

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121218\
 Data File : BN003881.D
 Acq On : 13 Dec 2018 08:49
 Operator : JU/SJ
 Sample : J6228-07
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E86

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-BN121218MA.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	3.017	3	5	26	rVB	280691	427455	52.07%	5.075%
2	4.946	326	333	341	rVB2	4449	10294	1.25%	0.122%
3	6.987	674	680	690	rVB2	44314	85988	10.48%	1.021%
4	7.164	705	710	718	rBB	36032	55402	6.75%	0.658%
5	7.358	738	743	750	rBB	46228	69737	8.50%	0.828%
6	7.828	818	823	830	rBB	71151	110891	13.51%	1.316%
7	8.522	936	941	949	rBB	46978	70937	8.64%	0.842%
8	8.999	1016	1022	1028	rBB	45708	73442	8.95%	0.872%
9	9.710	1138	1143	1149	rBB	35962	55171	6.72%	0.655%
10	10.240	1228	1233	1242	rBB	59953	93406	11.38%	1.109%
11	10.628	1292	1299	1307	rBB	121957	200049	24.37%	2.375%
12	10.769	1318	1323	1332	rBB	16522	25109	3.06%	0.298%
13	13.869	1845	1850	1854	rBV	134955	175017	21.32%	2.078%
14	13.916	1854	1858	1864	rVB	52158	68730	8.37%	0.816%
15	14.157	1893	1899	1908	rBV	155085	226651	27.61%	2.691%
16	14.469	1946	1952	1960	rBV2	205906	311653	37.97%	3.700%
17	14.663	1979	1985	1997	rBV	30934	51752	6.30%	0.614%
18	15.457	2114	2120	2129	rVB	222286	300419	36.60%	3.566%
19	15.581	2136	2141	2145	rBV	38323	46956	5.72%	0.557%
20	17.210	2412	2418	2424	rBV	300903	420966	51.28%	4.997%
21	17.310	2429	2435	2443	rVV	241269	342149	41.68%	4.062%
22	18.075	2562	2565	2574	rBV2	12962	16278	1.98%	0.193%
23	19.357	2780	2783	2790	rBV	8524	11982	1.46%	0.142%
24	19.504	2803	2808	2813	rBV2	36529	38532	4.69%	0.457%
25	19.604	2819	2825	2832	rBV	360755	470005	57.26%	5.580%
26	20.551	2982	2986	2991	rBV	38715	42946	5.23%	0.510%
