

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE052019\
 Data File : BE099666.D
 Acq On : 20 May 2019 18:37
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE052019

Quant Time: May 20 19:36:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE052019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 20 19:29:40 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	99	0.00
2	1,4-Dioxane	0.400	0.358	10.5	93	0.00
3	n-Nitrosodimethylamine	0.400	0.269	32.8#	76	0.00
4 S	2-Fluorophenol	0.400	0.343	14.2	97	0.00
5 S	Phenol-d6	0.400	0.348	13.0	104	0.00
6	bis(2-Chloroethyl)ether	0.400	0.348	13.0	98	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	98	0.00
8 S	Nitrobenzene-d5	0.400	0.340	15.0	97	0.00
9	Naphthalene	0.400	0.358	10.5	98	0.00
10	Hexachlorobutadiene	0.400	0.368	8.0	101	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.356	11.0	98	0.00
12	2-Methylnaphthalene	0.400	0.358	10.5	101	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	100	0.00
14 S	2,4,6-Tribromophenol	0.400	0.366	8.5	96	0.00
15 S	2-Fluorobiphenyl	0.400	0.344	14.0	97	0.00
16	Acenaphthylene	0.400	0.344	14.0	100	0.00
17	Acenaphthene	0.400	0.347	13.3	98	0.01
18	Fluorene	0.400	0.351	12.3	100	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	101	0.00
20	4-Bromophenyl-phenylether	0.400	0.341	14.8	99	0.00
21	Hexachlorobenzene	0.400	0.351	12.3	100	0.00
22	Pentachlorophenol	0.400	0.335	16.3	71	0.01
23	Phenanthrene	0.400	0.344	14.0	98	0.00
24	Anthracene	0.400	0.321	19.8	97	0.00
25 SURR	Fluoranthene-d10	0.400	0.345	13.8	100	0.00
26	Fluoranthene	0.400	0.350	12.5	101	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	102	0.00
28	Pvrene	0.400	0.384	4.0	100	0.00
29 S	Terphenyl-d14	0.400	0.375	6.3	99	0.00
30	Benzo(a)anthracene	0.400	0.368	8.0	98	0.00
31	Chrysene	0.400	0.391	2.3	101	0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.353	11.8	100	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.381	4.8	103	0.00
34 I	Perylene-d12	0.400	0.400	0.0	102	0.00
35	Benzo(b)fluoranthene	0.400	0.357	10.8	101	0.00
36	Benzo(k)fluoranthene	0.400	0.352	12.0	100	0.00
37 C	Benzo(a)pyrene	0.400	0.361	9.8	104	0.00
38	Dibenzo(a,h)anthracene	0.400	0.348	13.0	103	0.00
39	Benzo(a,h,i)perylene	0.400	0.356	11.0	102	0.00

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

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