

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082619\
 Data File : BN007410.D
 Acq On : 26 Aug 2019 15:25
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 26 16:00:42 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 24 00:38:17 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	111	0.00
2	1,4-Dioxane	0.665	0.680	-2.3	109	0.00
3	n-Nitrosodimethylamine	0.309	0.269	12.9	93	0.00
4 S	2-Fluorophenol	1.490	1.207	19.0	95	0.00
5 S	Phenol-d6	1.953	1.439	26.3#	94	0.00
6	bis(2-Chloroethyl)ether	1.445	1.368	5.3	106	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	112	-0.01
8 S	Nitrobenzene-d5	0.348	0.320	8.0	108	0.00
9	Naphthalene	1.143	1.093	4.4	105	0.00
10	Hexachlorobutadiene	0.234	0.217	7.3	100	0.00
11 SURR	2-Methylnaphthalene-d10	0.665	0.640	3.8	107	0.00
12	2-Methylnaphthalene	0.779	0.745	4.4	104	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	119	0.00
14 S	2,4,6-Tribromophenol	0.209	0.218	-4.3	126	0.00
15 S	2-Fluorobiphenyl	1.613	1.472	8.7	106	0.00
16	Acenaphthylene	2.367	2.070	12.5	103	0.00
17	Acenaphthene	1.385	1.287	7.1	108	0.00
18	Fluorene	1.751	1.712	2.2	113	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	126	0.00
20	4-Bromophenyl-phenylether	0.158	0.137	13.3	126	0.00
21	Hexachlorobenzene	0.268	0.260	3.0	121	0.00
22	Pentachlorophenol	0.083	0.080	3.6	129	0.00
23	Phenanthrene	1.203	1.169	2.8	122	0.00
24	Anthracene	1.210	1.087	10.2	112	0.00
25 SURR	Fluoranthene-d10	1.288	1.232	4.3	117	0.00
26	Fluoranthene	1.545	1.510	2.3	120	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	130	0.00
28	Pvrene	1.609	1.531	4.8	120	0.00
29 S	Terphenyl-d14	1.047	1.016	3.0	122	0.00
30	Benzo(a)anthracene	1.571	1.480	5.8	120	0.00
31	Chrysene	1.516	1.535	-1.3	126	0.00
32	Bis(2-ethylhexyl)phthalate	1.177	0.920	21.8	107	0.00
33	Indeno(1,2,3-cd)pyrene	1.827	1.921	-5.1	131	0.00
34 I	Perylene-d12	1.000	1.000	0.0	134	0.00
35	Benzo(b)fluoranthene	1.449	1.371	5.4	124	0.00
36	Benzo(k)fluoranthene	1.432	1.471	-2.7	132	0.00
37 C	Benzo(a)pyrene	1.379	1.338	3.0	126	0.00
38	Dibenzo(a,h)anthracene	1.376	1.377	-0.1	130	0.00
39	Benzo(a,h,i)perylene	1.371	1.413	-3.1	134	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082619\
Data File : BN007410.D
Acq On : 26 Aug 2019 15:25
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_N
LabSampleId :
SSTDCCC0.4

Quant Time: Aug 26 16:00:42 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Aug 24 00:38:17 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			