

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011022\
 Data File : BF126569.D
 Acq On : 10 Jan 2022 16:36
 Operator : JU/CG
 Sample : N1054-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20220020-CAPE-2-MODIFIED-ELUTRIATE

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-BF122921.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF126569.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.975	488	491	498	rVB	52238	65089	2.39%	0.444%
2	5.386	557	561	564	rBV	1733418	1555295	57.18%	10.620%
3	6.404	729	734	742	rBV	1341351	1318399	48.47%	9.003%
4	6.769	792	796	799	rBV	615441	494247	18.17%	3.375%
5	7.333	886	892	895	rBV	1765444	1892156	69.56%	12.920%
6	7.957	994	998	1007	rBV	303422	256568	9.43%	1.752%
7	8.051	1010	1014	1017	rBV	827961	681591	25.06%	4.654%
8	9.133	1192	1198	1201	rBV	2797861	2720187	100.00%	18.575%
9	9.810	1308	1313	1316	rBV	829377	711323	26.15%	4.857%
10	10.598	1442	1447	1451	rBV	1575584	1481880	54.48%	10.119%
11	11.292	1561	1565	1569	rVB	725015	634315	23.32%	4.331%
12	11.698	1626	1634	1637	rVB2	35292	37638	1.38%	0.257%
13	11.833	1654	1657	1665	rBV2	44098	54556	2.01%	0.373%
14	12.886	1831	1836	1839	rBV	1874247	1695681	62.34%	11.579%
15	13.792	1984	1990	2002	rBV2	70337	128041	4.71%	0.874%
16	13.939	2011	2015	2021	rVB	472242	404007	14.85%	2.759%
17	14.039	2026	2032	2039	rBV	82204	80624	2.96%	0.551%
18	15.398	2258	2263	2268	rBV	365265	433007	15.92%	2.957%

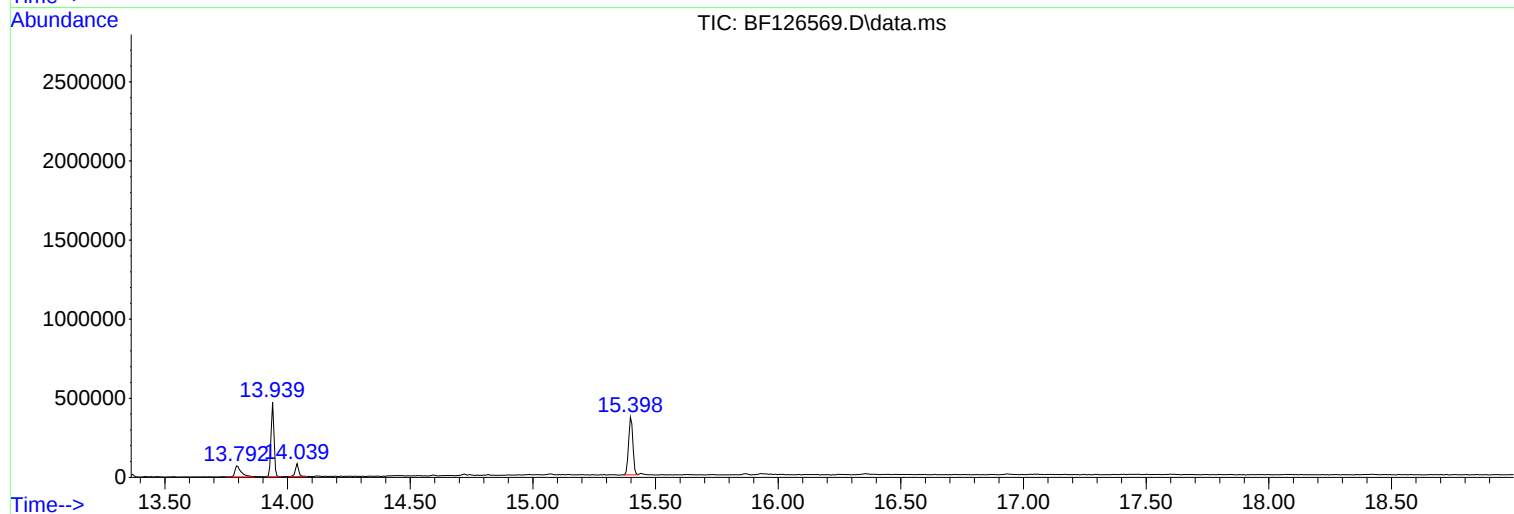
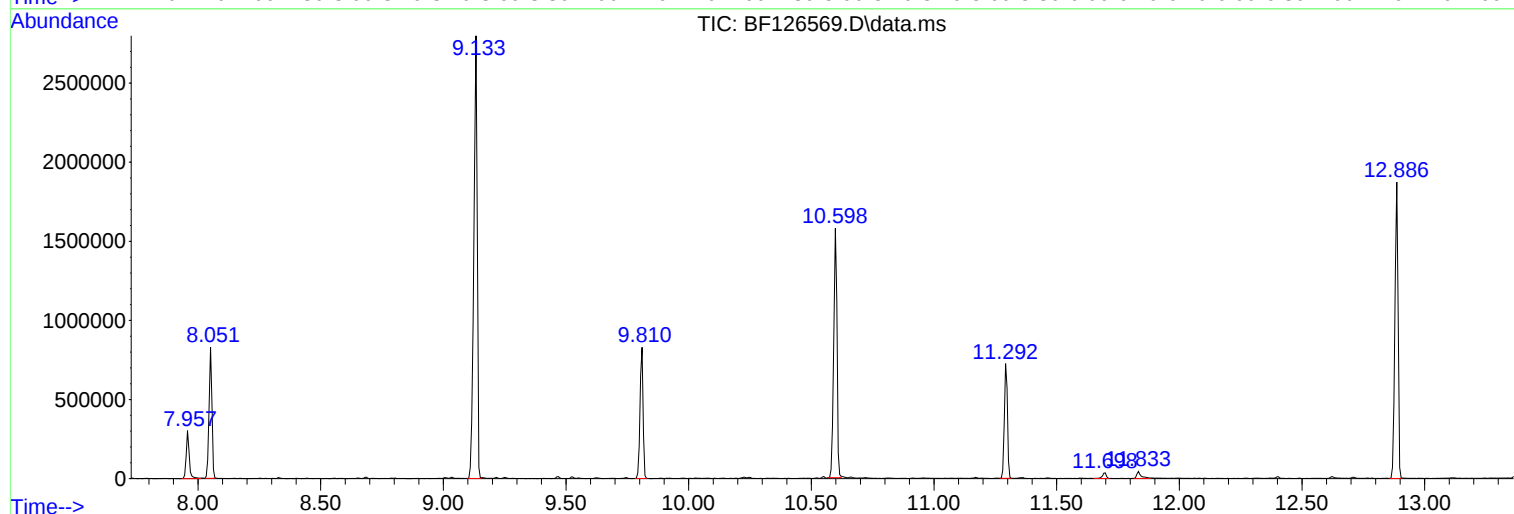
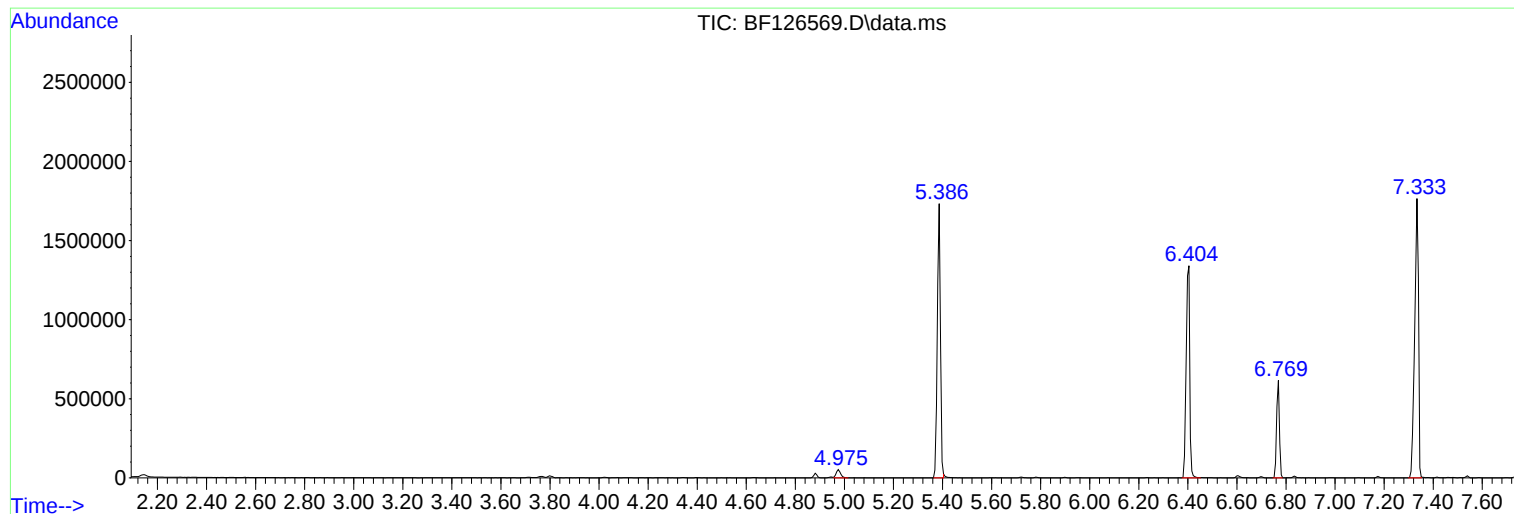
Sum of corrected areas: 14644604

