

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031519\
 Data File : BE098967.D
 Acq On : 15 Mar 2019 10:14
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC00.4

Quant Time: Mar 18 06:45:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 15 05:31:33 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	103	0.00
2	1,4-Dioxane	0.789	0.881	-11.7	119	0.00
3	n-Nitrosodimethylamine	1.006	0.749	25.5#	75	0.03
4 S	2-Fluorophenol	1.196	1.085	9.3	90	0.01
5 S	Phenol-d6	1.689	1.647	2.5	100	0.01
6	bis(2-Chloroethyl)ether	1.304	1.029	21.1	84	0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	115	0.00
8 S	Nitrobenzene-d5	0.419	0.357	14.8	109	0.01
9	Naphthalene	4.581	4.543	0.8	117	0.00
10	Hexachlorobutadiene	0.303	0.292	3.6	114	0.00
11 SURR	2-Methylnaphthalene-d10	0.739	0.713	3.5	113	0.00
12	2-Methylnaphthalene	0.743	0.677	8.9	108	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	119	0.00
14 S	2,4,6-Tribromophenol	0.202	0.143	29.2#	85	0.00
15 S	2-Fluorobiphenyl	1.656	1.621	2.1	118	0.00
16	Acenaphthylene	2.057	1.948	5.3	118	0.00
17	Acenaphthene	1.220	1.170	4.1	114	-0.01
18	Fluorene	1.760	1.623	7.8	112	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
20	4-Bromophenyl-phenylether	0.217	0.232	-6.9	116	0.01
21	Hexachlorobenzene	0.258	0.255	1.2	111	0.00
22	Pentachlorophenol	0.052	0.028	46.2#	77	0.01
23	Phenanthrene	1.137	1.056	7.1	101	0.00
24	Anthracene	0.972	0.790	18.7	91	0.00
25 SURR	Fluoranthene-d10	6.385	5.417	15.2	93	0.00
26	Fluoranthene	1.605	1.377	14.2	95	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	89	0.00
28	Pvrene	1.224	1.266	-3.4	93	0.00
29 S	Terphenyl-d14	0.972	0.926	4.7	86	0.00
30	Benzo(a)anthracene	1.248	1.141	8.6	83	0.00
31	Chrysene	1.336	1.223	8.5	84	0.00
32	Bis(2-ethylhexyl)phthalate	2.657	2.125	20.0	78	-0.01
33	Indeno(1,2,3-cd)pyrene	1.410	1.217	13.7	78	0.00
34 I	Perylene-d12	1.000	1.000	0.0	92	0.00
35	Benzo(b)fluoranthene	1.315	1.218	7.4	88	0.00
36	Benzo(k)fluoranthene	1.377	1.436	-4.3	101	0.00
37 C	Benzo(a)pyrene	1.261	1.204	4.5	91	0.00
38	Dibenzo(a,h)anthracene	1.193	0.851	28.7#	67	0.00
39	Benzo(a,h,i)perylene	1.248	1.073	14.0	81	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031519\
Data File : BE098967.D
Acq On : 15 Mar 2019 10:14
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Mar 18 06:45:02 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 15 05:31:33 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)

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