

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121117\
 Data File : BF101287.D
 Acq On : 11 Dec 2017 14:15
 Operator : SJ/JU
 Sample : I6746-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ENV-109-4(40-50)

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_F\METHODS\8270-BF113017.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	2.122	3	6	16	rVB2	35264	54959	2.75%	0.365%
2	5.063	502	506	520	rBV	91062	153951	7.70%	1.021%
3	5.457	568	573	588	rBV	1395506	1569010	78.52%	10.408%
4	6.475	740	746	758	rBV	1468202	1677257	83.94%	11.126%
5	6.604	763	768	771	rBV	1693530	1680061	84.08%	11.144%
6	6.828	803	806	814	rBB	373427	302241	15.13%	2.005%
7	6.987	828	833	836	rBV	1415346	1245111	62.31%	8.259%
8	7.398	897	903	906	rBV	968853	1009366	50.51%	6.695%
9	8.010	1004	1007	1016	rBB	25073	22763	1.14%	0.151%
10	8.110	1020	1024	1031	rBV	470990	397769	19.91%	2.639%
11	9.192	1202	1208	1211	rBV	2072715	1998184	100.00%	13.255%
12	9.581	1270	1274	1277	rBV	305004	235666	11.79%	1.563%
13	9.869	1319	1323	1327	rBV2	553157	454659	22.75%	3.016%
14	10.669	1454	1459	1462	rBV	1227243	1177493	58.93%	7.811%
15	11.363	1573	1577	1580	rBV	542172	463359	23.19%	3.074%
16	12.951	1842	1847	1850	rBV	1937747	1890044	94.59%	12.537%
17	13.827	1991	1996	2001	rBV	42613	47234	2.36%	0.313%
18	14.010	2023	2027	2031	rBV	457704	399517	19.99%	2.650%
19	14.833	2163	2167	2171	rVB	20269	21259	1.06%	0.141%
20	15.486	2271	2278	2284	rVB	226436	275412	13.78%	1.827%

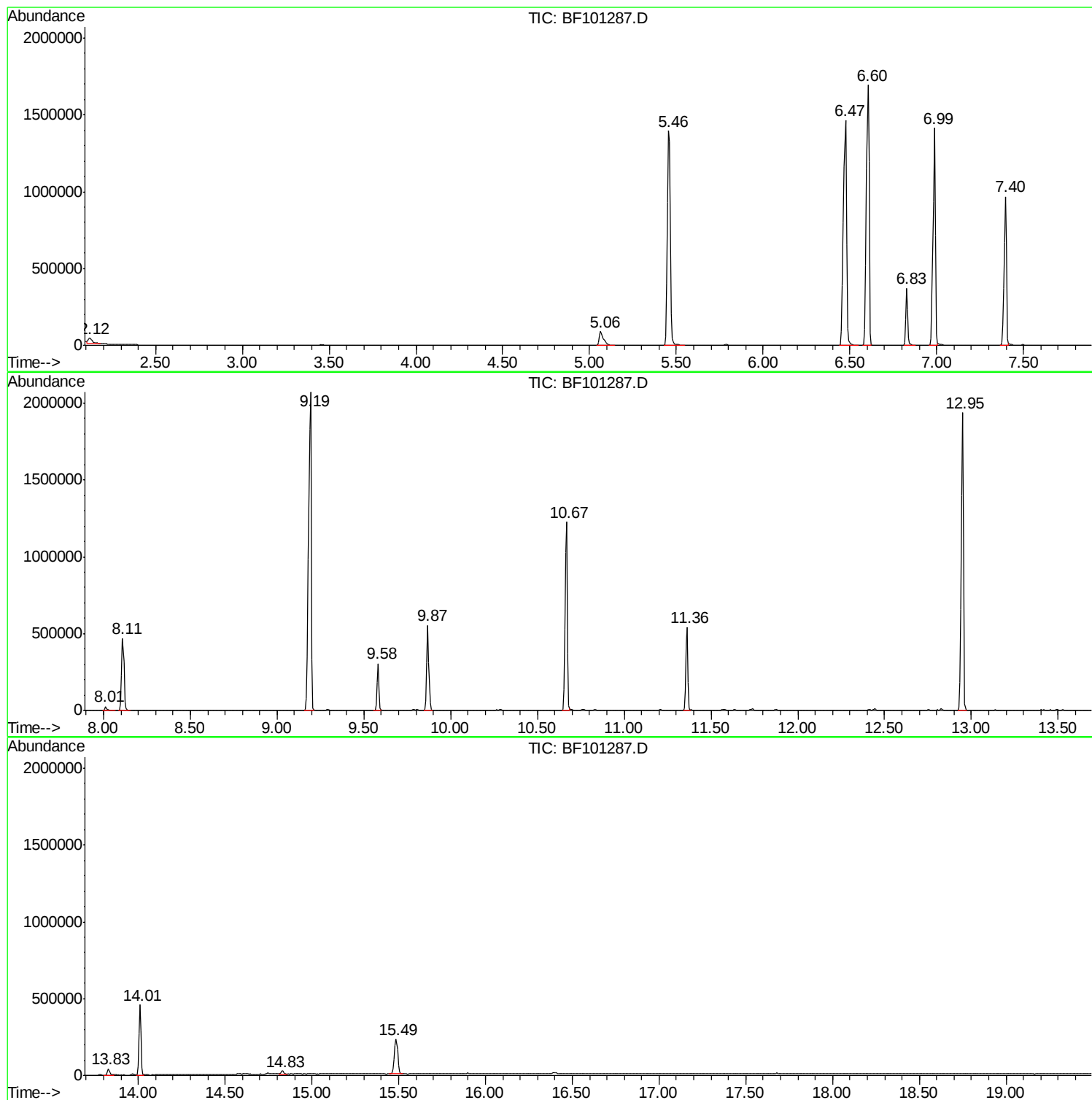
Sum of corrected areas: 15075315

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113017.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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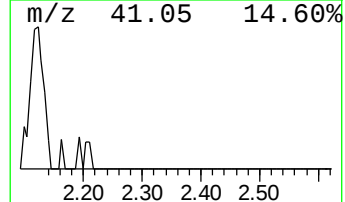