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122921.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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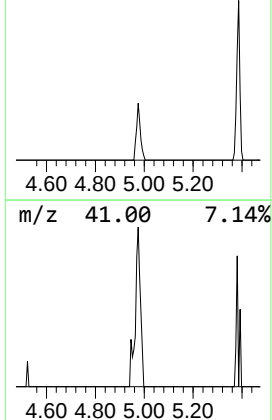
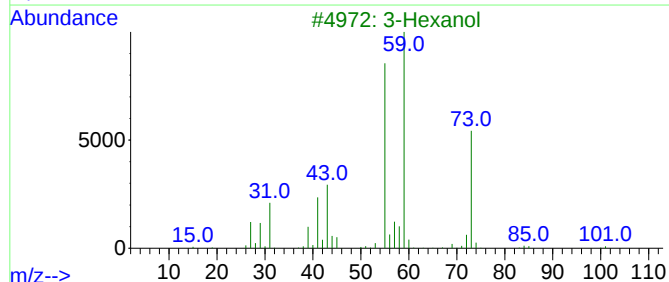
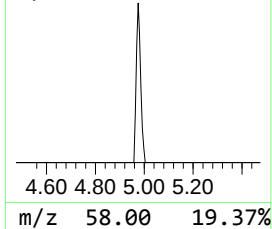
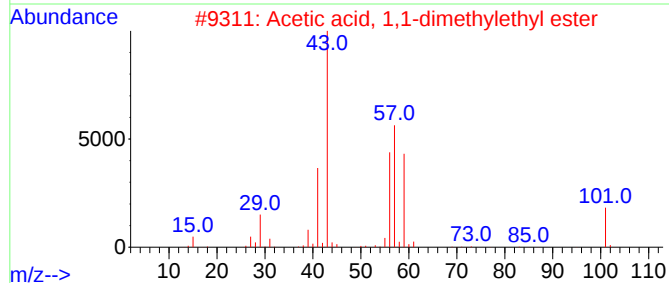
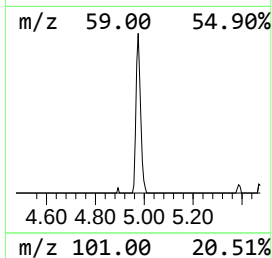
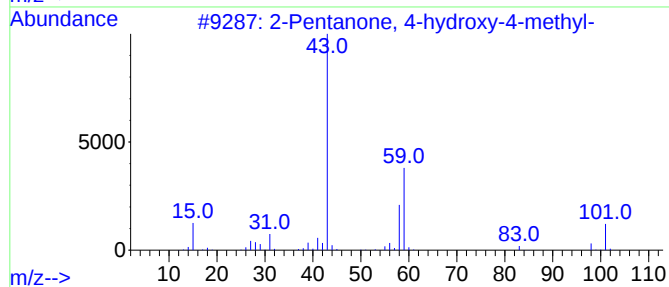
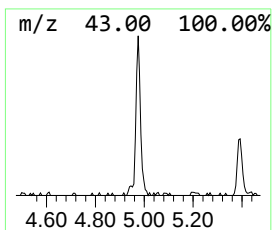
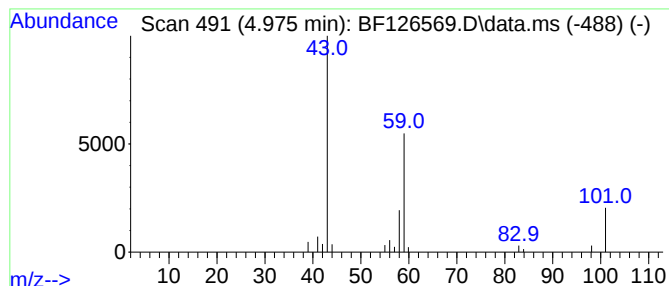
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122921.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.975	2.63 ng	65089	1,4-Dichlorobenzene-d4	6.769

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3		3-Hexanol	102	C6H14O	000623-37-0	23
4		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
5		4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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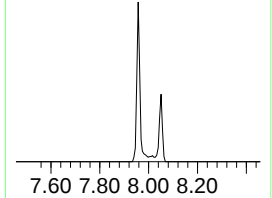
