

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE112119\
 Data File : BE100754.D
 Acq On : 21 Nov 2019 21:35
 Operator : JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE112119

Quant Time: Nov 22 06:44:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE112119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 22 06:40:44 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	101	0.00
2	1,4-Dioxane	0.400	0.397	0.8	100	0.00
3	n-Nitrosodimethylamine	0.400	0.406	-1.5	88	0.00
4 S	2-Fluorophenol	0.400	0.397	0.8	102	0.00
5 S	Phenol-d6	0.400	0.397	0.8	104	0.00
6	bis(2-Chloroethyl)ether	0.400	0.393	1.8	107	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	103	0.00
8 S	Nitrobenzene-d5	0.400	0.467	-16.8	140	0.00
9	Naphthalene	0.400	0.409	-2.2	104	0.00
10	Hexachlorobutadiene	0.400	0.412	-3.0	106	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.401	-0.3	106	0.00
12	2-Methylnaphthalene	0.400	0.397	0.8	104	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	102	0.00
14 S	2,4,6-Tribromophenol	0.400	0.419	-4.7	120	0.00
15 S	2-Fluorobiphenyl	0.400	0.416	-4.0	106	0.00
16	Acenaphthylene	0.400	0.379	5.3	102	0.00
17	Acenaphthene	0.400	0.398	0.5	104	0.00
18	Fluorene	0.400	0.393	1.8	104	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	103	0.00
20	4-Bromophenyl-phenylether	0.400	0.393	1.8	102	0.00
21	Hexachlorobenzene	0.400	0.405	-1.3	104	0.00
22	Pentachlorophenol	0.400	0.454	-13.5	116	0.00
23	Phenanthrene	0.400	0.402	-0.5	104	0.00
24	Anthracene	0.400	0.376	6.0	103	0.00
25 SURR	Fluoranthene-d10	0.400	0.377	5.8	100	0.00
26	Fluoranthene	0.400	0.380	5.0	101	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	101	0.00
28	Pvrene	0.400	0.386	3.5	100	0.00
29 S	Terphenyl-d14	0.400	0.384	4.0	99	0.00
30	Benzo(a)anthracene	0.400	0.379	5.3	97	-0.01
31	Chrysene	0.400	0.404	-1.0	101	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.413	-3.2	114	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.377	5.8	97	0.00
34 I	Perylene-d12	0.400	0.400	0.0	102	0.00
35	Benzo(b)fluoranthene	0.400	0.373	6.8	96	0.00
36	Benzo(k)fluoranthene	0.400	0.380	5.0	100	0.00
37 C	Benzo(a)pyrene	0.400	0.364	9.0	97	0.00
38	Dibenzo(a,h)anthracene	0.400	0.337	15.8	96	0.00
39	Benzo(a,h,i)perylene	0.400	0.382	4.5	98	0.00

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

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