

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN083118\
 Data File : BN002572.D
 Acq On : 31 Aug 2018 22:10
 Operator : JU/SJ
 Sample : J4665-15
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E0AB0

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SOM-EPA-BN081318MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.775	129	134	142	rBV	76689	114264	2.21%	0.208%
2	6.852	650	657	675	rBV	226111	428220	8.29%	0.781%
3	7.005	677	683	691	rBV	610776	977434	18.93%	1.782%
4	7.205	710	717	734	rVB	724256	1202271	23.29%	2.191%
5	7.663	789	795	803	rVB	598635	947863	18.36%	1.728%
6	8.375	909	916	927	rBV	607970	1067202	20.67%	1.945%
7	8.810	983	990	1000	rBV	813746	1408880	27.29%	2.568%
8	9.081	1031	1036	1046	rVB3	57090	109625	2.12%	0.200%
9	9.452	1091	1099	1105	rBV6	25759	67720	1.31%	0.123%
10	9.528	1105	1112	1119	rVV	723957	1177415	22.81%	2.146%
11	9.805	1152	1159	1167	rBV9	35527	94257	1.83%	0.172%
12	9.975	1182	1188	1194	rVB3	50657	104945	2.03%	0.191%
13	10.069	1195	1204	1214	rBV	1012595	1896506	36.73%	3.457%
14	10.175	1216	1222	1227	rVB3	73162	148782	2.88%	0.271%
15	10.434	1259	1266	1283	rVB	911014	1660991	32.17%	3.028%
16	10.593	1284	1293	1303	rBV3	168397	326153	6.32%	0.595%
17	10.693	1303	1310	1314	rBV7	43320	101100	1.96%	0.184%
18	11.010	1359	1364	1370	rBV2	62961	103851	2.01%	0.189%
19	11.175	1382	1392	1399	rBV	128882	291366	5.64%	0.531%
20	11.346	1416	1421	1429	rVB9	37702	97909	1.90%	0.178%
21	11.440	1430	1437	1444	rBV7	31675	82866	1.61%	0.151%
22	11.516	1444	1450	1453	rBV5	33938	76302	1.48%	0.139%
23	11.793	1493	1497	1504	rBV8	27303	60895	1.18%	0.111%
24	12.040	1535	1539	1543	rBV5	45164	77659	1.50%	0.142%
25	12.087	1543	1547	1555	rBV3	111492	255396	4.95%	0.466%
26	12.246	1571	1574	1584	rVB5	73283	140533	2.72%	0.256%
27	12.328	1584	1588	1594	rVB6	46554	91566	1.77%	0.167%
28	12.416	1595	1603	1612	rVB7	99244	250412	4.85%	0.456%
