

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA N\Data\BN030118\
 Data File : BN000225.D
 Acq On : 01 Mar 2018 19:39
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDC00.4

Quant Time: Mar 02 09:04:19 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	72	0.00
2	1,4-Dioxane	0.414	0.425	-2.7	78	0.00
3	n-Nitrosodimethylamine	0.282	0.295	-4.6	79	0.00
4 S	2-Fluorophenol	0.845	0.860	-1.8	71	0.00
5 S	Phenol-d6	1.073	1.091	-1.7	74	0.00
6	bis(2-Chloroethyl)ether	1.363	1.418	-4.0	72	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	74	0.00
8 S	Nitrobenzene-d5	0.237	0.222	6.3	69	0.00
9	Naphthalene	1.169	1.182	-1.1	73	0.00
10	Hexachlorobutadiene	0.198	0.200	-1.0	72	0.00
11 SURR	2-Methylnaphthalene-d10	0.602	0.619	-2.8	75	0.00
12	2-Methylnaphthalene	0.766	0.777	-1.4	74	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	79	0.00
14 S	2,4,6-Tribromophenol	0.127	0.118	7.1	74	0.00
15 S	2-Fluorobiphenyl	1.896	1.856	2.1	75	0.00
16	Acenaphthylene	1.885	1.645	12.7	70	0.00
17	Acenaphthene	1.487	1.420	4.5	75	0.00
18	Fluorene	1.854	1.779	4.0	76	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	80	0.00
20	4-Bromophenyl-phenylether	0.230	0.212	7.8	74	0.00
21	Hexachlorobenzene	0.308	0.290	5.8	75	0.00
22	Pentachlorophenol	0.082	0.057	30.5#	64	0.00
23	Phenanthrene	1.408	1.315	6.6	75	0.00
24	Anthracene	1.067	0.929	12.9	72	0.00
25 SURR	Fluoranthene-d10	1.069	0.968	9.4	74	0.00
26	Fluoranthene	1.460	1.299	11.0	72	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	74	0.00
28	Pvrene	1.570	1.564	0.4	73	0.00
29 S	Terphenyl-d14	0.703	0.719	-2.3	74	0.00
30	Benzo(a)anthracene	1.260	1.134	10.0	66	0.00
31	Chrysene	1.473	1.453	1.4	71	0.00
32	Bis(2-ethylhexyl)phthalate	0.469	0.419	10.7	71	0.00
33	Indeno(1,2,3-cd)pyrene	1.375	1.325	3.6	68	0.00
34 I	Perylene-d12	1.000	1.000	0.0	71	0.00
35	Benzo(b)fluoranthene	1.826	1.759	3.7	70	0.00
36	Benzo(k)fluoranthene	1.786	1.726	3.4	68	0.00
37 C	Benzo(a)pyrene	1.488	1.372	7.8	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	1.481	5.3	67	0.00

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA N\Data\BN030118\
Data File : BN000225.D
Acq On : 01 Mar 2018 19:39
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_N
LabSampleId :
SSTDCCC0.4

Quant Time: Mar 02 09:04:19 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			