

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE082818\
 Data File : BE097374.D
 Acq On : 28 Aug 2018 11:16
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 28 12:17:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE082218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 22 18:07:26 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	89	0.00
2	1,4-Dioxane	0.711	0.674	5.2	87	0.00
3	n-Nitrosodimethylamine	0.664	0.610	8.1	87	0.00
4 S	2-Fluorophenol	1.183	1.153	2.5	97	0.00
5 S	Phenol-d6	1.581	1.484	6.1	93	-0.01
6	bis(2-Chloroethyl)ether	1.455	1.393	4.3	91	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
8 S	Nitrobenzene-d5	0.287	0.293	-2.1	101	0.00
9	Naphthalene	3.641	3.794	-4.2	97	0.00
10	Hexachlorobutadiene	0.163	0.168	-3.1	96	0.00
11 SURR	2-Methylnaphthalene-d10	0.558	0.569	-2.0	97	0.00
12	2-Methylnaphthalene	0.574	0.583	-1.6	96	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	88	-0.01
14 S	2,4,6-Tribromophenol	0.135	0.133	1.5	105	0.00
15 S	2-Fluorobiphenyl	1.468	1.574	-7.2	99	0.00
16	Acenaphthylene	1.591	1.642	-3.2	102	0.00
17	Acenaphthene	1.061	1.112	-4.8	101	0.00
18	Fluorene	1.342	1.427	-6.3	102	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	91	0.00
20	4-Bromophenyl-phenylether	0.182	0.187	-2.7	104	0.00
21	Hexachlorobenzene	0.201	0.215	-7.0	106	0.00
22	Pentachlorophenol	0.055	0.049	10.9	106	0.00
23	Phenanthrene	0.952	0.998	-4.8	104	0.00
24	Anthracene	0.810	0.826	-2.0	104	0.00
25 SURR	Fluoranthene-d10	4.590	5.079	-10.7	111	0.00
26	Fluoranthene	1.193	1.326	-11.1	112	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	91	0.00
28	Pvrene	0.952	1.055	-10.8	109	0.00
29 S	Terphenyl-d14	0.738	0.787	-6.6	105	0.00
30	Benzo(a)anthracene	1.012	0.988	2.4	98	0.00
31	Chrysene	1.089	1.136	-4.3	101	0.00
32	Bis(2-ethylhexyl)phthalate	1.361	1.109	18.5	97	0.00
33	Indeno(1,2,3-cd)pyrene	1.546	1.175	24.0	87	0.00
34 I	Perylene-d12	1.000	1.000	0.0	90	0.00
35	Benzo(b)fluoranthene	1.144	1.035	9.5	97	0.00
36	Benzo(k)fluoranthene	1.264	1.199	5.1	100	0.00
37 C	Benzo(a)pyrene	1.032	0.997	3.4	101	0.00
38	Dibenzo(a,h)anthracene	1.356	1.030	24.0	85	0.00
39	Benzo(a,h,i)perylene	1.400	1.069	23.6	86	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE082818\
Data File : BE097374.D
Acq On : 28 Aug 2018 11:16
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Aug 28 12:17:07 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE082218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 22 18:07:26 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			