

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN063020\
 Data File : BN011080.D
 Acq On : 30 Jun 2020 14:29
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ICVBN063020

Quant Time: Jun 30 15:24:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN063020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 30 14:12:13 2020
 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	109	0.00
2	1,4-Dioxane	0.400	0.438	-9.5	116	0.00
3	n-Nitrosodimethylamine	0.400	0.428	-7.0	122	0.00
4 S	2-Fluorophenol	0.400	0.435	-8.7	115	-0.02
5 S	Phenol-d6	0.400	0.423	-5.7	110	0.00
6	bis(2-Chloroethyl)ether	0.400	0.411	-2.7	101	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	117	0.00
8 S	Nitrobenzene-d5	0.400	0.385	3.8	113	0.00
9	Naphthalene	0.400	0.403	-0.8	119	-0.01
10	Hexachlorobutadiene	0.400	0.383	4.3	103	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.361	9.8	101	0.00
12	2-Methylnaphthalene	0.400	0.370	7.5	107	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	110	0.00
14 S	2,4,6-Tribromophenol	0.400	0.390	2.5	106	-0.01
15 S	2-Fluorobiphenyl	0.400	0.413	-3.2	111	0.00
16	Acenaphthylene	0.400	0.399	0.3	109	0.00
17	Acenaphthene	0.400	0.402	-0.5	110	0.00
18	Fluorene	0.400	0.407	-1.7	110	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	110	-0.01
20	4,6-Dinitro-2-methylphenol	0.400	0.381	4.8	103	0.00
21	4-Bromophenyl-phenylether	0.400	0.390	2.5	105	0.00
22	Hexachlorobenzene	0.400	0.411	-2.7	111	0.00
23	Atrazine	0.400	0.399	0.3	106	0.00
24	Pentachlorophenol	0.400	0.418	-4.5	116	0.00
25	Phenanthrene	0.400	0.405	-1.3	109	0.00
26	Anthracene	0.400	0.399	0.3	107	-0.01
27 SURR	Fluoranthene-d10	0.400	0.402	-0.5	109	0.00
28	Fluoranthene	0.400	0.404	-1.0	110	0.00
29 I	Chrysene-d12	0.400	0.400	0.0	99	0.00
30	Pyrene	0.400	0.421	-5.2	109	0.00
31 S	Terphenyl-d14	0.400	0.425	-6.2	109	0.00
32	Benzo(a)anthracene	0.400	0.397	0.8	97	0.00
33	Chrysene	0.400	0.417	-4.2	104	0.00
34	Bis(2-ethylhexyl)phthalate	0.400	0.395	1.3	98	0.00
35	Indeno(1,2,3-cd)pyrene	0.400	0.468	-17.0	119	0.00
36 I	Perylene-d12	0.400	0.400	0.0	104	0.00
37	Benzo(b)fluoranthene	0.400	0.425	-6.2	107	0.00
38	Benzo(k)fluoranthene	0.400	0.416	-4.0	97	0.00
39 C	Benzo(a)pyrene	0.400	0.417	-4.2	108	0.00
40	Dibenzo(a,h)anthracene	0.400	0.445	-11.2	121	0.00
41	Benzo(g,h,i)perylene	0.400	0.435	-8.7	117	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN063020\
Data File : BN011080.D
Acq On : 30 Jun 2020 14:29
Operator : CG/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
ICVBN063020

Quant Time: Jun 30 15:24:38 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN063020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 30 14:12:13 2020
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	Amount	Calc.	%Dev	Area	% Dev(min)
----------	--------	-------	------	------	------------

(#) = Out of Range

SPCC's out = 0 CCC's out = 0