

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN082119\
 Data File : BN007295.D
 Acq On : 21 Aug 2019 18:10
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDC00.4

Quant Time: Aug 21 18:42:39 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 20 18:52:43 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	113	0.00
2	1,4-Dioxane	0.568	0.541	4.8	107	0.00
3	n-Nitrosodimethylamine	0.380	0.332	12.6	102	0.00
4 S	2-Fluorophenol	1.219	1.119	8.2	112	0.00
5 S	Phenol-d6	1.567	1.358	13.3	114	0.00
6	bis(2-Chloroethyl)ether	1.181	1.206	-2.1	112	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	117	0.00
8 S	Nitrobenzene-d5	0.307	0.295	3.9	111	0.00
9	Naphthalene	1.134	1.150	-1.4	114	0.00
10	Hexachlorobutadiene	0.250	0.252	-0.8	112	0.00
11 SURR	2-Methylnaphthalene-d10	0.691	0.692	-0.1	115	0.00
12	2-Methylnaphthalene	0.785	0.796	-1.4	117	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	114	0.00
14 S	2,4,6-Tribromophenol	0.225	0.188	16.4	103	0.00
15 S	2-Fluorobiphenyl	1.715	1.786	-4.1	114	0.00
16	Acenaphthylene	2.121	2.143	-1.0	117	0.00
17	Acenaphthene	1.352	1.395	-3.2	115	0.00
18	Fluorene	1.766	1.739	1.5	111	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
20	4-Bromophenyl-phenylether	0.229	0.238	-3.9	137	0.00
21	Hexachlorobenzene	0.265	0.281	-6.0	114	0.00
22	Pentachlorophenol	0.106	0.090	15.1	103	0.00
23	Phenanthrene	1.254	1.273	-1.5	110	0.00
24	Anthracene	1.129	1.116	1.2	110	0.00
25 SURR	Fluoranthene-d10	1.268	1.255	1.0	109	0.00
26	Fluoranthene	1.519	1.512	0.5	110	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	109	0.00
28	Pvrene	1.496	1.590	-6.3	112	0.00
29 S	Terphenyl-d14	1.001	1.021	-2.0	108	0.00
30	Benzo(a)anthracene	1.487	1.508	-1.4	110	0.00
31	Chrysene	1.518	1.601	-5.5	112	0.00
32	Bis(2-ethylhexyl)phthalate	0.594	0.561	5.6	116	0.00
33	Indeno(1,2,3-cd)pyrene	1.826	2.081	-14.0	120	0.00
34 I	Perylene-d12	1.000	1.000	0.0	121	0.00
35	Benzo(b)fluoranthene	1.491	1.398	6.2	112	0.00
36	Benzo(k)fluoranthene	1.461	1.503	-2.9	120	0.00
37 C	Benzo(a)pyrene	1.344	1.327	1.3	120	0.00
38	Dibenzo(a,h)anthracene	1.412	1.423	-0.8	119	0.00
39	Benzo(a,h,i)perylene	1.427	1.393	2.4	115	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN082119\
Data File : BN007295.D
Acq On : 21 Aug 2019 18:10
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_N
LabSampleId :
SSTDCCC0.4

Quant Time: Aug 21 18:42:39 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 20 18:52:43 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	AvgRF	CCRF	%Dev Area	% Dev(min)

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