

Data Path : U:\HPCHEM1\BNA E\DATA\BE012318\
 Data File : BE095097.D
 Acq On : 23 Jan 2018 15:00
 Operator : SJ/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICSVBE012318

Quant Time: Jan 23 15:53:52 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE012318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 23 15:47:25 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	96	0.00
2	1,4-Dioxane	0.393	0.417	-6.1	101	0.00
3	n-Nitrosodimethylamine	0.469	0.450	4.1	91	0.00
4 S	2-Fluorophenol	0.648	0.634	2.2	93	0.00
5 S	Phenol-d6	0.982	0.958	2.4	94	0.00
6	bis(2-Chloroethyl)ether	0.914	0.888	2.8	90	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	90	0.00
8 S	Nitrobenzene-d5	0.364	0.358	1.6	93	0.00
9	Naphthalene	4.577	4.626	-1.1	91	0.00
10	Hexachlorobutadiene	0.205	0.210	-2.4	93	0.00
11 SURR	2-Methylnaphthalene-d10	0.664	0.670	-0.9	92	0.00
12	2-Methylnaphthalene	0.785	0.791	-0.8	92	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	88	0.00
14 S	2,4,6-Tribromophenol	0.097	0.089	8.2	82	0.00
15 S	2-Fluorobiphenyl	1.758	1.823	-3.7	92	0.00
16	Acenaphthylene	2.141	2.096	2.1	89	0.00
17	Acenaphthene	1.393	1.384	0.6	89	-0.01
18	Fluorene	1.740	1.693	2.7	87	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	85	0.00
20	4-Bromophenyl-phenylether	0.235	0.232	1.3	85	0.00
21	Hexachlorobenzene	0.236	0.238	-0.8	87	0.00
22	Pentachlorophenol	0.055	0.040	27.3#	78	0.00
23	Phenanthrene	1.236	1.243	-0.6	86	0.00
24	Anthracene	1.102	1.082	1.8	86	0.00
25 SURR	Fluoranthene-d10	5.031	4.898	2.6	85	0.00
26	Fluoranthene	1.484	1.462	1.5	86	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	88	0.00
28	Pvrene	1.691	1.548	8.5	79	0.00
29 S	Terphenyl-d14	0.819	0.800	2.3	84	0.00
30	Benzo(a)anthracene	1.251	1.154	7.8	78	0.00
31	Chrysene	1.265	1.258	0.6	92	0.00
32	Bis(2-ethylhexyl)phthalate	1.740	1.395	19.8	73	0.00
33	Indeno(1,2,3-cd)pyrene	1.150	1.064	7.5	79	0.00
34 I	Perylene-d12	1.000	1.000	0.0	82	0.00
35	Benzo(b)fluoranthene	1.419	1.442	-1.6	84	0.00
36	Benzo(k)fluoranthene	1.452	1.415	2.5	82	0.00
37 C	Benzo(a)pyrene	1.301	1.251	3.8	79	0.00
38	Dibenzo(a,h)anthracene	1.095	1.027	6.2	77	0.00
39	Benzo(a,h,i)perylene	1.187	1.163	2.0	83	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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