

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE120419\
 Data File : BE100779.D
 Acq On : 4 Dec 2019 17:18
 Operator : JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE120419

Quant Time: Dec 04 18:00:15 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE120419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 04 17:18:11 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	103	0.00
2	1,4-Dioxane	0.729	0.724	0.7	103	-0.01
3	n-Nitrosodimethylamine	0.594	0.552	7.1	110	-0.01
4 S	2-Fluorophenol	0.949	0.939	1.1	109	0.00
5 S	Phenol-d6	1.433	1.360	5.1	106	0.00
6	bis(2-Chloroethyl)ether	1.385	1.333	3.8	105	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
8 S	Nitrobenzene-d5	0.190	0.196	-3.2	125	0.00
9	Naphthalene	4.250	4.099	3.6	106	0.00
10	Hexachlorobutadiene	0.166	0.163	1.8	107	0.00
11 SURR	2-Methylnaphthalene-d10	0.580	0.557	4.0	108	0.00
12	2-Methylnaphthalene	0.664	0.632	4.8	107	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
14 S	2,4,6-Tribromophenol	0.044	0.045	-2.3	124	0.00
15 S	2-Fluorobiphenyl	1.587	1.530	3.6	107	0.00
16	Acenaphthylene	1.655	1.438	13.1	107	0.00
17	Acenaphthene	1.140	1.051	7.8	108	0.00
18	Fluorene	1.471	1.347	8.4	110	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
20	4-Bromophenyl-phenylether	0.196	0.178	9.2	110	0.00
21	Hexachlorobenzene	0.213	0.202	5.2	110	0.00
22	Pentachlorophenol	0.015	0.004	73.3#	147	0.00
23	Phenanthrene	1.212	1.158	4.5	111	0.01
24	Anthracene	1.013	0.881	13.0	111	0.00
25 SURR	Fluoranthene-d10	4.728	4.161	12.0	107	0.00
26	Fluoranthene	1.446	1.267	12.4	106	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	99	0.00
28	Pvrene	1.300	1.283	1.3	105	0.00
29 S	Terphenyl-d14	0.958	0.962	-0.4	104	0.00
30	Benzo(a)anthracene	1.179	1.051	10.9	100	0.00
31	Chrysene	1.385	1.363	1.6	101	0.00
32	Bis(2-ethylhexyl)phthalate	1.069	0.815	23.8	100	0.00
33	Indeno(1,2,3-cd)pyrene	1.207	1.050	13.0	98	0.00
34 I	Perylene-d12	1.000	1.000	0.0	100	0.00
35	Benzo(b)fluoranthene	1.095	1.003	8.4	97	0.00
36	Benzo(k)fluoranthene	1.228	1.166	5.0	100	0.00
37 C	Benzo(a)pyrene	0.948	0.834	12.0	97	0.00
38	Dibenzo(a,h)anthracene	0.943	0.840	10.9	100	0.00
39	Benzo(a,h,i)perylene	1.009	0.927	8.1	98	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE120419\
Data File : BE100779.D
Acq On : 4 Dec 2019 17:18
Operator : JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
ICVBE120419

Quant Time: Dec 04 18:00:15 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE120419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 04 17:18:11 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			