

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101019\
 Data File : BE100631.D
 Acq On : 10 Oct 2019 15:11
 Operator : JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE101019

Quant Time: Oct 10 15:57:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 10 14:01:52 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	108	0.00
2	1,4-Dioxane	0.714	0.681	4.6	104	0.00
3	n-Nitrosodimethylamine	0.755	0.708	6.2	110	0.00
4 S	2-Fluorophenol	1.235	1.177	4.7	109	0.00
5 S	Phenol-d6	1.742	1.644	5.6	108	0.00
6	bis(2-Chloroethyl)ether	1.430	1.405	1.7	113	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
8 S	Nitrobenzene-d5	0.248	0.237	4.4	121	0.00
9	Naphthalene	4.206	4.115	2.2	112	0.00
10	Hexachlorobutadiene	0.183	0.179	2.2	114	0.00
11 SURR	2-Methylnaphthalene-d10	0.631	0.618	2.1	114	0.01
12	2-Methylnaphthalene	0.717	0.696	2.9	113	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	113	0.00
14 S	2,4,6-Tribromophenol	0.106	0.094	11.3	114	0.00
15 S	2-Fluorobiphenyl	1.457	1.438	1.3	114	0.00
16	Acenaphthylene	1.721	1.589	7.7	108	0.00
17	Acenaphthene	1.112	1.062	4.5	112	0.00
18	Fluorene	1.477	1.437	2.7	114	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	116	0.00
20	4-Bromophenyl-phenylether	0.204	0.196	3.9	116	0.00
21	Hexachlorobenzene	0.194	0.184	5.2	116	0.00
22	Pentachlorophenol	0.062	0.057	8.1	119	0.00
23	Phenanthrene	1.100	1.125	-2.3	118	0.00
24	Anthracene	1.021	0.975	4.5	115	0.00
25 SURR	Fluoranthene-d10	5.162	4.934	4.4	115	0.00
26	Fluoranthene	1.436	1.429	0.5	116	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	112	0.00
28	Pvrene	1.118	1.150	-2.9	116	0.00
29 S	Terphenyl-d14	0.875	0.867	0.9	117	0.00
30	Benzo(a)anthracene	1.305	1.260	3.4	109	0.00
31	Chrysene	1.304	1.269	2.7	110	0.00
32	Bis(2-ethylhexyl)phthalate	2.729	2.483	9.0	102	0.00
33	Indeno(1,2,3-cd)pyrene	1.603	1.363	15.0	98	0.00
34 I	Perylene-d12	1.000	1.000	0.0	100	0.00
35	Benzo(b)fluoranthene	1.166	1.094	6.2	101	0.00
36	Benzo(k)fluoranthene	1.200	1.150	4.2	102	0.00
37 C	Benzo(a)pyrene	1.086	0.992	8.7	98	0.00
38	Dibenzo(a,h)anthracene	1.122	1.011	9.9	99	0.00
39	Benzo(a,h,i)perylene	1.129	1.043	7.6	99	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101019\
Data File : BE100631.D
Acq On : 10 Oct 2019 15:11
Operator : JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
ICVBE101019

Quant Time: Oct 10 15:57:48 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 10 14:01:52 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			