

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE030719\
 Data File : BE098811.D
 Acq On : 7 Mar 2019 9:24
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 07 12:58:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 06 15:33:24 2019
 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	122	0.00
2	1,4-Dioxane	0.789	0.810	-2.7	130	0.00
3	n-Nitrosodimethylamine	1.006	0.835	17.0	99	0.00
4 S	2-Fluorophenol	1.196	1.169	2.3	115	0.00
5 S	Phenol-d6	1.689	1.615	4.4	117	0.00
6	bis(2-Chloroethyl)ether	1.304	1.136	12.9	110	0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	121	0.00
8 S	Nitrobenzene-d5	0.419	0.365	12.9	117	0.00
9	Naphthalene	4.581	4.569	0.3	123	0.00
10	Hexachlorobutadiene	0.303	0.313	-3.3	129	0.00
11 SURR	2-Methylnaphthalene-d10	0.739	0.715	3.2	119	0.00
12	2-Methylnaphthalene	0.743	0.707	4.8	118	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	126	0.00
14 S	2,4,6-Tribromophenol	0.202	0.197	2.5	124	0.00
15 S	2-Fluorobiphenyl	1.656	1.610	2.8	124	0.00
16	Acenaphthylene	2.057	1.983	3.6	127	0.00
17	Acenaphthene	1.220	1.190	2.5	124	0.00
18	Fluorene	1.760	1.683	4.4	124	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	120	0.00
20	4-Bromophenyl-phenylether	0.217	0.206	5.1	116	0.00
21	Hexachlorobenzene	0.258	0.251	2.7	123	0.00
22	Pentachlorophenol	0.052	0.044	15.4	138	0.00
23	Phenanthrene	1.137	1.114	2.0	120	0.00
24	Anthracene	0.972	0.880	9.5	115	0.00
25 SURR	Fluoranthene-d10	6.385	6.277	1.7	122	0.00
26	Fluoranthene	1.605	1.565	2.5	121	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	124	0.00
28	Pvrene	1.224	1.221	0.2	125	0.00
29 S	Terphenyl-d14	0.972	0.944	2.9	122	0.00
30	Benzo(a)anthracene	1.248	1.254	-0.5	128	0.00
31	Chrysene	1.336	1.314	1.6	125	0.00
32	Bis(2-ethylhexyl)phthalate	2.657	2.406	9.4	123	0.00
33	Indeno(1,2,3-cd)pyrene	1.410	1.452	-3.0	130	0.00
34 I	Perylene-d12	1.000	1.000	0.0	131	0.00
35	Benzo(b)fluoranthene	1.315	1.282	2.5	132	0.00
36	Benzo(k)fluoranthene	1.377	1.382	-0.4	138	0.00
37 C	Benzo(a)pyrene	1.261	1.229	2.5	132	0.00
38	Dibenzo(a,h)anthracene	1.193	1.164	2.4	130	0.00
39	Benzo(a,h,i)perylene	1.248	1.195	4.2	128	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE030719\
Data File : BE098811.D
Acq On : 7 Mar 2019 9:24
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Mar 07 12:58:00 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 06 15:33:24 2019
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			