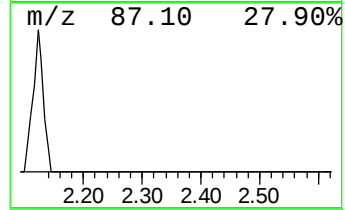
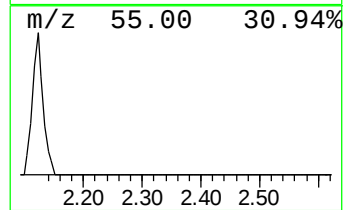
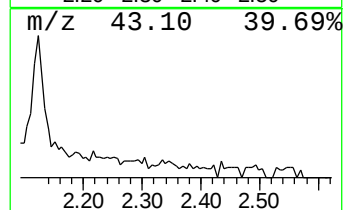
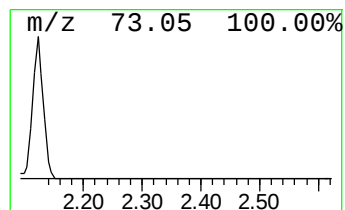
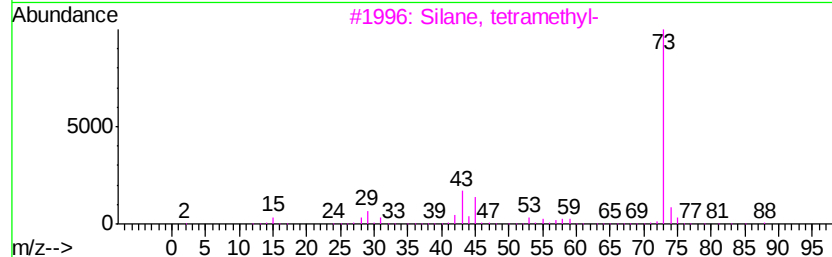
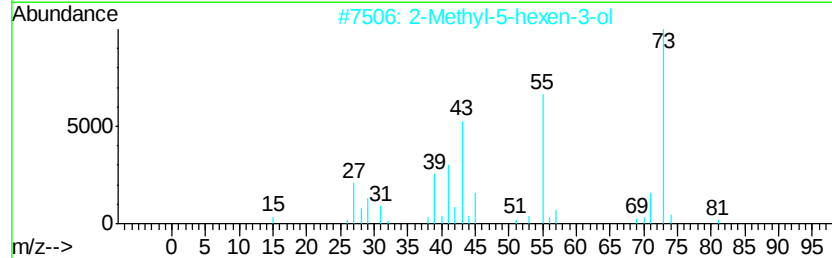
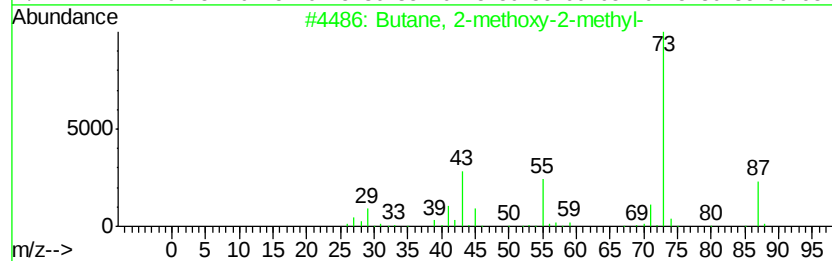
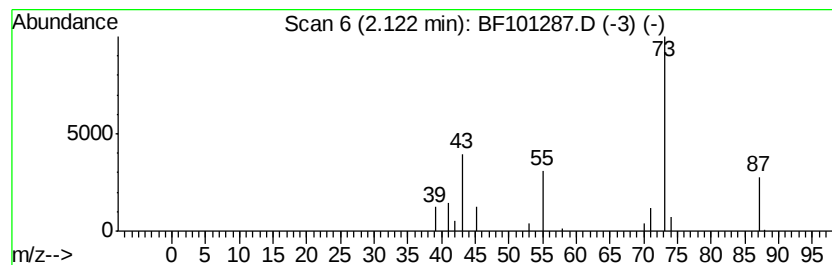
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.12	3.64 ng	54959	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-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	36
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
4			1,4-Butanediamine	88	C4H12N2	000110-60-1	9
5			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	9



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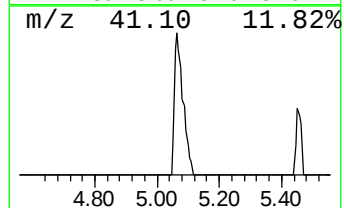
