

Data Path : Z:\HPCHEM1\BNA G\DATA\BG041118\
 Data File : BG033754.D
 Acq On : 11 Apr 2018 20:40
 Operator : SJ/JU
 Sample : J2320-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 040918-DBRI-E-U-S

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG031918.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	3.477	27	32	42	rBV	124992	255036	7.20%	1.789%
2	3.565	42	47	57	rVB	116070	178934	5.05%	1.255%
3	4.746	240	248	272	rBV	931925	1731332	48.85%	12.144%
4	6.285	504	510	522	rBV	101695	230383	6.50%	1.616%
5	7.971	791	797	810	rBV	55202	147093	4.15%	1.032%
6	8.365	856	864	881	rBV	223029	526516	14.86%	3.693%
7	8.859	941	948	957	rBV2	65940	159303	4.49%	1.117%
8	9.188	996	1004	1016	rBV2	381973	785217	22.16%	5.508%
9	10.057	1145	1152	1172	rBV	274117	676664	19.09%	4.746%
10	11.738	1431	1438	1459	rBV	107055	257336	7.26%	1.805%
11	14.100	1833	1840	1857	rBV	1029812	1716919	48.44%	12.043%
12	15.474	2068	2074	2081	rBV2	223968	397037	11.20%	2.785%
13	16.238	2198	2204	2210	rVB	52049	69555	1.96%	0.488%
14	16.943	2317	2324	2341	rBV	625228	1076974	30.39%	7.554%
15	17.878	2477	2483	2489	rBV2	53324	66321	1.87%	0.465%
16	18.201	2532	2538	2545	rBV	364246	632543	17.85%	4.437%
17	19.282	2717	2722	2726	rBV	35534	41633	1.17%	0.292%
18	19.887	2821	2825	2829	rVB	31385	38848	1.10%	0.272%
19	20.204	2871	2879	2884	rBV2	35532	58300	1.64%	0.409%
20	20.445	2916	2920	2925	rBV2	30181	36843	1.04%	0.258%
21	20.557	2932	2939	2944	rBV	34397	58660	1.66%	0.411%
22	20.721	2961	2967	2980	rBV	2519098	3544129	100.00%	24.859%
23	22.548	3268	3278	3285	rBV	392731	791509	22.33%	5.552%
24	24.176	3550	3555	3561	rVB4	17391	36190	1.02%	0.254%
25	26.303	3905	3917	3932	rVB2	209311	702473	19.82%	4.927%
26	29.194	4401	4409	4419	rVB7	11683	41228	1.16%	0.289%