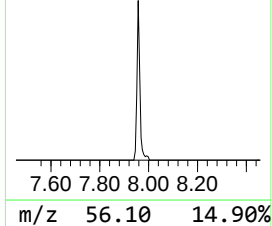
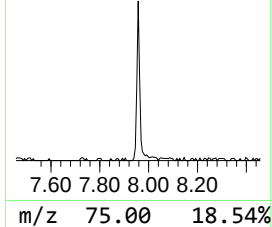
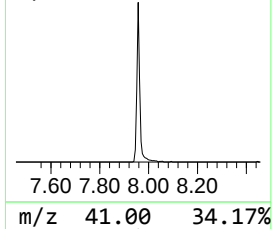
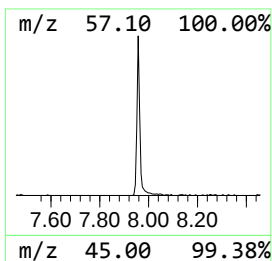
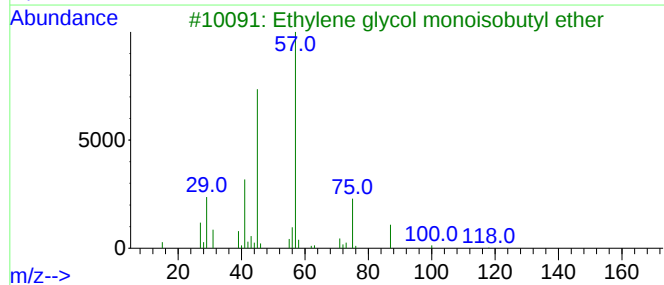
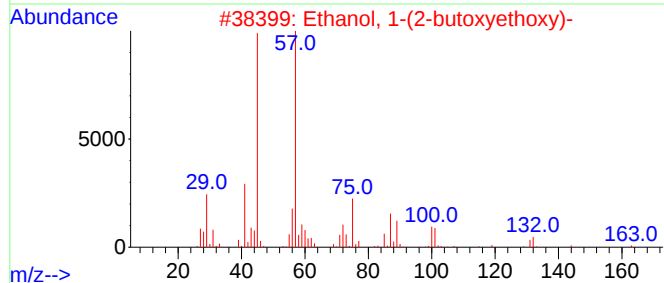
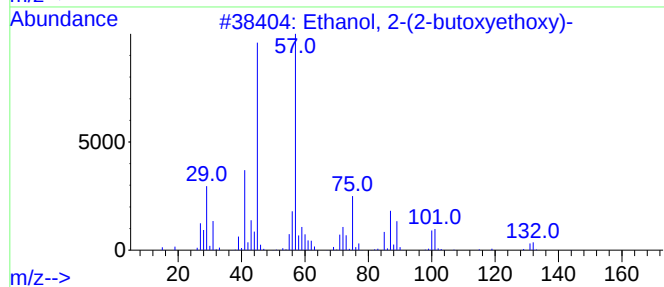
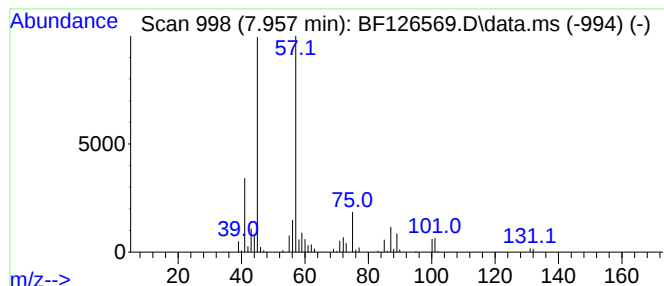
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.957	7.53 ng	256568	Naphthalene-d8	8.051

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			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
4			Di-sec-Butyl ether	130	C8H18O	006863-58-7	53
5			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	50



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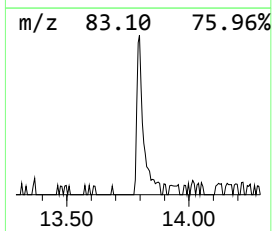
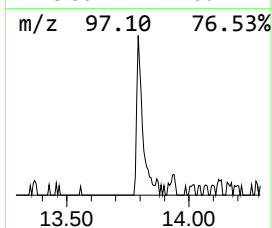
