

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112024\
Data File : BF140494.D
Acq On : 20 Nov 2024 12:15
Operator : RC/JU
Sample : P4868-03
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RW5-SP303-20241114

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-BF111324.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF140494.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.134	3	7	17	rVB	1820686	2812812	66.08%	11.417%
2	3.599	248	256	263	rBV	31949	57693	1.36%	0.234%
3	4.475	398	405	422	rBV	651925	973531	22.87%	3.951%
4	4.740	444	450	462	rVB	42303	65969	1.55%	0.268%
5	5.087	501	509	519	rBV	70629	105711	2.48%	0.429%
6	5.504	574	580	599	rBV	1102352	1497426	35.18%	6.078%
7	6.510	743	751	763	rBV	688226	1005366	23.62%	4.081%
8	6.875	807	813	820	rBV	504896	646310	15.18%	2.623%
9	7.434	902	908	918	rBV	1744484	2374092	55.78%	9.636%
10	7.616	931	939	952	rVB	28852	53963	1.27%	0.219%
11	8.151	1025	1030	1046	rBV	638873	840370	19.74%	3.411%
12	9.228	1207	1213	1218	rBV	3292251	4256471	100.00%	17.277%
13	9.904	1323	1328	1339	rBV	702047	934741	21.96%	3.794%
14	10.698	1457	1463	1485	rBV	1916681	2742795	64.44%	11.133%
15	11.398	1576	1582	1591	rBV	782855	1017631	23.91%	4.130%
16	12.986	1846	1852	1857	rBV	3171010	4183423	98.28%	16.980%
17	14.051	2028	2033	2043	rVB	518650	692667	16.27%	2.811%
18	15.568	2285	2291	2304	rBV2	207577	376222	8.84%	1.527%

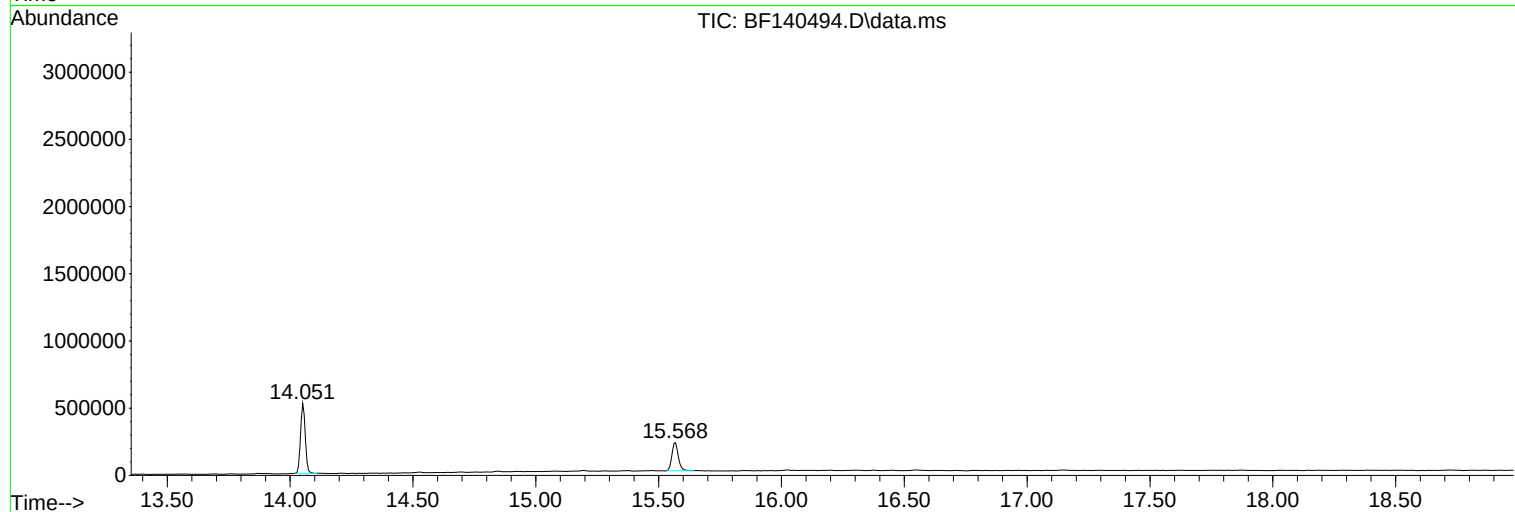
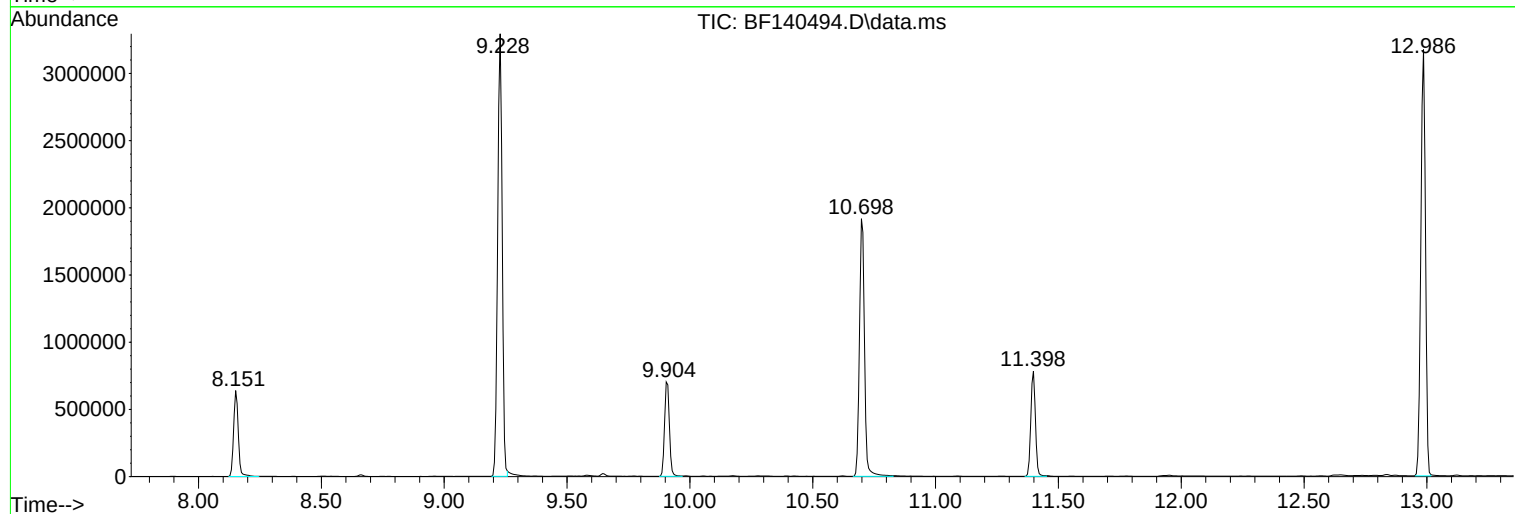
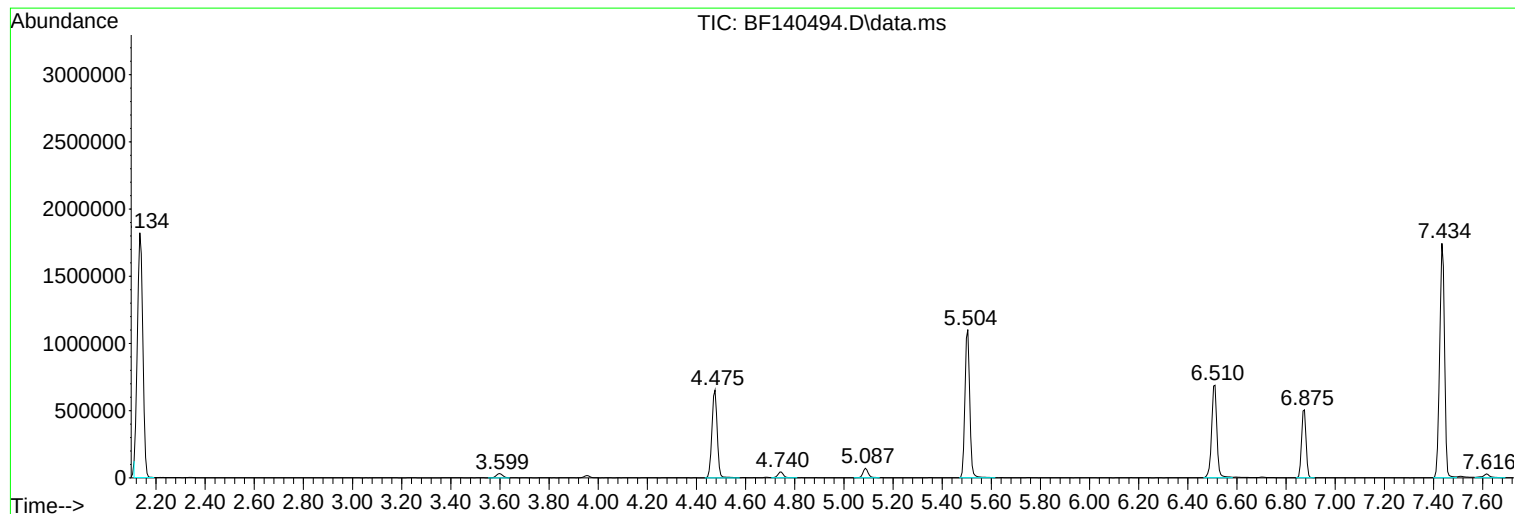
Sum of corrected areas: 24637193

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF111324.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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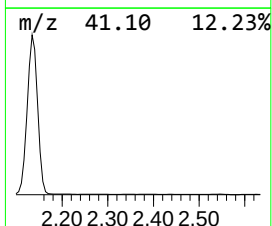