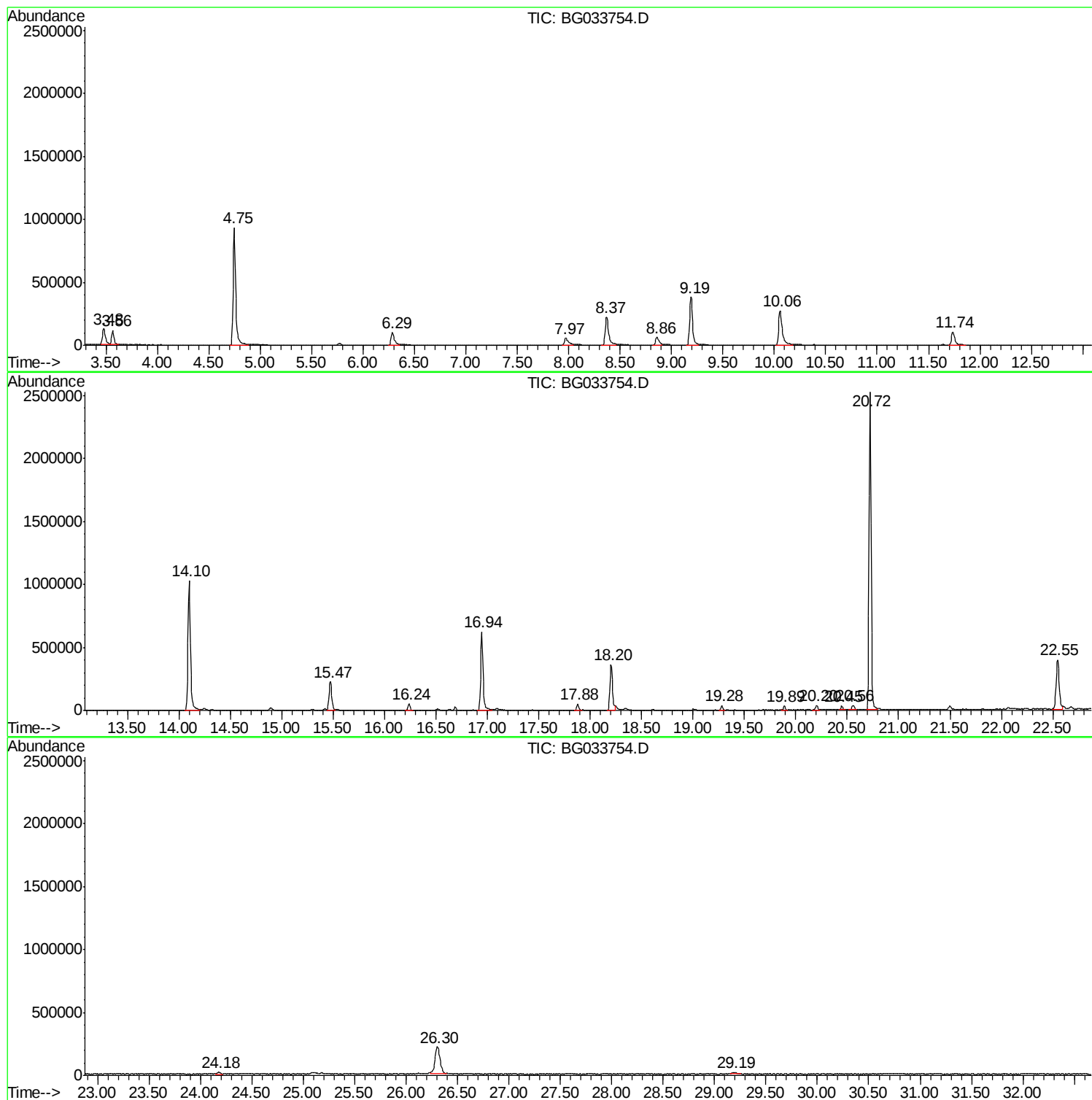
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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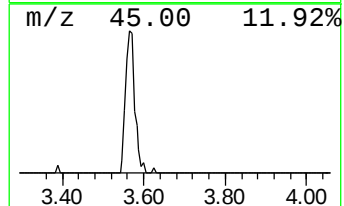
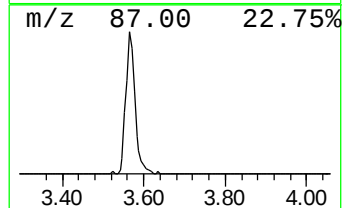
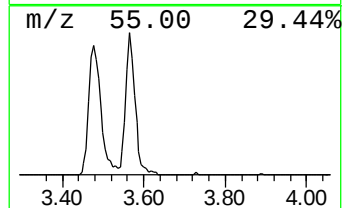
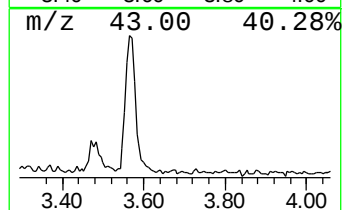
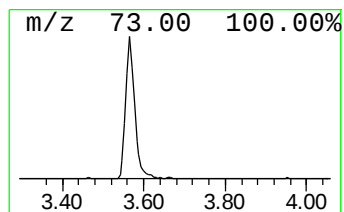
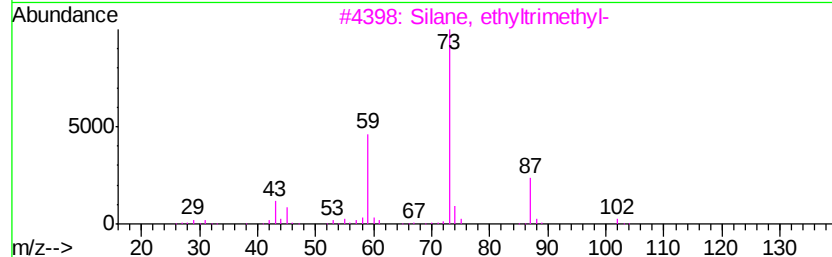
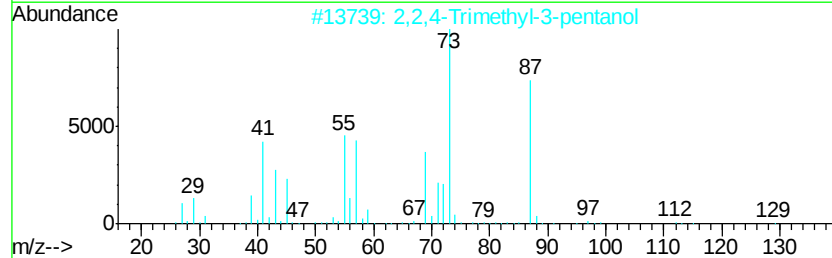
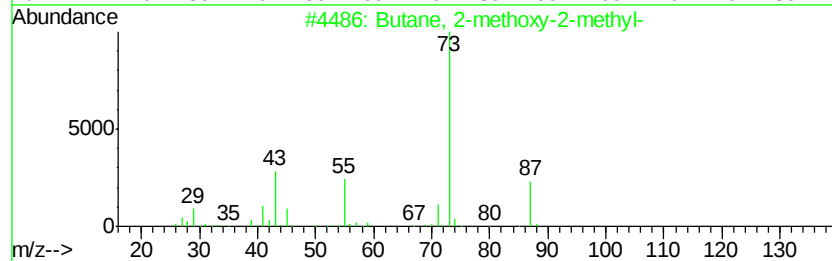
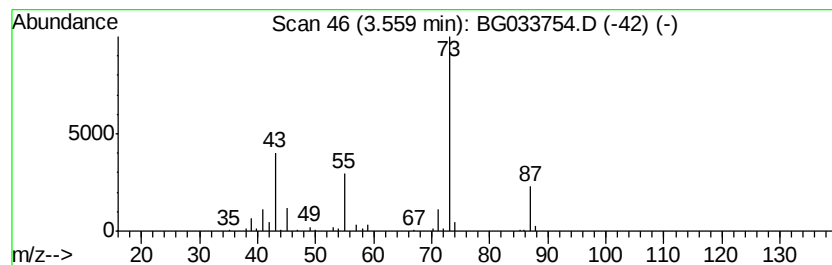
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.56	22.46 ng	178934	1,4-Dichlorobenzene-d4	8.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	17
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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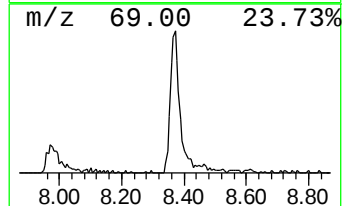
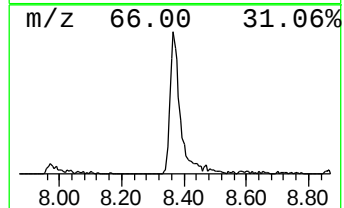
