

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121018\
 Data File : BE098299.D
 Acq On : 10 Dec 2018 15:32
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE121018

Quant Time: Dec 10 16:04:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 10 14:59:55 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	92	0.00
2	1,4-Dioxane	0.400	0.382	4.5	89	0.00
3	n-Nitrosodimethylamine	0.400	0.324	19.0	80	0.00
4 S	2-Fluorophenol	0.400	0.388	3.0	92	0.00
5 S	Phenol-d6	0.400	0.380	5.0	86	0.00
6	bis(2-Chloroethyl)ether	0.400	0.370	7.5	88	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	93	0.00
8 S	Nitrobenzene-d5	0.400	0.379	5.3	96	0.01
9	Naphthalene	0.400	0.401	-0.3	95	0.00
10	Hexachlorobutadiene	0.400	0.413	-3.2	99	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.407	-1.7	95	0.00
12	2-Methylnaphthalene	0.400	0.398	0.5	94	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	95	0.00
14 S	2,4,6-Tribromophenol	0.400	0.385	3.8	96	0.00
15 S	2-Fluorobiphenyl	0.400	0.367	8.3	95	0.00
16	Acenaphthylene	0.400	0.386	3.5	95	0.00
17	Acenaphthene	0.400	0.391	2.3	93	0.00
18	Fluorene	0.400	0.399	0.3	95	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	97	0.00
20	4-Bromophenyl-phenylether	0.400	0.379	5.3	96	0.00
21	Hexachlorobenzene	0.400	0.385	3.8	95	0.00
22	Pentachlorophenol	0.400	0.352	12.0	86	0.00
23	Phenanthrene	0.400	0.398	0.5	98	0.00
24	Anthracene	0.400	0.381	4.8	94	0.00
25 SURR	Fluoranthene-d10	0.400	0.424	-6.0	99	0.00
26	Fluoranthene	0.400	0.422	-5.5	98	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	99	0.00
28	Pvrene	0.400	0.407	-1.7	99	0.00
29 S	Terphenyl-d14	0.400	0.396	1.0	99	0.00
30	Benzo(a)anthracene	0.400	0.385	3.8	98	0.00
31	Chrysene	0.400	0.398	0.5	100	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.393	1.8	98	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.384	4.0	101	0.00
34 I	Perylene-d12	0.400	0.400	0.0	100	0.00
35	Benzo(b)fluoranthene	0.400	0.386	3.5	100	0.00
36	Benzo(k)fluoranthene	0.400	0.386	3.5	101	0.00
37 C	Benzo(a)pyrene	0.400	0.386	3.5	99	0.00
38	Dibenzo(a,h)anthracene	0.400	0.384	4.0	99	0.00
39	Benzo(a,h,i)perylene	0.400	0.392	2.0	101	0.00

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

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