

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE032019\
 Data File : BE099067.D
 Acq On : 20 Mar 2019 13:59
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 22 03:14:15 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 15 05:31:33 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	93	0.00
2	1,4-Dioxane	0.400	0.446	-11.5	104	0.00
3	n-Nitrosodimethylamine	0.400	0.317	20.8	73	0.03
4 S	2-Fluorophenol	0.400	0.386	3.5	87	0.01
5 S	Phenol-d6	0.400	0.420	-5.0	98	0.01
6	bis(2-Chloroethyl)ether	0.400	0.314	21.5	76	0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	109	-0.01
8 S	Nitrobenzene-d5	0.400	0.334	16.5	102	0.00
9	Naphthalene	0.400	0.379	5.3	106	-0.01
10	Hexachlorobutadiene	0.400	0.344	14.0	97	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.384	4.0	107	0.00
12	2-Methylnaphthalene	0.400	0.364	9.0	103	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	113	0.00
14 S	2,4,6-Tribromophenol	0.400	0.302	24.5	86	0.00
15 S	2-Fluorobiphenyl	0.400	0.383	4.3	109	-0.01
16	Acenaphthylene	0.400	0.370	7.5	110	-0.01
17	Acenaphthene	0.400	0.378	5.5	107	-0.01
18	Fluorene	0.400	0.357	10.8	103	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	100	-0.01
20	4-Bromophenyl-phenylether	0.400	0.364	9.0	92	0.00
21	Hexachlorobenzene	0.400	0.379	5.3	100	-0.01
22	Pentachlorophenol	0.400	0.403	-0.8	79	0.00
23	Phenanthrene	0.400	0.356	11.0	91	-0.01
24	Anthracene	0.400	0.326	18.5	86	0.00
25 SURR	Fluoranthene-d10	0.400	0.328	18.0	85	0.00
26	Fluoranthene	0.400	0.330	17.5	85	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	85	-0.01
28	Pvrene	0.400	0.400	0.0	86	0.00
29 S	Terphenyl-d14	0.400	0.374	6.5	80	0.00
30	Benzo(a)anthracene	0.400	0.336	16.0	73	0.00
31	Chrysene	0.400	0.369	7.8	81	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.358	10.5	83	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.271	32.3#	59	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	83	-0.01
35	Benzo(b)fluoranthene	0.400	0.347	13.3	75	0.00
36	Benzo(k)fluoranthene	0.400	0.403	-0.8	88	-0.01
37 C	Benzo(a)pyrene	0.400	0.380	5.0	82	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.244	39.0#	51	-0.01
39	Benzo(a,h,i)perylene	0.400	0.285	28.8#	61	0.00

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

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