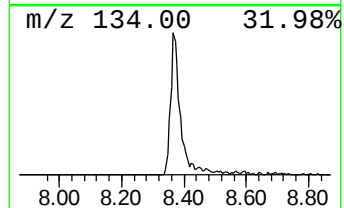
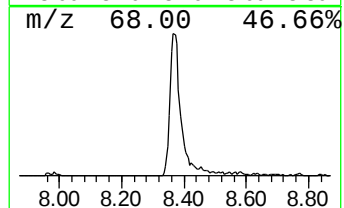
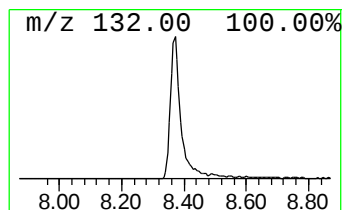
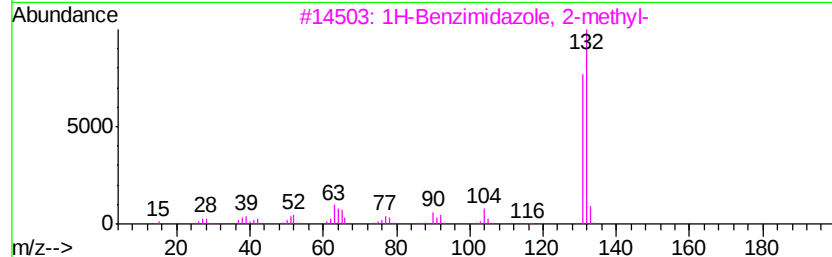
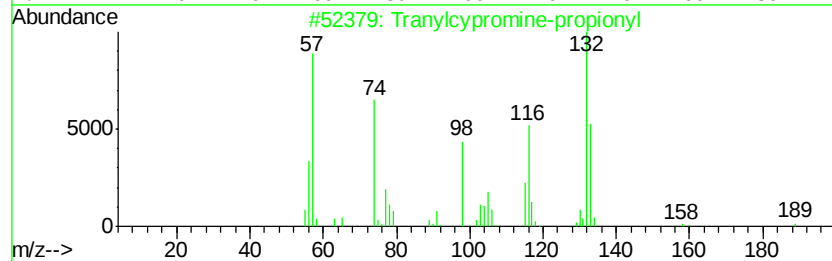
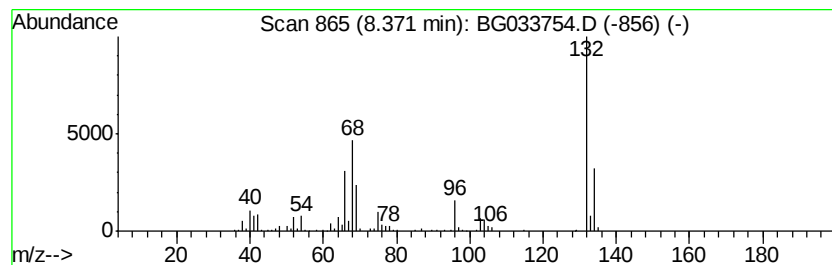
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Peak Number 4 unknown8.37 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.37	66.10 ng	526516	1,4-Dichlorobenzene-d4	8.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
2			5-Aminoindole	132	C8H8N2	005192-03-0	22
3			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
5			1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	12



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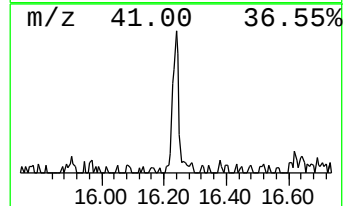
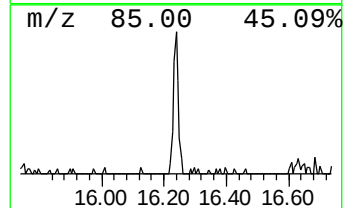
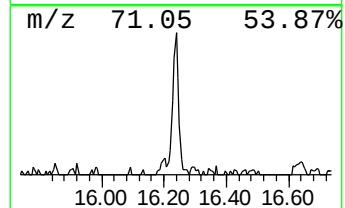
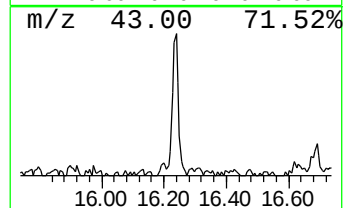