27	21.092	3073	3078	3095	rBV9	13428	54715	6.67%	0.650%
28	21.392	3124	3129	3135	rBV	559253	668988	81.50%	7.942%
29	21.957	3221	3225	3229	rBV	20165	23322	2.84%	0.277%
30	22.433	3302	3306	3312	rVB	23993	30493	3.71%	0.362%
31	22.563	3325	3328	3333	rVB	13037	15682	1.91%	0.186%
32	22.951	3389	3394	3401	rVB	53206	76726	9.35%	0.911%
33	23.557	3488	3497	3512	rVB	365929	741441	90.33%	8.802%
34	23.710	3515	3523	3532	rVB	433182	820850	100.00%	9.745%

LSC Area Percent Report

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35	23.863	3545	3549	3554	rBV3	10800	17453	2.13%	0.207%
36	24.198	3600	3606	3614	rBV	54443	106263	12.95%	1.261%
37	24.969	3731	3737	3745	rVB2	32273	68367	8.33%	0.812%
38	25.868	3883	3890	3899	rVB	33913	84391	10.28%	1.002%
39	26.233	3943	3952	3964	rVB2	192091	535507	65.24%	6.357%
40	27.263	4118	4127	4136	rBV2	78871	234677	28.59%	2.786%
41	27.351	4137	4142	4153	rVB4	30090	85963	10.47%	1.021%
42	27.810	4211	4220	4226	rBV7	85759	307155	37.42%	3.646%
43	28.568	4340	4349	4360	rVB3	40965	148110	18.04%	1.758%
44	29.004	4417	4423	4435	rVB5	40358	138180	16.83%	1.640%
45	29.204	4444	4457	4477	rVB8	81817	463375	56.45%	5.501%

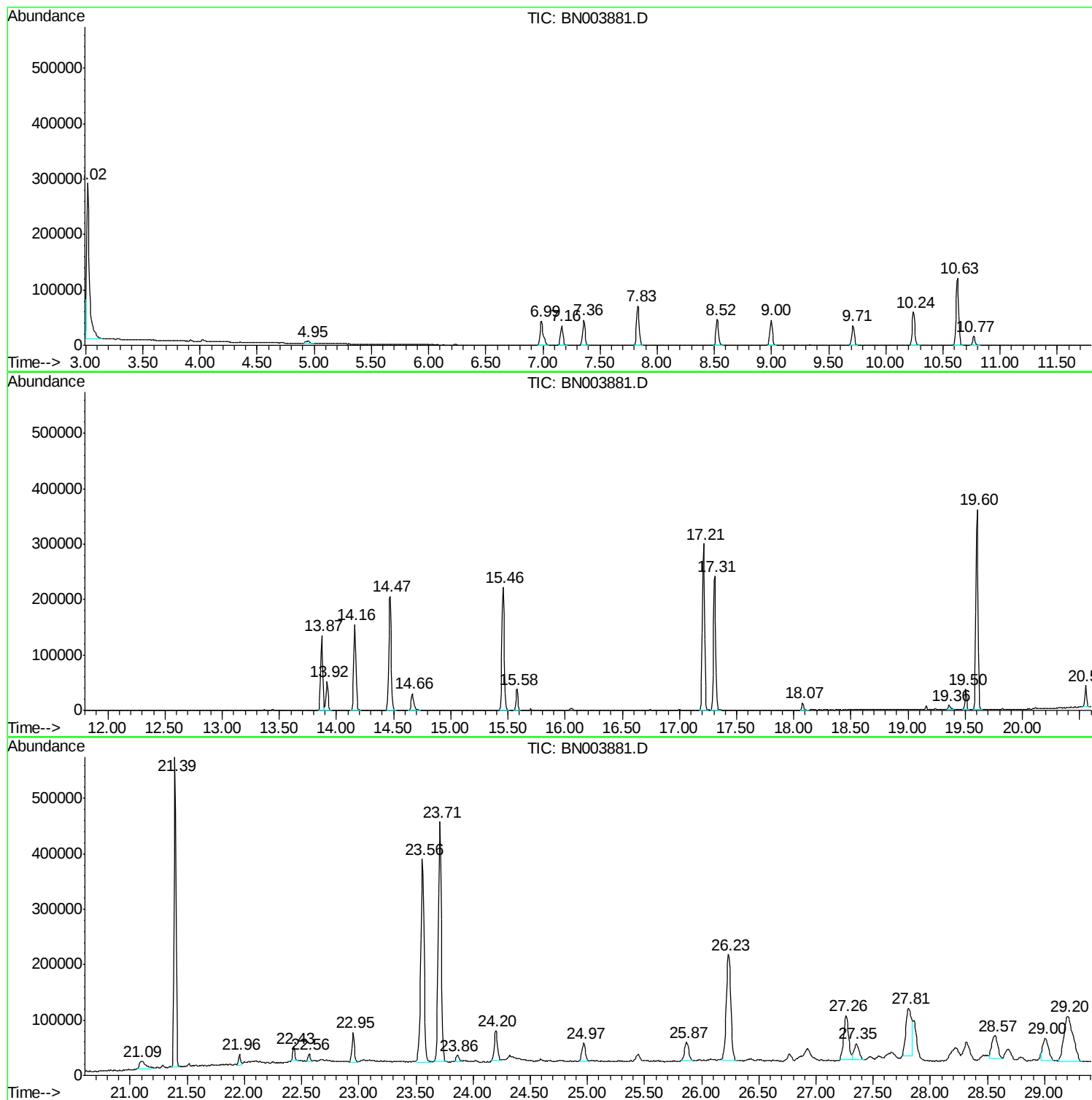
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN121218MA.M
Quant Title : SVOA CALIBRATION

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



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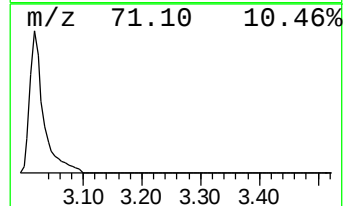
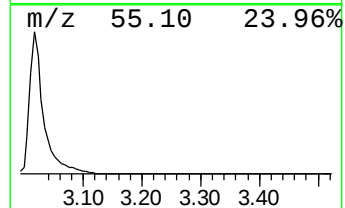
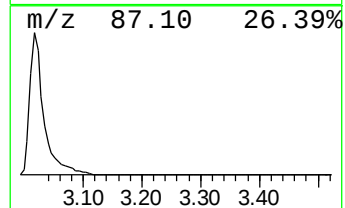
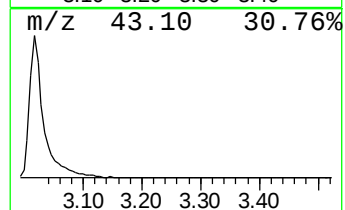
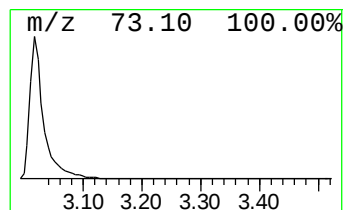
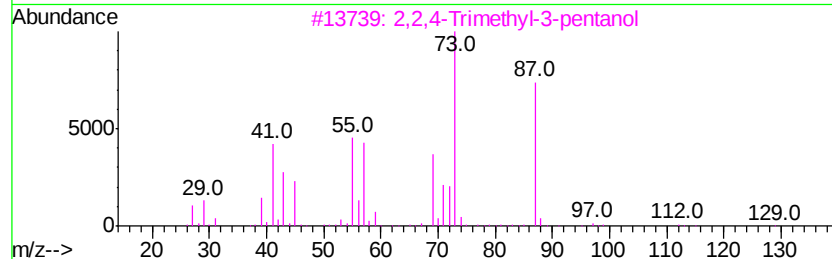
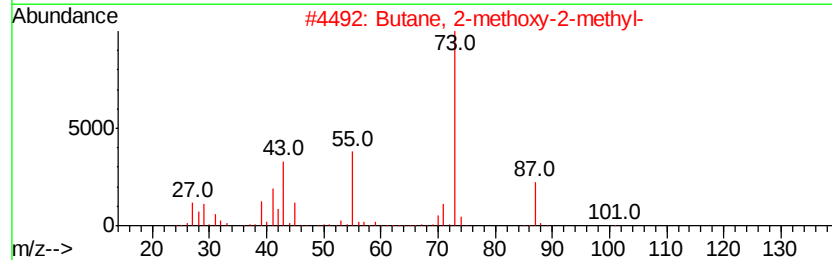
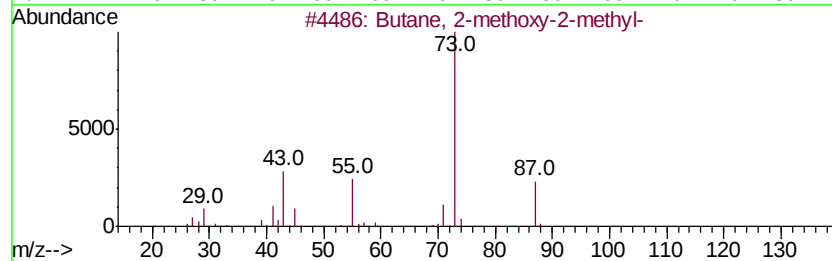
