

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE100818\
 Data File : Be097802.d
 Acq On : 8 Oct 2018 20:34
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC00.4

Quant Time: Oct 09 06:58:59 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	66	-0.01
2	1,4-Dioxane	0.797	0.741	7.0	66	0.00
3	n-Nitrosodimethylamine	0.725	0.626	13.7	61	0.00
4 S	2-Fluorophenol	0.946	0.877	7.3	65	-0.01
5 S	Phenol-d6	1.450	1.285	11.4	60	-0.01
6	bis(2-Chloroethyl)ether	1.501	1.406	6.3	65	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	69	0.00
8 S	Nitrobenzene-d5	0.130	0.219	-68.5#	125	0.00
9	Naphthalene	4.031	3.822	5.2	67	-0.01
10	Hexachlorobutadiene	0.214	0.200	6.5	67	-0.01
11 SURR	2-Methylnaphthalene-d10	0.636	0.600	5.7	67	-0.02
12	2-Methylnaphthalene	0.645	0.620	3.9	69	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	72	-0.02
14 S	2,4,6-Tribromophenol	0.072	0.080	-11.1	99	-0.01
15 S	2-Fluorobiphenyl	1.761	1.621	8.0	70	-0.02
16	Acenaphthylene	1.771	1.575	11.1	67	0.00
17	Acenaphthene	1.135	1.068	5.9	72	0.00
18	Fluorene	1.507	1.412	6.3	71	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	72	-0.01
20	4-Bromophenyl-phenylether	0.210	0.193	8.1	71	-0.01
21	Hexachlorobenzene	0.236	0.212	10.2	68	-0.01
22	Pentachlorophenol	0.037	0.052	-40.5#	114	-0.01
23	Phenanthrene	1.031	0.964	6.5	71	0.00
24	Anthracene	0.890	0.785	11.8	70	-0.01
25 SURR	Fluoranthene-d10	5.063	4.670	7.8	72	-0.01
26	Fluoranthene	1.313	1.205	8.2	71	-0.01
27 I	Chrysene-d12	1.000	1.000	0.0	65	-0.01
28	Pvrene	1.016	1.019	-0.3	72	-0.01
29 S	Terphenyl-d14	0.817	0.824	-0.9	73	-0.01
30	Benzo(a)anthracene	1.075	0.963	10.4	65	-0.01
31	Chrysene	1.202	1.153	4.1	67	-0.01
32	Bis(2-ethylhexyl)phthalate	0.787	0.746	5.2	71	-0.01
33	Indeno(1,2,3-cd)pyrene	1.044	0.901	13.7	61	-0.03
34 I	Perylene-d12	1.000	1.000	0.0	66	-0.02
35	Benzo(b)fluoranthene	1.098	0.949	13.6	64	-0.02
36	Benzo(k)fluoranthene	1.217	1.166	4.2	71	-0.01
37 C	Benzo(a)pyrene	0.994	0.857	13.8	64	-0.02
38	Dibenzo(a,h)anthracene	0.979	0.813	17.0	61	-0.03
39	Benzo(a,h,i)perylene	1.024	0.920	10.2	65	-0.03

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE100818\
Data File : Be097802.d
Acq On : 8 Oct 2018 20:34
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Oct 09 06:58:59 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	AvgRF	CCRF	%Dev Area	% Dev(min)

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