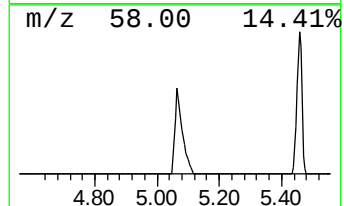
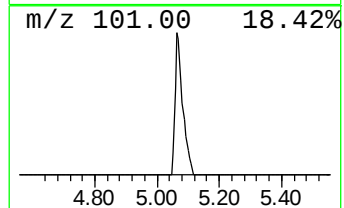
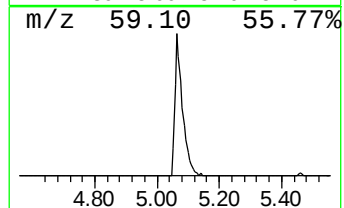
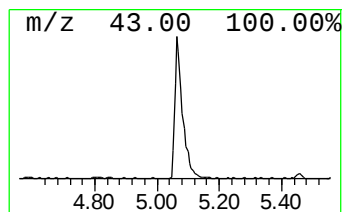
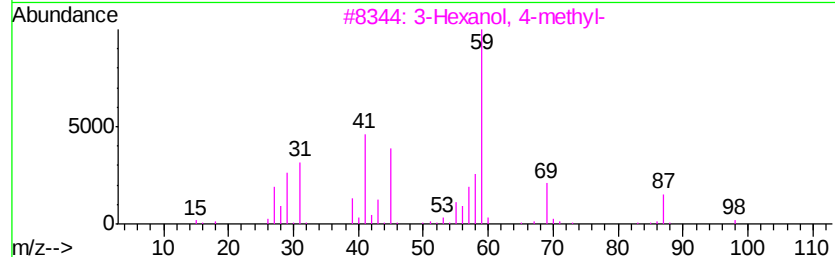
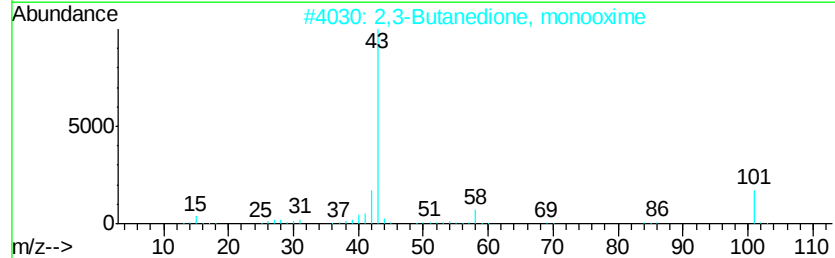
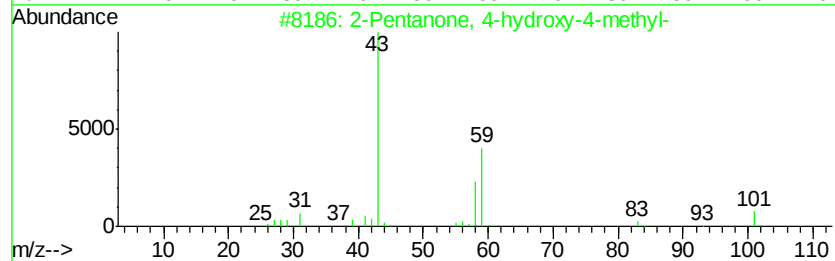
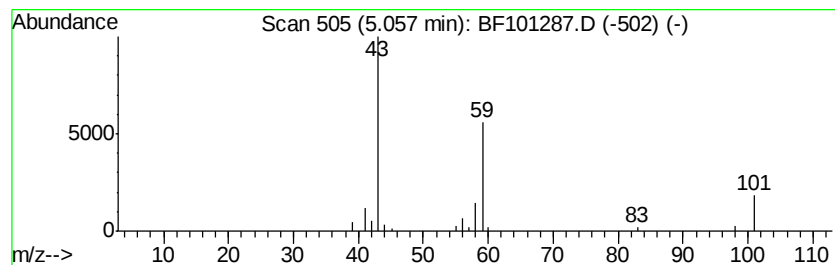
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.06	10.19 ng	153951	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	17
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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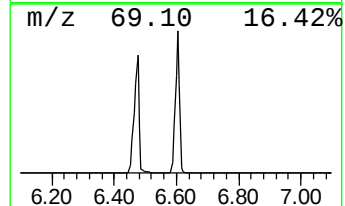
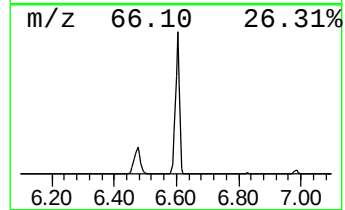
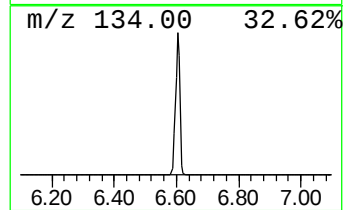
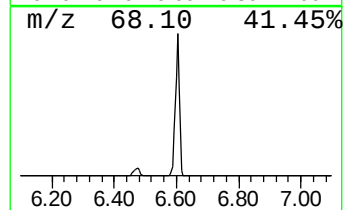
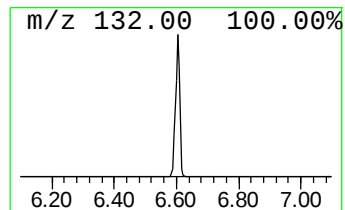
