

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081219\
 Data File : BF116186.D
 Acq On : 12 Aug 2019 12:38
 Operator : HP/JU
 Sample : PB122104BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB122104BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % 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_F\METHODS\8270-BF073019.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.181	8	16	34	rVB	173773	291339	6.13%	0.824%
2	5.469	571	575	610	rBV	3154900	3385464	71.24%	9.570%
3	6.475	741	746	764	rBV	3147989	3542150	74.53%	10.013%
4	6.610	764	769	772	rBV	3496461	3493737	73.51%	9.876%
5	6.834	803	807	819	rVB	916943	814619	17.14%	2.303%
6	6.987	829	833	853	rBV	2825270	2942220	61.91%	8.317%
7	7.398	898	903	923	rBV	2031402	2405744	50.62%	6.800%
8	8.110	1020	1024	1045	rBV	934932	1052538	22.15%	2.975%
9	9.186	1202	1207	1223	rBV	3914123	4492113	94.52%	12.698%
10	9.869	1318	1323	1341	rBV	1278304	1279514	26.92%	3.617%
11	10.663	1453	1458	1485	rBV	2336897	2830184	59.55%	8.000%
12	11.357	1572	1576	1601	rBV	1190070	1347022	28.34%	3.808%
13	12.939	1840	1845	1848	rBV	4458962	4752485	100.00%	13.434%
14	13.998	2021	2025	2042	rBV	953639	1137255	23.93%	3.215%
15	14.751	2149	2153	2157	rVB	64578	67983	1.43%	0.192%
16	15.080	2204	2209	2219	rVB	108363	134429	2.83%	0.380%
17	15.463	2266	2274	2296	rBV2	610723	1107509	23.30%	3.131%
18	15.880	2338	2345	2355	rBV	106324	178842	3.76%	0.506%
19	16.374	2423	2429	2435	rBV2	63216	121914	2.57%	0.345%

