

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE052319\
 Data File : BE099734.D
 Acq On : 23 May 2019 15:43
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: May 23 16:27:40 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	77	0.00
2	1,4-Dioxane	0.400	0.383	4.3	77	-0.01
3	n-Nitrosodimethylamine	0.400	0.451	-12.7	99	-0.02
4 S	2-Fluorophenol	0.400	0.449	-12.2	99	0.00
5 S	Phenol-d6	0.400	0.447	-11.7	104	0.00
6	bis(2-Chloroethyl)ether	0.400	0.412	-3.0	91	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	75	0.01
8 S	Nitrobenzene-d5	0.400	0.471	-17.7	103	0.00
9	Naphthalene	0.400	0.405	-1.3	85	0.01
10	Hexachlorobutadiene	0.400	0.406	-1.5	85	0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.415	-3.7	87	0.00
12	2-Methylnaphthalene	0.400	0.409	-2.2	88	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	71	0.00
14 S	2,4,6-Tribromophenol	0.400	0.528	-32.0#	118	-0.01
15 S	2-Fluorobiphenyl	0.400	0.428	-7.0	85	0.01
16	Acenaphthylene	0.400	0.432	-8.0	89	0.00
17	Acenaphthene	0.400	0.432	-8.0	86	0.00
18	Fluorene	0.400	0.412	-3.0	83	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	67	0.00
20	4-Bromophenyl-phenylether	0.400	0.406	-1.5	79	0.00
21	Hexachlorobenzene	0.400	0.410	-2.5	78	0.00
22	Pentachlorophenol	0.400	0.829	-107.2#	214	-0.01
23	Phenanthrene	0.400	0.405	-1.3	77	0.00
24	Anthracene	0.400	0.420	-5.0	85	0.00
25 SURR	Fluoranthene-d10	0.400	0.406	-1.5	78	0.00
26	Fluoranthene	0.400	0.403	-0.8	78	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	79	0.00
28	Pvrene	0.400	0.400	0.0	80	0.00
29 S	Terphenyl-d14	0.400	0.395	1.3	81	-0.01
30	Benzo(a)anthracene	0.400	0.449	-12.2	93	-0.01
31	Chrysene	0.400	0.402	-0.5	80	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.557	-39.3#	122	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.429	-7.2	90	0.00
34 I	Perylene-d12	0.400	0.400	0.0	76	0.00
35	Benzo(b)fluoranthene	0.400	0.414	-3.5	88	0.00
36	Benzo(k)fluoranthene	0.400	0.388	3.0	83	0.00
37 C	Benzo(a)pyrene	0.400	0.412	-3.0	89	0.00
38	Dibenzo(a,h)anthracene	0.400	0.408	-2.0	90	-0.01
39	Benzo(a,h,i)perylene	0.400	0.406	-1.5	88	0.00

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

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