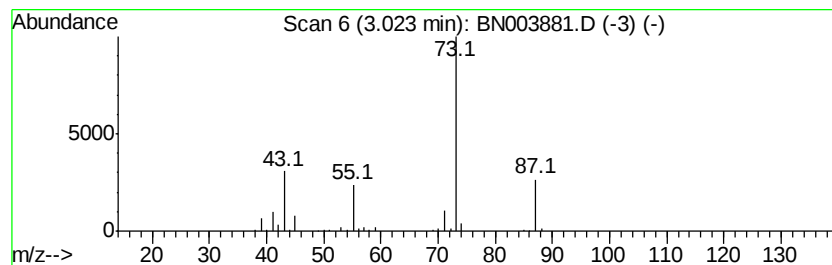
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.02	77.09 ng/ul	427455	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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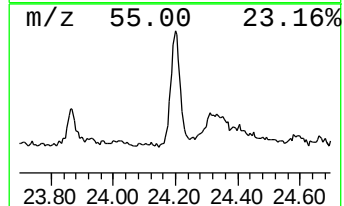
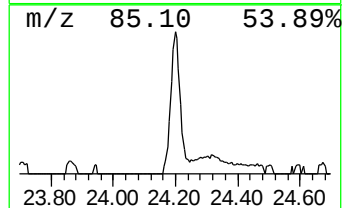
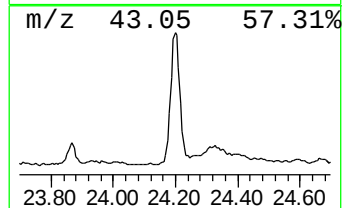
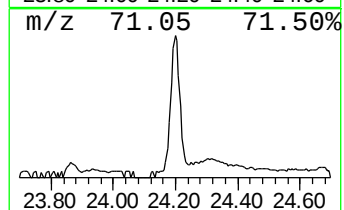
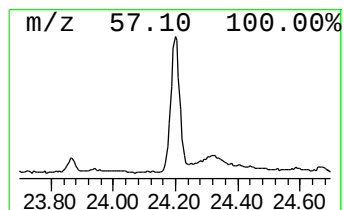
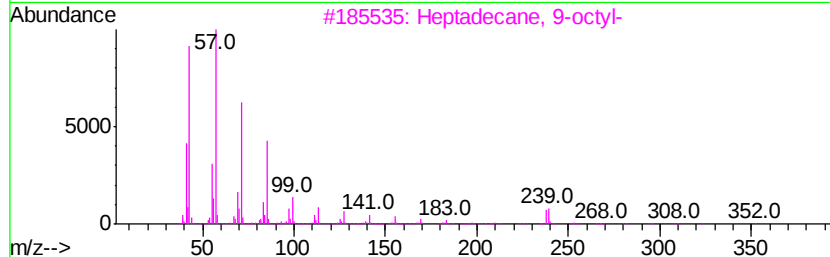
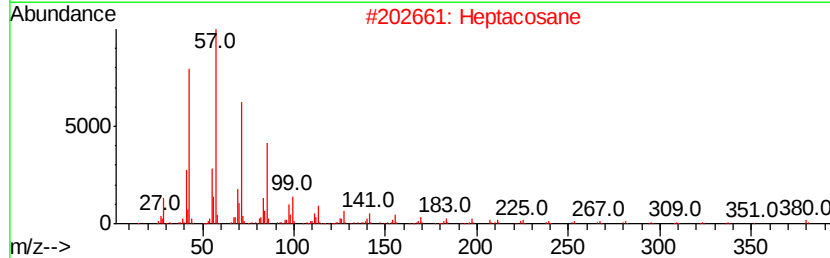
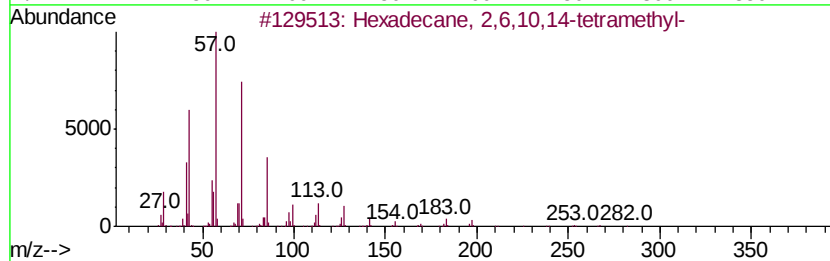
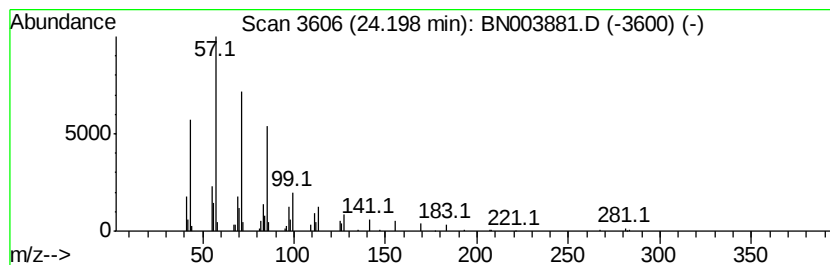
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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.20	2.59 ng/ul	106263	Perylene-d12	23.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	92
2		Heptacosane	380	C27H56	000593-49-7	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Heneicosane	296	C21H44	000629-94-7	91



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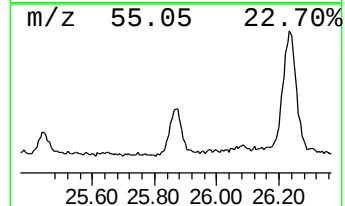
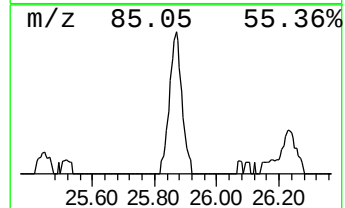
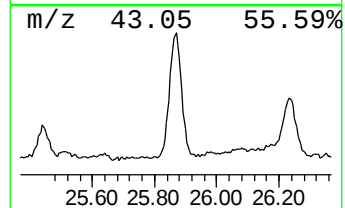
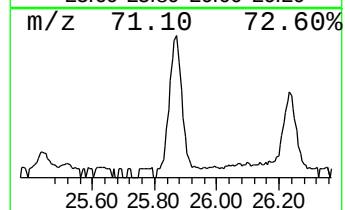
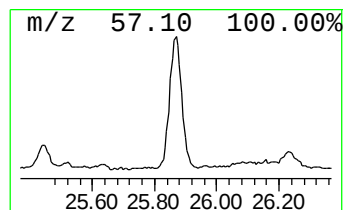
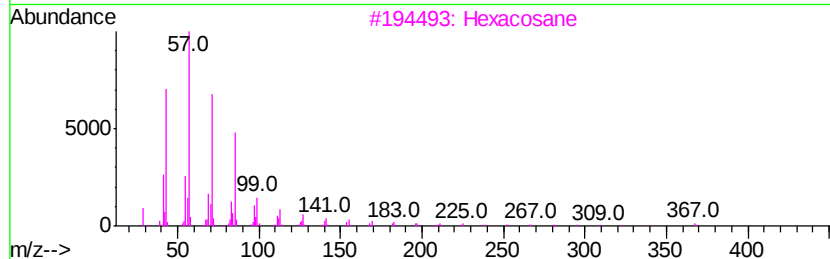
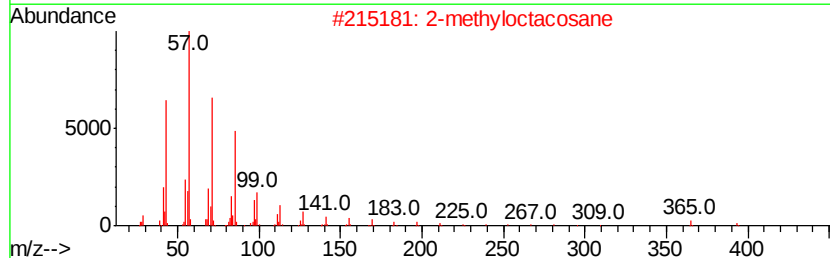
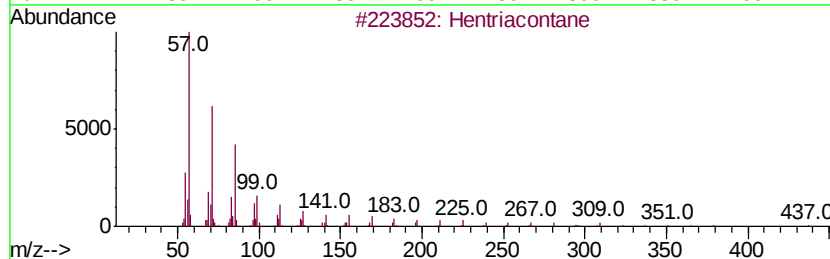
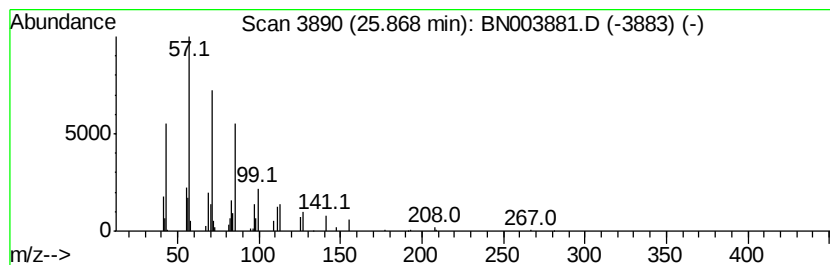
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Quant Title : SVOA CALIBRATION

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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.87	2.06 ng/ul	84391	Perylene-d12	23.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hentriacontane	437	C31H64	000630-04-6	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Octacosane	394	C28H58	000630-02-4	91