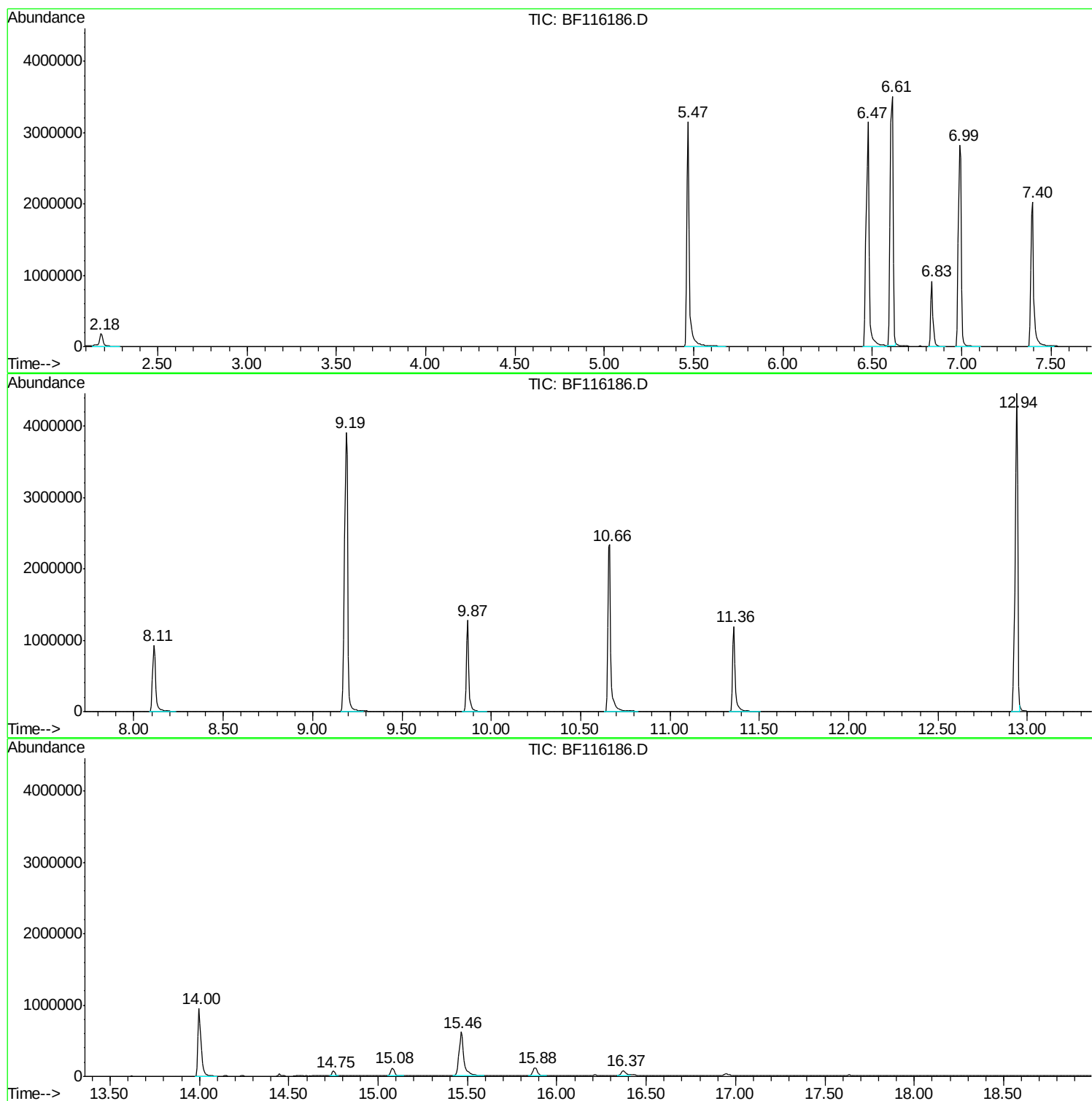
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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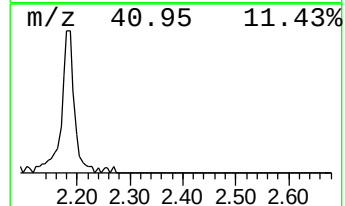
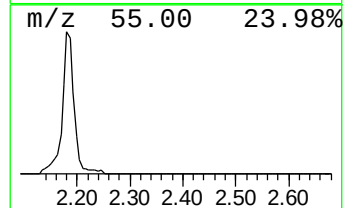
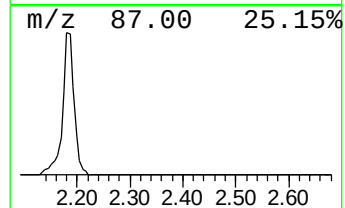
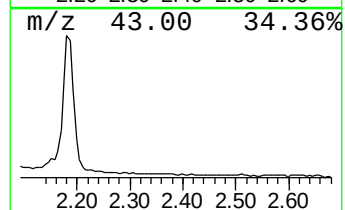
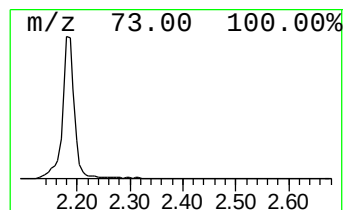
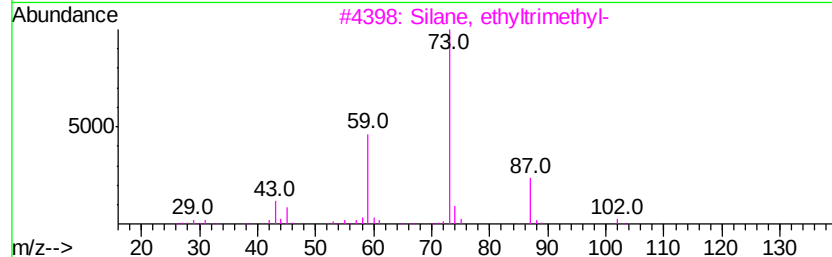
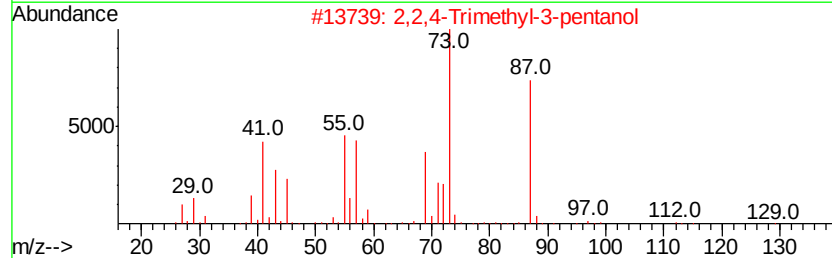
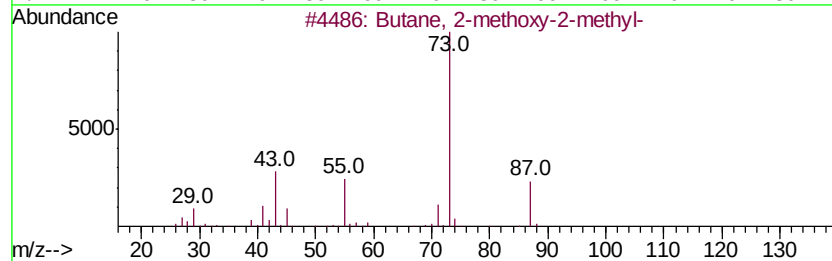
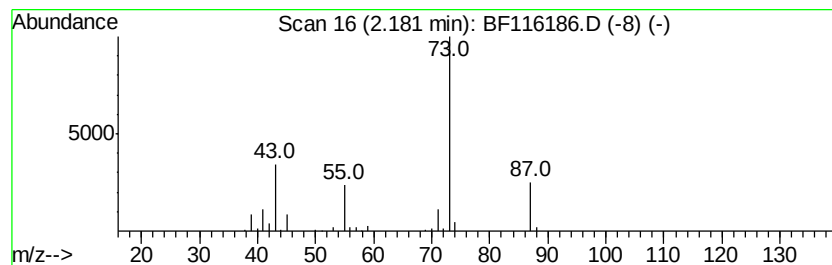
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.18	7.15 ng	291339	1,4-Dichlorobenzene-d4	6.83

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	10



