

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081319\
 Data File : BN007149.D
 Acq On : 14 Aug 2019 02:44
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 14 05:00:17 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	136	0.00
2	n-Nitrosodimethylamine	0.400	0.354	11.5	153	0.00
3 S	2-Fluorophenol	0.400	0.422	-5.5	149	0.00
4 S	Phenol-d6	0.400	0.425	-6.2	155	0.00
5	bis(2-Chloroethyl)ether	0.400	0.399	0.3	141	0.00
6 I	Naphthalene-d8	0.400	0.400	0.0	144	0.00
7 S	Nitrobenzene-d5	0.400	0.399	0.3	157	0.00
8	Naphthalene	0.400	0.375	6.3	144	0.00
9	Hexachlorobutadiene	0.400	0.380	5.0	144	-0.01
10 SURR	2-Methylnaphthalene-d10	0.400	0.369	7.8	143	0.00
11	2-Methylnaphthalene	0.400	0.373	6.8	145	0.00
12 I	Acenaphthene-d10	0.400	0.400	0.0	138	0.00
13 S	2,4,6-Tribromophenol	0.400	0.323	19.3	153	0.00
14 S	2-Fluorobiphenyl	0.400	0.196	51.0#	61	0.00
15	Acenaphthylene	0.400	0.439	-9.7	161	0.00
16	Acenaphthene	0.400	0.381	4.8	141	-0.01
17	Fluorene	0.400	0.341	14.8	142	-0.01
18 I	Phenanthrene-d10	0.400	0.400	0.0	124	0.00
19	4-Bromophenyl-phenylether	0.400	0.398	0.5	125	-0.01
20	Hexachlorobenzene	0.400	0.406	-1.5	128	0.00
21	Pentachlorophenol	0.400	0.362	9.5	130	0.00
22	Phenanthrene	0.400	0.389	2.8	126	0.00
23	Anthracene	0.400	0.391	2.3	131	-0.01
24 SURR	Fluoranthene-d10	0.400	0.354	11.5	115	0.00
25	Fluoranthene	0.400	0.365	8.8	118	0.00
26 I	Chrysene-d12	0.400	0.400	0.0	116	-0.02
27	Pyrene	0.400	0.406	-1.5	121	-0.01
28 S	Terphenyl-d14	0.400	0.393	1.8	118	0.00
29	Benzo(a)anthracene	0.400	0.389	2.8	115	-0.01
30	Chrysene	0.400	0.400	0.0	120	-0.01
31	Indeno(1,2,3-cd)pyrene	0.400	0.429	-7.2	132	-0.02
32 I	Perylene-d12	0.400	0.400	0.0	129	-0.02
33	Benzo(b)fluoranthene	0.400	0.377	5.8	124	-0.02
34	Benzo(k)fluoranthene	0.400	0.378	5.5	122	-0.02
35 C	Benzo(a)pyrene	0.400	0.389	2.8	130	-0.02
36	Dibenzo(a,h)anthracene	0.400	0.394	1.5	132	-0.03
37	Benzo(g,h,i)perylene	0.400	0.383	4.3	125	-0.02

(#) = Out of Range

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