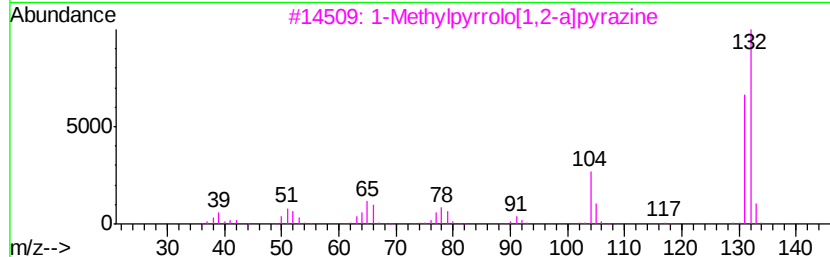
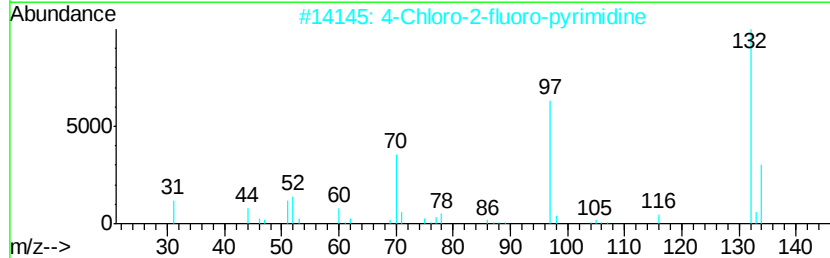
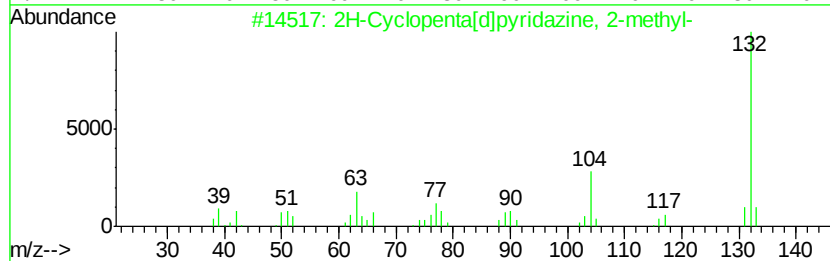
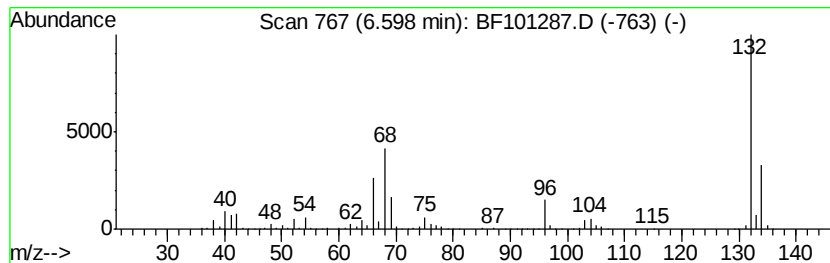
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Peak Number 3 unknown6.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	111.17 ng	1680060	1,4-Dichlorobenzene-d4	6.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
2		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
3		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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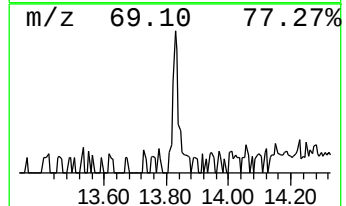
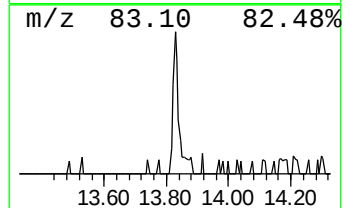
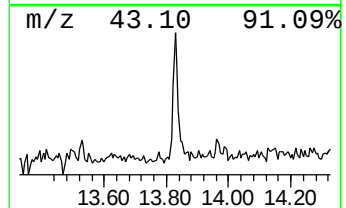
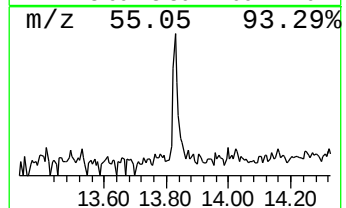
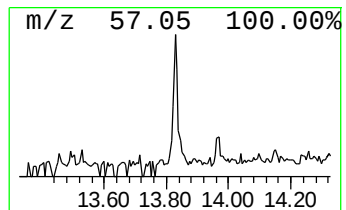
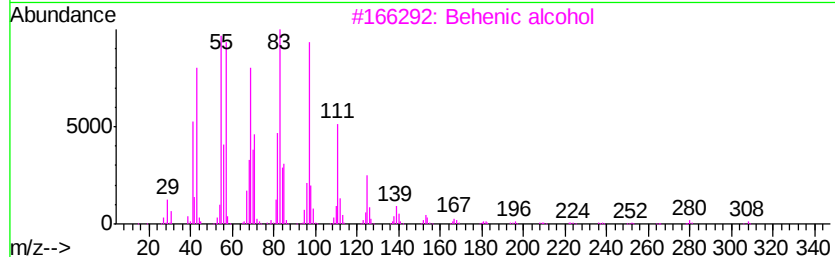
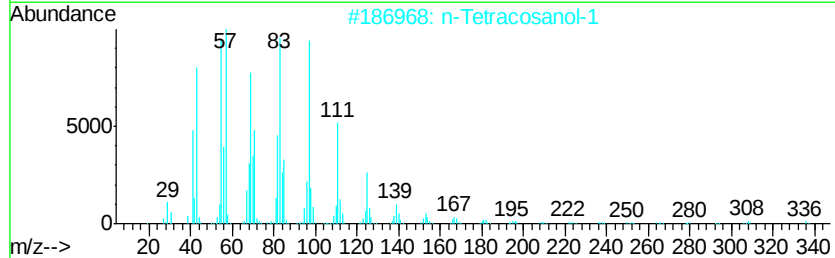
