

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE022619\
 Data File : BE098761.D
 Acq On : 26 Feb 2019 15:10
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE022619

Quant Time: Feb 26 16:22:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE022619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 26 15:44:42 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	95	-0.01
2	1,4-Dioxane	0.400	0.385	3.8	92	0.00
3	n-Nitrosodimethylamine	0.400	0.341	14.8	78	-0.01
4 S	2-Fluorophenol	0.400	0.376	6.0	91	0.00
5 S	Phenol-d6	0.400	0.372	7.0	91	0.00
6	bis(2-Chloroethyl)ether	0.400	0.356	11.0	85	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	93	0.00
8 S	Nitrobenzene-d5	0.400	0.383	4.3	93	0.00
9	Naphthalene	0.400	0.384	4.0	91	0.00
10	Hexachlorobutadiene	0.400	0.381	4.8	90	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.381	4.8	93	0.00
12	2-Methylnaphthalene	0.400	0.384	4.0	94	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	99	0.00
14 S	2,4,6-Tribromophenol	0.400	0.370	7.5	94	0.00
15 S	2-Fluorobiphenyl	0.400	0.367	8.3	91	0.01
16	Acenaphthylene	0.400	0.371	7.3	96	0.00
17	Acenaphthene	0.400	0.371	7.3	94	0.01
18	Fluorene	0.400	0.372	7.0	95	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	95	0.00
20	4-Bromophenyl-phenylether	0.400	0.376	6.0	96	0.00
21	Hexachlorobenzene	0.400	0.387	3.3	92	0.00
22	Pentachlorophenol	0.400	0.522	-30.5#	131	0.00
23	Phenanthrene	0.400	0.387	3.3	95	0.00
24	Anthracene	0.400	0.377	5.8	94	0.00
25 SURR	Fluoranthene-d10	0.400	0.384	4.0	94	0.00
26	Fluoranthene	0.400	0.382	4.5	95	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	95	0.00
28	Pyrene	0.400	0.375	6.3	94	0.00
29 S	Terphenyl-d14	0.400	0.373	6.8	93	0.00
30	Benzo(a)anthracene	0.400	0.402	-0.5	105	0.00
31	Chrysene	0.400	0.373	6.8	89	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.341	14.8	90	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.512	-28.0#	128	0.00
34 I	Perylene-d12	0.400	0.400	0.0	92	0.00
35	Benzo(b)fluoranthene	0.400	0.401	-0.3	94	0.00
36	Benzo(k)fluoranthene	0.400	0.419	-4.7	99	0.00
37 C	Benzo(a)pyrene	0.400	0.411	-2.7	98	0.00
38	Dibenzo(a,h)anthracene	0.400	0.602	-50.5#	171	0.00
39	Benzo(g,h,i)perylene	0.400	0.475	-18.7	111	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0