

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081319\
 Data File : BN007147.D
 Acq On : 14 Aug 2019 00:20
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 14 05:00:06 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN081019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 13 12:12:02 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	153	0.00
2	n-Nitrosodimethylamine	0.400	0.346	13.5	169	0.00
3 S	2-Fluorophenol	0.400	0.417	-4.2	167	0.00
4 S	Phenol-d6	0.400	0.429	-7.2	177	0.00
5	bis(2-Chloroethyl)ether	0.400	0.390	2.5	156	0.00
6 I	Naphthalene-d8	0.400	0.400	0.0	164	0.00
7 S	Nitrobenzene-d5	0.400	0.398	0.5	178	0.00
8	Naphthalene	0.400	0.374	6.5	164	0.00
9	Hexachlorobutadiene	0.400	0.380	5.0	164	-0.01
10 SURR	2-Methylnaphthalene-d10	0.400	0.366	8.5	162	0.00
11	2-Methylnaphthalene	0.400	0.371	7.3	164	0.00
12 I	Acenaphthene-d10	0.400	0.400	0.0	159	0.00
13 S	2,4,6-Tribromophenol	0.400	0.319	20.3	174	0.00
14 S	2-Fluorobiphenyl	0.400	0.212	47.0#	76	0.00
15	Acenaphthylene	0.400	0.462	-15.5	195	0.00
16	Acenaphthene	0.400	0.374	6.5	159	-0.01
17	Fluorene	0.400	0.343	14.2	164	0.00
18 I	Phenanthrene-d10	0.400	0.400	0.0	141	0.00
19	4-Bromophenyl-phenylether	0.400	0.396	1.0	141	-0.01
20	Hexachlorobenzene	0.400	0.405	-1.3	144	0.00
21	Pentachlorophenol	0.400	0.357	10.8	145	0.00
22	Phenanthrene	0.400	0.388	3.0	141	0.00
23	Anthracene	0.400	0.392	2.0	148	-0.01
24 SURR	Fluoranthene-d10	0.400	0.355	11.3	131	0.00
25	Fluoranthene	0.400	0.364	9.0	133	0.00
26 I	Chrysene-d12	0.400	0.400	0.0	127	-0.01
27	Pyrene	0.400	0.418	-4.5	137	0.00
28 S	Terphenyl-d14	0.400	0.407	-1.7	134	-0.01
29	Benzo(a)anthracene	0.400	0.409	-2.2	133	-0.01
30	Chrysene	0.400	0.400	0.0	132	-0.01
31	Indeno(1,2,3-cd)pyrene	0.400	0.436	-9.0	147	-0.02
32 I	Perylene-d12	0.400	0.400	0.0	144	-0.01
33	Benzo(b)fluoranthene	0.400	0.380	5.0	139	-0.01
34	Benzo(k)fluoranthene	0.400	0.376	6.0	136	-0.01
35 C	Benzo(a)pyrene	0.400	0.389	2.8	146	-0.02
36	Dibenzo(a,h)anthracene	0.400	0.395	1.3	148	-0.02
37	Benzo(g,h,i)perylene	0.400	0.383	4.3	140	-0.02

(#) = Out of Range

SPCC's out = 0 CCC's out = 0