29	12.516	1612	1620	1624	rBV6	65424	144615	2.80%	0.264%
30	12.569	1624	1629	1634	rBV6	52599	101258	1.96%	0.185%
31	12.628	1634	1639	1645	rVB6	46036	104610	2.03%	0.191%
32	12.698	1647	1651	1656	rVB6	30952	52181	1.01%	0.095%
33	12.757	1658	1661	1665	rBV5	36217	52312	1.01%	0.095%
34	13.110	1716	1721	1725	rBV4	91637	202114	3.91%	0.368%

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35	13.322	1754	1757	1762	rBV7	49641	70794	1.37%	0.129%
36	13.546	1791	1795	1801	rVV4	106184	172301	3.34%	0.314%
37	13.710	1818	1823	1829	rBV	1936870	2739846	53.07%	4.994%
38	13.987	1864	1870	1886	rVB	2303100	3592280	69.58%	6.548%
39	14.245	1911	1914	1917	rBV2	77026	103227	2.00%	0.188%
40	14.298	1917	1923	1931	rVV2	1375185	2158460	41.81%	3.934%
41	14.440	1945	1947	1951	rVB4	53537	60277	1.17%	0.110%
42	14.534	1959	1963	1970	rBV	204645	393521	7.62%	0.717%
43	15.292	2086	2092	2099	rBV	2863030	4219863	81.74%	7.692%
44	15.363	2099	2104	2109	rBV2	225018	392650	7.61%	0.716%
45	15.428	2110	2115	2121	rVB	1110901	1533400	29.70%	2.795%
46	15.551	2131	2136	2141	rVB	584168	864486	16.74%	1.576%
47	16.628	2316	2319	2326	rBV5	125407	220719	4.28%	0.402%
48	17.045	2385	2390	2396	rBV	1718589	2452008	47.49%	4.470%
49	17.145	2402	2407	2415	rVB2	3043009	4458794	86.37%	8.127%
50	17.569	2475	2479	2485	rVB	312064	425266	8.24%	0.775%
51	17.969	2543	2547	2555	rBV	653008	1001444	19.40%	1.825%
52	18.186	2580	2584	2591	rBV	521982	745830	14.45%	1.359%
53	19.251	2763	2765	2773	rVB	327514	418516	8.11%	0.763%
54	19.451	2794	2799	2806	rVB	3566980	5162674	100.00%	9.410%
55	21.245	3100	3104	3112	rVB	2029302	2645483	51.24%	4.822%
56	23.327	3452	3458	3470	rVB	2232567	4397795	85.18%	8.016%
57	23.463	3476	3481	3490	rVB	1125908	2193066	42.48%	3.998%
58	23.980	3565	3569	3577	rVB	223185	418626	8.11%	0.763%

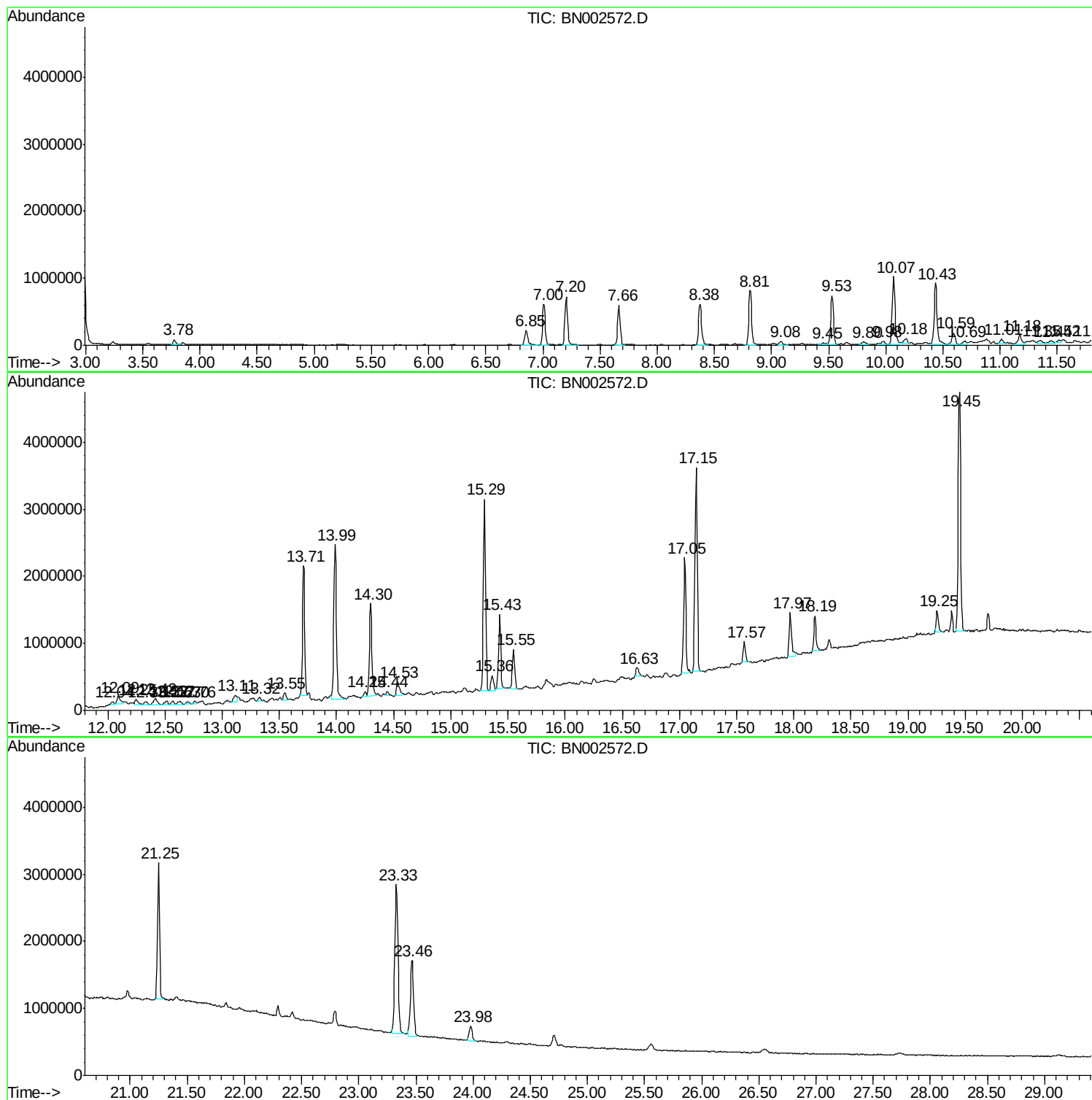
