

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE100318\
 Data File : BE097718.D
 Acq On : 3 Oct 2018 19:16
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 04 04:43:12 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE100318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 03 12:17:19 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	93	0.00
2	1,4-Dioxane	0.797	0.775	2.8	97	0.00
3	n-Nitrosodimethylamine	0.725	0.671	7.4	91	0.00
4 S	2-Fluorophenol	0.946	0.919	2.9	95	0.00
5 S	Phenol-d6	1.450	1.357	6.4	89	0.00
6	bis(2-Chloroethyl)ether	1.501	1.398	6.9	90	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
8 S	Nitrobenzene-d5	0.130	0.149	-14.6	114	0.00
9	Naphthalene	4.031	3.821	5.2	90	0.00
10	Hexachlorobutadiene	0.214	0.201	6.1	91	0.00
11 SURR	2-Methylnaphthalene-d10	0.636	0.596	6.3	89	0.00
12	2-Methylnaphthalene	0.645	0.618	4.2	93	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	94	0.00
14 S	2,4,6-Tribromophenol	0.072	0.078	-8.3	124	0.00
15 S	2-Fluorobiphenyl	1.761	1.647	6.5	93	-0.01
16	Acenaphthylene	1.771	1.579	10.8	88	0.00
17	Acenaphthene	1.135	1.080	4.8	94	0.00
18	Fluorene	1.507	1.475	2.1	96	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	92	0.00
20	4-Bromophenyl-phenylether	0.210	0.200	4.8	94	0.00
21	Hexachlorobenzene	0.236	0.230	2.5	95	0.00
22	Pentachlorophenol	0.037	0.040	-8.1	112	0.00
23	Phenanthrene	1.031	0.968	6.1	91	0.00
24	Anthracene	0.890	0.812	8.8	92	0.00
25 SURR	Fluoranthene-d10	5.063	4.664	7.9	92	0.00
26	Fluoranthene	1.313	1.208	8.0	91	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	91	0.00
28	Pvrene	1.016	0.915	9.9	91	0.00
29 S	Terphenyl-d14	0.817	0.734	10.2	92	0.00
30	Benzo(a)anthracene	1.075	0.961	10.6	91	0.00
31	Chrysene	1.202	1.122	6.7	92	0.00
32	Bis(2-ethylhexyl)phthalate	0.787	0.755	4.1	101	0.00
33	Indeno(1,2,3-cd)pyrene	1.044	1.021	2.2	97	0.00
34 I	Perylene-d12	1.000	1.000	0.0	100	0.00
35	Benzo(b)fluoranthene	1.098	0.896	18.4	92	0.00
36	Benzo(k)fluoranthene	1.217	1.068	12.2	98	0.00
37 C	Benzo(a)pyrene	0.994	0.843	15.2	95	0.00
38	Dibenzo(a,h)anthracene	0.979	0.853	12.9	98	0.00
39	Benzo(a,h,i)perylene	1.024	0.895	12.6	96	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE100318\
Data File : BE097718.D
Acq On : 3 Oct 2018 19:16
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Oct 04 04:43:12 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE100318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 03 12:17:19 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	AvgRF	CCRF	%Dev Area	% Dev(min)

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