

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE071618\
 Data File : BE096941.D
 Acq On : 16 Jul 2018 16:24
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE071618

Quant Time: Jul 17 15:09:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 17 15:02:56 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	99	0.00
2	1,4-Dioxane	0.400	0.345	13.8	89	0.00
3	n-Nitrosodimethylamine	0.400	0.347	13.3	92	0.00
4 S	2-Fluorophenol	0.400	0.334	16.5	94	0.00
5 S	Phenol-d6	0.400	0.338	15.5	92	0.00
6	bis(2-Chloroethyl)ether	0.400	0.350	12.5	90	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	98	0.00
8 S	Nitrobenzene-d5	0.400	0.345	13.8	91	0.00
9	Naphthalene	0.400	0.354	11.5	91	0.00
10	Hexachlorobutadiene	0.400	0.353	11.8	91	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.345	13.8	89	0.00
12	2-Methylnaphthalene	0.400	0.345	13.8	89	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	97	0.00
14 S	2,4,6-Tribromophenol	0.400	0.361	9.8	89	0.00
15 S	2-Fluorobiphenyl	0.400	0.362	9.5	90	0.00
16	Acenaphthylene	0.400	0.343	14.2	87	0.00
17	Acenaphthene	0.400	0.352	12.0	89	0.00
18	Fluorene	0.400	0.347	13.3	89	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	100	-0.01
20	4-Bromophenyl-phenylether	0.400	0.342	14.5	90	-0.01
21	Hexachlorobenzene	0.400	0.350	12.5	90	-0.01
22	Pentachlorophenol	0.400	0.365	8.8	88	0.00
23	Phenanthrene	0.400	0.346	13.5	90	0.00
24	Anthracene	0.400	0.339	15.3	91	0.00
25 SURR	Fluoranthene-d10	0.400	0.328	18.0	89	0.00
26	Fluoranthene	0.400	0.329	17.8	88	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	96	0.00
28	Pvrene	0.400	0.351	12.3	88	0.00
29 S	Terphenyl-d14	0.400	0.355	11.3	90	0.00
30	Benzo(a)anthracene	0.400	0.335	16.3	88	-0.01
31	Chrysene	0.400	0.351	12.3	86	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.315	21.3	87	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.323	19.3	87	0.00
34 I	Perylene-d12	0.400	0.400	0.0	99	0.00
35	Benzo(b)fluoranthene	0.400	0.325	18.8	87	0.00
36	Benzo(k)fluoranthene	0.400	0.326	18.5	85	0.00
37 C	Benzo(a)pyrene	0.400	0.345	13.8	85	0.00
38	Dibenzo(a,h)anthracene	0.400	0.356	11.0	90	0.00
39	Benzo(a,h,i)perylene	0.400	0.338	15.5	93	0.00

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

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