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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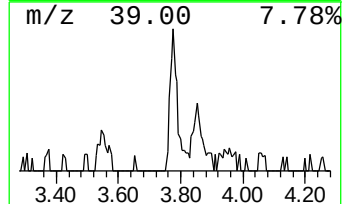
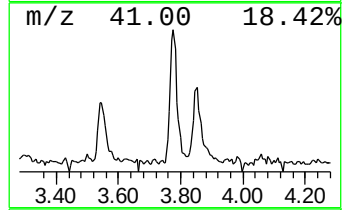
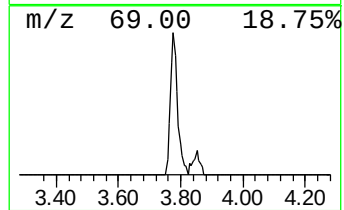
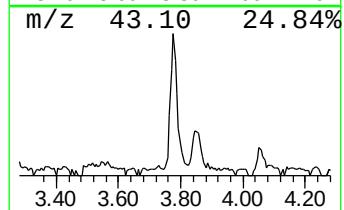
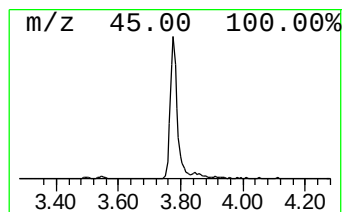
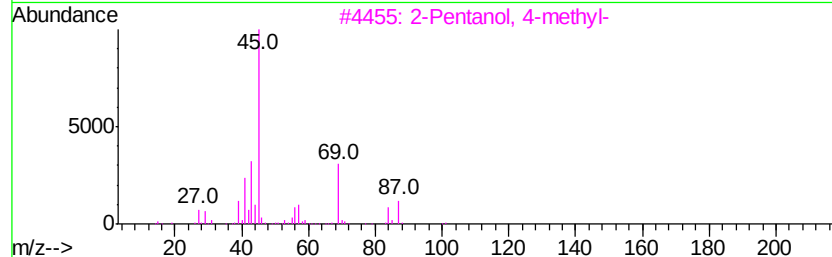
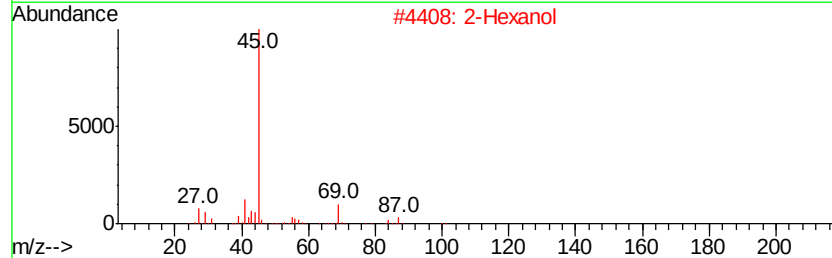
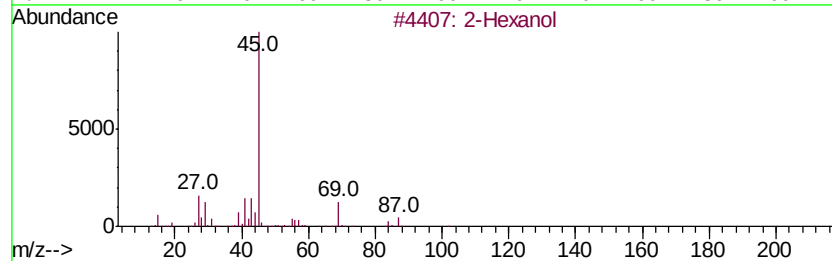
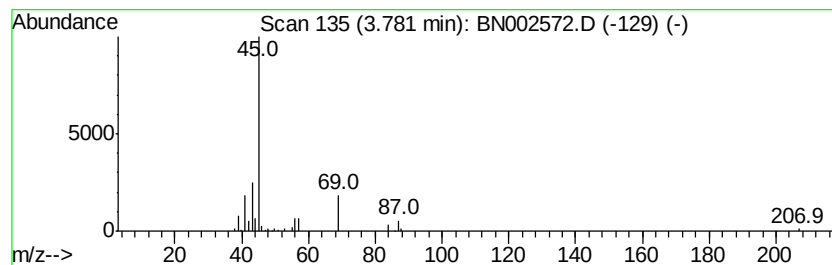
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Hexanol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.78	2.41 ng/ul	114264	1,4-Dichlorobenzene-d4	7.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol			102	C6H14O	000626-93-7	72
2	2-Hexanol			102	C6H14O	000626-93-7	56
3	2-Pentanol, 4-methyl-			102	C6H14O	000108-11-2	56
4	2-Hexanol			102	C6H14O	000626-93-7	56
5	2-Nonanol			144	C9H20O	000628-99-9	56



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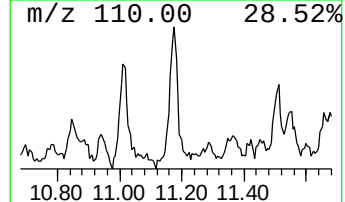
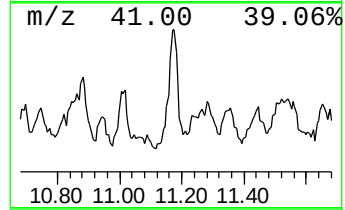
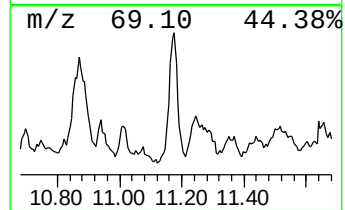
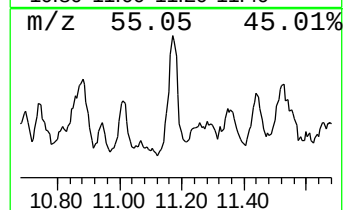
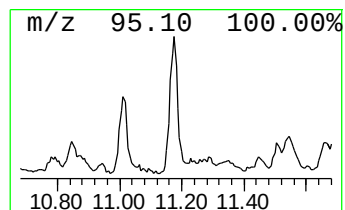
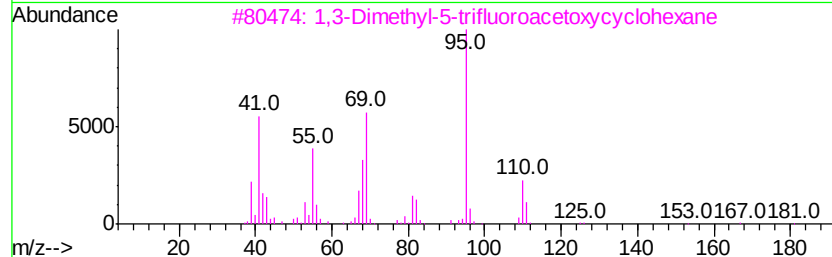
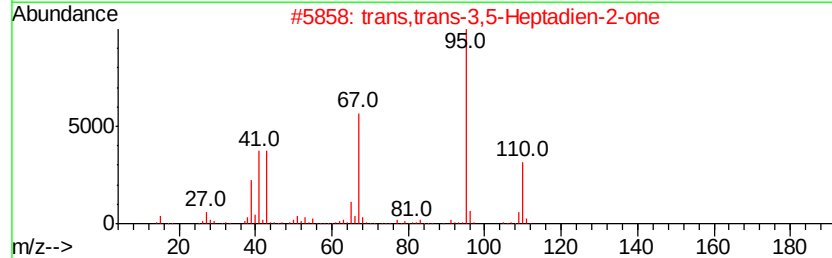
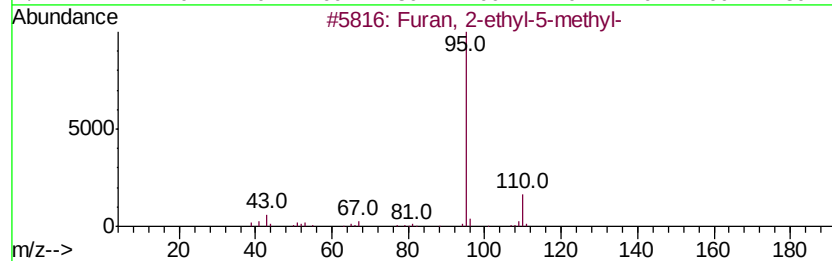
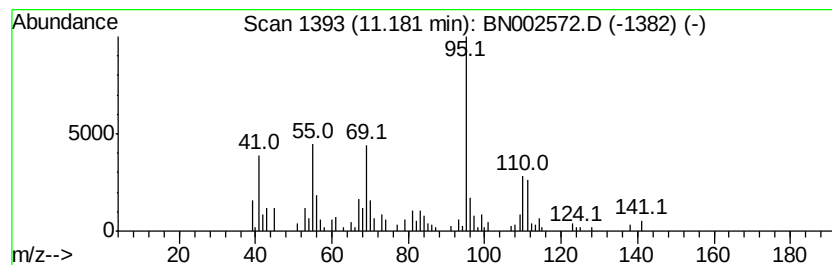
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Furan, 2-ethyl-5-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.18	3.51 ng/ul	291366	Napthalene-d8	10.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Furan, 2-ethyl-5-methyl-	110 C7H100	001703-52-2 52
2		trans,trans-3,5-Heptadien-2-one	110 C7H100	018402-90-9 50
3		1,3-Dimethyl-5-trifluoroacetoxycyclohexane	224 C10H15F3O2	1000216-63-7 42
4		Furan, 2-ethyl-5-methyl-	110 C7H100	001703-52-2 42
5		1,7,7-Trimethylbicyclo[2.2.1]hept-2-ene	154 C10H180	010385-78-1 42



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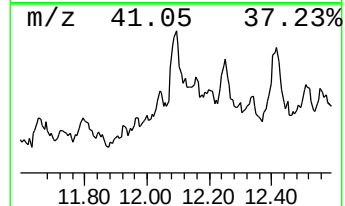
