

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012120\
 Data File : BN009407.D
 Acq On : 21 Jan 2020 10:29
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 21 14:29:39 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN011420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 14 09:33:55 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	144	0.00
2	1,4-Dioxane	0.417	0.420	-0.7	138	0.00
3	n-Nitrosodimethylamine	0.644	0.583	9.5	140	-0.02
4 S	2-Fluorophenol	0.921	0.868	5.8	132	0.00
5 S	Phenol-d6	1.065	1.037	2.6	138	0.00
6	bis(2-Chloroethyl)ether	1.097	1.043	4.9	149	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	138	0.00
8 S	Nitrobenzene-d5	0.319	0.302	5.3	138	-0.01
9	Naphthalene	1.082	1.040	3.9	132	-0.01
10	Hexachlorobutadiene	0.237	0.237	0.0	135	-0.01
11 SURR	2-Methylnaphthalene-d10	0.773	0.743	3.9	133	0.00
12	2-Methylnaphthalene	0.754	0.720	4.5	133	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	145	0.00
14 S	2,4,6-Tribromophenol	0.151	0.135	10.6	138	0.00
15 S	2-Fluorobiphenyl	1.497	1.445	3.5	139	0.00
16	Acenaphthylene	1.725	1.582	8.3	139	0.00
17	Acenaphthene	1.213	1.160	4.4	138	0.00
18	Fluorene	1.581	1.513	4.3	138	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	153#	0.00
20	4-Bromophenyl-phenylether	0.241	0.217	10.0	141	0.00
21	Hexachlorobenzene	0.256	0.233	9.0	137	0.00
22	Pentachlorophenol	0.072	0.059	18.1	155#	0.00
23	Phenanthrene	1.202	1.114	7.3	140	0.00
24	Anthracene	1.009	0.902	10.6	141	0.00
25 SURR	Fluoranthene-d10	1.372	1.263	7.9	144	0.00
26	Fluoranthene	1.410	1.292	8.4	143	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	151#	-0.01
28	Pvrene	1.516	1.445	4.7	142	0.00
29 S	Terphenyl-d14	0.948	0.881	7.1	140	0.00
30	Benzo(a)anthracene	1.315	1.255	4.6	150	0.00
31	Chrysene	1.504	1.420	5.6	143	0.00
32	Bis(2-ethylhexyl)phthalate	0.581	0.518	10.8	146	0.00
33	Indeno(1,2,3-cd)pyrene	1.564	1.550	0.9	146	0.00
34 I	Perylene-d12	1.000	1.000	0.0	156#	0.00
35	Benzo(b)fluoranthene	1.471	1.322	10.1	146	0.00
36	Benzo(k)fluoranthene	1.624	1.569	3.4	144	0.00
37 C	Benzo(a)pyrene	1.295	1.217	6.0	148	0.00
38	Dibenzo(a,h)anthracene	1.303	1.196	8.2	145	0.00
39	Benzo(a,h,i)perylene	1.434	1.331	7.2	144	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012120\
Data File : BN009407.D
Acq On : 21 Jan 2020 10:29
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_N
LabSampleId :
SSTDCCC0.4

Quant Time: Jan 21 14:29:39 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN011420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 14 09:33:55 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	AvgRF	CCRF	%Dev Area	% Dev(min)

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