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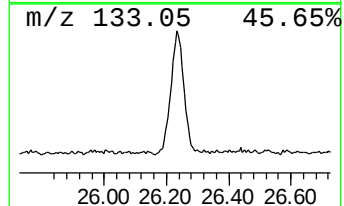
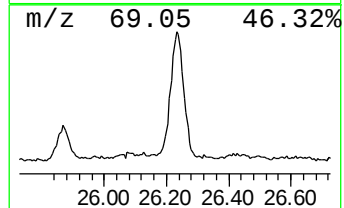
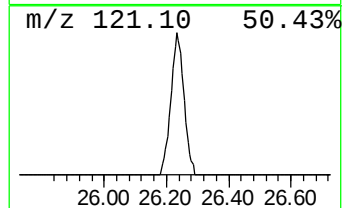
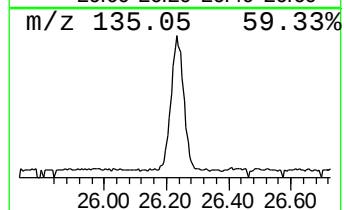
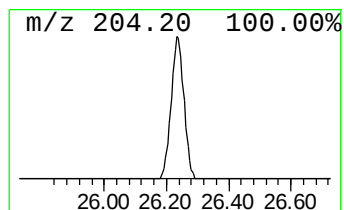
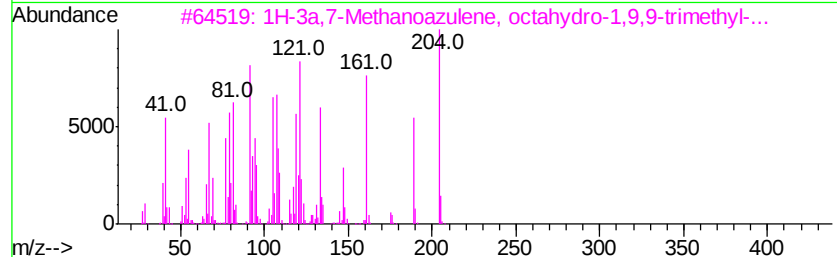
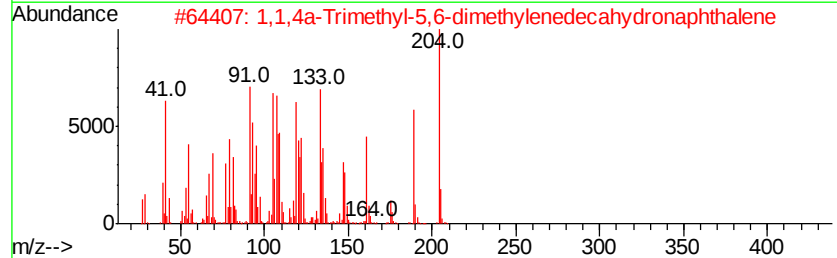
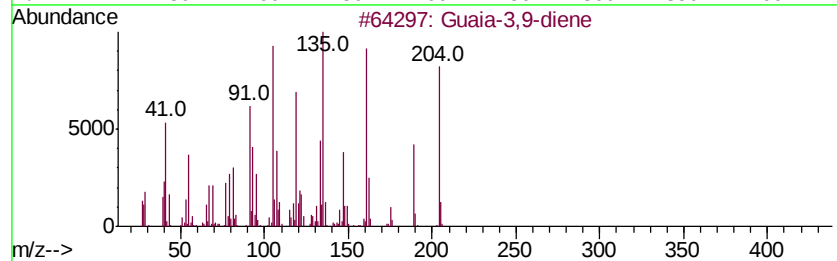
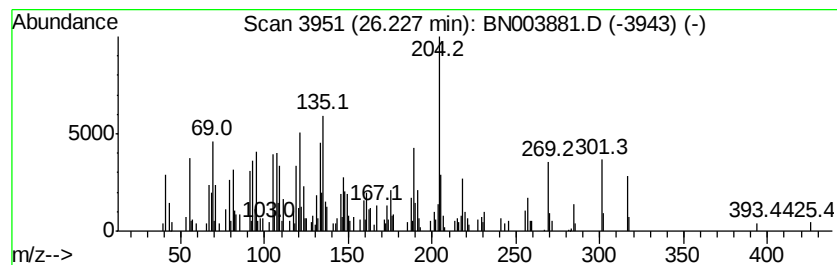
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Peak Number 4 Guaia-3,9-diene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.23	13.05 ng/ul	535507	Perylene-d12	23.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Guaia-3,9-diene	204	C15H24	000489-83-8	53
2		1,1,4a-Trimethyl-5,6-dimethylene...	204	C15H24	1000193-60-8	50
3		1H-3a,7-Methanoazulene, octahydr...	204	C15H24	000508-55-4	47
4		Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	004630-07-3	46
5		2,5,5,8a-Tetramethyl-6,7,8,8a-te...	204	C14H20O	124957-09-1	45



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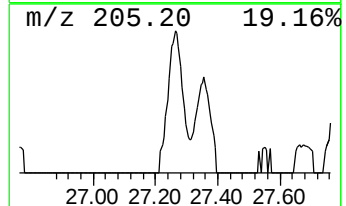
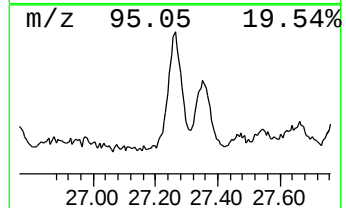
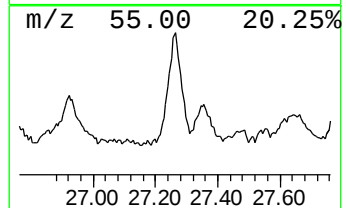
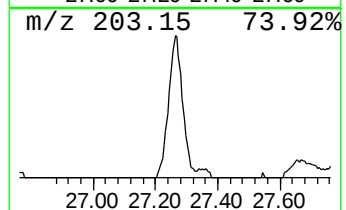
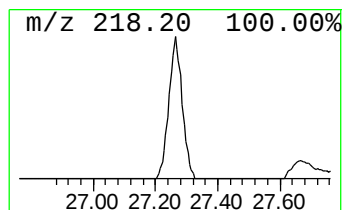
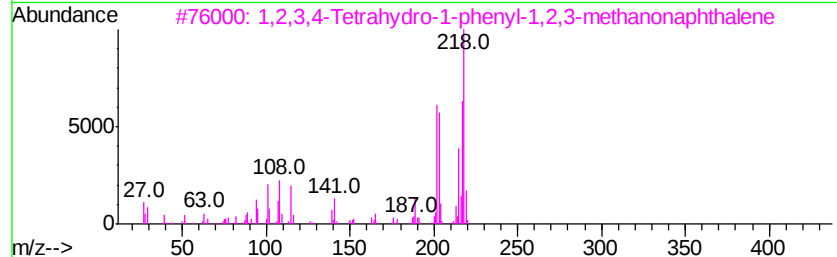
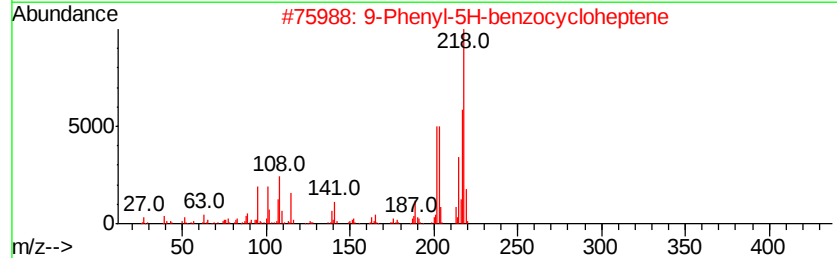
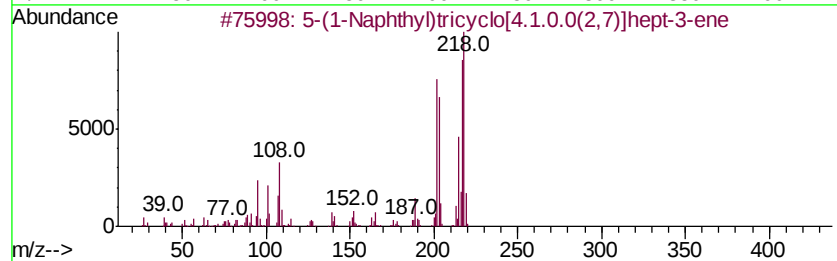
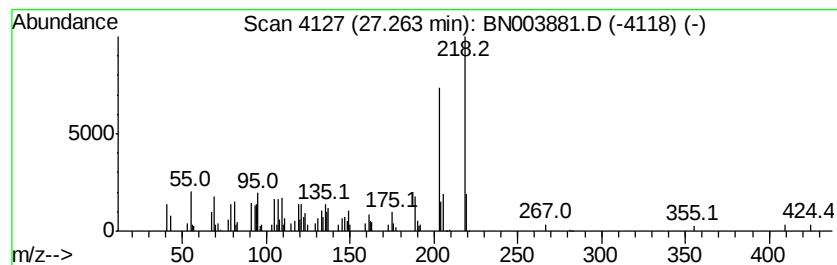
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 5-(1-Naphthyl)tricyclo[4.1.... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.26	5.72 ng/ul	234677	Perylene-d12	23.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-(1-Naphthyl)tricyclo[4.1.0.0(2...	218	C17H14	107729-05-5	87		