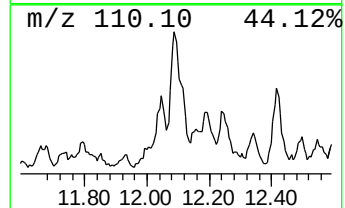
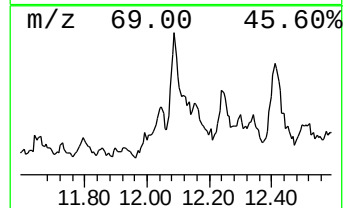
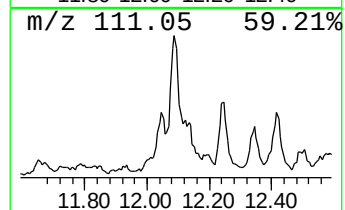
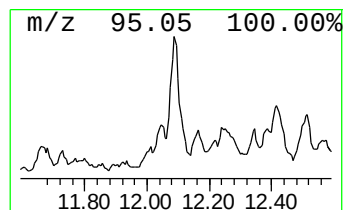
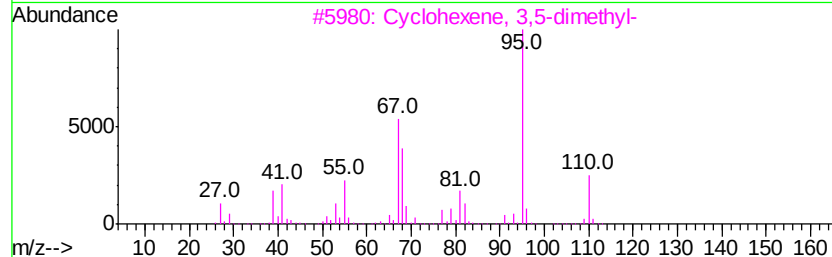
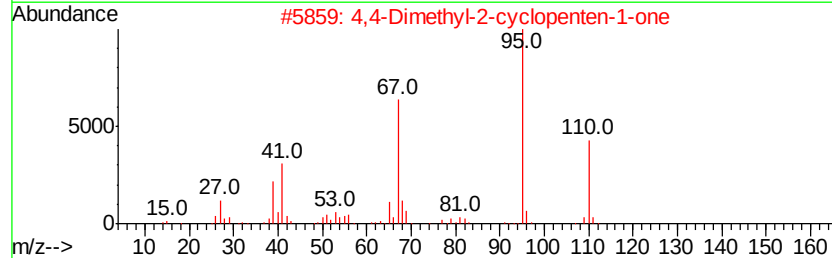
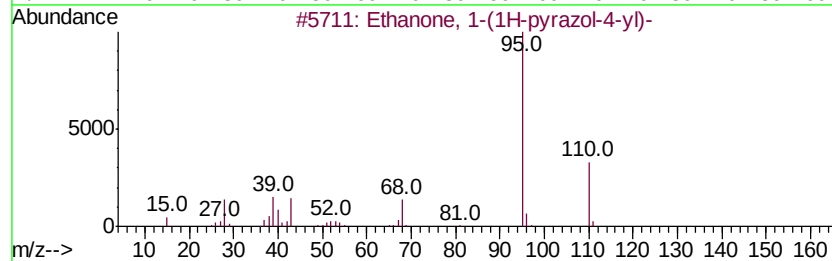
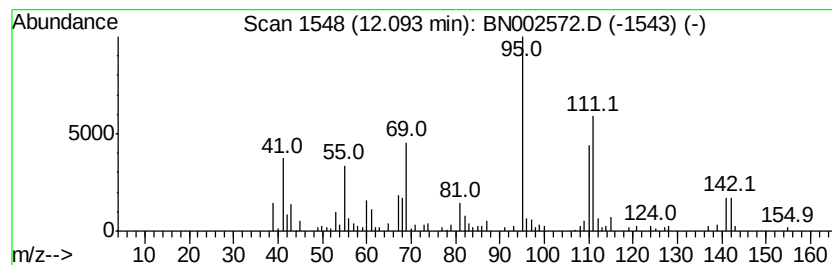
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST11.L
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Peak Number 3 Ethanone, 1-(1H-pyrazol-4-yl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.09	3.08 ng/ul	255396	Naphthalene-d8	10.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanone, 1-(1H-pyrazol-4-yl)-	110	C5H6N2O	025016-16-4	50
2		4,4-Dimethyl-2-cyclopenten-1-one	110	C7H10O	022748-16-9	47
3		Cyclohexene, 3,5-dimethyl-	110	C8H14	000823-17-6	47
4		4-Methyl-1,3-heptadiene	110	C8H14	017603-57-5	43
5		Ethanone, 1-(2-furanyl)-	110	C6H6O2	001192-62-7	43



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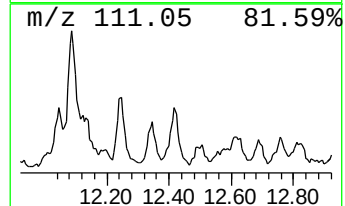
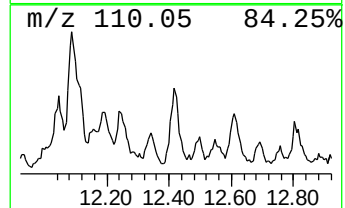
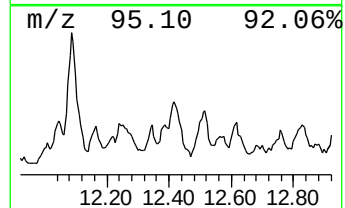
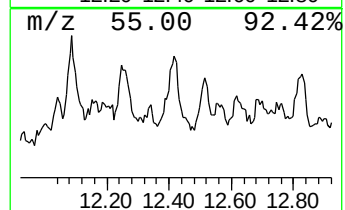
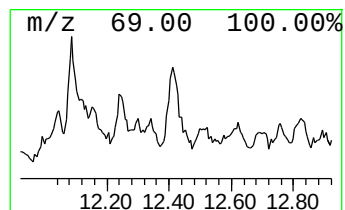
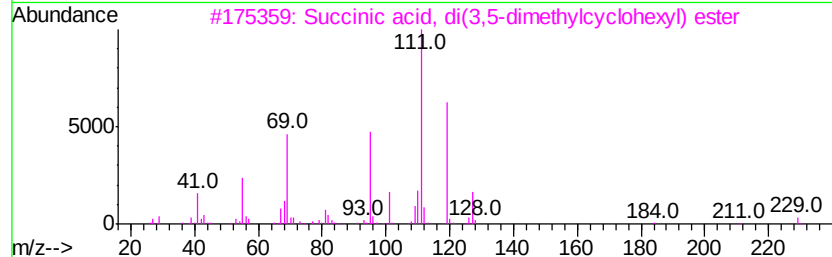
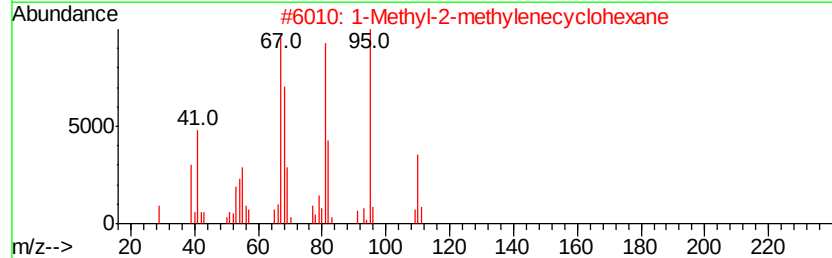
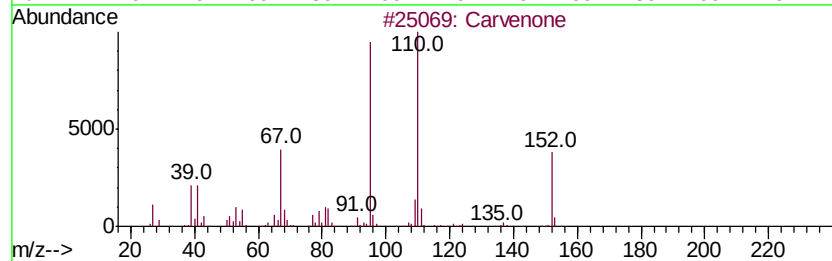
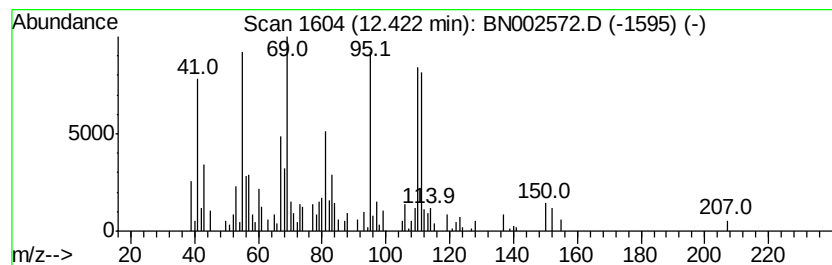
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.42	2.32 ng/ul	250412	Acenaphthene-d10	14.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Carvenone		152 C10H16O	000499-74-1 45
2	1-Methyl-2-methylenecyclohexane		110 C8H14	002808-75-5 43
3	Succinic acid, di(3,5-dimethylcv...		338 C20H34O4	1000330-05-9 43
4	Carvenone		152 C10H16O	000499-74-1 38
5	2,4-Hexadiene, 2,5-dimethyl-		110 C8H14	000764-13-6 38



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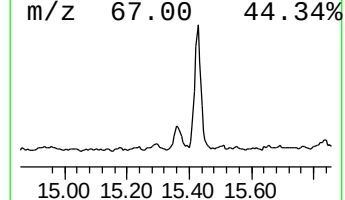