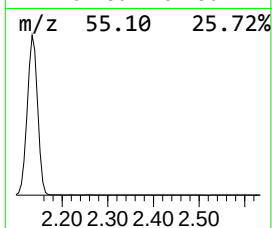
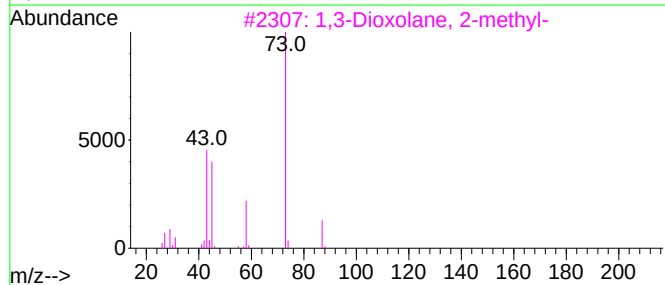
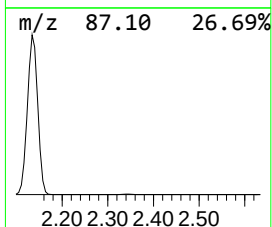
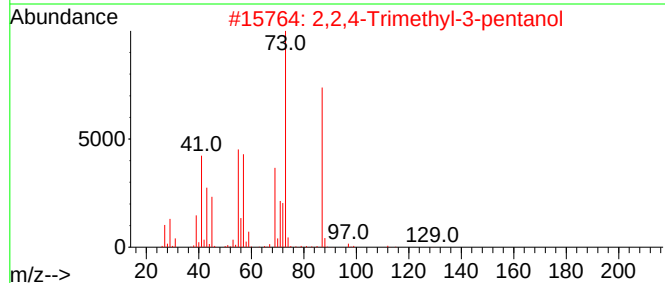
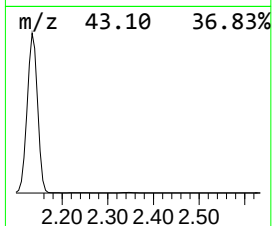
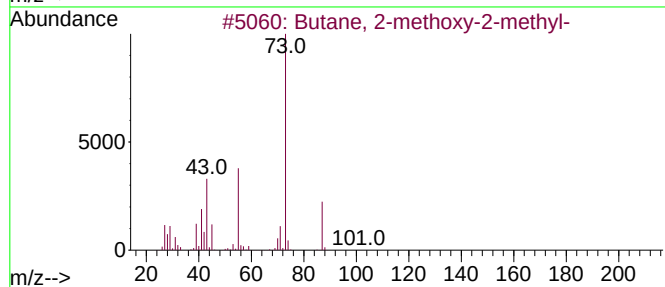
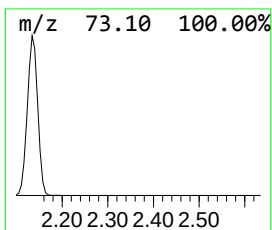
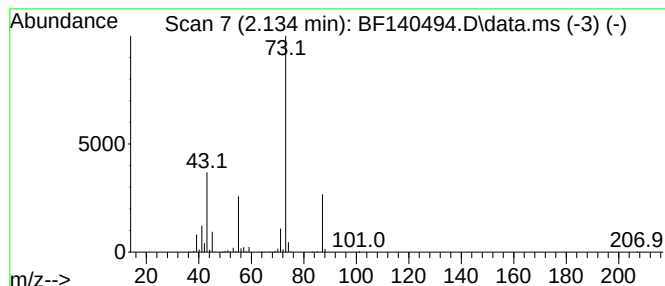
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF111324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.134	87.04 ng	2812810	1,4-Dichlorobenzene-d4	6.875

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	25
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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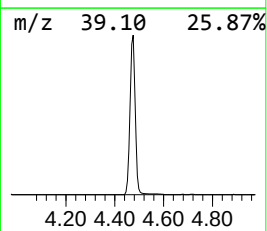
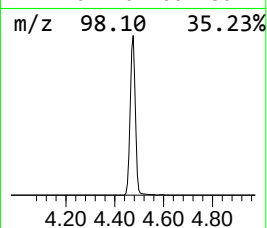
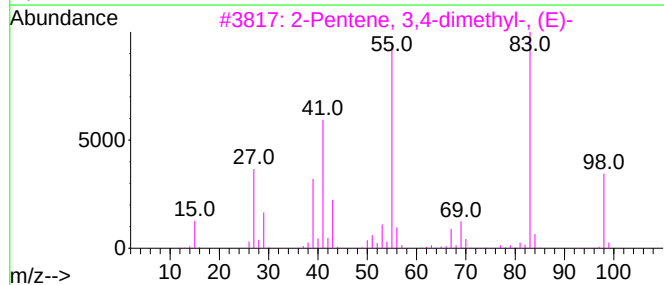
