

Data Path : Z:\HPCHEM1\BNA_E\Data\BE091117\
 Data File : BE093973.D
 Acq On : 11 Sep 2017 21:28
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 12 08:26:40 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 05 13:34:56 2017
 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	106	0.00
2	1,4-Dioxane	0.551	0.595	-8.0	116	0.00
3	n-Nitrosodimethylamine	0.646	0.621	3.9	102	0.00
4 S	2-Fluorophenol	1.294	1.278	1.2	106	0.00
5 S	Phenol-d6	1.941	1.887	2.8	106	0.00
6	bis(2-Chloroethyl)ether	1.587	1.577	0.6	106	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
8 S	Nitrobenzene-d5	0.338	0.334	1.2	106	0.00
9	Naphthalene	4.585	4.575	0.2	106	0.00
10	Hexachlorobutadiene	0.167	0.160	4.2	102	0.00
11 SURR	2-Methylnaphthalene-d10	0.640	0.635	0.8	104	0.00
12	2-Methylnaphthalene	0.762	0.751	1.4	104	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
14 S	2,4,6-Tribromophenol	0.124	0.126	-1.6	103	0.00
15 S	2-Fluorobiphenyl	1.582	1.577	0.3	102	0.00
16	Acenaphthylene	1.997	1.971	1.3	107	0.00
17	Acenaphthene	1.349	1.349	0.0	104	0.00
18	Fluorene	1.741	1.711	1.7	105	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
20	4-Bromophenyl-phenylether	0.219	0.255	-16.4	96	0.00
21	Hexachlorobenzene	0.236	0.219	7.2	85	0.00
22	Pentachlorophenol	0.063	0.074	-17.5	114	0.00
23	Phenanthrene	1.293	1.415	-9.4	94	0.00
24	Anthracene	1.218	1.269	-4.2	100	0.00
25 SURR	Fluoranthene-d10	5.960	5.950	0.2	100	0.00
26	Fluoranthene	1.783	1.787	-0.2	102	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	98	-0.01
28	Pyrene	1.376	1.388	-0.9	100	0.00
29 S	Terphenyl-d14	0.736	0.739	-0.4	99	0.00
30	Benzo(a)anthracene	1.229	1.164	5.3	93	0.00
31	Chrysene	1.300	1.363	-4.8	104	0.00
32	Bis(2-ethylhexyl)phthalate	2.983	2.659	10.9	92	0.00
33	Indeno(1,2,3-cd)pyrene	0.867	0.761	12.2	88	0.00
34 I	Perylene-d12	1.000	1.000	0.0	96	0.00
35	Benzo(b)fluoranthene	1.274	1.279	-0.4	96	0.00
36	Benzo(k)fluoranthene	1.561	1.554	0.4	96	0.00
37 C	Benzo(a)pyrene	1.297	1.315	-1.4	98	0.00
38	Dibenzo(a,h)anthracene	0.860	0.813	5.5	91	0.00
39	Benzo(g,h,i)perylene	0.907	0.869	4.2	92	0.00

Data Path : Z:\HPCHEM1\BNA_E\Data\BE091117\
Data File : BE093973.D
Acq On : 11 Sep 2017 21:28
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Sep 12 08:26:40 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 05 13:34:56 2017
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		