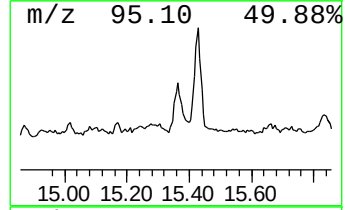
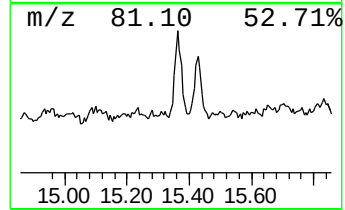
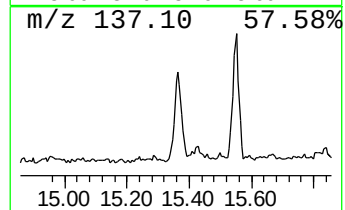
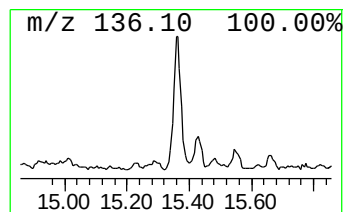
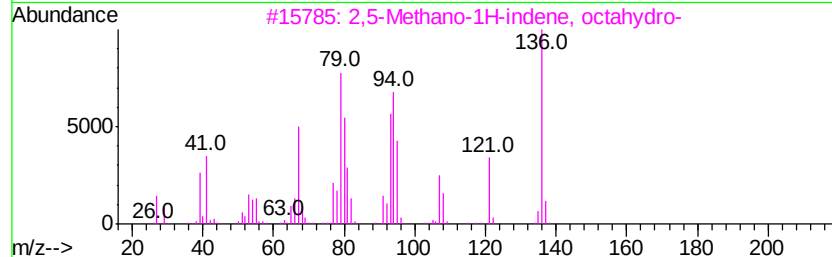
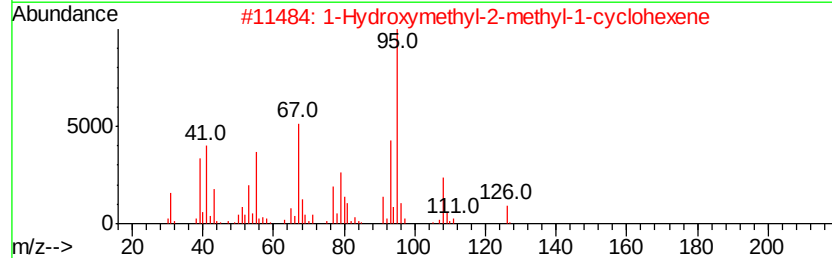
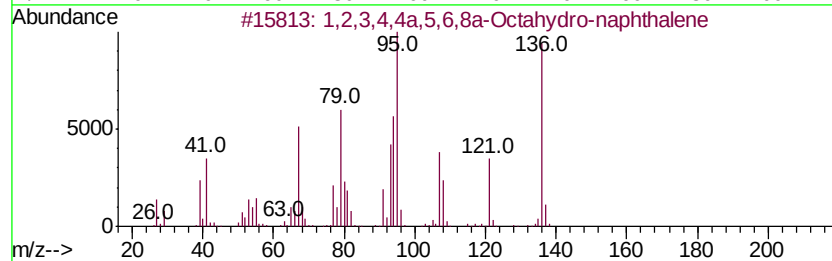
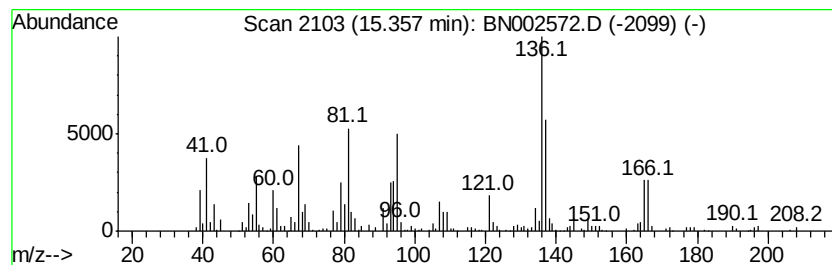
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1,2,3,4,4a,5,6,8a-Octahydro... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.36	3.64 ng/ul	392650	Acenaphthene-d10	14.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,4,4a,5,6,8a-Octahydro-naph...	136	C10H16	031244-58-3	58
2	1-Hydroxymethyl-2-methyl-1-cyclo...	126	C8H14O	029474-11-1	47
3	2,5-Methano-1H-indene, octahydro-	136	C10H16	019026-94-9	46
4	Phenol, 3-(dimethylamino)-	137	C8H11NO	000099-07-0	45
5	Phenol, 3-(dimethylamino)-	137	C8H11NO	000099-07-0	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN083118\
Data File : BN002572.D
Acq On : 31 Aug 2018 22:10
Operator : JU/SJ
Sample : J4665-15
Misc :
ALS Vial : 22 Sample Multiplier: 1

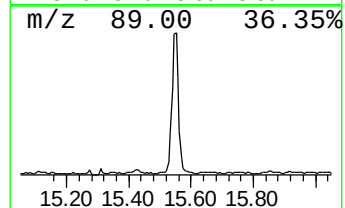
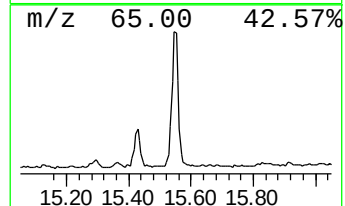
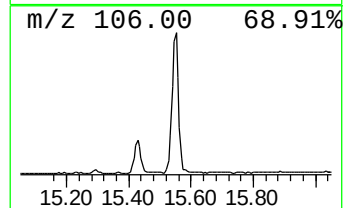
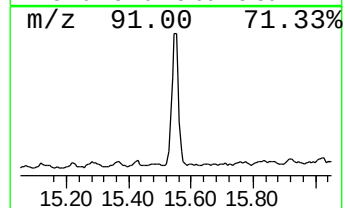
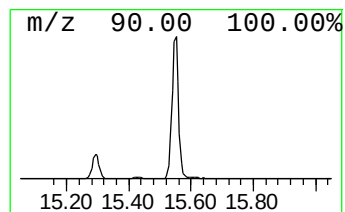
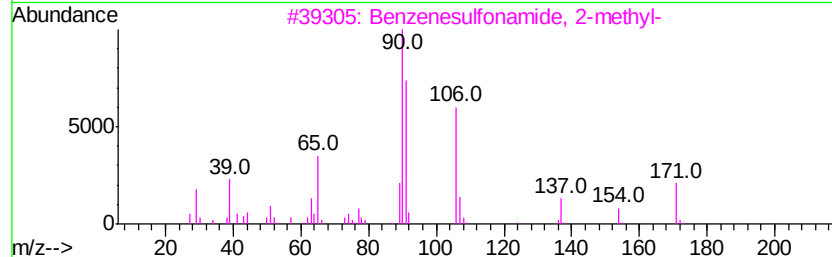
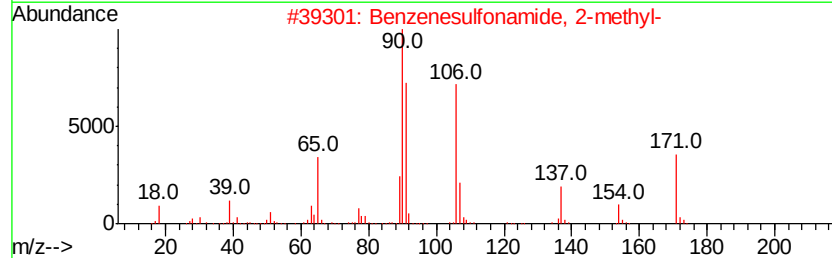
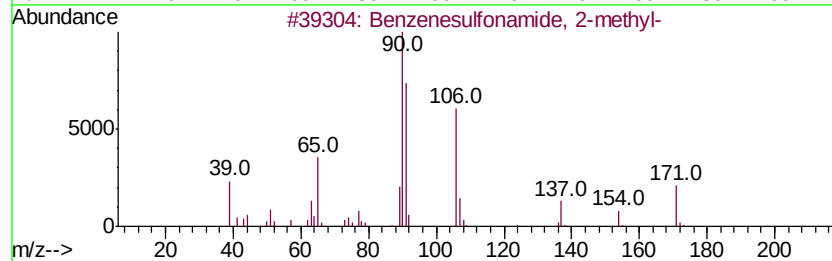
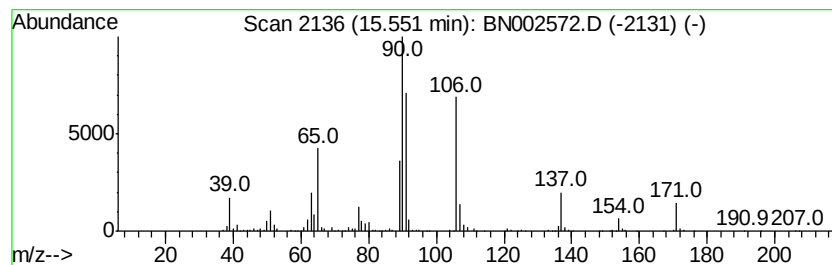
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzenesulfonamide, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	8.01 ng/ul	864486	Acenaphthene-d10	14.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzenesulfonamide, 2-methyl-	171 C7H9NO2S	000088-19-7	95
2	Benzenesulfonamide, 2-methyl-	171 C7H9NO2S	000088-19-7	95
3	Benzenesulfonamide, 2-methyl-	171 C7H9NO2S	000088-19-7	94
4	Benzenesulfonamide, 2-methyl-	171 C7H9NO2S	000088-19-7	94