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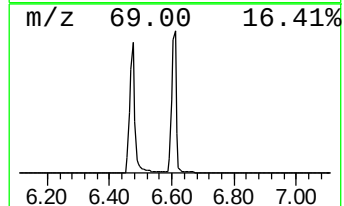
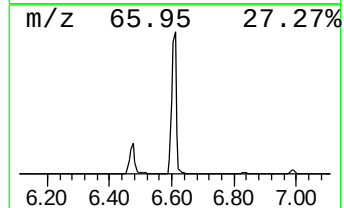
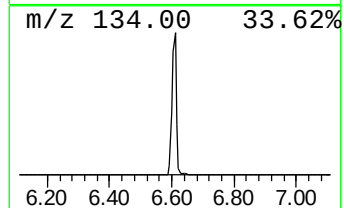
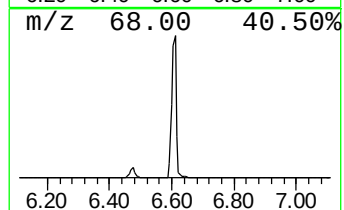
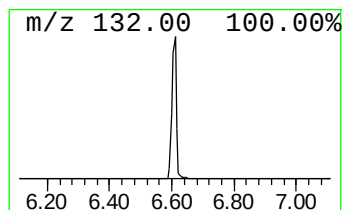
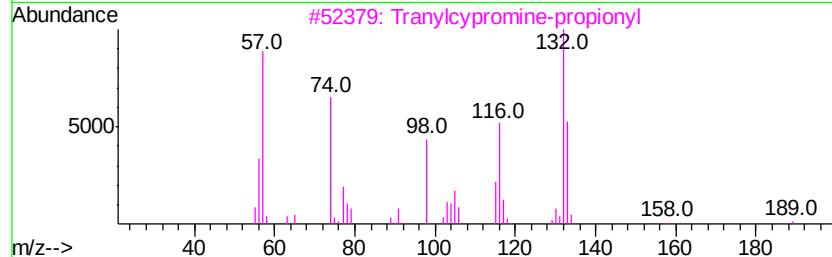
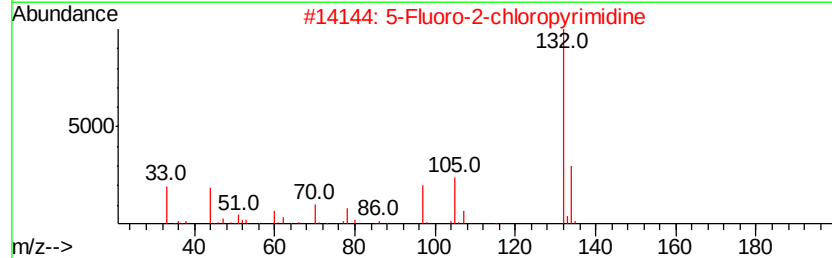
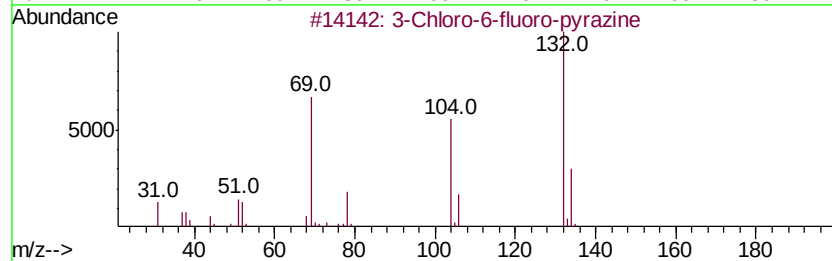
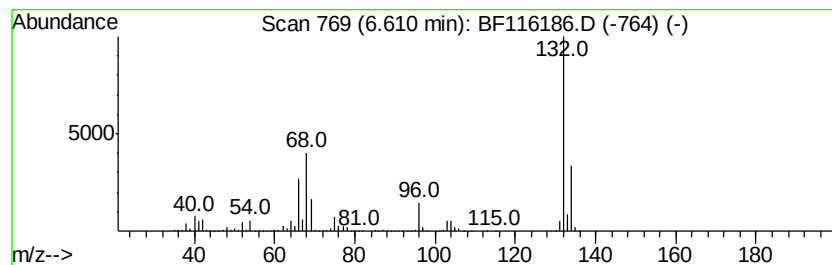
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	85.78 ng	3493740	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9



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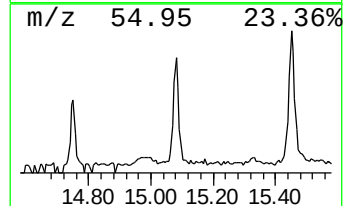
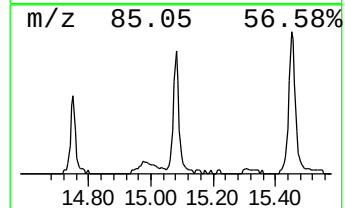
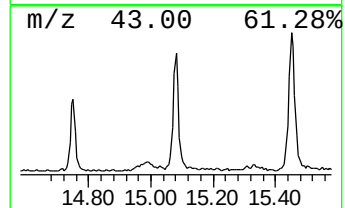
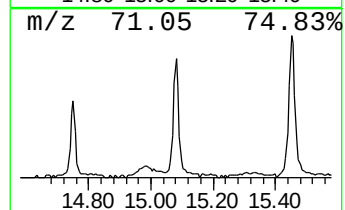
