

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110921\
 Data File : BF126095.D
 Acq On : 9 Nov 2021 22:36
 Operator : JU/CG
 Sample : M4583-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VWE-B7-110621-10-20

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF126095.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.146	3	10	15	rBV	1529265	1968335	31.78%	5.120%
2	2.246	23	27	33	rVB	54324	75292	1.22%	0.196%
3	5.069	504	507	518	rVB	266076	355760	5.74%	0.925%
4	5.475	570	576	579	rBV	4562966	4756903	76.81%	12.373%
5	6.481	741	747	750	rBV	4656316	4996221	80.68%	12.996%
6	6.681	779	781	791	rVB2	68752	81161	1.31%	0.211%
7	6.775	791	797	805	rBV	79886	76304	1.23%	0.198%
8	6.851	806	810	813	rBV	1612703	1190795	19.23%	3.097%
9	7.416	900	906	909	rBV	3675493	3273011	52.85%	8.513%
10	7.451	909	912	916	rVB	119177	99640	1.61%	0.259%
11	8.033	1008	1011	1023	rBV	466844	482143	7.79%	1.254%
12	8.133	1023	1028	1031	rBV	1861626	1534817	24.78%	3.992%
13	9.210	1206	1211	1214	rBV	7622636	6192783	100.00%	16.108%
14	9.886	1322	1326	1329	rBV	2051982	1579119	25.50%	4.107%
15	10.674	1456	1460	1463	rBV	3551279	3243768	52.38%	8.437%
16	11.369	1574	1578	1581	rBV	1667507	1321759	21.34%	3.438%
17	11.439	1586	1590	1595	rVB3	68968	73377	1.18%	0.191%
18	12.580	1781	1784	1789	rVB	172774	137039	2.21%	0.356%
19	12.810	1820	1823	1826	rVB	121187	97990	1.58%	0.255%
20	12.957	1844	1848	1851	rBV	5163118	4308042	69.57%	11.206%
