

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071919\
 Data File : BE100469.D
 Acq On : 19 Jul 2019 13:58
 Operator : HP/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE071919

Quant Time: Jul 19 14:35:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 19 13:59:01 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	117	0.00
2	1,4-Dioxane	0.400	0.403	-0.8	105	-0.01
3	n-Nitrosodimethylamine	0.400	0.380	5.0	114	-0.01
4 S	2-Fluorophenol	0.400	0.383	4.3	110	0.00
5 S	Phenol-d6	0.400	0.377	5.8	112	0.00
6	bis(2-Chloroethyl)ether	0.400	0.398	0.5	117	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	117	0.00
8 S	Nitrobenzene-d5	0.400	0.400	0.0	127	0.00
9	Naphthalene	0.400	0.394	1.5	117	0.00
10	Hexachlorobutadiene	0.400	0.387	3.3	114	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.384	4.0	117	0.00
12	2-Methylnaphthalene	0.400	0.391	2.3	118	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	121	0.00
14 S	2,4,6-Tribromophenol	0.400	0.376	6.0	125	0.00
15 S	2-Fluorobiphenyl	0.400	0.386	3.5	118	0.00
16	Acenaphthylene	0.400	0.377	5.8	118	0.00
17	Acenaphthene	0.400	0.393	1.8	121	0.00
18	Fluorene	0.400	0.382	4.5	119	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	119	0.00
20	4-Bromophenyl-phenylether	0.400	0.377	5.8	117	0.00
21	Hexachlorobenzene	0.400	0.381	4.8	118	-0.01
22	Pentachlorophenol	0.400	0.439	-9.7	126	0.00
23	Phenanthrene	0.400	0.380	5.0	120	0.00
24	Anthracene	0.400	0.384	4.0	123	0.00
25 SURR	Fluoranthene-d10	0.400	0.367	8.3	113	0.00
26	Fluoranthene	0.400	0.364	9.0	112	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	99	0.00
28	Pvrene	0.400	0.408	-2.0	110	0.00
29 S	Terphenyl-d14	0.400	0.400	0.0	109	0.00
30	Benzo(a)anthracene	0.400	0.387	3.3	99	0.00
31	Chrysene	0.400	0.393	1.8	98	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.371	7.3	93	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.390	2.5	95	0.00
34 I	Perylene-d12	0.400	0.400	0.0	96	0.00
35	Benzo(b)fluoranthene	0.400	0.380	5.0	95	0.00
36	Benzo(k)fluoranthene	0.400	0.382	4.5	93	0.00
37 C	Benzo(a)pyrene	0.400	0.386	3.5	96	0.00
38	Dibenzo(a,h)anthracene	0.400	0.386	3.5	95	0.00
39	Benzo(a,h,i)perylene	0.400	0.391	2.3	96	-0.01

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Compound	Amount	Calc.	%Dev Area	% Dev(min)

(#) = Out of Range	SPCC's out = 0 CCC's out = 0			