

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071219\
 Data File : BE100423.D
 Acq On : 12 Jul 2019 17:17
 Operator : HP/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE071219

Quant Time: Jul 12 18:04:23 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 17:51:04 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	-0.01
2	1,4-Dioxane	0.705	0.762	-8.1	122	-0.01
3	n-Nitrosodimethylamine	0.433	0.322	25.6#	102	-0.01
4 S	2-Fluorophenol	1.235	1.284	-4.0	110	0.00
5 S	Phenol-d6	1.817	1.826	-0.5	106	0.00
6	bis(2-Chloroethyl)ether	1.355	1.393	-2.8	110	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	110	0.00
8 S	Nitrobenzene-d5	0.201	0.216	-7.5	108	0.00
9	Naphthalene	4.295	4.664	-8.6	110	0.00
10	Hexachlorobutadiene	0.208	0.215	-3.4	108	0.00
11 SURR	2-Methylnaphthalene-d10	0.667	0.722	-8.2	111	0.00
12	2-Methylnaphthalene	0.746	0.811	-8.7	111	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
14 S	2,4,6-Tribromophenol	0.152	0.149	2.0	115	0.00
15 S	2-Fluorobiphenyl	1.473	1.598	-8.5	112	0.00
16	Acenaphthylene	2.086	2.227	-6.8	109	0.00
17	Acenaphthene	1.229	1.332	-8.4	111	0.01
18	Fluorene	1.647	1.784	-8.3	108	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
20	4-Bromophenyl-phenylether	0.228	0.245	-7.5	112	0.00
21	Hexachlorobenzene	0.235	0.252	-7.2	112	0.00
22	Pentachlorophenol	0.068	0.070	-2.9	117	0.00
23	Phenanthrene	1.096	1.231	-12.3	113	0.00
24	Anthracene	1.097	1.193	-8.8	110	0.01
25 SURR	Fluoranthene-d10	5.869	6.294	-7.2	106	0.00
26	Fluoranthene	1.575	1.717	-9.0	107	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	104	0.00
28	Pvrene	1.257	1.437	-14.3	105	0.00
29 S	Terphenyl-d14	0.883	0.994	-12.6	105	0.00
30	Benzo(a)anthracene	1.428	1.548	-8.4	101	0.00
31	Chrysene	1.368	1.498	-9.5	102	0.00
32	Bis(2-ethylhexyl)phthalate	2.190	2.066	5.7	106	0.00
33	Indeno(1,2,3-cd)pyrene	1.684	1.708	-1.4	98	0.00
34 I	Perylene-d12	1.000	1.000	0.0	101	0.00
35	Benzo(b)fluoranthene	1.260	1.338	-6.2	96	0.00
36	Benzo(k)fluoranthene	1.285	1.474	-14.7	108	0.00
37 C	Benzo(a)pyrene	1.191	1.268	-6.5	98	0.00
38	Dibenzo(a,h)anthracene	1.184	1.242	-4.9	98	0.00
39	Benzo(a,h,i)perylene	1.175	1.269	-8.0	100	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071219\
Data File : BE100423.D
Acq On : 12 Jul 2019 17:17
Operator : HP/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
ICVBE071219

Quant Time: Jul 12 18:04:23 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 12 17:51:04 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			