

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE101618\
 Data File : BE097912.D
 Acq On : 16 Oct 2018 15:42
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE101618

Quant Time: Oct 16 17:37:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 16 15:36:44 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	101	0.00
2	1,4-Dioxane	0.697	0.670	3.9	105	0.00
3	n-Nitrosodimethylamine	0.779	0.740	5.0	108	0.00
4 S	2-Fluorophenol	1.261	1.213	3.8	103	0.00
5 S	Phenol-d6	1.864	1.816	2.6	101	0.00
6	bis(2-Chloroethyl)ether	1.583	1.511	4.5	103	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
8 S	Nitrobenzene-d5	0.379	0.381	-0.5	107	0.00
9	Naphthalene	3.983	3.923	1.5	106	-0.01
10	Hexachlorobutadiene	0.215	0.208	3.3	105	0.00
11 SURR	2-Methylnaphthalene-d10	0.665	0.660	0.8	106	0.00
12	2-Methylnaphthalene	0.682	0.680	0.3	108	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	108	0.00
14 S	2,4,6-Tribromophenol	0.215	0.193	10.2	110	0.00
15 S	2-Fluorobiphenyl	1.635	1.548	5.3	108	0.00
16	Acenaphthylene	1.848	1.755	5.0	108	0.00
17	Acenaphthene	1.129	1.093	3.2	110	0.00
18	Fluorene	1.554	1.519	2.3	111	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	118	0.00
20	4-Bromophenyl-phenylether	0.219	0.205	6.4	113	0.00
21	Hexachlorobenzene	0.215	0.206	4.2	115	0.00
22	Pentachlorophenol	0.062	0.046	25.8#	116	0.00
23	Phenanthrene	1.016	0.977	3.8	117	0.00
24	Anthracene	0.960	0.905	5.7	116	0.00
25 SURR	Fluoranthene-d10	5.421	5.144	5.1	120	0.00
26	Fluoranthene	1.334	1.262	5.4	119	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	106	0.00
28	Pvrene	1.154	1.205	-4.4	117	0.00
29 S	Terphenyl-d14	0.917	0.909	0.9	115	0.00
30	Benzo(a)anthracene	1.227	1.192	2.9	109	0.00
31	Chrysene	1.162	1.091	6.1	108	0.00
32	Bis(2-ethylhexyl)phthalate	2.865	2.463	14.0	93	0.00
33	Indeno(1,2,3-cd)pyrene	1.249	1.150	7.9	98	0.00
34 I	Perylene-d12	1.000	1.000	0.0	96	0.00
35	Benzo(b)fluoranthene	1.151	1.069	7.1	96	0.00
36	Benzo(k)fluoranthene	1.199	1.128	5.9	100	0.00
37 C	Benzo(a)pyrene	1.122	1.074	4.3	98	0.00
38	Dibenzo(a,h)anthracene	1.099	1.056	3.9	99	0.00
39	Benzo(a,h,i)perylene	1.126	1.067	5.2	100	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE101618\
Data File : BE097912.D
Acq On : 16 Oct 2018 15:42
Operator : SJ/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
ICVBE101618

Quant Time: Oct 16 17:37:14 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 16 15:36:44 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			