

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031119\
 Data File : BE098888.D
 Acq On : 12 Mar 2019 8:24
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 12 16:39:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 06 15:33:24 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	115	0.00
2	1,4-Dioxane	0.400	0.320	20.0	99	0.00
3	n-Nitrosodimethylamine	0.400	0.178	55.5#	50	0.01
4 S	2-Fluorophenol	0.400	0.329	17.8	91	0.01
5 S	Phenol-d6	0.400	0.369	7.8	106	0.00
6	bis(2-Chloroethyl)ether	0.400	0.321	19.8	96	0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	119	0.00
8 S	Nitrobenzene-d5	0.400	0.336	16.0	111	0.00
9	Naphthalene	0.400	0.399	0.3	122	-0.01
10	Hexachlorobutadiene	0.400	0.381	4.8	117	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.399	0.3	121	0.00
12	2-Methylnaphthalene	0.400	0.389	2.8	120	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	127	0.00
14 S	2,4,6-Tribromophenol	0.400	0.329	17.8	105	-0.01
15 S	2-Fluorobiphenyl	0.400	0.390	2.5	125	0.00
16	Acenaphthylene	0.400	0.370	7.5	123	0.00
17	Acenaphthene	0.400	0.381	4.8	122	-0.01
18	Fluorene	0.400	0.371	7.3	121	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	116	0.00
20	4-Bromophenyl-phenylether	0.400	0.363	9.3	107	0.00
21	Hexachlorobenzene	0.400	0.397	0.8	121	0.00
22	Pentachlorophenol	0.400	0.435	-8.7	112	0.00
23	Phenanthrene	0.400	0.384	4.0	114	0.00
24	Anthracene	0.400	0.376	6.0	115	0.00
25 SURR	Fluoranthene-d10	0.400	0.361	9.8	108	0.00
26	Fluoranthene	0.400	0.360	10.0	108	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	239	-0.01
28	Pvrene	0.400	0.182	54.5#	109	0.00
29 S	Terphenyl-d14	0.400	0.335	16.3	202	0.00
30	Benzo(a)anthracene	0.400	0.306	23.5	187	0.00
31	Chrysene	0.400	0.343	14.2	210	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.326	18.5	213	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.347	13.3	212	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	236	-0.01
35	Benzo(b)fluoranthene	0.400	0.358	10.5	218	0.00
36	Benzo(k)fluoranthene	0.400	0.378	5.5	234	-0.01
37 C	Benzo(a)pyrene	0.400	0.339	15.3	206	-0.02
38	Dibenzo(a,h)anthracene	0.400	0.363	9.3	217	-0.02
39	Benzo(a,h,i)perylene	0.400	0.360	10.0	217	-0.02

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

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