2	9-Phenyl-5H-benzocycloheptene	218	C17H14	138089-71-1	80		
3	1,2,3,4-Tetrahydro-1-phenyl-1,2,...	218	C17H14	138089-69-7	72		
4	3a-Methyl-3,3a-dihydroperrolo[3,...	218	C11H10N2OS	017163-68-7	64		
5	8-Amino-6,7-dimethoxy-4-methylqu...	218	C12H14N2O2	1000213-07-2	59		



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121218\
Data File : BN003881.D
Acq On : 13 Dec 2018 08:49
Operator : JU/SJ
Sample : J6228-07
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampled :
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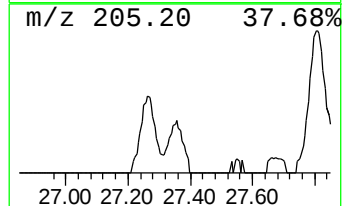
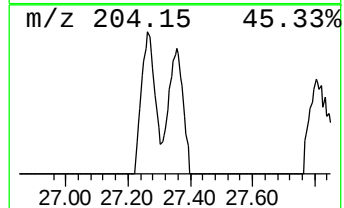
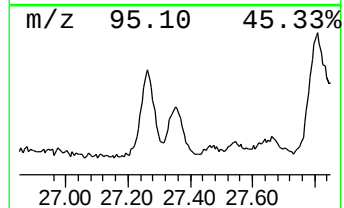
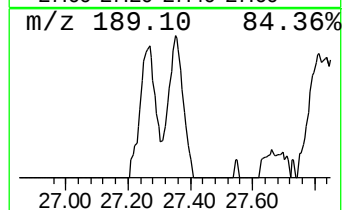
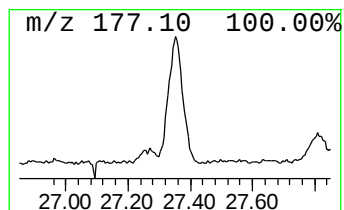
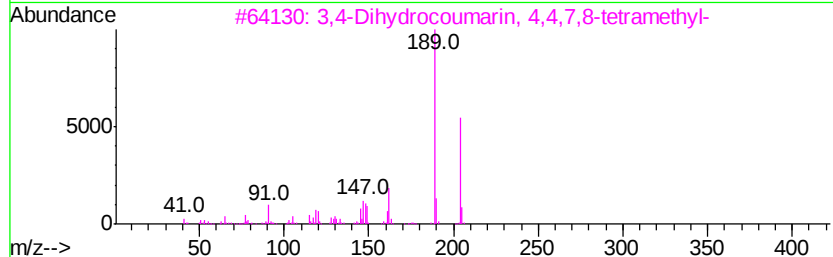
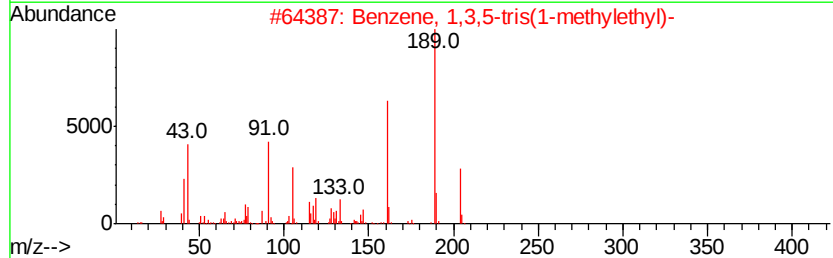
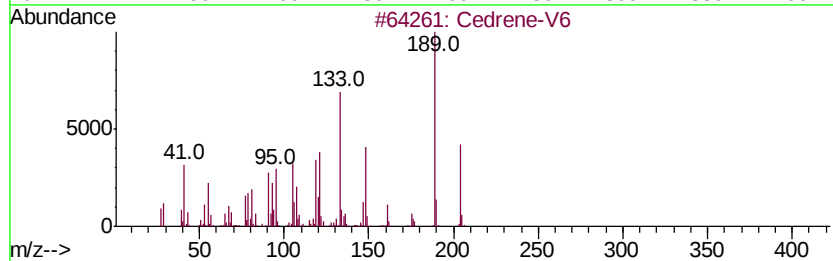
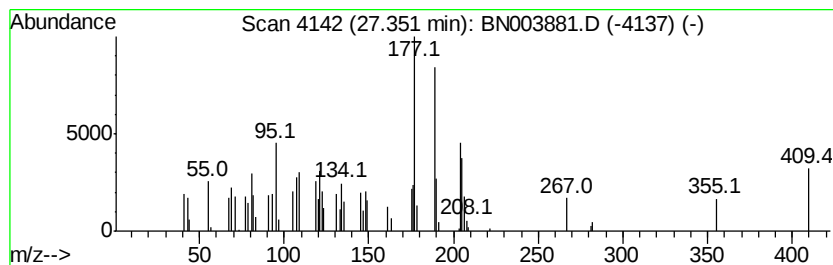
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.35	2.09 ng/ul	85963	Perylene-d12	23.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cedrene-V6	204	C15H24	1000162-76-8	35
2		Benzene, 1,3,5-tris(1-methylethyl)-	204	C15H24	000717-74-8	30
3		3,4-Dihydrocoumarin, 4,4,7,8-tet...	204	C13H16O2	040614-36-6	27
4		7-Chloro-3-methyl-quinoline-8-ca...	249	C13H12ClN02	1000373-12-1	27
5		7-Chloro-3-methyl-quinoline-8-ca...	277	C15H16ClN02	1000373-13-5	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121218\
Data File : BN003881.D
Acq On : 13 Dec 2018 08:49
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Sample : J6228-07
Misc :
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Instrument :
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ClientSampleId :
