

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE030619\
 Data File : BE098798.D
 Acq On : 6 Mar 2019 14:55
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE030619

Quant Time: Mar 06 15:38:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 06 15:33:24 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	101	0.01
2	1,4-Dioxane	0.400	0.377	5.8	98	0.00
3	n-Nitrosodimethylamine	0.400	0.356	11.0	88	-0.01
4 S	2-Fluorophenol	0.400	0.388	3.0	95	0.00
5 S	Phenol-d6	0.400	0.403	-0.8	102	0.01
6	bis(2-Chloroethyl)ether	0.400	0.383	4.3	100	0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	95	0.01
8 S	Nitrobenzene-d5	0.400	0.339	15.3	90	0.01
9	Naphthalene	0.400	0.407	-1.7	99	0.01
10	Hexachlorobutadiene	0.400	0.404	-1.0	99	0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.408	-2.0	99	0.01
12	2-Methylnaphthalene	0.400	0.393	1.8	97	0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	101	0.01
14 S	2,4,6-Tribromophenol	0.400	0.401	-0.3	102	0.00
15 S	2-Fluorobiphenyl	0.400	0.384	4.0	98	0.00
16	Acenaphthylene	0.400	0.383	4.3	102	0.01
17	Acenaphthene	0.400	0.396	1.0	101	0.00
18	Fluorene	0.400	0.377	5.8	98	0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	99	0.00
20	4-Bromophenyl-phenylether	0.400	0.368	8.0	93	0.01
21	Hexachlorobenzene	0.400	0.375	6.3	98	0.00
22	Pentachlorophenol	0.400	0.448	-12.0	104	0.00
23	Phenanthrene	0.400	0.383	4.3	98	0.00
24	Anthracene	0.400	0.368	8.0	97	0.01
25 SURR	Fluoranthene-d10	0.400	0.379	5.3	98	0.00
26	Fluoranthene	0.400	0.373	6.8	96	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	101	0.00
28	Pvrene	0.400	0.381	4.8	97	0.01
29 S	Terphenyl-d14	0.400	0.377	5.8	96	0.00
30	Benzo(a)anthracene	0.400	0.372	7.0	96	0.01
31	Chrysene	0.400	0.394	1.5	102	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.360	10.0	99	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.399	0.3	103	0.02
34 I	Perylene-d12	0.400	0.400	0.0	103	0.00
35	Benzo(b)fluoranthene	0.400	0.376	6.0	100	0.00
36	Benzo(k)fluoranthene	0.400	0.401	-0.3	108	0.00
37 C	Benzo(a)pyrene	0.400	0.381	4.8	101	0.00
38	Dibenzo(a,h)anthracene	0.400	0.388	3.0	101	0.00
39	Benzo(a,h,i)perylene	0.400	0.388	3.0	102	0.02

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

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