

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011819\
 Data File : BE098663.D
 Acq On : 18 Jan 2019 8:34
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 18 11:37:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE011519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 15 14:30:55 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	81	0.00
2	1,4-Dioxane	0.400	0.398	0.5	87	0.00
3	n-Nitrosodimethylamine	0.400	0.360	10.0	72	0.00
4 S	2-Fluorophenol	0.400	0.396	1.0	82	0.00
5 S	Phenol-d6	0.400	0.396	1.0	80	0.00
6	bis(2-Chloroethyl)ether	0.400	0.367	8.3	80	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	83	0.00
8 S	Nitrobenzene-d5	0.400	0.431	-7.7	93	0.01
9	Naphthalene	0.400	0.365	8.8	81	0.00
10	Hexachlorobutadiene	0.400	0.397	0.8	87	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.366	8.5	82	0.00
12	2-Methylnaphthalene	0.400	0.361	9.8	81	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	89	0.00
14 S	2,4,6-Tribromophenol	0.400	0.386	3.5	87	0.00
15 S	2-Fluorobiphenyl	0.400	0.338	15.5	83	0.01
16	Acenaphthylene	0.400	0.352	12.0	84	0.00
17	Acenaphthene	0.400	0.352	12.0	82	0.00
18	Fluorene	0.400	0.363	9.3	84	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	88	0.00
20	4-Bromophenyl-phenylether	0.400	0.347	13.3	83	0.00
21	Hexachlorobenzene	0.400	0.358	10.5	83	0.00
22	Pentachlorophenol	0.400	0.491	-22.7	114	0.00
23	Phenanthrene	0.400	0.356	11.0	84	0.00
24	Anthracene	0.400	0.332	17.0	81	0.00
25 SURR	Fluoranthene-d10	0.400	0.361	9.8	86	0.00
26	Fluoranthene	0.400	0.355	11.3	85	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	90	0.00
28	Pvrene	0.400	0.354	11.5	83	0.00
29 S	Terphenyl-d14	0.400	0.354	11.5	87	0.00
30	Benzo(a)anthracene	0.400	0.358	10.5	86	0.00
31	Chrysene	0.400	0.356	11.0	85	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.325	18.8	83	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.373	6.8	87	0.00
34 I	Perylene-d12	0.400	0.400	0.0	91	0.00
35	Benzo(b)fluoranthene	0.400	0.353	11.8	86	0.00
36	Benzo(k)fluoranthene	0.400	0.379	5.3	89	0.00
37 C	Benzo(a)pyrene	0.400	0.364	9.0	87	0.00
38	Dibenzo(a,h)anthracene	0.400	0.387	3.3	93	0.00
39	Benzo(a,h,i)perylene	0.400	0.364	9.0	86	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011819\
Data File : BE098663.D
Acq On : 18 Jan 2019 8:34
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Jan 18 11:37:44 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE011519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 15 14:30:55 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				