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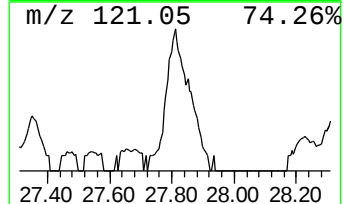
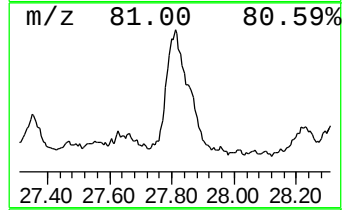
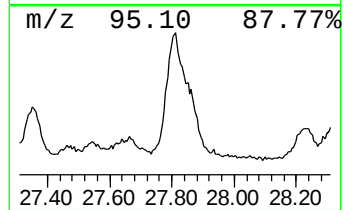
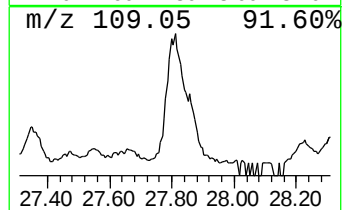
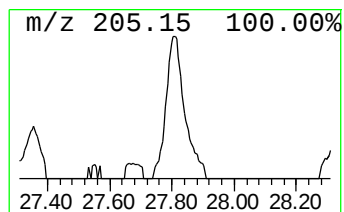
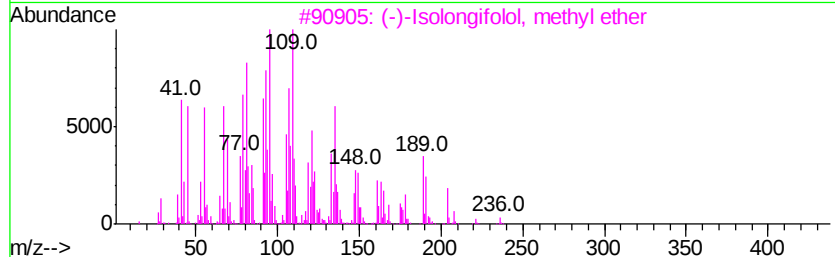
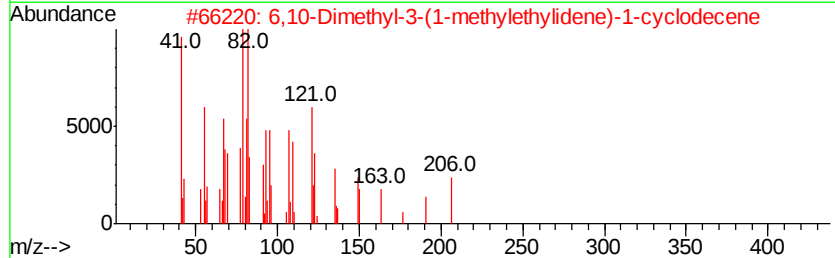
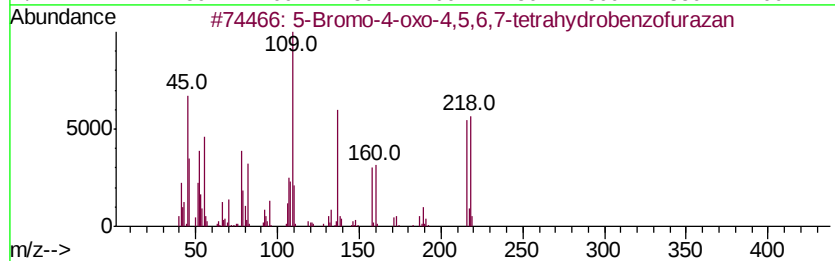
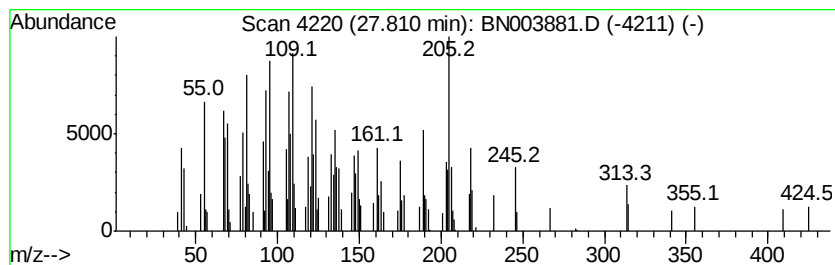
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.81	7.48 ng/ul	307155	Perylene-d12	23.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	42
2		6,10-Dimethyl-3-(1-methylethylidene...	206	C15H26	069239-71-0	38
3		(-)-Isolongifolol, methyl ether	236	C16H28O	1000333-80-8	30
4		(3H,6H)Thieno[3,4-c]isoxazole, 3...	205	C11H11NOS	1000115-54-7	30
5		1-(2-Pyridyl)piperazine	163	C9H13N3	034803-66-2	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121218\
Data File : BN003881.D
Acq On : 13 Dec 2018 08:49
Operator : JU/SJ
Sample : J6228-07
Misc :
ALS Vial : 30 Sample Multiplier: 1

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ClientSampleId :
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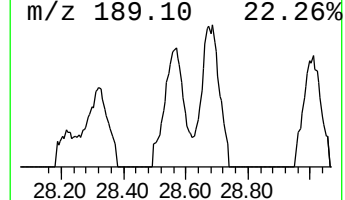
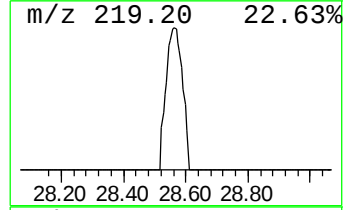
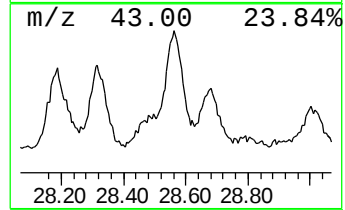
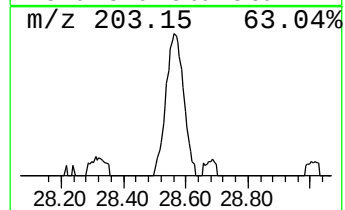
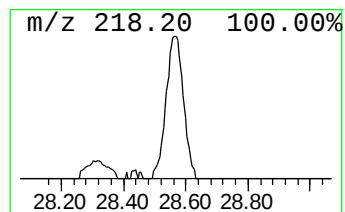
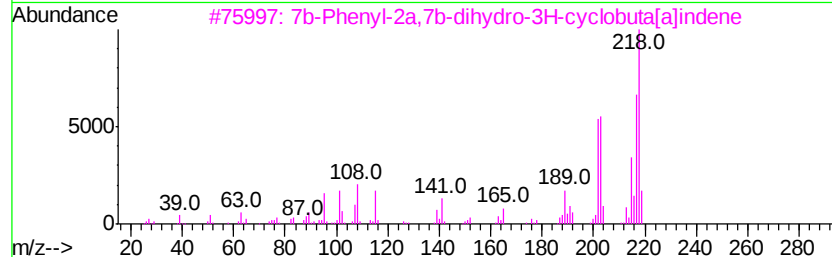
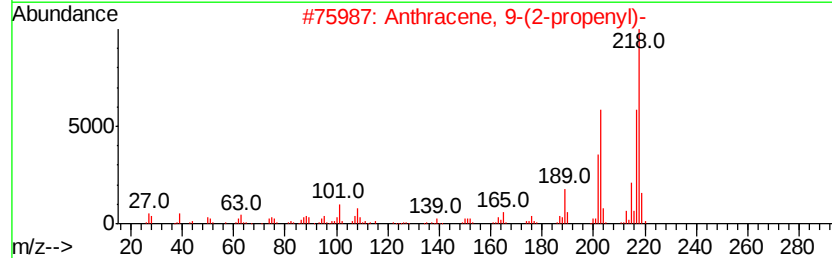
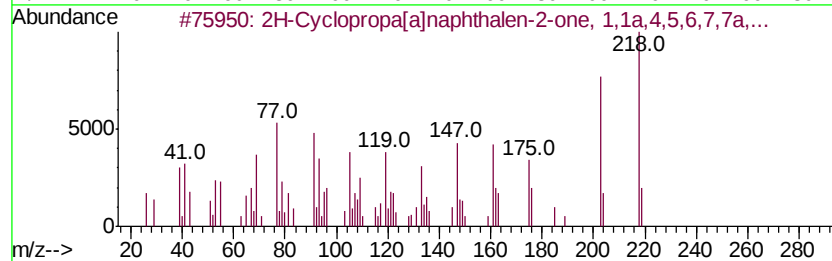
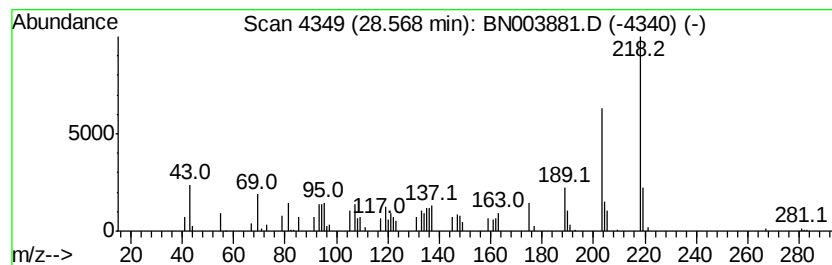
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2H-Cyclopropa[a]naphthalen-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.57	3.61 ng/ul	148110	Perylene-d12	23.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-Cyclopropa[a]naphthalen-2-one...	218	C15H22O	006831-17-0	76