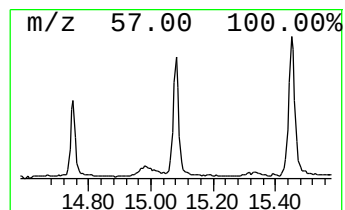
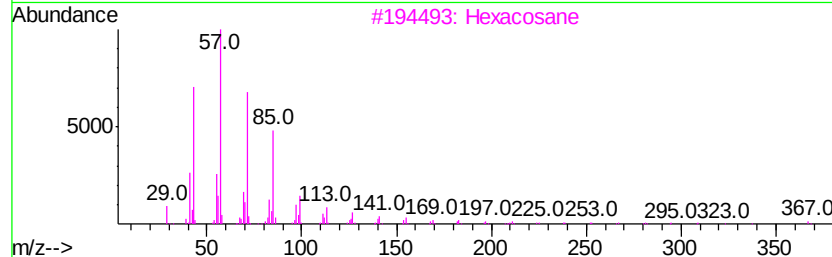
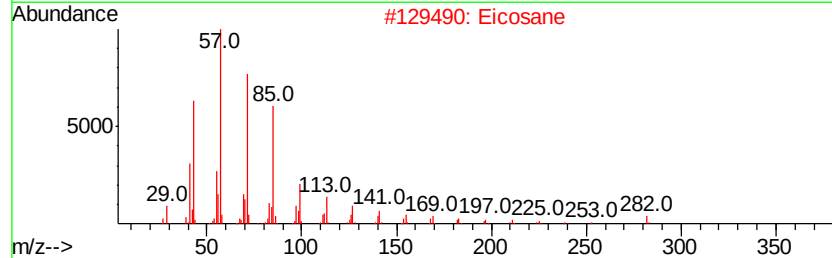
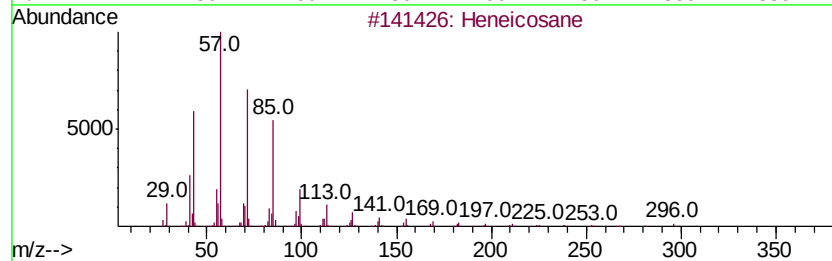
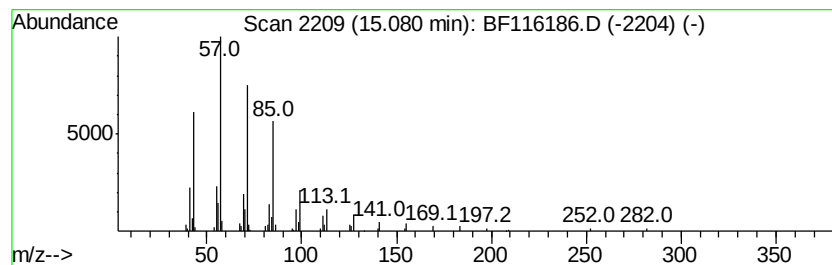
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Peak Number 4 Heneicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.08	2.43 ng	134429	Perylene-d12	15.46

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Eicosane		282	C20H42	000112-95-8	91
3	Hexacosane		366	C26H54	000630-01-3	90
4	2-methyloctacosane		408	C29H60	1000376-72-8	90
5	Octacosane		394	C28H58	000630-02-4	90



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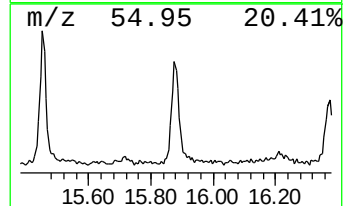
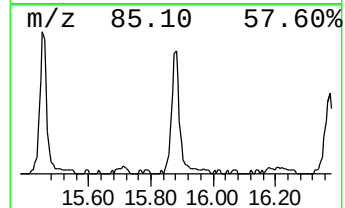