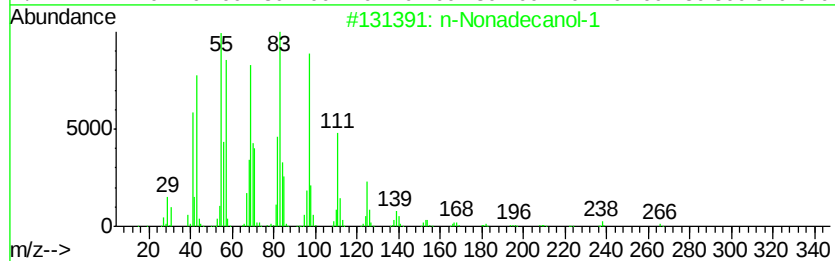
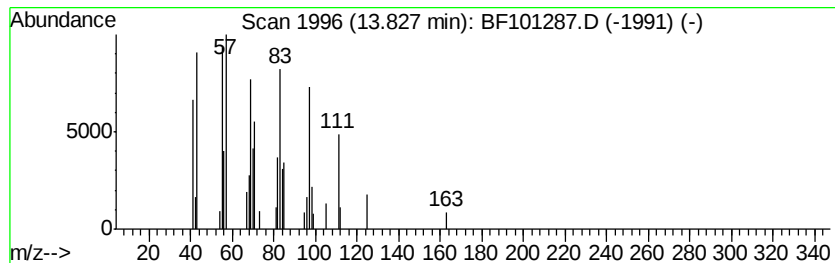
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Peak Number 5 n-Nonadecanol-1 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	2.36 ng	47234	Chrysene-d12	14.01

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Nonadecanol-1		284	C19H40O	001454-84-8	91
2	n-Tetracosanol-1		354	C24H50O	000506-51-4	91
3	Behenic alcohol		326	C22H46O	000661-19-8	91
4	Trifluoroacetoxy hexadecane		338	C18H33F3O2	006222-03-3	91
5	3-Eicosene, (E)-		280	C20H40	074685-33-9	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.12	3.6	ng	54959	1	6.83	302241	20.0
2-Pentanone, 4-hy...	5.06	10.2	ng	153951	1	6.83	302241	20.0
unknown6.60	6.60	111.2	ng	1680060	1	6.83	302241	20.0
n-Nonadecanol-1	13.83	2.4	ng	47234	5	14.01	399517	20.0