21	13.439	1927	1930	1936	rBV4	55839	107537	1.74%	0.280%
22	13.862	1998	2002	2011	rVB2	220336	291643	4.71%	0.759%
23	14.004	2021	2026	2030	rBV	1150482	1231304	19.88%	3.203%
24	14.486	2105	2108	2111	rBV3	50580	73975	1.19%	0.192%
25	15.474	2271	2276	2280	rBV	777511	896658	14.48%	2.332%

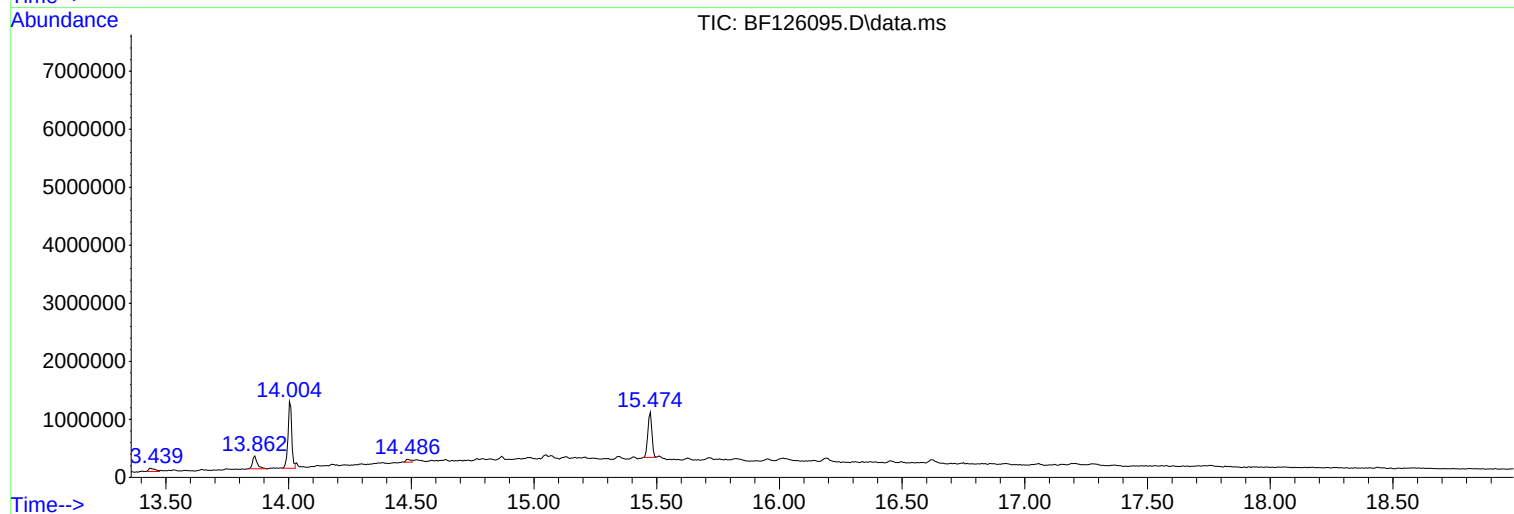
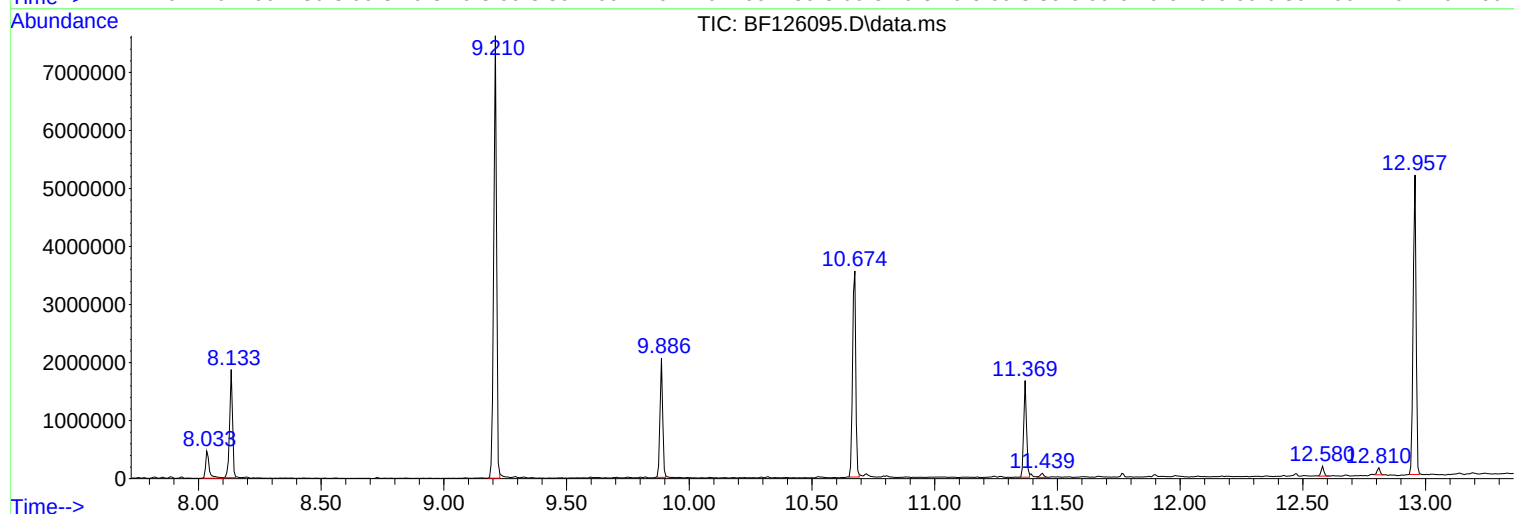
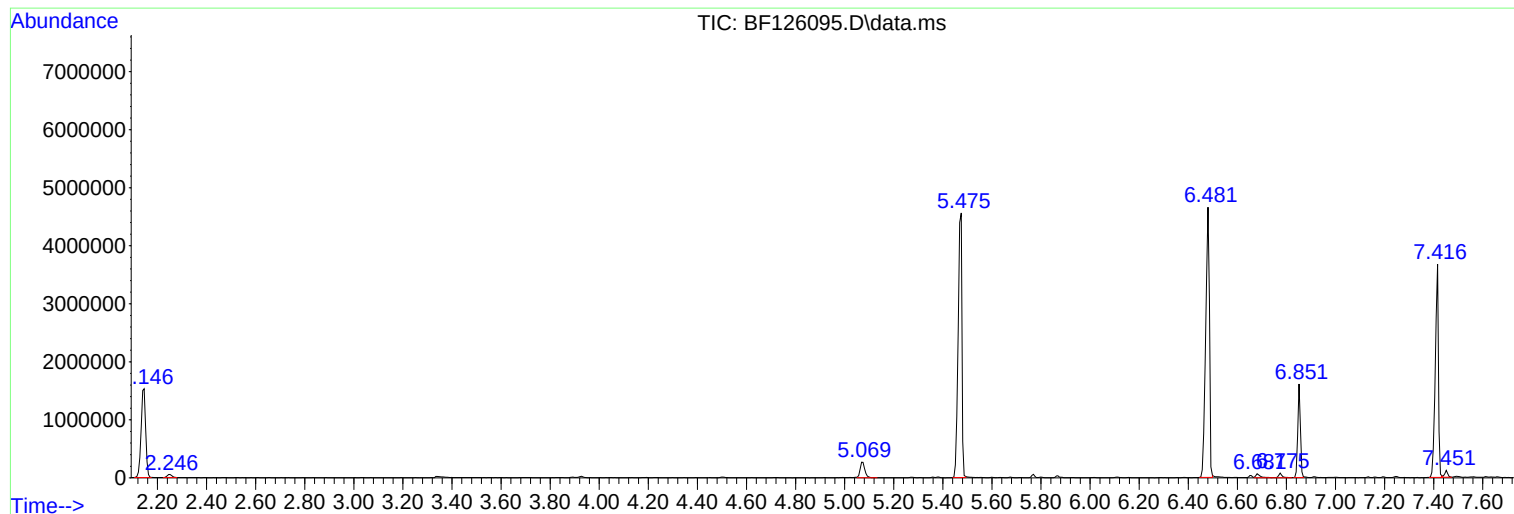
Sum of corrected areas: 38445376

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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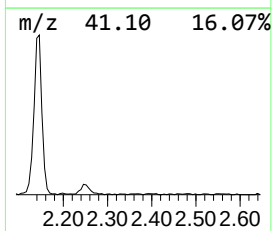
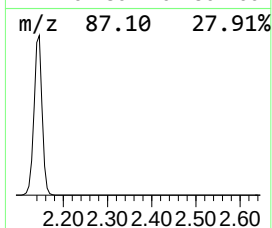
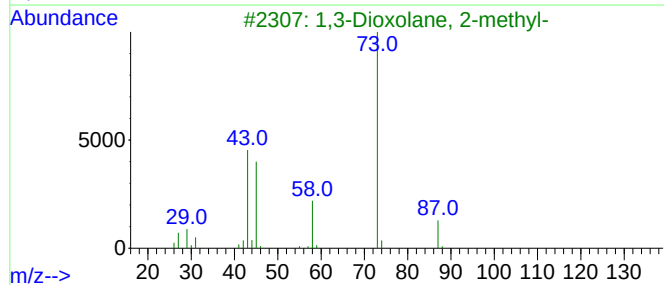
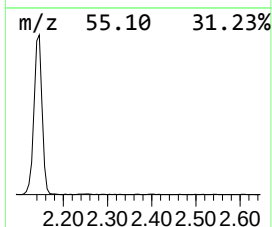
