

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE091718\
 Data File : BE097497.D
 Acq On : 17 Sep 2018 15:52
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 18 05:59:24 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE091718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 17 18:57:11 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)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	95	0.00
2	1,4-Dioxane	0.692	0.738	-6.6	107	0.00
3	n-Nitrosodimethylamine	0.578	0.525	9.2	106	0.00
4 S	2-Fluorophenol	1.355	1.322	2.4	105	0.00
5 S	Phenol-d6	1.719	1.625	5.5	112	0.00
6	bis(2-Chloroethyl)ether	1.374	1.422	-3.5	106	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
8 S	Nitrobenzene-d5	0.322	0.303	5.9	110	0.00
9	Naphthalene	3.631	3.714	-2.3	107	0.00
10	Hexachlorobutadiene	0.172	0.174	-1.2	105	0.00
11 SURR	2-Methylnaphthalene-d10	0.572	0.574	-0.3	106	0.00
12	2-Methylnaphthalene	0.581	0.597	-2.8	110	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
14 S	2,4,6-Tribromophenol	0.211	0.176	16.6	100	0.00
15 S	2-Fluorobiphenyl	1.397	1.458	-4.4	109	0.00
16	Acenaphthylene	1.740	1.666	4.3	104	0.00
17	Acenaphthene	1.074	1.122	-4.5	109	0.00
18	Fluorene	1.369	1.420	-3.7	109	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	97	-0.01
20	4-Bromophenyl-phenylether	0.193	0.190	1.6	108	0.00
21	Hexachlorobenzene	0.205	0.208	-1.5	107	0.01
22	Pentachlorophenol	0.036	0.028	22.2	128	0.00
23	Phenanthrene	0.941	0.981	-4.3	109	0.00
24	Anthracene	0.869	0.850	2.2	105	0.00
25 SURR	Fluoranthene-d10	4.806	4.812	-0.1	107	0.00
26	Fluoranthene	1.227	1.229	-0.2	106	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	94	0.00
28	Pyrene	1.029	1.067	-3.7	106	0.00
29 S	Terphenyl-d14	0.765	0.793	-3.7	109	0.00
30	Benzo(a)anthracene	1.052	1.031	2.0	104	0.00
31	Chrysene	1.079	1.157	-7.2	105	0.00
32	Bis(2-ethylhexyl)phthalate	2.235	2.166	3.1	96	0.00
33	Indeno(1,2,3-cd)pyrene	1.286	1.363	-6.0	107	0.00
34 I	Perylene-d12	1.000	1.000	0.0	96	0.00
35	Benzo(b)fluoranthene	1.023	1.019	0.4	106	0.00
36	Benzo(k)fluoranthene	1.115	1.163	-4.3	107	0.00
37 C	Benzo(a)pyrene	1.016	1.047	-3.1	107	0.00
38	Dibenzo(a,h)anthracene	1.060	1.101	-3.9	109	0.00
39	Benzo(g,h,i)perylene	1.082	1.101	-1.8	105	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE091718\
Data File : BE097497.D
Acq On : 17 Sep 2018 15:52
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Sep 18 05:59:24 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE091718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 17 18:57:11 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		