

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072220\
 Data File : BN011222.D
 Acq On : 21 Jul 2020 17:07
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 21 17:37:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN071120.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 21 14:35:20 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	91	0.00
2	1,4-Dioxane	0.400	0.348	13.0	73	0.00
3	n-Nitrosodimethylamine	0.400	0.365	8.8	87	0.00
4 S	2-Fluorophenol	0.400	0.363	9.3	79	0.00
5 S	Phenol-d6	0.400	0.399	0.3	88	0.00
6	bis(2-Chloroethyl)ether	0.400	0.409	-2.2	99	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	100	0.00
8 S	Nitrobenzene-d5	0.400	0.363	9.3	90	0.01
9	Naphthalene	0.400	0.374	6.5	94	0.00
10	Hexachlorobutadiene	0.400	0.363	9.3	88	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.411	-2.7	98	0.00
12	2-Methylnaphthalene	0.400	0.408	-2.0	98	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	118	0.00
14 S	2,4,6-Tribromophenol	0.400	0.371	7.3	107	0.00
15 S	2-Fluorobiphenyl	0.400	0.353	11.8	97	0.00
16	Acenaphthylene	0.400	0.360	10.0	109	0.01
17	Acenaphthene	0.400	0.372	7.0	110	0.00
18	Fluorene	0.400	0.378	5.5	98	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	135	0.00
20	4,6-Dinitro-2-methylphenol	0.400	0.376	6.0	113	0.00
21	4-Bromophenyl-phenylether	0.400	0.363	9.3	122	0.00
22	Hexachlorobenzene	0.400	0.341	14.8	113	0.00
23	Atrazine	0.400	0.383	4.3	129	-0.01
24	Pentachlorophenol	0.400	0.346	13.5	126	0.00
25	Phenanthrene	0.400	0.369	7.8	125	0.00
26	Anthracene	0.400	0.356	11.0	121	0.00
27 SURR	Fluoranthene-d10	0.400	0.367	8.3	122	0.00
28	Fluoranthene	0.400	0.361	9.8	116	0.00
29 I	Chrysene-d12	0.400	0.400	0.0	110	-0.01
30	Pyrene	0.400	0.453	-13.2	110	0.00
31 S	Terphenyl-d14	0.400	0.455	-13.7	118	0.00
32	Benzo(a)anthracene	0.400	0.370	7.5	104	0.00
33	Chrysene	0.400	0.407	-1.7	112	0.00
34	Bis(2-ethylhexyl)phthalate	0.400	0.477	-19.2	132	0.00
35	Indeno(1,2,3-cd)pyrene	0.400	0.352	12.0	84	0.00
36 I	Perylene-d12	0.400	0.400	0.0	104	0.00
37	Benzo(b)fluoranthene	0.400	0.349	12.8	101	0.00
38	Benzo(k)fluoranthene	0.400	0.380	5.0	99	-0.01
39 C	Benzo(a)pyrene	0.400	0.341	14.8	88	0.00
40	Dibenzo(a,h)anthracene	0.400	0.310	22.5	79	-0.01
41	Benzo(g,h,i)perylene	0.400	0.323	19.3	82	-0.01

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072220\
Data File : BN011222.D
Acq On : 21 Jul 2020 17:07
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_N
LabSampleId :
SSTDCCC0.4

Quant Time: Jul 21 17:37:40 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN071120.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 21 14:35:20 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