

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE032219\
 Data File : BE099082.D
 Acq On : 22 Mar 2019 16:10
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE032219

Quant Time: Mar 23 00:49:46 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE032219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 23 00:46:37 2019
 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	97	0.00
2	1,4-Dioxane	0.400	0.335	16.3	94	0.00
3	n-Nitrosodimethylamine	0.400	0.253	36.8#	67	-0.01
4 S	2-Fluorophenol	0.400	0.334	16.5	91	0.00
5 S	Phenol-d6	0.400	0.335	16.3	93	0.00
6	bis(2-Chloroethyl)ether	0.400	0.354	11.5	96	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	94	0.00
8 S	Nitrobenzene-d5	0.400	0.418	-4.5	124	0.00
9	Naphthalene	0.400	0.353	11.8	94	0.00
10	Hexachlorobutadiene	0.400	0.359	10.3	94	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.358	10.5	96	-0.01
12	2-Methylnaphthalene	0.400	0.340	15.0	93	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	96	0.00
14 S	2,4,6-Tribromophenol	0.400	0.362	9.5	111	0.00
15 S	2-Fluorobiphenyl	0.400	0.360	10.0	94	0.00
16	Acenaphthylene	0.400	0.332	17.0	94	0.00
17	Acenaphthene	0.400	0.344	14.0	96	0.00
18	Fluorene	0.400	0.344	14.0	96	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	96	0.00
20	4-Bromophenyl-phenylether	0.400	0.342	14.5	96	0.00
21	Hexachlorobenzene	0.400	0.344	14.0	98	0.00
22	Pentachlorophenol	0.400	0.475	-18.7	108	0.00
23	Phenanthrene	0.400	0.348	13.0	99	0.00
24	Anthracene	0.400	0.331	17.3	95	0.00
25 SURR	Fluoranthene-d10	0.400	0.352	12.0	97	0.00
26	Fluoranthene	0.400	0.340	15.0	98	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	100	0.00
28	Pyrene	0.400	0.331	17.3	98	0.00
29 S	Terphenyl-d14	0.400	0.350	12.5	99	0.00
30	Benzo(a)anthracene	0.400	0.337	15.8	99	0.00
31	Chrysene	0.400	0.347	13.3	103	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.270	32.5#	95	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.343	14.2	102	0.00
34 I	Perylene-d12	0.400	0.400	0.0	105	0.00
35	Benzo(b)fluoranthene	0.400	0.329	17.8	104	0.00
36	Benzo(k)fluoranthene	0.400	0.331	17.3	104	0.00
37 C	Benzo(a)pyrene	0.400	0.322	19.5	103	0.00
38	Dibenzo(a,h)anthracene	0.400	0.326	18.5	104	-0.01
39	Benzo(g,h,i)perylene	0.400	0.330	17.5	104	0.00

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

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