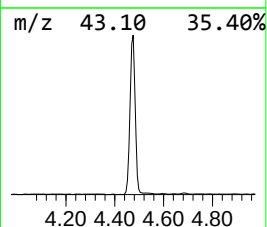
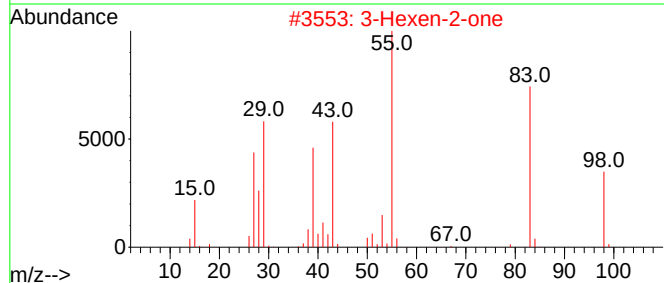
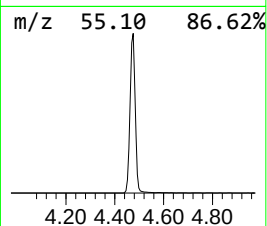
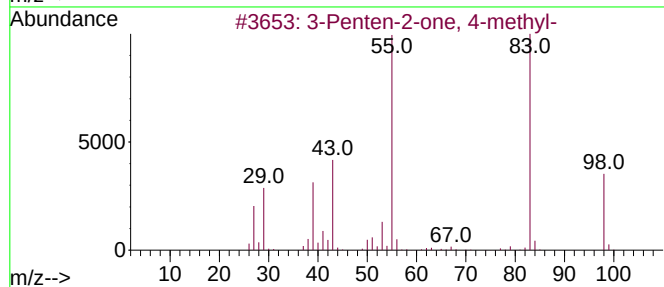
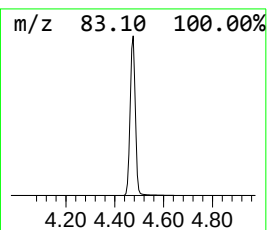
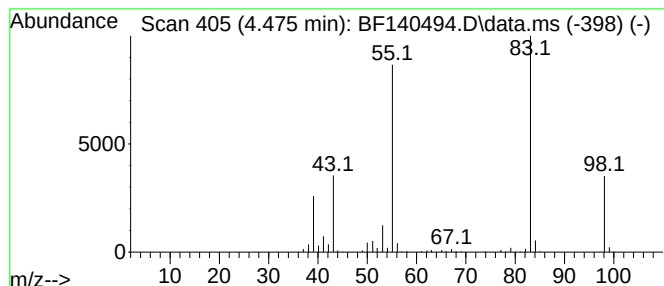
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF111324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.475	30.13 ng	973531	1,4-Dichlorobenzene-d4	6.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	94
2			3-Hexen-2-one	98	C6H10O	000763-93-9	91
3			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4			2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5			2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	80



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Misc :
ALS Vial : 8 Sample Multiplier: 1