2		Anthracene, 9-(2-propenyl)-	218	C17H14	023707-65-5	64
3		7b-Phenyl-2a,7b-dihydro-3H-cyclo...	218	C17H14	138089-70-0	64
4		(1S,6R,9S)-5,5,9,10-Tetramethylt...	218	C16H26	1000298-97-8	59
5		1,2,3,4-Tetrahydro-1-phenyl-1,2,...	218	C17H14	138089-69-7	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121218\
Data File : BN003881.D
Acq On : 13 Dec 2018 08:49
Operator : JU/SJ
Sample : J6228-07
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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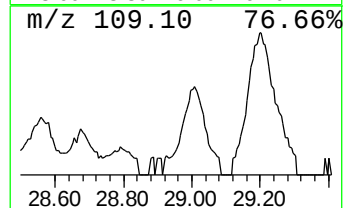
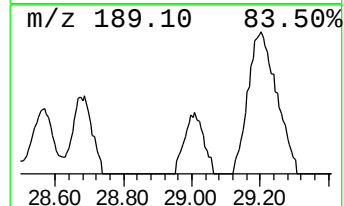
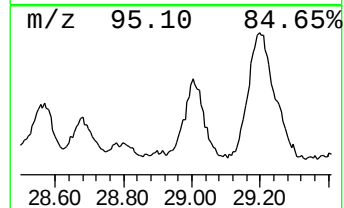
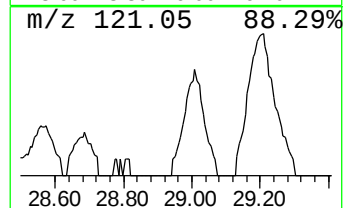
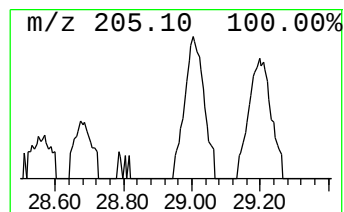
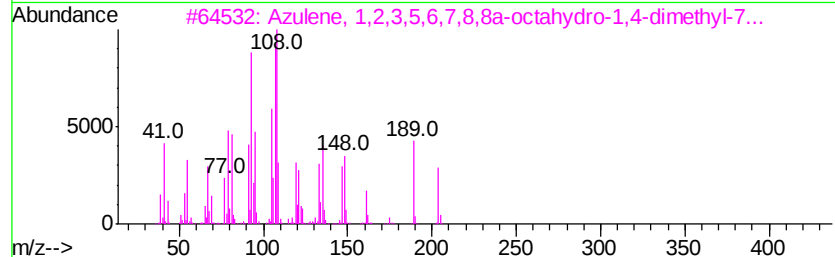
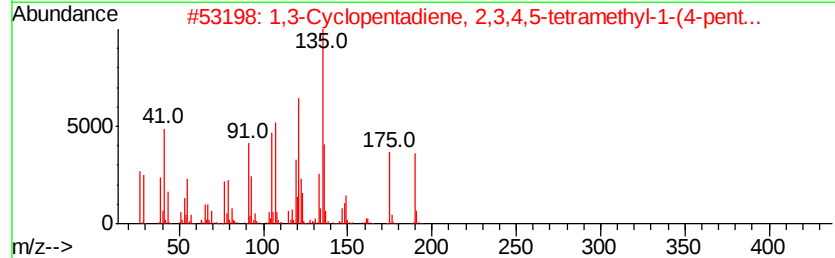
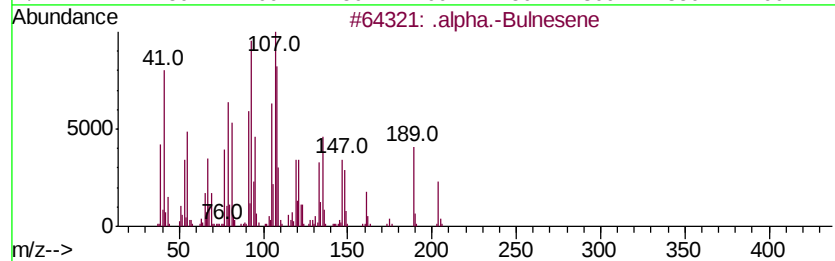
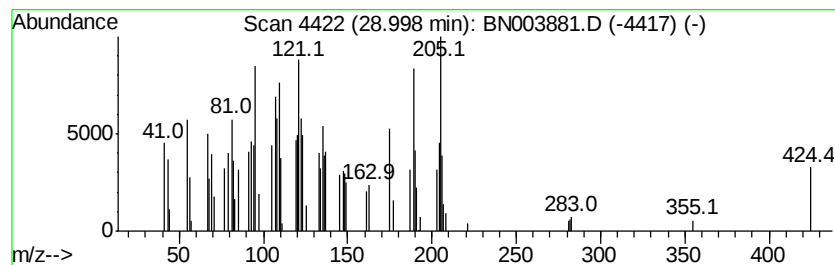
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-03 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.00	3.37 ng/ul	138180	Perylene-d12	23.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.alpha.-Bulnesene	204	C15H24	1000374-19-9	42
2		1,3-Cyclopentadiene, 2,3,4,5-tet...	190	C14H22	1000161-28-7	42
3		Azulene, 1,2,3,5,6,7,8,8a-octahv...	204	C15H24	003691-11-0	38
4		Thuiopsene-(I2)	204	C15H24	1000152-20-4	38
5		1-(2-Pyridyl)piperazine	163	C9H13N3	034803-66-2	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121218\
Data File : BN003881.D
Acq On : 13 Dec 2018 08:49
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Sample : J6228-07
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampled :
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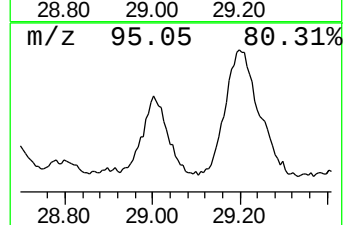
