

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE072318\
 Data File : BE097062.D
 Acq On : 24 Jul 2018 7:37
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 24 08:13:49 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 20 11:34:20 2018
 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	78	0.00
2	1,4-Dioxane	0.400	0.427	-6.7	86	0.00
3	n-Nitrosodimethylamine	0.400	0.434	-8.5	90	0.00
4 S	2-Fluorophenol	0.400	0.508	-27.0#	112	0.00
5 S	Phenol-d6	0.400	0.504	-26.0#	107	0.00
6	bis(2-Chloroethyl)ether	0.400	0.410	-2.5	82	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	72	0.00
8 S	Nitrobenzene-d5	0.400	0.485	-21.2	94	0.00
9	Naphthalene	0.400	0.397	0.8	75	0.00
10	Hexachlorobutadiene	0.400	0.431	-7.7	82	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.398	0.5	76	0.00
12	2-Methylnaphthalene	0.400	0.399	0.3	75	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	88	0.00
14 S	2,4,6-Tribromophenol	0.400	0.615	-53.7#	155	0.00
15 S	2-Fluorobiphenyl	0.400	0.360	10.0	81	0.00
16	Acenaphthylene	0.400	0.379	5.3	87	0.00
17	Acenaphthene	0.400	0.372	7.0	86	0.00
18	Fluorene	0.400	0.429	-7.2	100	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	104	0.00
20	4-Bromophenyl-phenylether	0.400	0.390	2.5	107	0.00
21	Hexachlorobenzene	0.400	0.380	5.0	102	0.00
22	Pentachlorophenol	0.400	0.772	-93.0#	233	0.00
23	Phenanthrene	0.400	0.394	1.5	107	0.00
24	Anthracene	0.400	0.404	-1.0	112	-0.01
25 SURR	Fluoranthene-d10	0.400	0.458	-14.5	129	0.00
26	Fluoranthene	0.400	0.423	-5.7	117	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	131	0.00
28	Pvrene	0.400	0.363	9.3	123	0.00
29 S	Terphenyl-d14	0.400	0.387	3.3	133	0.00
30	Benzo(a)anthracene	0.400	0.425	-6.2	152	0.00
31	Chrysene	0.400	0.399	0.3	134	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.786	-96.5#	296	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.558	-39.5#	205	0.00
34 I	Perylene-d12	0.400	0.400	0.0	154	0.00
35	Benzo(b)fluoranthene	0.400	0.381	4.8	158	0.00
36	Benzo(k)fluoranthene	0.400	0.370	7.5	149	0.00
37 C	Benzo(a)pyrene	0.400	0.432	-8.0	179	0.00
38	Dibenzo(a,h)anthracene	0.400	0.488	-22.0	217	0.00
39	Benzo(a,h,i)perylene	0.400	0.462	-15.5	197	0.00

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

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