

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN102918\
 Data File : BN003329.D
 Acq On : 29 Oct 2018 09:58
 Operator : MJ/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 30 06:44:09 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN102618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 29 16:15:49 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	99	0.00
2	1,4-Dioxane	0.518	0.457	11.8	94	0.00
3	n-Nitrosodimethylamine	0.238	0.236	0.8	102	0.00
4 S	2-Fluorophenol	0.991	0.935	5.7	97	0.00
5 S	Phenol-d6	1.251	1.181	5.6	98	0.00
6	bis(2-Chloroethyl)ether	1.116	1.079	3.3	96	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	97	-0.01
8 S	Nitrobenzene-d5	0.175	0.219	-25.1#	133	0.00
9	Naphthalene	0.979	0.909	7.2	93	-0.01
10	Hexachlorobutadiene	0.179	0.168	6.1	94	0.00
11 SURR	2-Methylnaphthalene-d10	0.651	0.592	9.1	92	0.00
12	2-Methylnaphthalene	0.684	0.625	8.6	92	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	97	-0.01
14 S	2,4,6-Tribromophenol	0.159	0.147	7.5	106	0.00
15 S	2-Fluorobiphenyl	1.721	1.528	11.2	92	-0.01
16	Acenaphthylene	1.783	1.552	13.0	90	-0.01
17	Acenaphthene	1.236	1.116	9.7	91	0.00
18	Fluorene	1.546	1.355	12.4	88	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	97	0.00
20	4-Bromophenyl-phenylether	0.232	0.207	10.8	91	-0.01
21	Hexachlorobenzene	0.244	0.223	8.6	91	0.00
22	Pentachlorophenol	0.066	0.074	-12.1	118	0.00
23	Phenanthrene	1.097	0.982	10.5	91	0.00
24	Anthracene	0.941	0.806	14.3	88	-0.01
25 SURR	Fluoranthene-d10	1.163	1.034	11.1	91	0.00
26	Fluoranthene	1.260	1.117	11.3	91	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	96	0.00
28	Pvrene	1.106	0.974	11.9	90	0.00
29 S	Terphenyl-d14	0.930	0.793	14.7	90	0.00
30	Benzo(a)anthracene	1.048	0.927	11.5	95	-0.01
31	Chrysene	1.131	1.023	9.5	93	-0.01
32	Bis(2-ethylhexyl)phthalate	0.310	0.290	6.5	103	-0.01
33	Indeno(1,2,3-cd)pyrene	1.170	1.130	3.4	99	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	102	0.00
35	Benzo(b)fluoranthene	1.248	1.080	13.5	95	0.00
36	Benzo(k)fluoranthene	1.236	1.061	14.2	94	0.00
37 C	Benzo(a)pyrene	1.136	0.997	12.2	95	0.00
38	Dibenzo(a,h)anthracene	1.058	0.933	11.8	99	-0.01
39	Benzo(a,h,i)perylene	1.058	0.958	9.5	102	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN102918\
Data File : BN003329.D
Acq On : 29 Oct 2018 09:58
Operator : MJ/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_N
LabSampleId :
SSTDCCC0.4

Quant Time: Oct 30 06:44:09 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN102618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 29 16:15:49 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			