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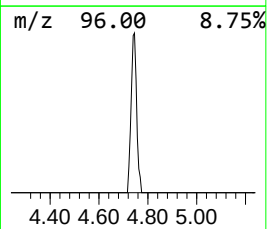
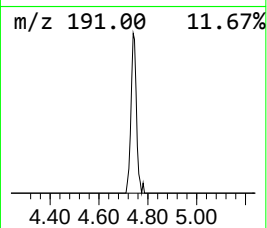
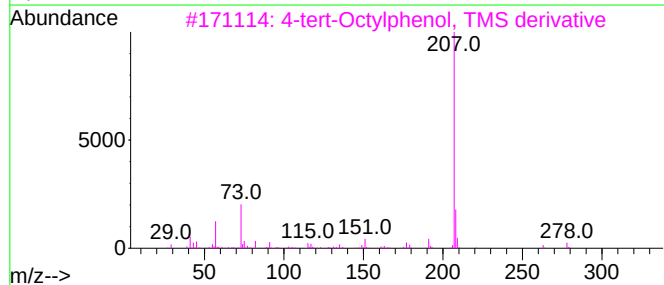
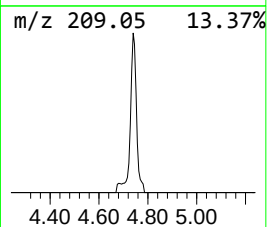
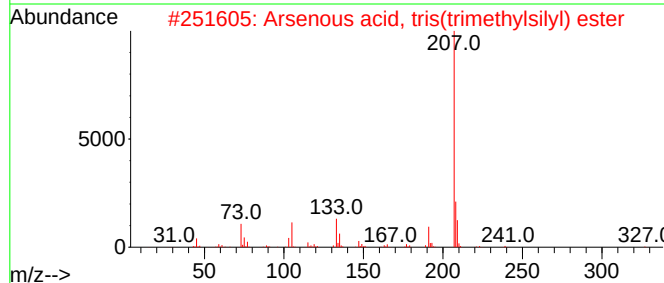
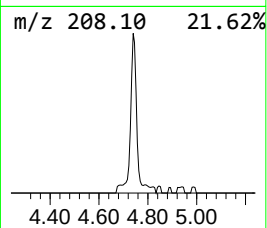
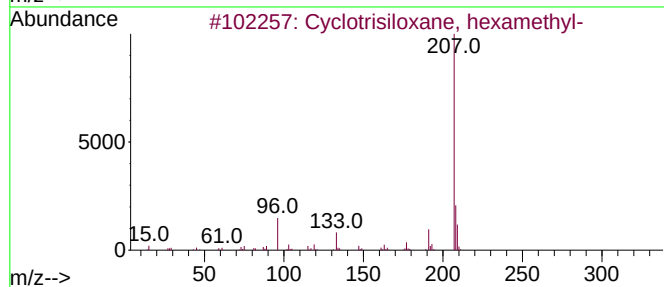
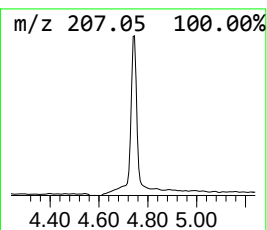
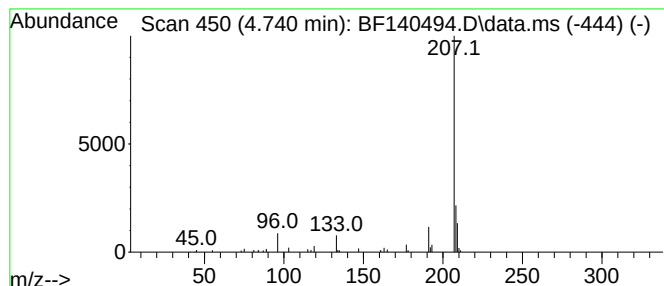
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclotrisiloxane, hexamethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.740	2.04 ng	65969	1,4-Dichlorobenzene-d4	6.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	90
2			Arsenous acid, tris(trimethylsil...	342	C9H27AsO3Si3	055429-29-3	83
3			4-tert-Octylphenol, TMS derivative	278	C17H30OSi	078721-87-6	39
4			Benzo[h]quinoline, 2,4-dimethyl-	207	C15H13N	000605-67-4	39
