

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070119\
 Data File : BN006652.D
 Acq On : 01 Jul 2019 08:39
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 02 05:03:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN062819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 28 13:47:56 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	108	0.00
2	1,4-Dioxane	0.554	0.586	-5.8	122	0.00
3	n-Nitrosodimethylamine	0.435	0.411	5.5	115	0.00
4 S	2-Fluorophenol	0.981	1.006	-2.5	115	0.00
5 S	Phenol-d6	1.299	1.282	1.3	112	0.00
6	bis(2-Chloroethyl)ether	1.158	1.166	-0.7	104	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
8 S	Nitrobenzene-d5	0.269	0.273	-1.5	106	0.00
9	Naphthalene	1.051	1.098	-4.5	99	-0.01
10	Hexachlorobutadiene	0.216	0.237	-9.7	104	-0.01
11 SURR	2-Methylnaphthalene-d10	0.630	0.618	1.9	93	0.00
12	2-Methylnaphthalene	0.726	0.724	0.3	94	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	85	-0.01
14 S	2,4,6-Tribromophenol	0.188	0.178	5.3	90	0.00
15 S	2-Fluorobiphenyl	1.482	1.563	-5.5	90	-0.01
16	Acenaphthylene	1.907	1.943	-1.9	88	0.00
17	Acenaphthene	1.256	1.329	-5.8	88	0.00
18	Fluorene	1.585	1.597	-0.8	84	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	80	0.00
20	4-Bromophenyl-phenylether	0.222	0.235	-5.9	85	-0.01
21	Hexachlorobenzene	0.280	0.305	-8.9	84	0.00
22	Pentachlorophenol	0.045	0.038	15.6	86	0.00
23	Phenanthrene	1.137	1.174	-3.3	82	0.00
24	Anthracene	0.987	0.997	-1.0	83	0.00
25 SURR	Fluoranthene-d10	1.175	1.205	-2.6	81	0.00
26	Fluoranthene	1.382	1.443	-4.4	81	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	74	0.00
28	Pvrene	1.400	1.571	-12.2	81	0.00
29 S	Terphenyl-d14	0.838	0.910	-8.6	80	0.00
30	Benzo(a)anthracene	1.308	1.385	-5.9	80	-0.01
31	Chrysene	1.410	1.512	-7.2	75	-0.01
32	Bis(2-ethylhexyl)phthalate	0.658	0.771	-17.2	97	0.00
33	Indeno(1,2,3-cd)pyrene	1.427	1.902	-33.3#	85	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	76	-0.01
35	Benzo(b)fluoranthene	1.361	1.373	-0.9	76	-0.01
36	Benzo(k)fluoranthene	1.450	1.457	-0.5	75	-0.01
37 C	Benzo(a)pyrene	1.272	1.362	-7.1	78	-0.01
38	Dibenzo(a,h)anthracene	1.111	1.317	-18.5	84	-0.01
39	Benzo(a,h,i)perylene	1.109	1.353	-22.0	87	-0.01

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070119\
Data File : BN006652.D
Acq On : 01 Jul 2019 08:39
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_N
LabSampleId :
SSTDCCC0.4

Quant Time: Jul 02 05:03:21 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN062819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 28 13:47:56 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			