

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE061818\
 Data File : BE096806.D
 Acq On : 18 Jun 2018 13:17
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE061818

Quant Time: Jun 18 14:42:29 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 18 13:28:13 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	93	0.00
2	1,4-Dioxane	0.400	0.399	0.3	94	0.00
3	n-Nitrosodimethylamine	0.400	0.381	4.8	95	0.00
4 S	2-Fluorophenol	0.400	0.391	2.3	101	0.00
5 S	Phenol-d6	0.400	0.392	2.0	104	0.00
6	bis(2-Chloroethyl)ether	0.400	0.370	7.5	93	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	92	0.00
8 S	Nitrobenzene-d5	0.400	0.391	2.3	101	0.00
9	Naphthalene	0.400	0.378	5.5	93	0.00
10	Hexachlorobutadiene	0.400	0.382	4.5	92	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.375	6.3	93	0.00
12	2-Methylnaphthalene	0.400	0.372	7.0	93	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	91	0.00
14 S	2,4,6-Tribromophenol	0.400	0.381	4.8	109	0.01
15 S	2-Fluorobiphenyl	0.400	0.381	4.8	94	0.00
16	Acenaphthylene	0.400	0.358	10.5	94	0.00
17	Acenaphthene	0.400	0.373	6.8	94	0.00
18	Fluorene	0.400	0.369	7.8	94	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	93	0.00
20	4-Bromophenyl-phenylether	0.400	0.370	7.5	94	0.00
21	Hexachlorobenzene	0.400	0.369	7.8	93	0.00
22	Pentachlorophenol	0.400	0.379	5.3	109	0.00
23	Phenanthrene	0.400	0.374	6.5	94	0.00
24	Anthracene	0.400	0.365	8.8	98	0.00
25 SURR	Fluoranthene-d10	0.400	0.362	9.5	96	0.00
26	Fluoranthene	0.400	0.358	10.5	94	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	94	0.00
28	Pvrene	0.400	0.378	5.5	96	0.00
29 S	Terphenyl-d14	0.400	0.381	4.8	95	0.00
30	Benzo(a)anthracene	0.400	0.351	12.3	96	0.00
31	Chrysene	0.400	0.376	6.0	95	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.380	5.0	106	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.378	5.5	104	0.00
34 I	Perylene-d12	0.400	0.400	0.0	102	0.00
35	Benzo(b)fluoranthene	0.400	0.339	15.3	98	0.00
36	Benzo(k)fluoranthene	0.400	0.352	12.0	101	0.00
37 C	Benzo(a)pyrene	0.400	0.325	18.8	96	0.00
38	Dibenzo(a,h)anthracene	0.400	0.359	10.3	104	0.00
39	Benzo(a,h,i)perylene	0.400	0.350	12.5	100	0.00

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Compound	Amount	Calc.	%Dev	Area	% Dev(min)

(#) = Out of Range	SPCC's out = 0 CCC's out = 0				