

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE060719\
 Data File : BE099853.D
 Acq On : 7 Jun 2019 15:25
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE060719

Quant Time: Jun 07 18:36:37 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE060719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 07 15:25:21 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	115	0.00
2	1,4-Dioxane	0.400	0.395	1.3	128	0.00
3	n-Nitrosodimethylamine	0.400	0.386	3.5	118	0.00
4 S	2-Fluorophenol	0.400	0.381	4.8	115	0.01
5 S	Phenol-d6	0.400	0.386	3.5	110	0.01
6	bis(2-Chloroethyl)ether	0.400	0.392	2.0	122	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	118	0.00
8 S	Nitrobenzene-d5	0.400	0.389	2.8	120	0.01
9	Naphthalene	0.400	0.395	1.3	119	0.01
10	Hexachlorobutadiene	0.400	0.405	-1.3	120	0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.405	-1.3	120	0.01
12	2-Methylnaphthalene	0.400	0.370	7.5	110	0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	107	0.01
14 S	2,4,6-Tribromophenol	0.400	0.390	2.5	98	0.00
15 S	2-Fluorobiphenyl	0.400	0.413	-3.2	115	0.00
16	Acenaphthylene	0.400	0.395	1.3	108	0.00
17	Acenaphthene	0.400	0.395	1.3	109	0.00
18	Fluorene	0.400	0.399	0.3	106	0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	103	0.00
20	4-Bromophenyl-phenylether	0.400	0.393	1.8	106	0.01
21	Hexachlorobenzene	0.400	0.390	2.5	102	0.01
22	Pentachlorophenol	0.400	0.518	-29.5#	456	-0.01
23	Phenanthrene	0.400	0.391	2.3	100	0.00
24	Anthracene	0.400	0.381	4.8	99	0.00
25 SURR	Fluoranthene-d10	0.400	0.400	0.0	101	0.00
26	Fluoranthene	0.400	0.400	0.0	101	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	119	0.00
28	Pvrene	0.400	0.384	4.0	100	0.00
29 S	Terphenyl-d14	0.400	0.378	5.5	101	0.00
30	Benzo(a)anthracene	0.400	0.411	-2.7	113	0.00
31	Chrysene	0.400	0.381	4.8	103	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.409	-2.2	113	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.396	1.0	110	0.00
34 I	Perylene-d12	0.400	0.400	0.0	107	0.00
35	Benzo(b)fluoranthene	0.400	0.418	-4.5	114	0.00
36	Benzo(k)fluoranthene	0.400	0.398	0.5	105	0.00
37 C	Benzo(a)pyrene	0.400	0.412	-3.0	111	0.00
38	Dibenzo(a,h)anthracene	0.400	0.402	-0.5	108	0.00
39	Benzo(a,h,i)perylene	0.400	0.404	-1.0	109	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE060719\
Data File : BE099853.D
Acq On : 7 Jun 2019 15:25
Operator : JU/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
ICVBE060719

Quant Time: Jun 07 18:36:37 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE060719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 07 15:25:21 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				