5			1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	38



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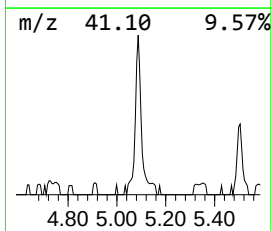
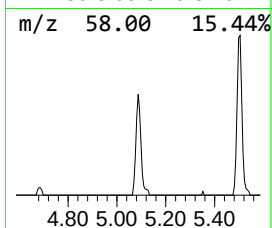
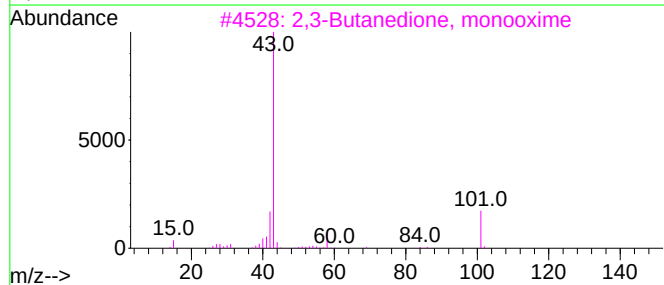
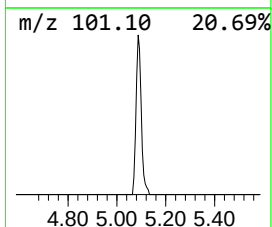
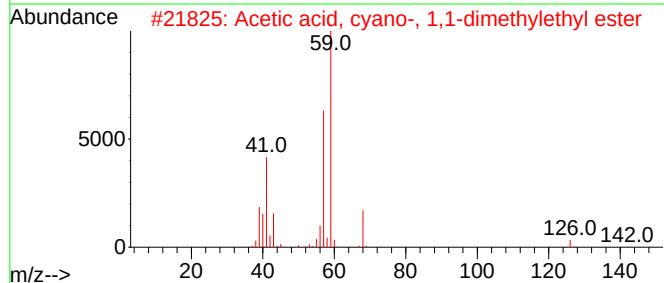
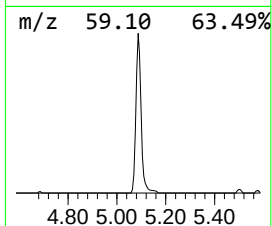
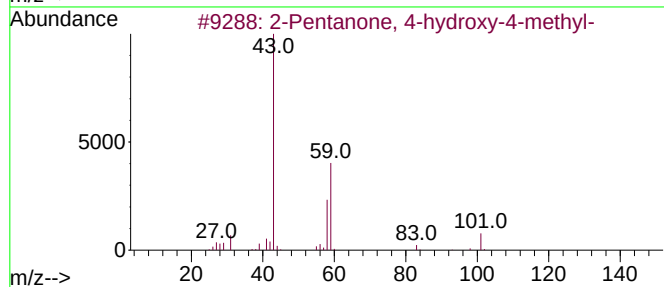
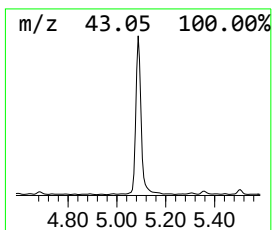
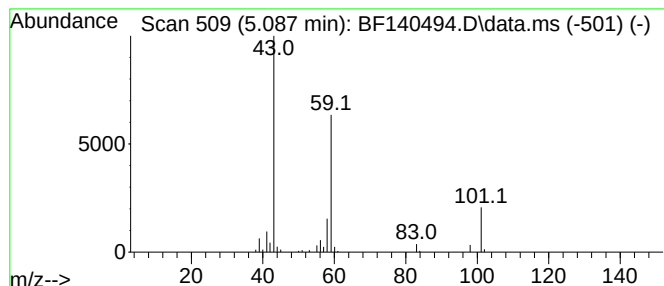
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.087	3.27 ng	105711	1,4-Dichlorobenzene-d4	6.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.134	87.0	ng	2812810	1	6.875	646310	20.0
3-Penten-2-one,...	4.475	30.1	ng	973531	1	6.875	646310	20.0
Cyclotrisiloxan...	4.740	2.0	ng	65969	1	6.875	646310	20.0
2-Pentanone, 4-...	5.087	3.3	ng	105711	1	6.875	646310	20.0