

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN073018\
 Data File : BN002190.D
 Acq On : 30 Jul 2018 17:50
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDC00.4

Quant Time: Jul 31 01:11:09 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 16:17:29 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	87	0.00
2	1,4-Dioxane	0.450	0.435	3.3	93	0.00
3	n-Nitrosodimethylamine	0.302	0.247	18.2	80	0.00
4 S	2-Fluorophenol	0.933	0.875	6.2	86	0.00
5 S	Phenol-d6	1.135	1.058	6.8	87	0.00
6	bis(2-Chloroethyl)ether	1.091	1.014	7.1	81	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	90	0.00
8 S	Nitrobenzene-d5	0.301	0.288	4.3	88	0.00
9	Naphthalene	0.902	0.882	2.2	90	0.00
10	Hexachlorobutadiene	0.195	0.190	2.6	90	0.00
11 SURR	2-Methylnaphthalene-d10	0.589	0.567	3.7	90	0.00
12	2-Methylnaphthalene	0.610	0.592	3.0	89	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	85	0.00
14 S	2,4,6-Tribromophenol	0.171	0.143	16.4	72	0.00
15 S	2-Fluorobiphenyl	1.536	1.512	1.6	88	0.00
16	Acenaphthylene	1.709	1.605	6.1	83	0.01
17	Acenaphthene	1.099	1.067	2.9	85	0.00
18	Fluorene	1.352	1.285	5.0	81	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	71	0.00
20	4-Bromophenyl-phenylether	0.226	0.223	1.3	76	0.00
21	Hexachlorobenzene	0.250	0.255	-2.0	77	0.00
22	Pentachlorophenol	0.046	0.039	15.2	63	0.00
23	Phenanthrene	0.995	0.952	4.3	70	0.00
24	Anthracene	0.823	0.747	9.2	68	0.00
25 SURR	Fluoranthene-d10	0.988	0.858	13.2	60	0.00
26	Fluoranthene	1.083	0.933	13.9	59	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	54	0.00
28	Pvrene	1.183	1.212	-2.5	59	0.00
29 S	Terphenyl-d14	0.855	0.843	1.4	57	0.00
30	Benzo(a)anthracene	1.006	0.929	7.7	52	0.00
31	Chrysene	1.126	1.094	2.8	54	0.00
32	Bis(2-ethylhexyl)phthalate	0.460	0.386	16.1	46#	-0.01
33	Indeno(1,2,3-cd)pyrene	1.061	1.155	-8.9	69	0.00
34 I	Perylene-d12	1.000	1.000	0.0	60	0.00
35	Benzo(b)fluoranthene	1.176	1.080	8.2	56	0.00
36	Benzo(k)fluoranthene	1.221	1.175	3.8	61	0.00
37 C	Benzo(a)pyrene	1.066	1.036	2.8	61	0.00
38	Dibenzo(a,h)anthracene	1.037	1.048	-1.1	69	0.00
39	Benzo(a,h,i)perylene	1.060	1.083	-2.2	69	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN073018\
Data File : BN002190.D
Acq On : 30 Jul 2018 17:50
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
LabSampleId :
SSTDCCC0.4

Quant Time: Jul 31 01:11:09 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 30 16:17:29 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			