

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031419\
 Data File : BE098938.D
 Acq On : 14 Mar 2019 10:12
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 15 05:33:16 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	94	0.00
2	1,4-Dioxane	0.400	0.389	2.8	94	0.00
3	n-Nitrosodimethylamine	0.400	0.323	19.3	75	0.04
4 S	2-Fluorophenol	0.400	0.360	10.0	82	0.02
5 S	Phenol-d6	0.400	0.348	13.0	82	0.02
6	bis(2-Chloroethyl)ether	0.400	0.323	19.3	79	0.02
7 I	Naphthalene-d8	0.400	0.400	0.0	107	0.00
8 S	Nitrobenzene-d5	0.400	0.331	17.3	99	0.01
9	Naphthalene	0.400	0.388	3.0	107	0.00
10	Hexachlorobutadiene	0.400	0.391	2.3	108	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.389	2.8	106	0.00
12	2-Methylnaphthalene	0.400	0.362	9.5	100	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	113	0.00
14 S	2,4,6-Tribromophenol	0.400	0.301	24.8	86	0.00
15 S	2-Fluorobiphenyl	0.400	0.370	7.5	106	0.00
16	Acenaphthylene	0.400	0.369	7.8	110	0.00
17	Acenaphthene	0.400	0.389	2.8	111	-0.01
18	Fluorene	0.400	0.372	7.0	108	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	103	0.00
20	4-Bromophenyl-phenylether	0.400	0.342	14.5	89	0.00
21	Hexachlorobenzene	0.400	0.387	3.3	105	0.00
22	Pentachlorophenol	0.400	0.377	5.8	66	0.03
23	Phenanthrene	0.400	0.385	3.8	101	0.00
24	Anthracene	0.400	0.333	16.8	90	0.00
25 SURR	Fluoranthene-d10	0.400	0.350	12.5	93	0.00
26	Fluoranthene	0.400	0.349	12.8	93	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	90	0.00
28	Pvrene	0.400	0.408	-2.0	92	0.00
29 S	Terphenyl-d14	0.400	0.376	6.0	85	0.00
30	Benzo(a)anthracene	0.400	0.365	8.8	84	0.00
31	Chrysene	0.400	0.377	5.8	87	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.309	22.8	76	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.355	11.3	81	0.00
34 I	Perylene-d12	0.400	0.400	0.0	92	0.00
35	Benzo(b)fluoranthene	0.400	0.361	9.8	86	0.00
36	Benzo(k)fluoranthene	0.400	0.401	-0.3	97	0.00
37 C	Benzo(a)pyrene	0.400	0.376	6.0	89	0.00
38	Dibenzo(a,h)anthracene	0.400	0.327	18.3	76	0.00
39	Benzo(a,h,i)perylene	0.400	0.346	13.5	81	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031419\
Data File : BE098938.D
Acq On : 14 Mar 2019 10:12
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Mar 15 05:33:16 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				