

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031119\
 Data File : BE098857.D
 Acq On : 11 Mar 2019 11:08
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 12 16:30:39 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 06 15:33:24 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	110	0.00
2	1,4-Dioxane	0.400	0.402	-0.5	113	0.00
3	n-Nitrosodimethylamine	0.400	0.223	44.3#	60	0.02
4 S	2-Fluorophenol	0.400	0.313	21.8	83	0.01
5 S	Phenol-d6	0.400	0.371	7.3	102	0.00
6	bis(2-Chloroethyl)ether	0.400	0.341	14.8	97	0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	114	0.00
8 S	Nitrobenzene-d5	0.400	0.348	13.0	110	0.00
9	Naphthalene	0.400	0.403	-0.8	118	0.00
10	Hexachlorobutadiene	0.400	0.384	4.0	113	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.392	2.0	114	0.00
12	2-Methylnaphthalene	0.400	0.383	4.3	113	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	118	0.00
14 S	2,4,6-Tribromophenol	0.400	0.319	20.3	95	0.00
15 S	2-Fluorobiphenyl	0.400	0.411	-2.7	123	0.00
16	Acenaphthylene	0.400	0.384	4.0	119	0.00
17	Acenaphthene	0.400	0.384	4.0	114	0.00
18	Fluorene	0.400	0.378	5.5	114	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	110	0.00
20	4-Bromophenyl-phenylether	0.400	0.360	10.0	100	0.00
21	Hexachlorobenzene	0.400	0.390	2.5	112	0.00
22	Pentachlorophenol	0.400	0.382	4.5	73	0.01
23	Phenanthrene	0.400	0.380	5.0	107	0.00
24	Anthracene	0.400	0.355	11.3	103	0.00
25 SURR	Fluoranthene-d10	0.400	0.354	11.5	101	0.00
26	Fluoranthene	0.400	0.358	10.5	102	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	99	0.00
28	Pvrene	0.400	0.407	-1.7	101	0.00
29 S	Terphenyl-d14	0.400	0.383	4.3	95	0.00
30	Benzo(a)anthracene	0.400	0.382	4.5	96	0.00
31	Chrysene	0.400	0.385	3.8	98	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.309	22.8	83	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.397	0.8	100	0.00
34 I	Perylene-d12	0.400	0.400	0.0	100	0.00
35	Benzo(b)fluoranthene	0.400	0.383	4.3	99	0.00
36	Benzo(k)fluoranthene	0.400	0.409	-2.2	108	0.00
37 C	Benzo(a)pyrene	0.400	0.389	2.8	101	0.00
38	Dibenzo(a,h)anthracene	0.400	0.376	6.0	96	0.00
39	Benzo(a,h,i)perylene	0.400	0.394	1.5	101	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031119\
Data File : BE098857.D
Acq On : 11 Mar 2019 11:08
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Mar 12 16:30:39 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 06 15:33:24 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	Amount	Calc.	%Dev	Area	% Dev(min)

(#) = Out of Range	SPCC's out = 0 CCC's out = 0				