

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE032219\
 Data File : BE099084.D
 Acq On : 22 Mar 2019 17:27
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 23 01:14:23 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE032219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 23 00:46:37 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	83	0.00
2	1,4-Dioxane	0.621	0.623	-0.3	97	0.00
3	n-Nitrosodimethylamine	0.320	0.244	23.8	69	0.00
4 S	2-Fluorophenol	1.191	1.019	14.4	80	0.00
5 S	Phenol-d6	1.715	1.518	11.5	84	0.00
6	bis(2-Chloroethyl)ether	1.394	1.326	4.9	89	-0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	83	0.00
8 S	Nitrobenzene-d5	0.175	0.193	-10.3	116	0.00
9	Naphthalene	4.714	4.254	9.8	85	0.00
10	Hexachlorobutadiene	0.231	0.213	7.8	86	0.00
11 SURR	2-Methylnaphthalene-d10	0.720	0.654	9.2	86	-0.01
12	2-Methylnaphthalene	0.807	0.729	9.7	87	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	87	0.00
14 S	2,4,6-Tribromophenol	0.118	0.106	10.2	101	0.00
15 S	2-Fluorobiphenyl	1.591	1.409	11.4	84	0.00
16	Acenaphthylene	2.135	1.731	18.9	83	-0.01
17	Acenaphthene	1.266	1.097	13.3	88	0.00
18	Fluorene	1.847	1.610	12.8	88	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
20	4-Bromophenyl-phenylether	0.221	0.198	10.4	93	0.00
21	Hexachlorobenzene	0.206	0.179	13.1	92	0.00
22	Pentachlorophenol	0.061	0.052	14.8	111	-0.01
23	Phenanthrene	1.188	1.036	12.8	92	0.00
24	Anthracene	1.124	0.958	14.8	91	0.00
25 SURR	Fluoranthene-d10	5.336	4.796	10.1	92	0.00
26	Fluoranthene	1.671	1.461	12.6	93	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	87	0.00
28	Pyrene	1.357	1.210	10.8	92	0.00
29 S	Terphenyl-d14	0.846	0.777	8.2	91	0.00
30	Benzo(a)anthracene	1.409	1.239	12.1	90	0.00
31	Chrysene	1.408	1.285	8.7	95	0.00
32	Bis(2-ethylhexyl)phthalate	1.829	1.205	34.1#	81	0.00
33	Indeno(1,2,3-cd)pyrene	1.514	1.300	14.1	89	0.00
34 I	Perylene-d12	1.000	1.000	0.0	87	0.00
35	Benzo(b)fluoranthene	1.384	1.193	13.8	91	0.00
36	Benzo(k)fluoranthene	1.356	1.195	11.9	92	0.00
37 C	Benzo(a)pyrene	1.279	1.087	15.0	90	0.00
38	Dibenzo(a,h)anthracene	1.244	1.070	14.0	91	0.00
39	Benzo(g,h,i)perylene	1.278	1.111	13.1	91	-0.01

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