5	Bicyclo[4.1.0]hepta-1,3,5-triene	90 C7H6	004646-69-9	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN083118\
Data File : BN002572.D
Acq On : 31 Aug 2018 22:10
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Sample : J4665-15
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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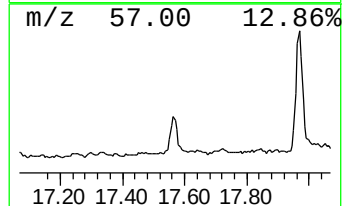
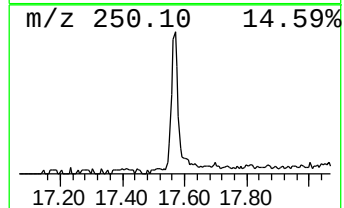
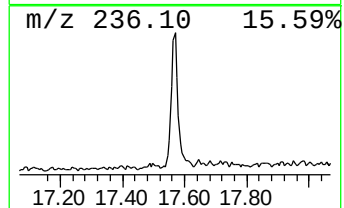
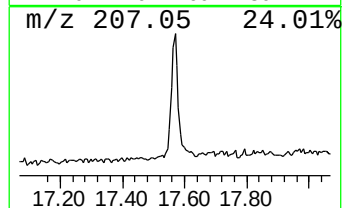
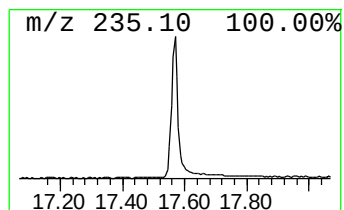
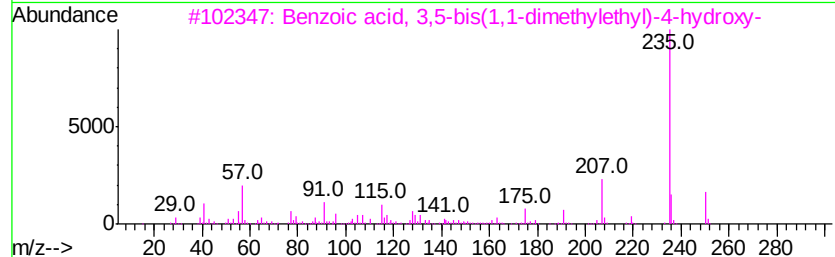
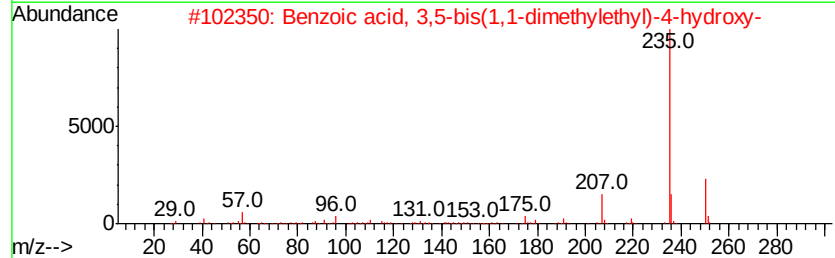
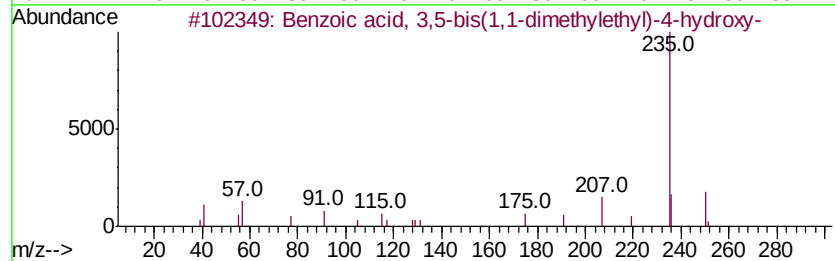
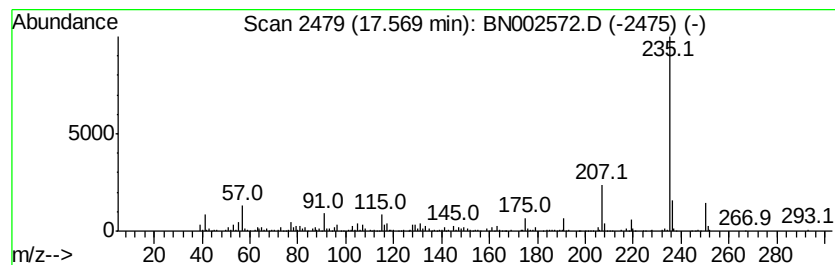
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzoic acid, 3,5-bis(1,1-d... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.57	3.47 ng/ul	425266	Phenanthrene-d10	17.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	98
2		Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	91
3		Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	87
4		6,7,8,9-Tetrahydro-1,2,3-trimeth...	250	C15H22O3	1000242-60-6	64
5		Hexobarbital-2(or 4)-methyl deri...	250	C13H18N2O3	1000123-51-5	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN083118\
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 Sample : J4665-15
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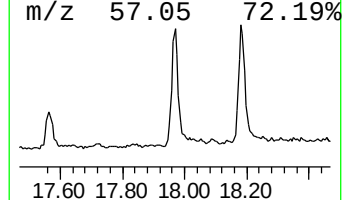
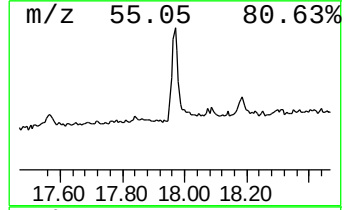
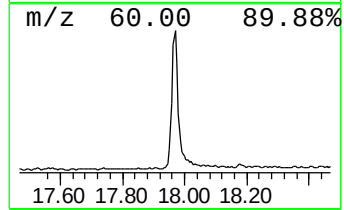
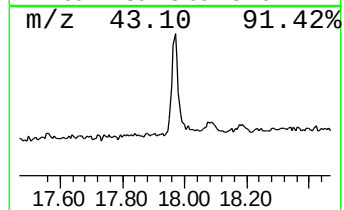
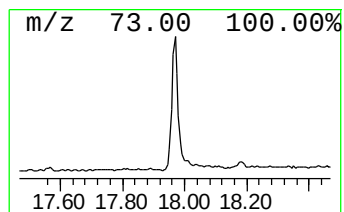
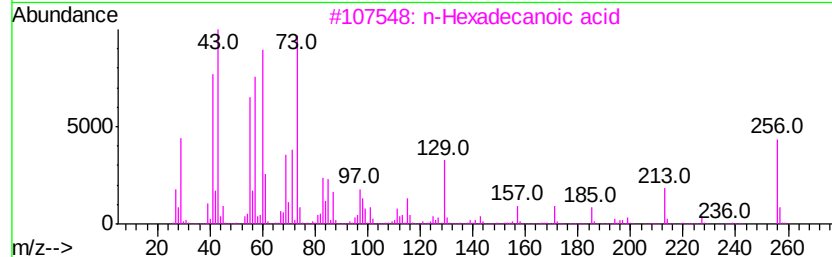
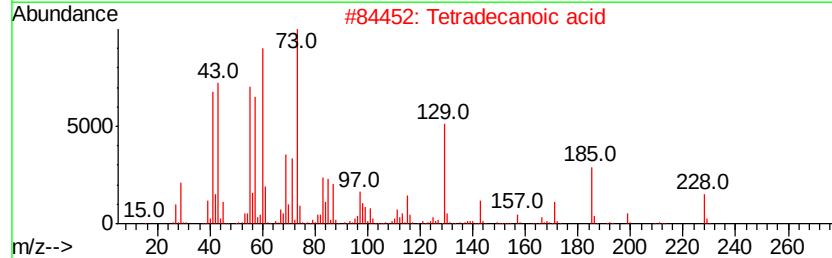
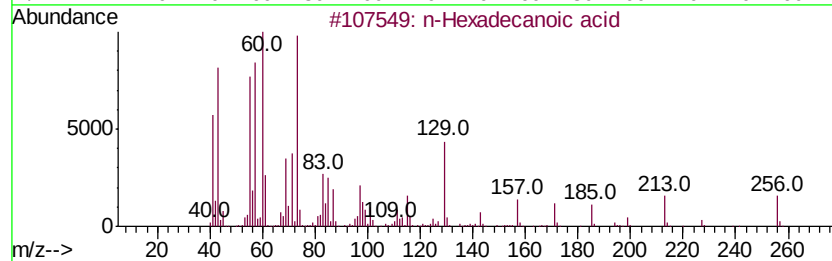
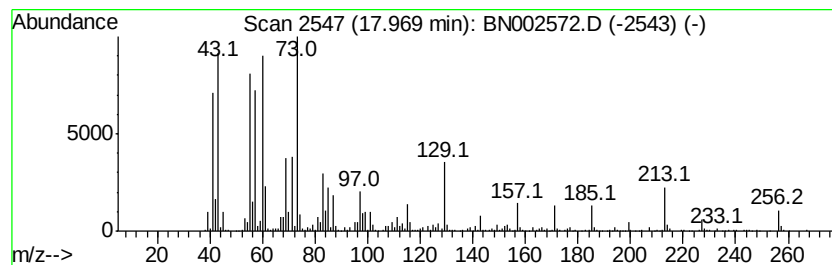