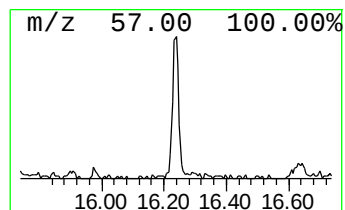
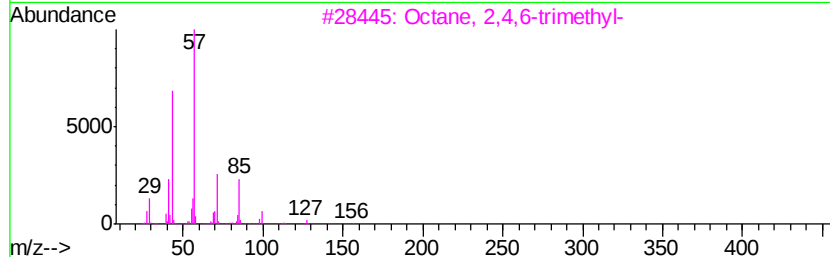
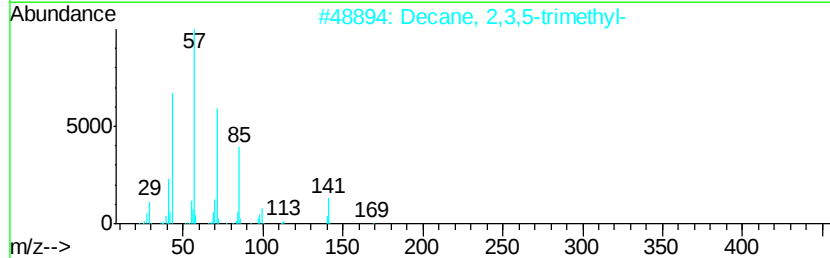
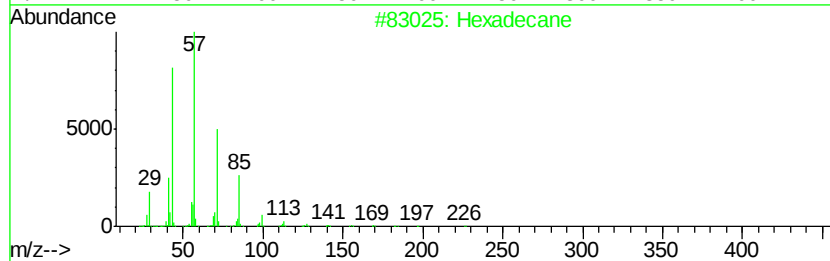
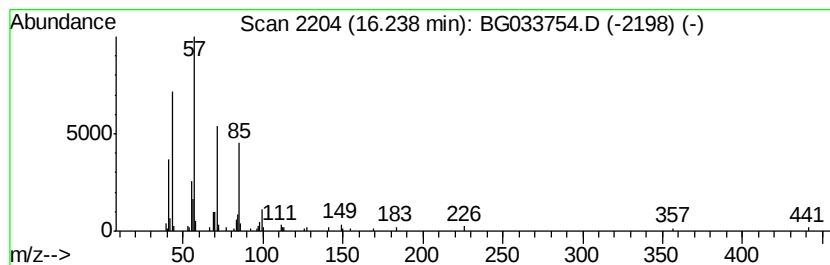
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Peak Number 6 Hexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.24	3.50 ng	69555	Acenaphthene-d10	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	93
2		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	86
3		Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	86
4		Pentadecane	212	C15H32	000629-62-9	72
5		Decane, 3-bromo-	220	C10H21Br	030571-71-2	53



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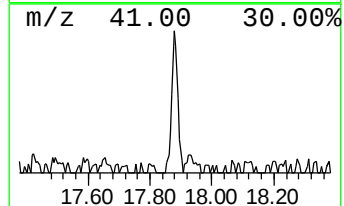
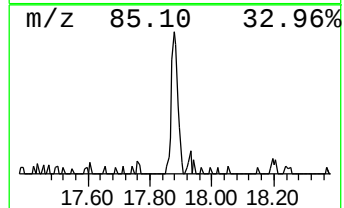
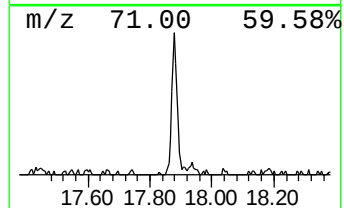
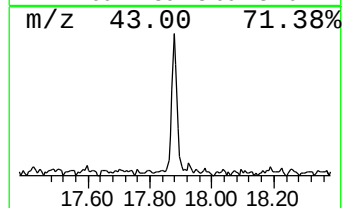
