

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE111218\
 Data File : BE098189.D
 Acq On : 12 Nov 2018 14:13
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE111218

Quant Time: Nov 12 14:43:48 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	103	0.00
2	1,4-Dioxane	0.650	0.656	-0.9	117	0.00
3	n-Nitrosodimethylamine	0.785	0.704	10.3	99	0.00
4 S	2-Fluorophenol	1.129	1.122	0.6	104	0.00
5 S	Phenol-d6	1.814	1.683	7.2	98	0.00
6	bis(2-Chloroethyl)ether	1.464	1.338	8.6	100	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
8 S	Nitrobenzene-d5	0.399	0.400	-0.3	108	0.01
9	Naphthalene	4.068	3.950	2.9	102	0.00
10	Hexachlorobutadiene	0.253	0.235	7.1	98	0.00
11 SURR	2-Methylnaphthalene-d10	0.702	0.679	3.3	102	0.00
12	2-Methylnaphthalene	0.706	0.693	1.8	104	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
14 S	2,4,6-Tribromophenol	0.174	0.148	14.9	102	0.01
15 S	2-Fluorobiphenyl	1.675	1.594	4.8	105	0.00
16	Acenaphthylene	1.834	1.806	1.5	106	0.00
17	Acenaphthene	1.120	1.096	2.1	105	0.00
18	Fluorene	1.434	1.393	2.9	104	0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
20	4-Bromophenyl-phenylether	0.244	0.241	1.2	108	0.01
21	Hexachlorobenzene	0.254	0.245	3.5	102	0.01
22	Pentachlorophenol	0.044	0.040	9.1	101	0.00
23	Phenanthrene	0.996	0.965	3.1	103	0.01
24	Anthracene	0.928	0.886	4.5	102	0.01
25 SURR	Fluoranthene-d10	4.995	4.679	6.3	100	0.00
26	Fluoranthene	1.234	1.180	4.4	102	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	100	0.00
28	Pvrene	1.101	1.014	7.9	101	0.00
29 S	Terphenyl-d14	0.908	0.811	10.7	99	0.00
30	Benzo(a)anthracene	1.137	1.050	7.7	102	0.00
31	Chrysene	1.154	1.070	7.3	101	0.00
32	Bis(2-ethylhexyl)phthalate	2.360	2.034	13.8	98	0.00
33	Indeno(1,2,3-cd)pyrene	1.058	1.002	5.3	100	0.00
34 I	Perylene-d12	1.000	1.000	0.0	100	0.00
35	Benzo(b)fluoranthene	1.135	1.044	8.0	99	0.00
36	Benzo(k)fluoranthene	1.197	1.157	3.3	103	0.00
37 C	Benzo(a)pyrene	1.088	1.029	5.4	101	0.00
38	Dibenzo(a,h)anthracene	0.927	0.857	7.6	98	0.01
39	Benzo(a,h,i)perylene	0.968	0.916	5.4	103	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE111218\
Data File : BE098189.D
Acq On : 12 Nov 2018 14:13
Operator : SJ/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
ICVBE111218

Quant Time: Nov 12 14:43:48 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			