

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE111318\
 Data File : BE098212.D
 Acq On : 13 Nov 2018 16:58
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 13 17:33:48 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE111218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 12 13:32:15 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	107	0.00
2	1,4-Dioxane	0.650	0.591	9.1	110	0.00
3	n-Nitrosodimethylamine	0.785	0.723	7.9	106	0.00
4 S	2-Fluorophenol	1.129	1.147	-1.6	111	0.00
5 S	Phenol-d6	1.814	1.921	-5.9	116	0.00
6	bis(2-Chloroethyl)ether	1.464	1.431	2.3	112	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	109	0.00
8 S	Nitrobenzene-d5	0.399	0.395	1.0	114	0.00
9	Naphthalene	4.068	3.932	3.3	109	0.00
10	Hexachlorobutadiene	0.253	0.240	5.1	107	0.00
11 SURR	2-Methylnaphthalene-d10	0.702	0.692	1.4	111	0.00
12	2-Methylnaphthalene	0.706	0.713	-1.0	114	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	116	0.00
14 S	2,4,6-Tribromophenol	0.174	0.162	6.9	125	0.00
15 S	2-Fluorobiphenyl	1.675	1.501	10.4	110	0.00
16	Acenaphthylene	1.834	1.788	2.5	117	0.00
17	Acenaphthene	1.120	1.075	4.0	115	0.00
18	Fluorene	1.434	1.415	1.3	118	0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	116	0.00
20	4-Bromophenyl-phenylether	0.244	0.226	7.4	115	0.00
21	Hexachlorobenzene	0.254	0.240	5.5	113	0.01
22	Pentachlorophenol	0.044	0.056	-27.3#	160#	0.00
23	Phenanthrene	0.996	0.964	3.2	117	0.01
24	Anthracene	0.928	0.880	5.2	115	0.00
25 SURR	Fluoranthene-d10	4.995	4.747	5.0	116	0.00
26	Fluoranthene	1.234	1.185	4.0	117	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	114	0.00
28	Pvrene	1.101	1.028	6.6	117	0.00
29 S	Terphenyl-d14	0.908	0.822	9.5	115	0.00
30	Benzo(a)anthracene	1.137	1.065	6.3	118	0.00
31	Chrysene	1.154	1.065	7.7	115	0.00
32	Bis(2-ethylhexyl)phthalate	2.360	2.397	-1.6	133	0.00
33	Indeno(1,2,3-cd)pyrene	1.058	1.015	4.1	116	0.00
34 I	Perylene-d12	1.000	1.000	0.0	113	0.00
35	Benzo(b)fluoranthene	1.135	1.072	5.6	116	0.00
36	Benzo(k)fluoranthene	1.197	1.159	3.2	118	0.00
37 C	Benzo(a)pyrene	1.088	1.036	4.8	115	0.00
38	Dibenzo(a,h)anthracene	0.927	0.943	-1.7	123	0.00
39	Benzo(a,h,i)perylene	0.968	0.919	5.1	117	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE111318\
Data File : BE098212.D
Acq On : 13 Nov 2018 16:58
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Nov 13 17:33:48 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE111218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 12 13:32:15 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			