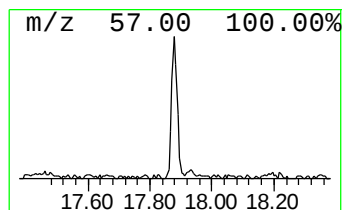
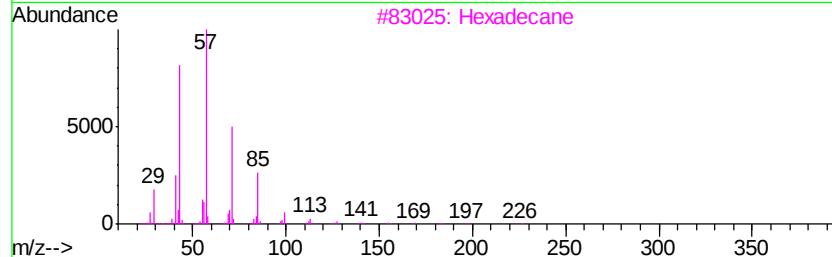
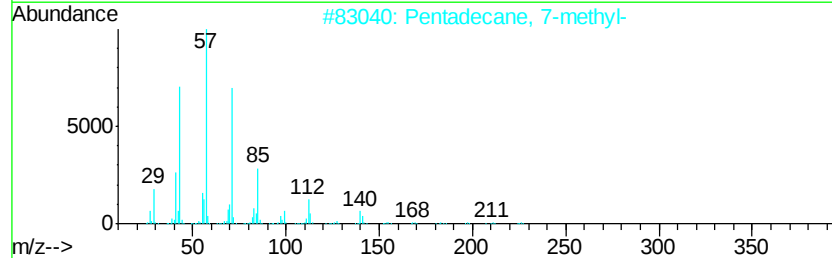
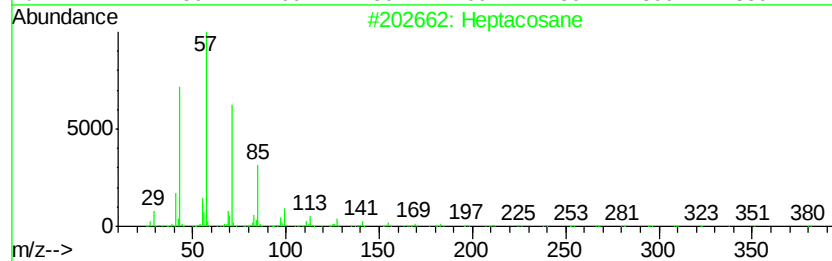
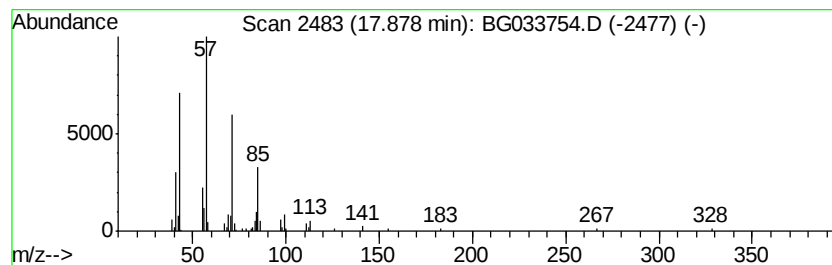
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Peak Number 7 Heptacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.88	2.10 ng	66321	Phenanthrene-d10	18.20

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	80
2	Pentadecane, 7-methyl-		226	C16H34	006165-40-8	78
3	Hexadecane		226	C16H34	000544-76-3	72
4	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	64
5	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	64



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.56	22.5	ng	178934	1	8.86	159303	20.0
unknown8.37	8.37	66.1	ng	526516	1	8.86	159303	20.0
Hexadecane	16.24	3.5	ng	69555	3	15.47	397037	20.0
Heptacosane	17.88	2.1	ng	66321	4	18.20	632543	20.0