

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE092719\
 Data File : BE100578.D
 Acq On : 27 Sep 2019 18:34
 Operator : JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE092719

Quant Time: Sep 27 19:09:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE092719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 27 18:32:20 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	99	0.00
2	1,4-Dioxane	0.721	0.679	5.8	100	0.00
3	n-Nitrosodimethylamine	0.384	0.346	9.9	99	0.00
4 S	2-Fluorophenol	1.179	1.095	7.1	96	0.00
5 S	Phenol-d6	1.601	1.470	8.2	96	0.00
6	bis(2-Chloroethyl)ether	1.383	1.312	5.1	98	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
8 S	Nitrobenzene-d5	0.289	0.263	9.0	98	0.00
9	Naphthalene	4.156	4.050	2.6	101	0.00
10	Hexachlorobutadiene	0.127	0.123	3.1	101	0.00
11 SURR	2-Methylnaphthalene-d10	0.623	0.594	4.7	100	0.00
12	2-Methylnaphthalene	0.707	0.678	4.1	100	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
14 S	2,4,6-Tribromophenol	0.115	0.106	7.8	99	0.01
15 S	2-Fluorobiphenyl	1.385	1.349	2.6	100	0.00
16	Acenaphthylene	1.782	1.716	3.7	100	0.00
17	Acenaphthene	1.145	1.088	5.0	99	0.00
18	Fluorene	1.488	1.461	1.8	100	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
20	4-Bromophenyl-phenylether	0.186	0.173	7.0	103	0.00
21	Hexachlorobenzene	0.153	0.147	3.9	101	0.00
22	Pentachlorophenol	0.055	0.041	25.5#	103	0.00
23	Phenanthrene	1.128	1.108	1.8	102	0.00
24	Anthracene	1.021	0.945	7.4	102	0.00
25 SURR	Fluoranthene-d10	5.084	4.720	7.2	98	0.00
26	Fluoranthene	1.418	1.357	4.3	98	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	99	0.01
28	Pvrene	1.211	1.227	-1.3	97	0.00
29 S	Terphenyl-d14	0.988	0.935	5.4	97	0.00
30	Benzo(a)anthracene	1.178	1.133	3.8	97	0.00
31	Chrysene	1.201	1.206	-0.4	99	0.01
32	Bis(2-ethylhexyl)phthalate	3.127	2.840	9.2	97	0.00
33	Indeno(1,2,3-cd)pyrene	1.481	1.455	1.8	97	0.00
34 I	Perylene-d12	1.000	1.000	0.0	97	0.00
35	Benzo(b)fluoranthene	1.094	1.039	5.0	97	0.00
36	Benzo(k)fluoranthene	1.193	1.125	5.7	97	0.00
37 C	Benzo(a)pyrene	1.045	0.992	5.1	96	0.00
38	Dibenzo(a,h)anthracene	1.065	1.019	4.3	98	0.00
39	Benzo(a,h,i)perylene	1.088	1.043	4.1	96	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE092719\
Data File : BE100578.D
Acq On : 27 Sep 2019 18:34
Operator : JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
ICVBE092719

Quant Time: Sep 27 19:09:32 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE092719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 27 18:32:20 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			