

Data Path : Z:\HPCHEM1\BNA E\Data\BE032318\
 Data File : BE095821.D
 Acq On : 23 Mar 2018 12:21
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE032318

Quant Time: Mar 23 12:57:01 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE032318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 23 12:41:11 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	101	-0.01
2	1,4-Dioxane	0.400	0.404	-1.0	108	0.00
3	n-Nitrosodimethylamine	0.400	0.366	8.5	89	0.00
4 S	2-Fluorophenol	0.400	0.387	3.3	94	-0.01
5 S	Phenol-d6	0.400	0.381	4.8	92	0.00
6	bis(2-Chloroethyl)ether	0.400	0.373	6.8	91	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	96	0.00
8 S	Nitrobenzene-d5	0.400	0.400	0.0	95	0.00
9	Naphthalene	0.400	0.394	1.5	92	0.00
10	Hexachlorobutadiene	0.400	0.399	0.3	92	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.407	-1.7	94	0.00
12	2-Methylnaphthalene	0.400	0.407	-1.7	95	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	99	0.00
14 S	2,4,6-Tribromophenol	0.400	0.374	6.5	95	0.00
15 S	2-Fluorobiphenyl	0.400	0.391	2.3	95	0.00
16	Acenaphthylene	0.400	0.393	1.8	97	0.00
17	Acenaphthene	0.400	0.392	2.0	97	0.00
18	Fluorene	0.400	0.392	2.0	96	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	100	0.00
20	4-Bromophenyl-phenylether	0.400	0.394	1.5	97	0.00
21	Hexachlorobenzene	0.400	0.392	2.0	96	0.00
22	Pentachlorophenol	0.400	0.354	11.5	88	0.00
23	Phenanthrene	0.400	0.400	0.0	97	0.00
24	Anthracene	0.400	0.394	1.5	97	0.00
25 SURR	Fluoranthene-d10	0.400	0.401	-0.3	98	0.00
26	Fluoranthene	0.400	0.401	-0.3	99	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	100	0.00
28	Pvrene	0.400	0.442	-10.5	100	0.00
29 S	Terphenyl-d14	0.400	0.399	0.3	98	0.00
30	Benzo(a)anthracene	0.400	0.405	-1.3	99	0.00
31	Chrysene	0.400	0.404	-1.0	99	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.359	10.3	95	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.407	-1.7	100	0.00
34 I	Perylene-d12	0.400	0.400	0.0	101	0.00
35	Benzo(b)fluoranthene	0.400	0.404	-1.0	101	0.00
36	Benzo(k)fluoranthene	0.400	0.408	-2.0	102	0.00
37 C	Benzo(a)pyrene	0.400	0.402	-0.5	100	0.00
38	Dibenzo(a,h)anthracene	0.400	0.407	-1.7	100	0.00
39	Benzo(a,h,i)perylene	0.400	0.393	1.8	98	0.00

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

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