

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE081018\
 Data File : BE097255.D
 Acq On : 10 Aug 2018 9:01
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 10 15:56:29 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE072418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 31 10:49:28 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	85	0.00
2	1,4-Dioxane	0.636	0.522	17.9	72	0.00
3	n-Nitrosodimethylamine	0.711	0.547	23.1	70	0.00
4 S	2-Fluorophenol	1.124	1.164	-3.6	93	0.00
5 S	Phenol-d6	1.649	1.780	-7.9	98	0.00
6	bis(2-Chloroethyl)ether	1.506	1.379	8.4	84	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
8 S	Nitrobenzene-d5	0.347	0.359	-3.5	96	0.00
9	Naphthalene	3.590	3.459	3.6	89	-0.01
10	Hexachlorobutadiene	0.161	0.159	1.2	91	0.00
11 SURR	2-Methylnaphthalene-d10	0.599	0.549	8.3	85	0.00
12	2-Methylnaphthalene	0.616	0.573	7.0	86	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	80	0.00
14 S	2,4,6-Tribromophenol	0.132	0.127	3.8	85	0.00
15 S	2-Fluorobiphenyl	1.381	1.364	1.2	81	0.00
16	Acenaphthylene	1.795	1.730	3.6	79	0.00
17	Acenaphthene	1.044	0.992	5.0	78	-0.01
18	Fluorene	1.379	1.207	12.5	73	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	65	0.00
20	4-Bromophenyl-phenylether	0.194	0.201	-3.6	70	0.00
21	Hexachlorobenzene	0.209	0.227	-8.6	73	0.00
22	Pentachlorophenol	0.057	0.048	15.8	69	0.01
23	Phenanthrene	0.953	0.919	3.6	66	-0.01
24	Anthracene	0.853	0.832	2.5	68	0.00
25 SURR	Fluoranthene-d10	4.003	3.935	1.7	70	0.00
26	Fluoranthene	1.070	1.048	2.1	68	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	69	-0.01
28	Pyrene	1.097	1.069	2.6	71	0.00
29 S	Terphenyl-d14	0.768	0.720	6.3	69	0.00
30	Benzo(a)anthracene	0.995	0.977	1.8	76	0.00
31	Chrysene	1.067	1.027	3.7	69	0.00
32	Bis(2-ethylhexyl)phthalate	1.243	1.519	-22.2	103	0.00
33	Indeno(1,2,3-cd)pyrene	0.966	1.228	-27.1#	94	0.00
34 I	Perylene-d12	1.000	1.000	0.0	85	0.00
35	Benzo(b)fluoranthene	1.015	0.916	9.8	81	0.00
36	Benzo(k)fluoranthene	1.185	1.098	7.3	85	0.00
37 C	Benzo(a)pyrene	0.955	0.946	0.9	89	0.00
38	Dibenzo(a,h)anthracene	0.953	1.001	-5.0	94	0.00
39	Benzo(g,h,i)perylene	1.011	0.987	2.4	87	0.01

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE081018\
Data File : BE097255.D
Acq On : 10 Aug 2018 9:01
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Aug 10 15:56:29 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE072418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 31 10:49:28 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		