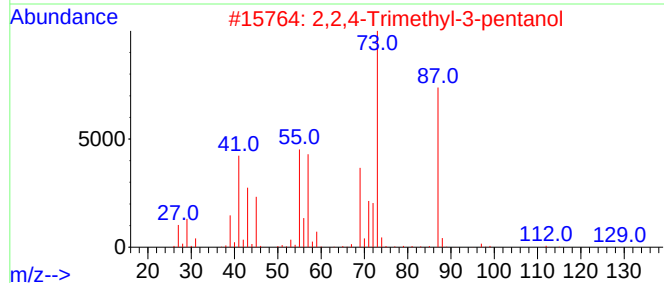
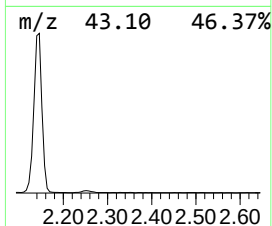
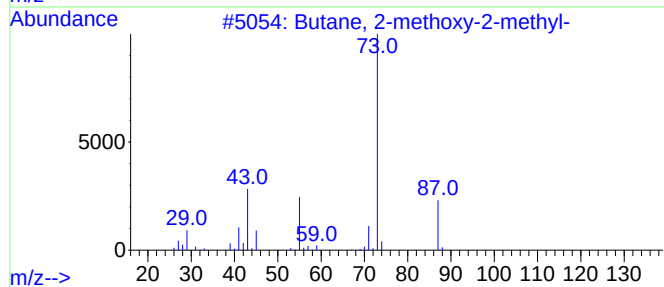
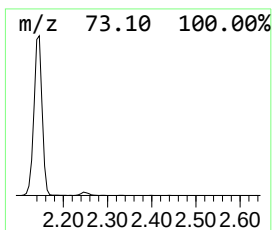
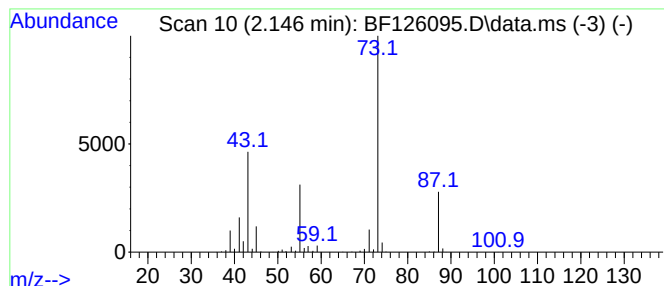
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.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.
2.146	33.06 ng	1968340	1,4-Dichlorobenzene-d4	6.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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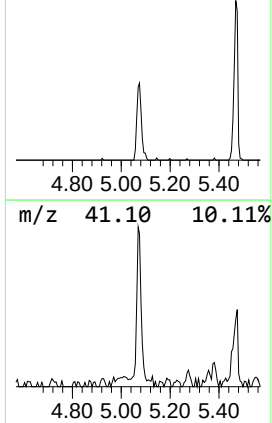
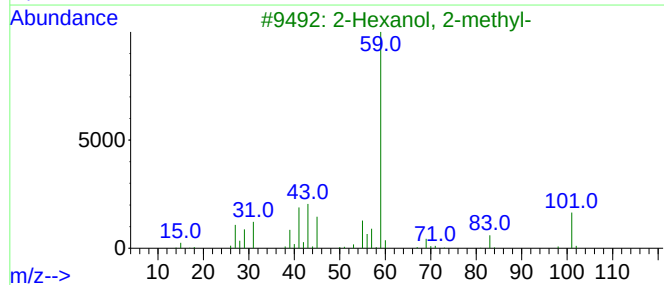
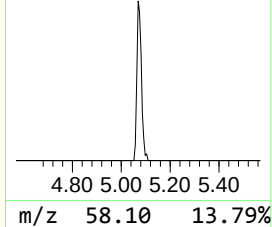
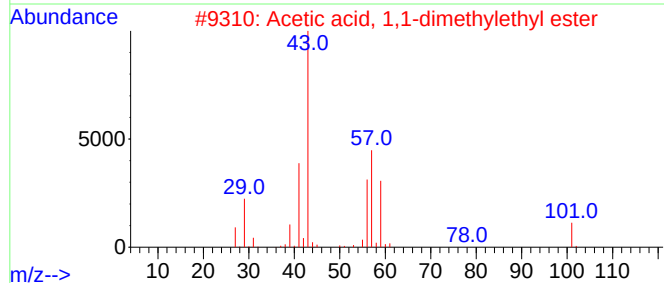
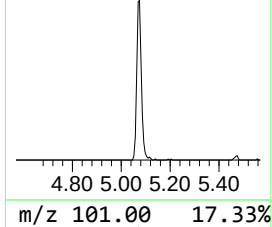
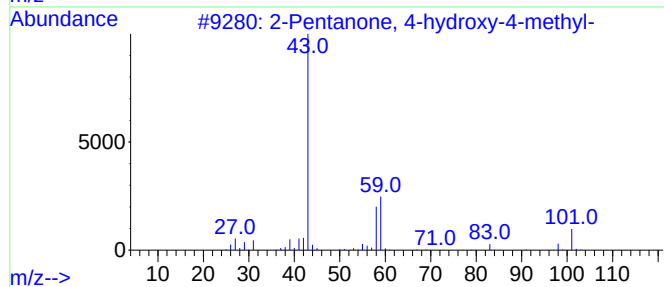
