

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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	115	0.00
2	1,4-Dioxane	0.596	0.556	6.7	128	0.00
3	n-Nitrosodimethylamine	0.382	0.368	3.7	118	0.00
4 S	2-Fluorophenol	1.339	1.275	4.8	115	0.01
5 S	Phenol-d6	1.650	1.591	3.6	110	0.01
6	bis(2-Chloroethyl)ether	1.213	1.188	2.1	122	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	118	0.00
8 S	Nitrobenzene-d5	0.403	0.392	2.7	120	0.01
9	Naphthalene	4.347	4.297	1.2	119	0.01
10	Hexachlorobutadiene	0.320	0.324	-1.3	120	0.01
11 SURR	2-Methylnaphthalene-d10	0.724	0.732	-1.1	120	0.01
12	2-Methylnaphthalene	0.739	0.684	7.4	110	0.01
13 I	Acenaphthene-d10	1.000	1.000	0.0	107	0.01
14 S	2,4,6-Tribromophenol	0.396	0.385	2.8	98	0.00
15 S	2-Fluorobiphenyl	1.446	1.492	-3.2	115	0.00
16	Acenaphthylene	1.981	1.956	1.3	108	0.00
17	Acenaphthene	1.086	1.072	1.3	109	0.00
18	Fluorene	1.661	1.657	0.2	106	0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
20	4-Bromophenyl-phenylether	0.229	0.225	1.7	106	0.01
21	Hexachlorobenzene	0.224	0.218	2.7	102	0.01
22	Pentachlorophenol	0.069	0.059	14.5	456#	-0.01
23	Phenanthrene	0.992	0.970	2.2	100	0.00
24	Anthracene	0.951	0.907	4.6	99	0.00
25 SURR	Fluoranthene-d10	6.199	6.192	0.1	101	0.00
26	Fluoranthene	1.751	1.752	-0.1	101	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	119	0.00
28	Pvrene	1.059	1.017	4.0	100	0.00
29 S	Terphenyl-d14	0.746	0.704	5.6	101	0.00
30	Benzo(a)anthracene	1.271	1.307	-2.8	113	0.00
31	Chrysene	1.248	1.187	4.9	103	0.00
32	Bis(2-ethylhexyl)phthalate	2.073	2.118	-2.2	113	0.00
33	Indeno(1,2,3-cd)pyrene	1.520	1.505	1.0	110	0.00
34 I	Perylene-d12	1.000	1.000	0.0	107	0.00
35	Benzo(b)fluoranthene	1.111	1.160	-4.4	114	0.00
36	Benzo(k)fluoranthene	1.111	1.106	0.5	105	0.00
37 C	Benzo(a)pyrene	1.078	1.109	-2.9	111	0.00
38	Dibenzo(a,h)anthracene	1.102	1.109	-0.6	108	0.00
39	Benzo(a,h,i)perylene	1.120	1.130	-0.9	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	AvgRF	CCRF	%Dev Area	% Dev(min)

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