

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE092718\
 Data File : BE097696.D
 Acq On : 27 Sep 2018 16:17
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE092718

Quant Time: Sep 27 17:01:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE092718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 27 15:45:00 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	94	0.00
2	1,4-Dioxane	0.400	0.374	6.5	95	0.00
3	n-Nitrosodimethylamine	0.400	0.349	12.8	87	0.00
4 S	2-Fluorophenol	0.400	0.345	13.8	87	0.00
5 S	Phenol-d6	0.400	0.349	12.8	90	0.01
6	bis(2-Chloroethyl)ether	0.400	0.382	4.5	95	0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	98	0.00
8 S	Nitrobenzene-d5	0.400	0.361	9.8	90	0.00
9	Naphthalene	0.400	0.385	3.8	97	0.00
10	Hexachlorobutadiene	0.400	0.390	2.5	100	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.378	5.5	97	0.00
12	2-Methylnaphthalene	0.400	0.375	6.3	96	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	96	0.00
14 S	2,4,6-Tribromophenol	0.400	0.323	19.3	91	0.00
15 S	2-Fluorobiphenyl	0.400	0.387	3.3	98	0.00
16	Acenaphthylene	0.400	0.375	6.3	97	0.00
17	Acenaphthene	0.400	0.397	0.8	98	0.00
18	Fluorene	0.400	0.384	4.0	95	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	98	0.00
20	4-Bromophenyl-phenylether	0.400	0.376	6.0	96	0.00
21	Hexachlorobenzene	0.400	0.387	3.3	100	0.00
22	Pentachlorophenol	0.400	0.337	15.8	85	0.01
23	Phenanthrene	0.400	0.379	5.3	97	0.00
24	Anthracene	0.400	0.354	11.5	94	0.00
25 SURR	Fluoranthene-d10	0.400	0.379	5.3	98	0.00
26	Fluoranthene	0.400	0.380	5.0	98	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	97	0.00
28	Pvrene	0.400	0.390	2.5	97	0.00
29 S	Terphenyl-d14	0.400	0.388	3.0	98	0.00
30	Benzo(a)anthracene	0.400	0.359	10.3	93	0.00
31	Chrysene	0.400	0.408	-2.0	102	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.323	19.3	89	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.377	5.8	96	0.00
34 I	Perylene-d12	0.400	0.400	0.0	96	0.00
35	Benzo(b)fluoranthene	0.400	0.386	3.5	99	0.00
36	Benzo(k)fluoranthene	0.400	0.388	3.0	96	0.00
37 C	Benzo(a)pyrene	0.400	0.369	7.8	95	0.00
38	Dibenzo(a,h)anthracene	0.400	0.372	7.0	96	0.00
39	Benzo(a,h,i)perylene	0.400	0.379	5.3	96	0.00

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

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