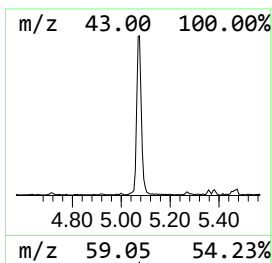
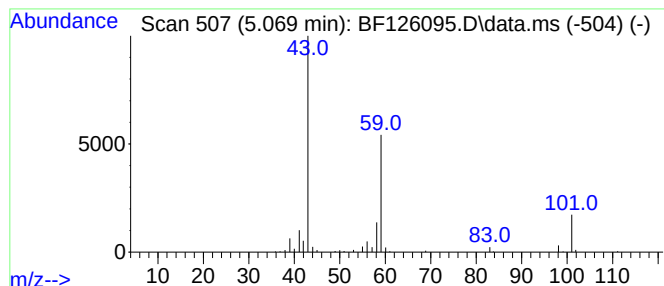
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.069	5.98 ng	355760	1,4-Dichlorobenzene-d4	6.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12



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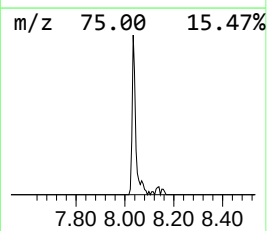
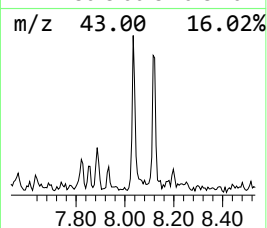
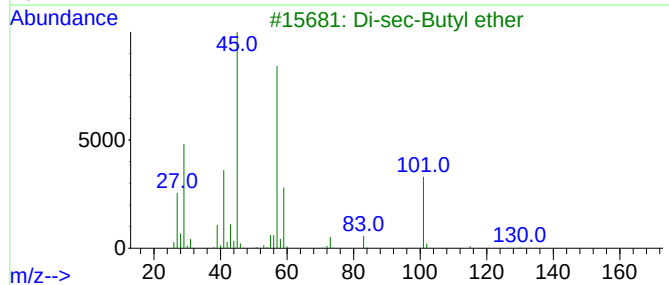
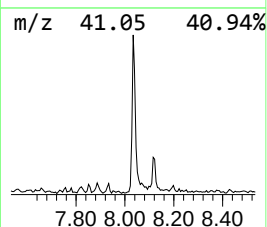
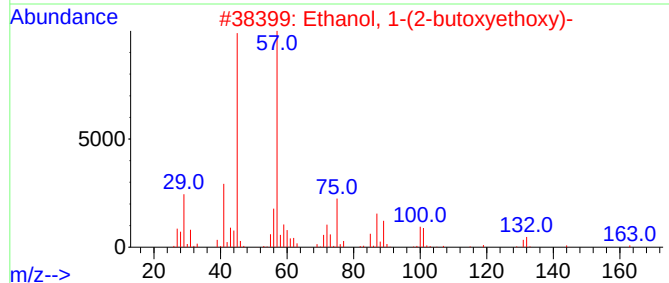
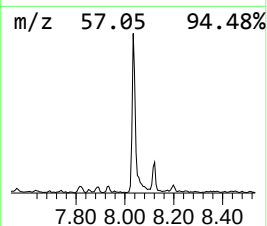
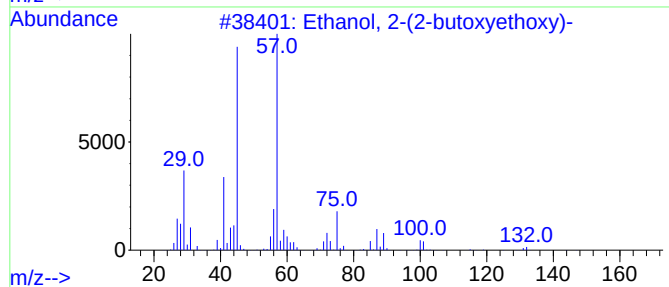
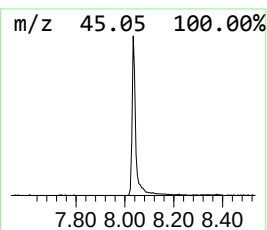
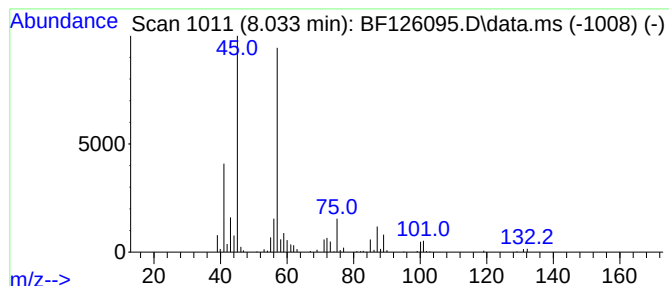