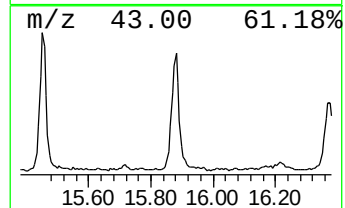
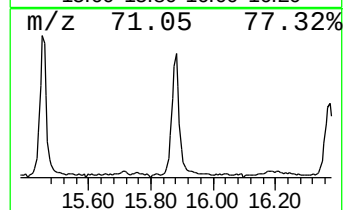
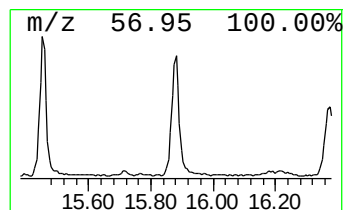
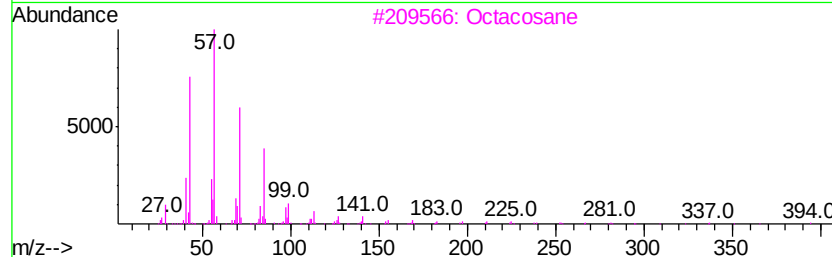
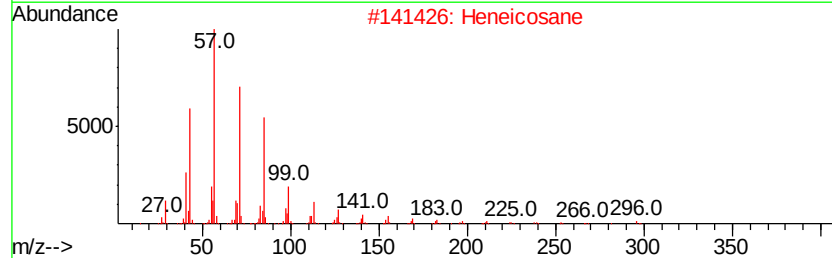
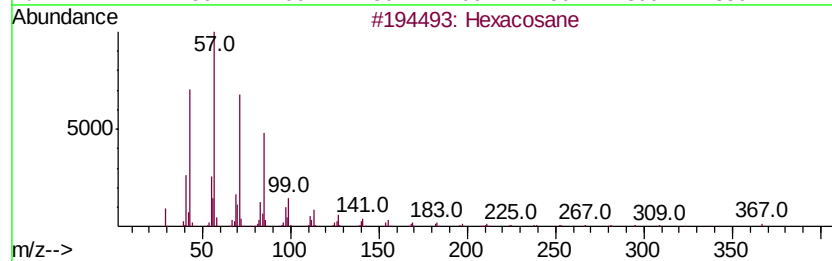
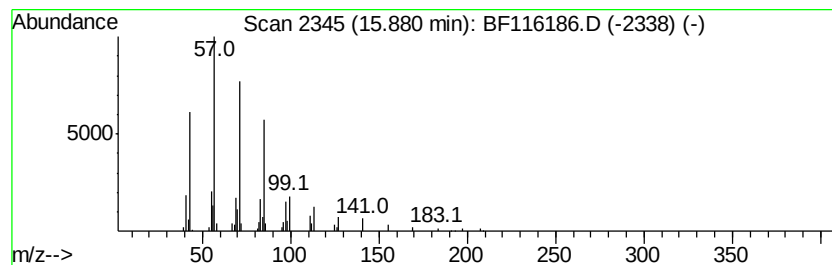
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Hexacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.88	3.23 ng	178842	Perylene-d12	15.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



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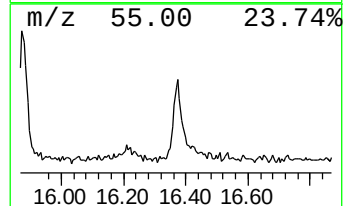
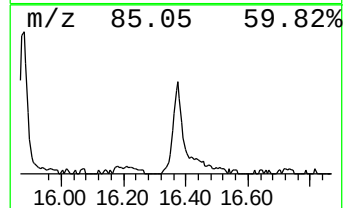
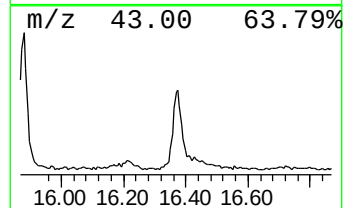
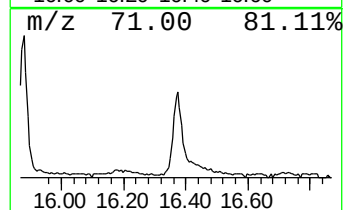
