

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011519\
 Data File : BE098627.D
 Acq On : 15 Jan 2019 14:34
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICSVBE011519

Quant Time: Jan 15 15:06:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE011519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 15 14:30:55 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	94	0.00
2	1,4-Dioxane	0.789	0.687	12.9	95	0.00
3	n-Nitrosodimethylamine	0.579	0.485	16.2	83	0.00
4 S	2-Fluorophenol	1.067	1.015	4.9	92	0.00
5 S	Phenol-d6	1.519	1.516	0.2	94	0.00
6	bis(2-Chloroethyl)ether	1.149	1.031	10.3	91	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
8 S	Nitrobenzene-d5	0.317	0.322	-1.6	97	0.00
9	Naphthalene	3.984	3.792	4.8	93	0.00
10	Hexachlorobutadiene	0.285	0.276	3.2	94	0.00
11 SURR	2-Methylnaphthalene-d10	0.658	0.616	6.4	94	0.00
12	2-Methylnaphthalene	0.660	0.613	7.1	93	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
14 S	2,4,6-Tribromophenol	0.174	0.170	2.3	96	0.00
15 S	2-Fluorobiphenyl	1.638	1.507	8.0	98	0.00
16	Acenaphthylene	1.732	1.615	6.8	96	0.00
17	Acenaphthene	1.087	1.019	6.3	94	0.00
18	Fluorene	1.414	1.330	5.9	94	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	93	0.00
20	4-Bromophenyl-phenylether	0.212	0.198	6.6	94	0.00
21	Hexachlorobenzene	0.256	0.247	3.5	94	0.00
22	Pentachlorophenol	0.075	0.062	17.3	99	0.00
23	Phenanthrene	0.948	0.908	4.2	95	0.00
24	Anthracene	0.812	0.750	7.6	95	0.00
25 SURR	Fluoranthene-d10	4.955	4.712	4.9	96	0.00
26	Fluoranthene	1.220	1.145	6.1	95	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	96	0.00
28	Pvrene	1.101	1.035	6.0	95	0.00
29 S	Terphenyl-d14	0.885	0.808	8.7	96	0.00
30	Benzo(a)anthracene	1.075	1.016	5.5	97	0.00
31	Chrysene	1.115	1.075	3.6	98	0.00
32	Bis(2-ethylhexyl)phthalate	2.281	1.976	13.4	94	0.00
33	Indeno(1,2,3-cd)pyrene	1.150	1.130	1.7	98	0.00
34 I	Perylene-d12	1.000	1.000	0.0	98	0.00
35	Benzo(b)fluoranthene	1.091	1.017	6.8	98	0.00
36	Benzo(k)fluoranthene	1.193	1.129	5.4	96	0.00
37 C	Benzo(a)pyrene	1.072	1.001	6.6	97	0.00
38	Dibenzo(a,h)anthracene	0.949	0.965	-1.7	106	0.00
39	Benzo(a,h,i)perylene	1.040	1.009	3.0	98	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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