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Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.033	6.28 ng	482143	Naphthalene-d8	8.133

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3			Di-sec-Butyl ether	130	C8H18O	006863-58-7	59
4			3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	53
5			Tetraethylene glycol	194	C8H18O5	000112-60-7	50



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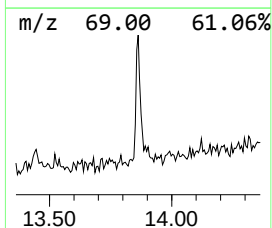
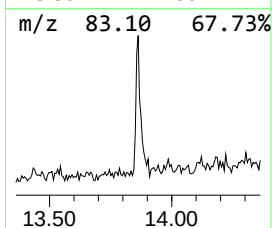
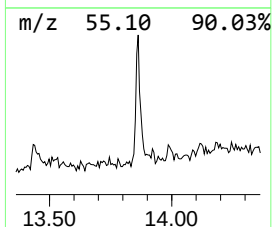
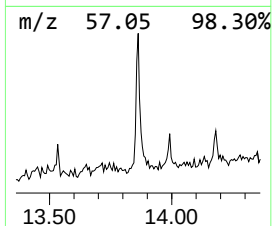
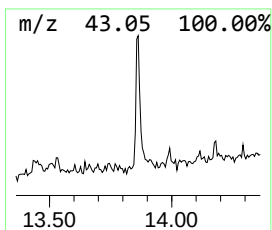
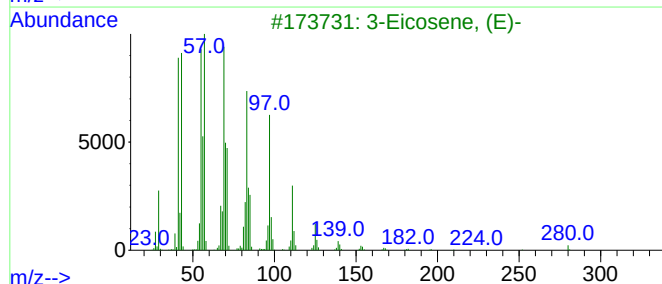
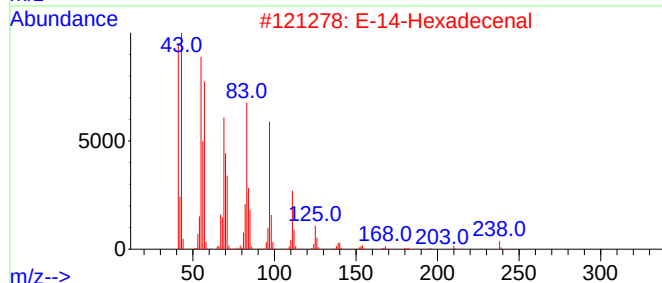
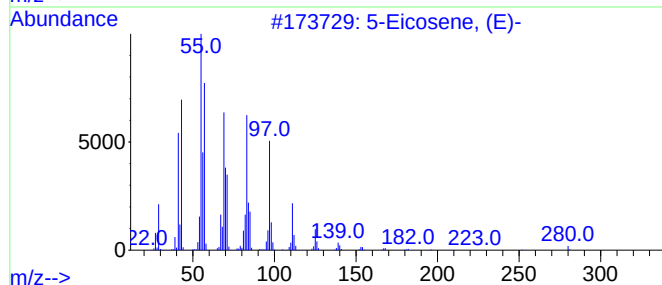
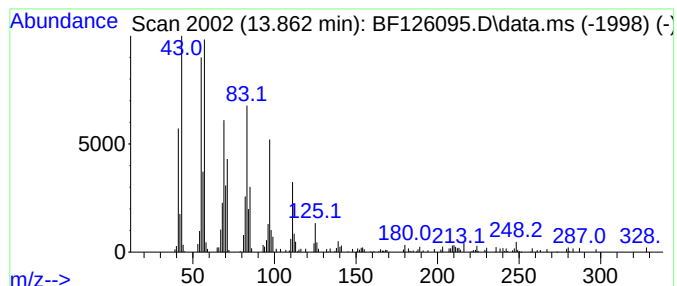
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Peak Number 5 5-Eicosene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.863	4.74 ng	291643	Chrysene-d12	14.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Eicosene, (E)-	280	C20H40	074685-30-6	95
2			E-14-Hexadecenal	238	C16H30O	330207-53-9	94
3			3-Eicosene, (E)-	280	C20H40	074685-33-9	94
4			1-Docosene	308	C22H44	001599-67-3	93
5			1-Eicosene	280	C20H40	003452-07-1	93



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.146	33.1	ng	1968340	1	6.851	1190800	20.0
2-Pentanone, 4-...	5.069	6.0	ng	355760	1	6.851	1190800	20.0
Ethanol, 2-(2-b...	8.033	6.3	ng	482143	2	8.133	1534820	20.0
5-Eicosene, (E)-	13.863	4.7	ng	291643	5	14.004	1231300	20.0