromophenol	0.154	0.136	11.7	95	0.00
15 S	2-Fluorobiphenyl	1.531	1.578	-3.1	98	0.00
16	Acenaphthylene	1.569	1.452	7.5	96	0.00
17	Acenaphthene	1.118	1.076	3.8	94	0.00
18	Fluorene	1.485	1.453	2.2	97	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	97	0.00
20	4-Bromophenyl-phenylether	0.243	0.237	2.5	96	0.00
21	Hexachlorobenzene	0.249	0.255	-2.4	99	0.00
22	Pentachlorophenol	0.085	0.072	15.3	95	0.00
23	Phenanthrene	1.099	1.086	1.2	97	0.00
24	Anthracene	0.940	0.869	7.6	95	0.00
25 SURR	Fluoranthene-d10	1.206	1.140	5.5	96	0.00
26	Fluoranthene	1.347	1.309	2.8	98	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	97	0.00
28	Pvrene	1.245	1.223	1.8	96	0.00
29 S	Terphenyl-d14	0.864	0.870	-0.7	98	0.00
30	Benzo(a)anthracene	1.274	1.187	6.8	94	-0.01
31	Chrysene	1.366	1.387	-1.5	98	-0.01
32	Bis(2-ethylhexyl)phthalate	0.430	0.311	27.7#	78	0.00
33	Indeno(1,2,3-cd)pyrene	1.463	1.503	-2.7	92	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	93	0.00
35	Benzo(b)fluoranthene	1.359	1.285	5.4	97	0.00
36	Benzo(k)fluoranthene	1.383	1.376	0.5	95	0.00
37 C	Benzo(a)pyrene	1.147	1.065	7.1	95	0.00
38	Dibenzo(a,h)anthracene	1.240	1.232	0.6	92	-0.01
39	Benzo(a,h,i)perylene	1.248	1.210	3.0	95	-0.01

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