

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE072318\
 Data File : BE097046.D
 Acq On : 23 Jul 2018 20:58
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 24 05:28:06 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	70	0.00
2	1,4-Dioxane	0.400	0.438	-9.5	79	0.00
3	n-Nitrosodimethylamine	0.400	0.449	-12.2	83	0.00
4 S	2-Fluorophenol	0.400	0.494	-23.5	98	0.00
5 S	Phenol-d6	0.400	0.474	-18.5	90	0.00
6	bis(2-Chloroethyl)ether	0.400	0.410	-2.5	74	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	64	0.00
8 S	Nitrobenzene-d5	0.400	0.481	-20.2	83	0.00
9	Naphthalene	0.400	0.397	0.8	67	0.00
10	Hexachlorobutadiene	0.400	0.434	-8.5	73	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.396	1.0	67	0.00
12	2-Methylnaphthalene	0.400	0.401	-0.3	68	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	76	0.00
14 S	2,4,6-Tribromophenol	0.400	0.594	-48.5#	129	0.00
15 S	2-Fluorobiphenyl	0.400	0.364	9.0	71	0.00
16	Acenaphthylene	0.400	0.392	2.0	78	0.00
17	Acenaphthene	0.400	0.375	6.3	75	0.00
18	Fluorene	0.400	0.431	-7.7	87	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	90	0.00
20	4-Bromophenyl-phenylether	0.400	0.403	-0.8	96	0.00
21	Hexachlorobenzene	0.400	0.386	3.5	90	0.00
22	Pentachlorophenol	0.400	0.782	-95.5#	206	0.00
23	Phenanthrene	0.400	0.394	1.5	93	0.00
24	Anthracene	0.400	0.401	-0.3	97	0.00
25 SURR	Fluoranthene-d10	0.400	0.458	-14.5	112	0.00
26	Fluoranthene	0.400	0.423	-5.7	102	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	110	0.00
28	Pvrene	0.400	0.376	6.0	107	0.00
29 S	Terphenyl-d14	0.400	0.393	1.8	113	0.00
30	Benzo(a)anthracene	0.400	0.408	-2.0	123	0.00
31	Chrysene	0.400	0.398	0.5	112	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.578	-44.5#	183	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.546	-36.5#	168	0.00
34 I	Perylene-d12	0.400	0.400	0.0	125	0.00
35	Benzo(b)fluoranthene	0.400	0.380	5.0	128	0.00
36	Benzo(k)fluoranthene	0.400	0.352	12.0	115	0.00
37 C	Benzo(a)pyrene	0.400	0.410	-2.5	136	0.00
38	Dibenzo(a,h)anthracene	0.400	0.498	-24.5	180	0.00
39	Benzo(a,h,i)perylene	0.400	0.472	-18.0	163	0.00

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

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