

Data Path : Z:\HPCHEM1\BNA E\Data\BE072117\
 Data File : BE093511.D
 Acq On : 21 Jul 2017 16:12
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE072117

Quant Time: Jul 21 17:20:38 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE072117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 21 16:13:20 2017
 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	0.00
2	1,4-Dioxane	0.400	0.372	7.0	95	0.00
3	n-Nitrosodimethylamine	0.400	0.400	0.0	109	0.00
4 S	2-Fluorophenol	0.400	0.395	1.3	101	0.00
5 S	Phenol-d6	0.400	0.380	5.0	100	0.00
6	bis(2-Chloroethyl)ether	0.400	0.402	-0.5	103	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	102	0.00
8 S	Nitrobenzene-d5	0.400	0.381	4.8	102	0.00
9	Naphthalene	0.400	0.387	3.3	102	0.00
10	Hexachlorobutadiene	0.400	0.392	2.0	102	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.388	3.0	101	0.00
12	2-Methylnaphthalene	0.400	0.382	4.5	101	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	103	0.00
14 S	2,4,6-Tribromophenol	0.400	0.355	11.3	88	0.00
15 S	2-Fluorobiphenyl	0.400	0.393	1.8	99	0.00
16	Acenaphthylene	0.400	0.374	6.5	100	0.00
17	Acenaphthene	0.400	0.384	4.0	99	0.00
18	Fluorene	0.400	0.382	4.5	98	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	109	0.00
20	4-Bromophenyl-phenylether	0.400	0.309	22.8	96	0.00
21	Hexachlorobenzene	0.400	0.323	19.3	97	0.00
22	Pentachlorophenol	0.400	0.395	1.3	96	0.00
23	Phenanthrene	0.400	0.369	7.8	105	0.00
24	Anthracene	0.400	0.377	5.8	107	0.00
25 SURR	Fluoranthene-d10	0.400	0.348	13.0	98	0.00
26	Fluoranthene	0.400	0.344	14.0	98	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	99	0.00
28	Pvrene	0.400	0.396	1.0	98	0.00
29 S	Terphenyl-d14	0.400	0.399	0.3	98	0.00
30	Benzo(a)anthracene	0.400	0.386	3.5	97	0.00
31	Chrysene	0.400	0.402	-0.5	99	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.344	14.0	87	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.389	2.8	98	0.00
34 I	Perylene-d12	0.400	0.400	0.0	100	0.00
35	Benzo(b)fluoranthene	0.400	0.394	1.5	98	0.00
36	Benzo(k)fluoranthene	0.400	0.394	1.5	100	0.00
37 C	Benzo(a)pyrene	0.400	0.382	4.5	97	0.00
38	Dibenzo(a,h)anthracene	0.400	0.389	2.8	99	0.00
39	Benzo(a,h,i)perylene	0.400	0.389	2.8	97	0.00

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

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