

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE100818\
 Data File : Be097810.d
 Acq On : 9 Oct 2018 1:30
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 09 07:00:19 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE100318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 08 06:37:06 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	76	-0.01
2	1,4-Dioxane	0.797	0.740	7.2	75	0.00
3	n-Nitrosodimethylamine	0.725	0.615	15.2	68	0.00
4 S	2-Fluorophenol	0.946	0.816	13.7	69	-0.01
5 S	Phenol-d6	1.450	1.223	15.7	65	0.00
6	bis(2-Chloroethyl)ether	1.501	1.394	7.1	73	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	78	-0.01
8 S	Nitrobenzene-d5	0.130	0.214	-64.6#	137	-0.01
9	Naphthalene	4.031	3.828	5.0	76	-0.01
10	Hexachlorobutadiene	0.214	0.211	1.4	79	-0.01
11 SURR	2-Methylnaphthalene-d10	0.636	0.621	2.4	78	-0.02
12	2-Methylnaphthalene	0.645	0.608	5.7	76	-0.02
13 I	Acenaphthene-d10	1.000	1.000	0.0	82	-0.02
14 S	2,4,6-Tribromophenol	0.072	0.083	-15.3	117	-0.01
15 S	2-Fluorobiphenyl	1.761	1.602	9.0	79	-0.02
16	Acenaphthylene	1.771	1.549	12.5	76	0.00
17	Acenaphthene	1.135	1.058	6.8	81	0.00
18	Fluorene	1.507	1.394	7.5	80	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	82	-0.01
20	4-Bromophenyl-phenylether	0.210	0.197	6.2	82	-0.01
21	Hexachlorobenzene	0.236	0.210	11.0	77	-0.01
22	Pentachlorophenol	0.037	0.045	-21.6	111	-0.01
23	Phenanthrene	1.031	0.953	7.6	79	-0.01
24	Anthracene	0.890	0.770	13.5	77	-0.01
25 SURR	Fluoranthene-d10	5.063	4.644	8.3	81	-0.01
26	Fluoranthene	1.313	1.187	9.6	79	-0.01
27 I	Chrysene-d12	1.000	1.000	0.0	73	-0.01
28	Pvrene	1.016	1.024	-0.8	81	-0.01
29 S	Terphenyl-d14	0.817	0.823	-0.7	82	-0.01
30	Benzo(a)anthracene	1.075	0.964	10.3	73	-0.01
31	Chrysene	1.202	1.147	4.6	75	-0.01
32	Bis(2-ethylhexyl)phthalate	0.787	0.782	0.6	84	-0.01
33	Indeno(1,2,3-cd)pyrene	1.044	0.915	12.4	69	-0.03
34 I	Perylene-d12	1.000	1.000	0.0	74	-0.02
35	Benzo(b)fluoranthene	1.098	0.960	12.6	73	-0.02
36	Benzo(k)fluoranthene	1.217	1.166	4.2	79	-0.02
37 C	Benzo(a)pyrene	0.994	0.832	16.3	69	-0.02
38	Dibenzo(a,h)anthracene	0.979	0.803	18.0	68	-0.03
39	Benzo(a,h,i)perylene	1.024	0.923	9.9	73	-0.03

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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