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 Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 8 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.97	8.17 ng/ul	1001440	Phenanthrene-d10	17.05	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	95
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
4	Tridecanoic acid	214	C13H26O2	000638-53-9	89
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN083118\
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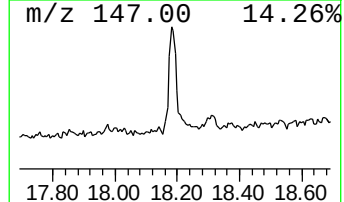
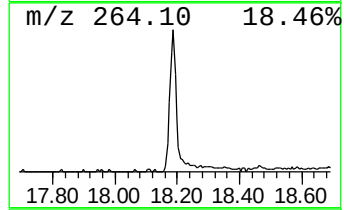
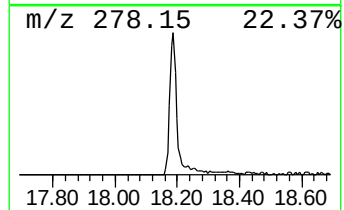
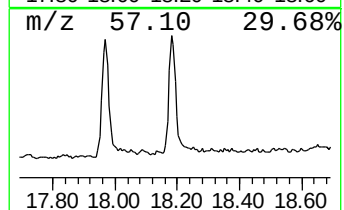
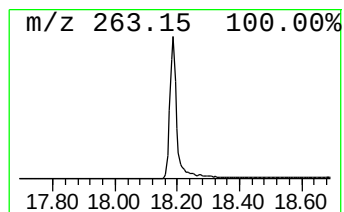
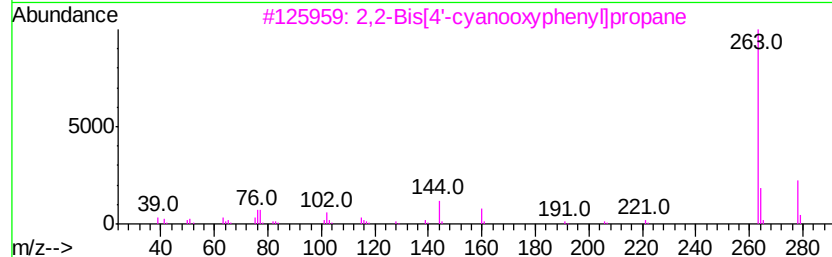
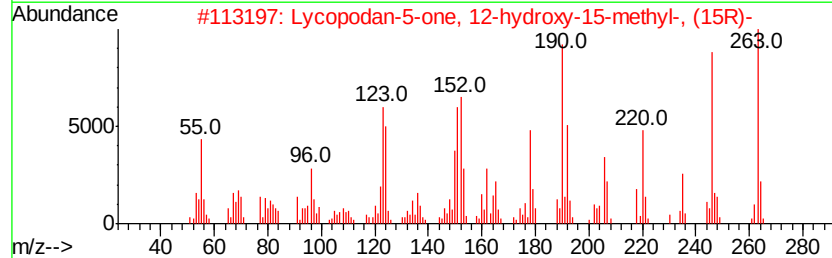
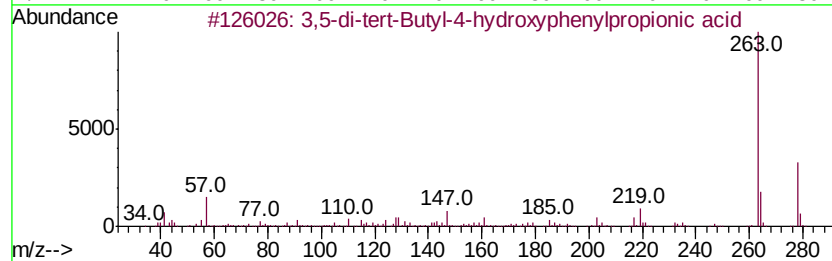
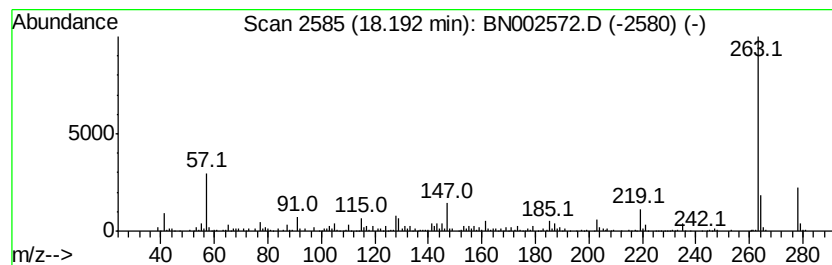
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.19	6.08 ng/ul	745830	Phenanthrene-d10	17.05		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-di-tert-Butyl-4-hydroxyphenyl...	278	C17H26O3	020170-32-5	94	
2	Lycopodan-5-one, 12-hydroxy-15-m...	263	C16H25NO2	006900-92-1	92	
3	2,2-Bis[4'-cyanoxyphenyl]propane	278	C17H14N2O2	1000322-13-3	72	
4	3,5-di-tert-Butyl-4-hydroxyphenyl...	278	C17H26O3	020170-32-5	70	
5	Ethanone, [1-(2,5-diphenyl-2H-1,...	263	C16H13N3O	322647-34-7	52	



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN083118\
Data File : BN002572.D
Acq On : 31 Aug 2018 22:10
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Misc :
ALS Vial : 22 Sample Multiplier: 1

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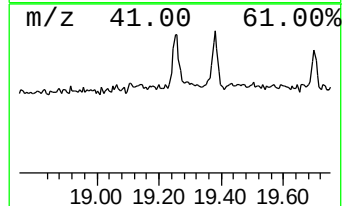
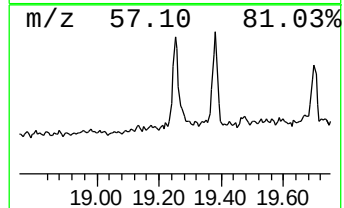
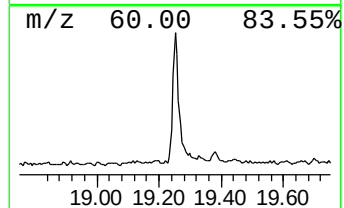
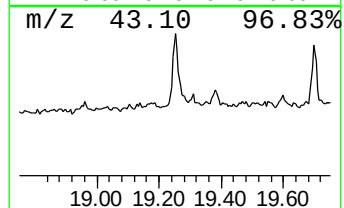
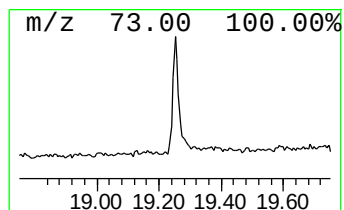
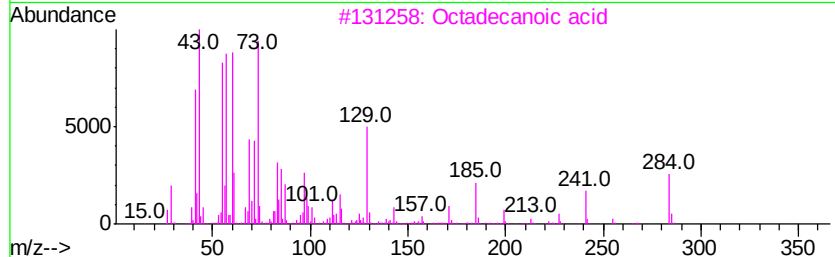
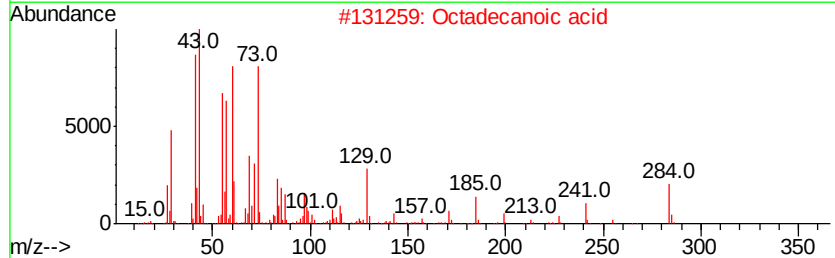
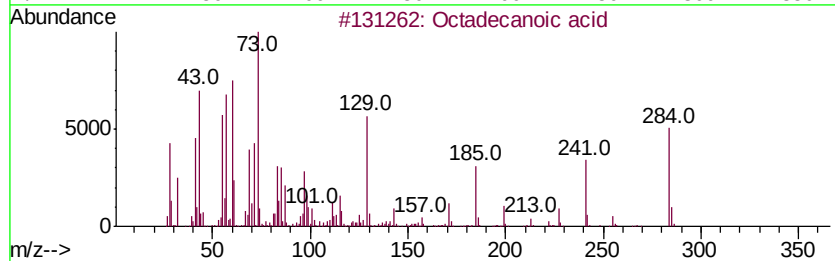
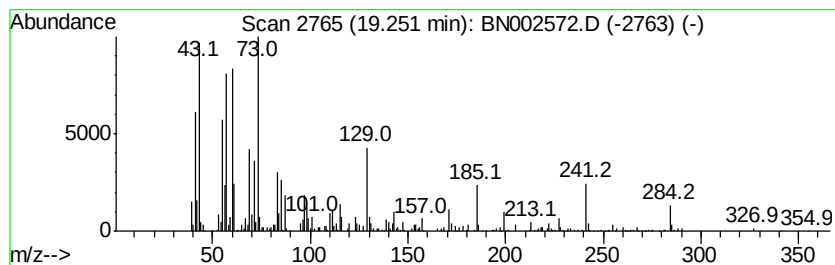
