

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA N\Data\BN030118\
 Data File : BN000216.D
 Acq On : 01 Mar 2018 14:28
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 01 17:46:19 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-SIM-BN022718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 15:07:57 2018
 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	66	0.00
2	1,4-Dioxane	0.414	0.440	-6.3	75	0.00
3	n-Nitrosodimethylamine	0.282	0.225	20.2	55	0.00
4 S	2-Fluorophenol	0.845	0.897	-6.2	69	0.00
5 S	Phenol-d6	1.073	1.129	-5.2	70	0.00
6	bis(2-Chloroethyl)ether	1.363	1.404	-3.0	66	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	67	0.00
8 S	Nitrobenzene-d5	0.237	0.228	3.8	64	0.00
9	Naphthalene	1.169	1.186	-1.5	67	0.00
10	Hexachlorobutadiene	0.198	0.202	-2.0	66	0.00
11 SURR	2-Methylnaphthalene-d10	0.602	0.618	-2.7	68	0.00
12	2-Methylnaphthalene	0.766	0.774	-1.0	67	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	73	0.00
14 S	2,4,6-Tribromophenol	0.127	0.125	1.6	73	0.00
15 S	2-Fluorobiphenyl	1.896	1.818	4.1	68	0.00
16	Acenaphthylene	1.885	1.639	13.1	65	0.00
17	Acenaphthene	1.487	1.409	5.2	69	0.00
18	Fluorene	1.854	1.748	5.7	69	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	76	0.00
20	4-Bromophenyl-phenylether	0.230	0.212	7.8	70	0.00
21	Hexachlorobenzene	0.308	0.287	6.8	70	0.00
22	Pentachlorophenol	0.082	0.070	14.6	73	0.00
23	Phenanthrene	1.408	1.315	6.6	70	0.00
24	Anthracene	1.067	0.944	11.5	68	0.00
25 SURR	Fluoranthene-d10	1.069	0.991	7.3	71	0.00
26	Fluoranthene	1.460	1.318	9.7	69	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	76	0.00
28	Pvrene	1.570	1.460	7.0	70	0.00
29 S	Terphenyl-d14	0.703	0.669	4.8	71	0.00
30	Benzo(a)anthracene	1.260	1.175	6.7	71	0.00
31	Chrysene	1.473	1.425	3.3	71	0.00
32	Bis(2-ethylhexyl)phthalate	0.469	0.411	12.4	72	0.00
33	Indeno(1,2,3-cd)pyrene	1.375	1.288	6.3	68	0.00
34 I	Perylene-d12	1.000	1.000	0.0	72	0.00
35	Benzo(b)fluoranthene	1.826	1.710	6.4	68	0.00
36	Benzo(k)fluoranthene	1.786	1.699	4.9	68	0.00
37 C	Benzo(a)pyrene	1.488	1.364	8.3	67	0.00
38	Dibenzo(a,h)anthracene	1.325	1.261	4.8	69	0.00
39	Benzo(a,h,i)perylene	1.564	1.484	5.1	68	0.00

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA N\Data\BN030118\
Data File : BN000216.D
Acq On : 01 Mar 2018 14:28
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_N
LabSampleId :
SSTDCCC0.4

Quant Time: Mar 01 17:46:19 2018
Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-SIM-BN022718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 27 15:07:57 2018
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			