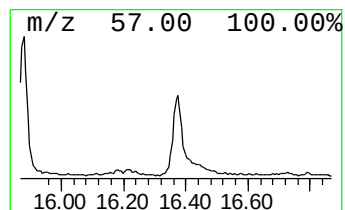
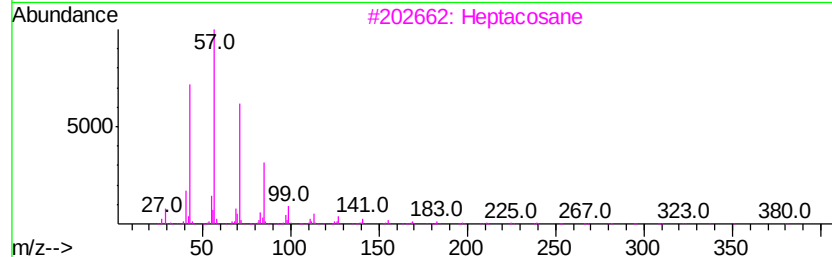
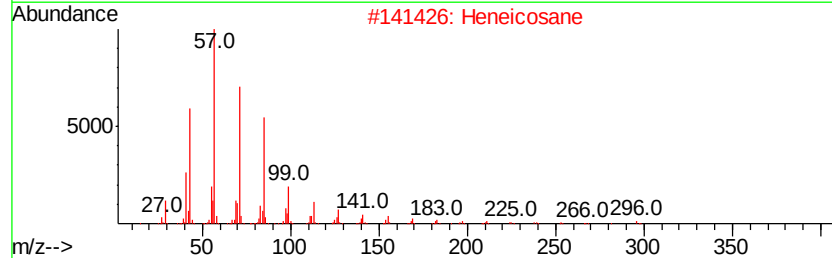
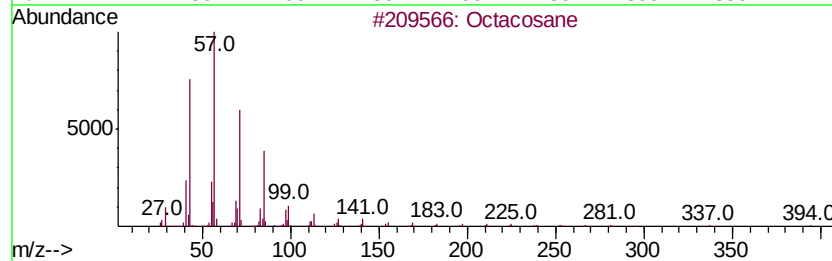
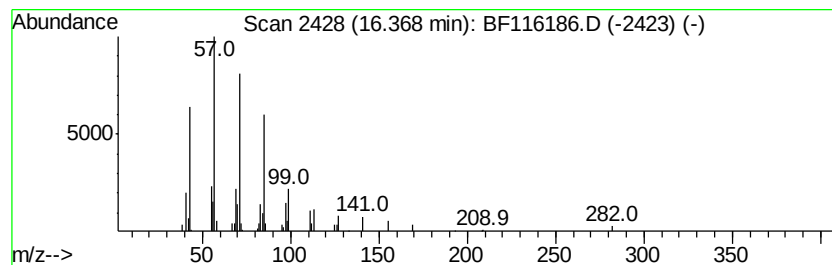
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.37	2.20 ng	121914	Perylene-d12	15.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		triacontane	422	C30H62	000638-68-6	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	90



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.18	7.2	ng	291339	1	6.83	814619	20.0
unknown6.61	6.61	85.8	ng	3493740	1	6.83	814619	20.0
Heneicosane	15.08	2.4	ng	134429	6	15.46	1107510	20.0
Hexacosane	15.88	3.2	ng	178842	6	15.46	1107510	20.0
Octacosane	16.37	2.2	ng	121914	6	15.46	1107510	20.0