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.25	3.16 ng/ul	418516	Chrysene-d12	21.25

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	98
3			Octadecanoic acid	284	C18H36O2	000057-11-4	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	95



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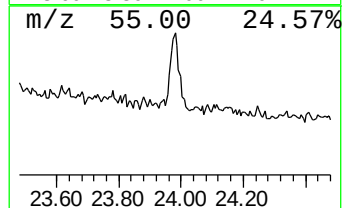
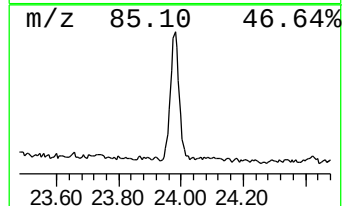
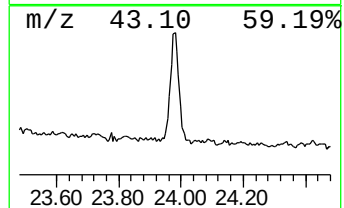
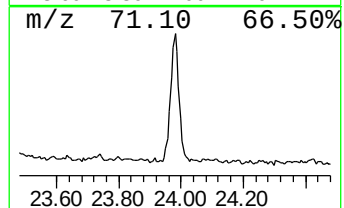
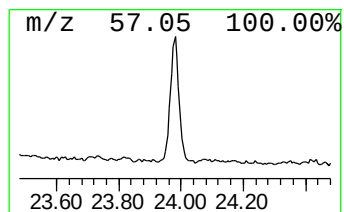
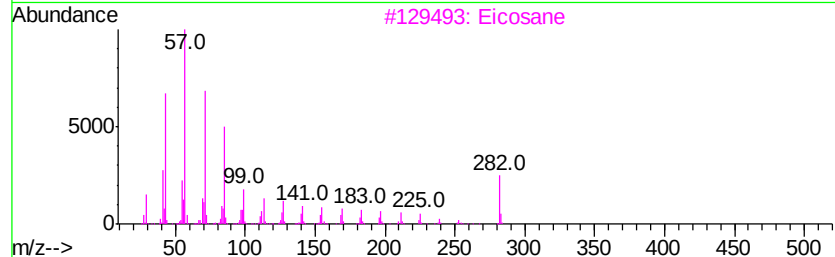
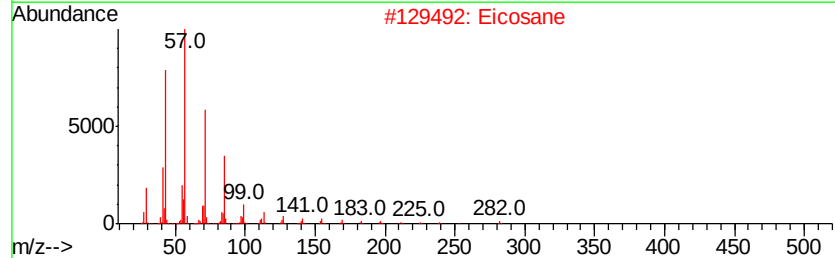
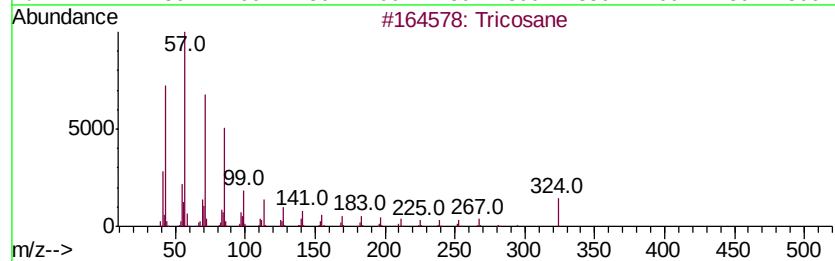
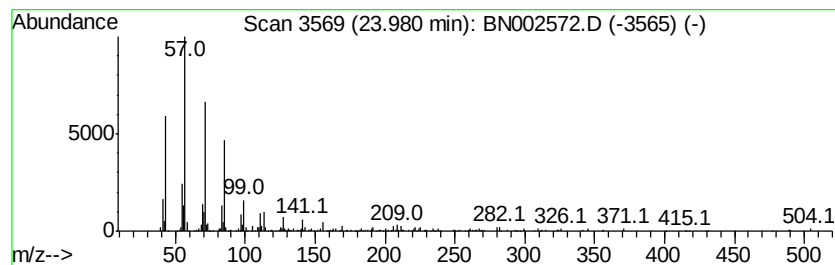
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.98	3.82 ng/ul	418626	Perylene-d12	23.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	93
2		Eicosane	282	C20H42	000112-95-8	93
3		Eicosane	282	C20H42	000112-95-8	93
4		Heneicosane	296	C21H44	000629-94-7	90
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN083118\
Data File : BN002572.D
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Operator : JU/SJ
Sample : J4665-15
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Hexanol	3.78	2.4	ng/ul	114264	1	7.66	947863	20.0
Furan, 2-ethyl-5-...	11.18	3.5	ng/ul	291366	2	10.43	1660990	20.0
Ethanone, 1-(1H-p...	12.09	3.1	ng/ul	255396	2	10.43	1660990	20.0
unknown-01	12.42	2.3	ng/ul	250412	3	14.30	2158460	20.0
1,2,3,4,4a,5,6,8a...	15.36	3.6	ng/ul	392650	3	14.30	2158460	20.0
Benzenesulfonamid...	15.55	8.0	ng/ul	864486	3	14.30	2158460	20.0
Benzoic acid, 3,5...	17.57	3.5	ng/ul	425266	4	17.05	2452010	20.0
n-Hexadecanoic acid	17.97	8.2	ng/ul	1001440	4	17.05	2452010	20.0
3,5-di-tert-Butyl...	18.19	6.1	ng/ul	745830	4	17.05	2452010	20.0
Octadecanoic acid	19.25	3.2	ng/ul	418516	5	21.25	2645480	20.0
(DEL) Alkane: Str...	23.98	3.8	ng/ul	418626	6	23.47	2193070	20.0