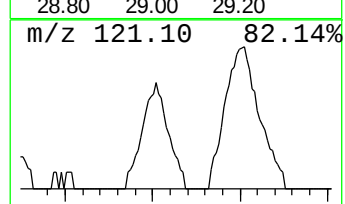
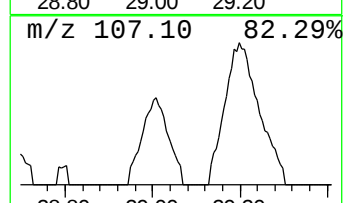
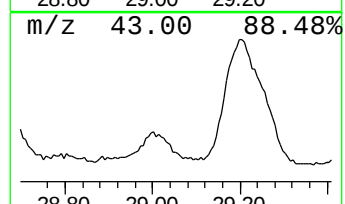
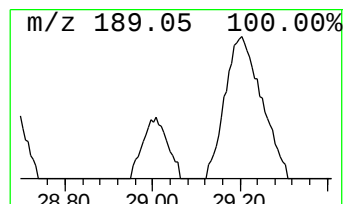
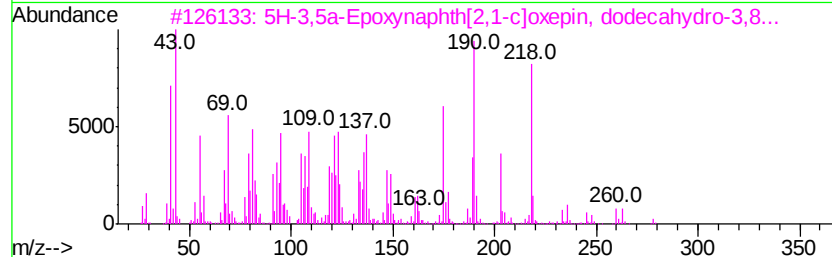
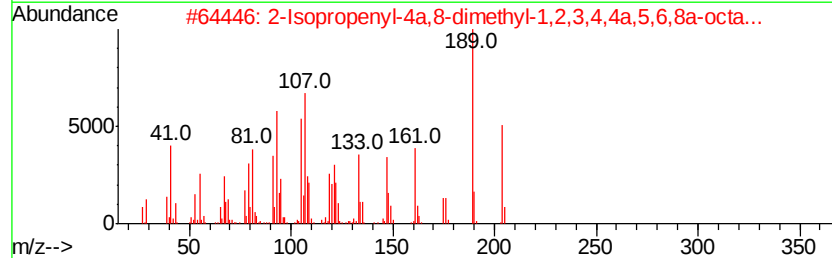
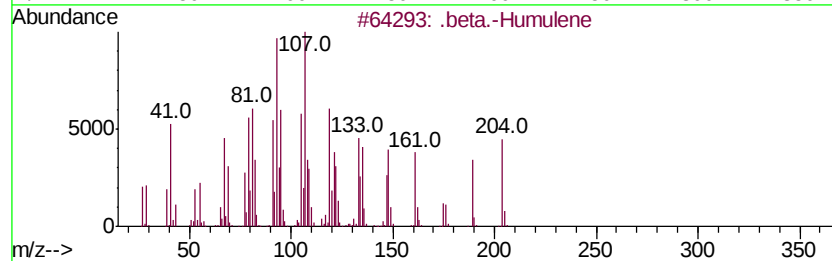
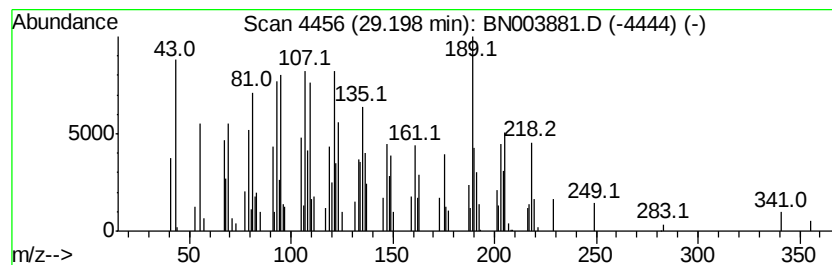
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 .beta.-Humulene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.20	11.29 ng/ul	463375	Perylene-d12	23.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-Humulene	204	C15H24	000116-04-1	64
2		2-Isopropenyl-4a,8-dimethyl-1,2,...	204	C15H24	1000193-57-0	49
3		5H-3,5a-Epoxy-naphth[2,1-c]oxepin...	278	C18H30O2	001153-34-0	38
4		(-)-Isolongifolol, methyl ether	236	C16H28O	1000333-80-8	38
5		Acetic acid, 3-hydroxy-6-isoprop...	278	C17H26O3	1000185-44-8	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121218\
Data File : BN003881.D
Acq On : 13 Dec 2018 08:49
Operator : JU/SJ
Sample : J6228-07
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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
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(DEL) Alkane: Str...	24.20	2.6	ng/ul	106263	6	23.71	820850	20.0
(DEL) Alkane: Str...	25.87	2.1	ng/ul	84391	6	23.71	820850	20.0
Guaia-3,9-diene	26.23	13.1	ng/ul	535507	6	23.71	820850	20.0
5-(1-Naphthyl)tri...	27.26	5.7	ng/ul	234677	6	23.71	820850	20.0
unknown-01	27.35	2.1	ng/ul	85963	6	23.71	820850	20.0
unknown-02	27.81	7.5	ng/ul	307155	6	23.71	820850	20.0
2H-Cyclopropa[a]n...	28.57	3.6	ng/ul	148110	6	23.71	820850	20.0
unknown-03	29.00	3.4	ng/ul	138180	6	23.71	820850	20.0
.beta.-Humulene	29.20	11.3	ng/ul	463375	6	23.71	820850	20.0