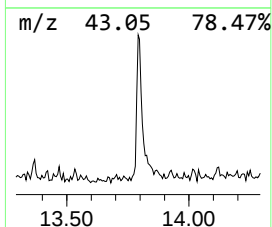
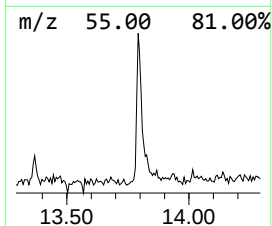
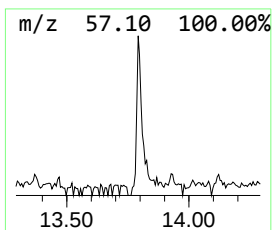
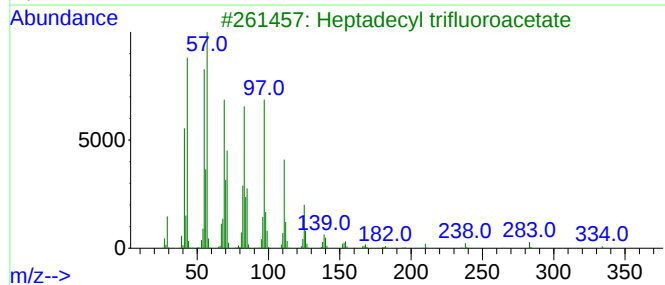
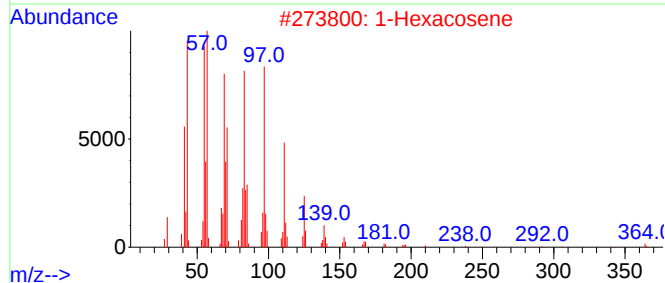
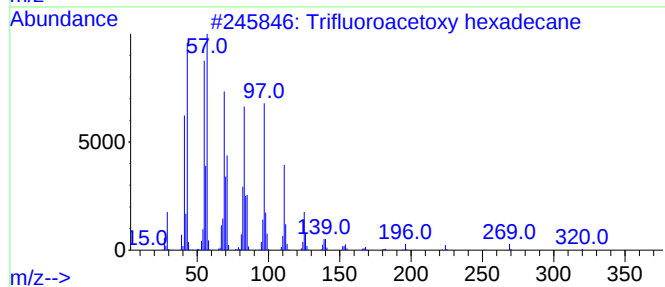
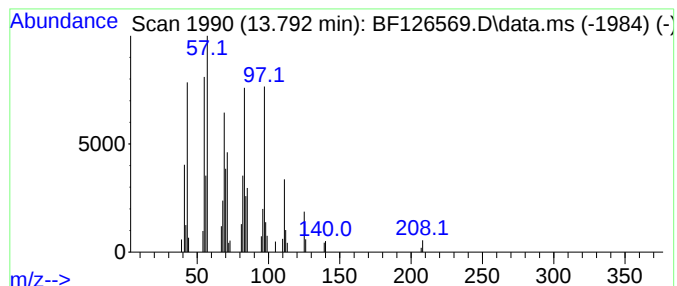
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Peak Number 3 Trifluoroacetoxy hexadecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.792	6.34 ng	128041	Chrysene-d12	13.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
2			1-Hexacosene	364	C26H52	018835-33-1	91
3			Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	91
4			1-Docosene	308	C22H44	001599-67-3	91
5			1-Tricosene	322	C23H46	018835-32-0	91



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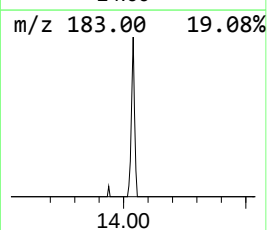
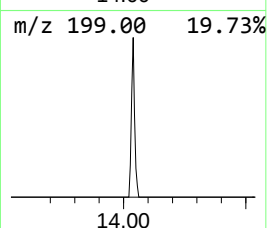
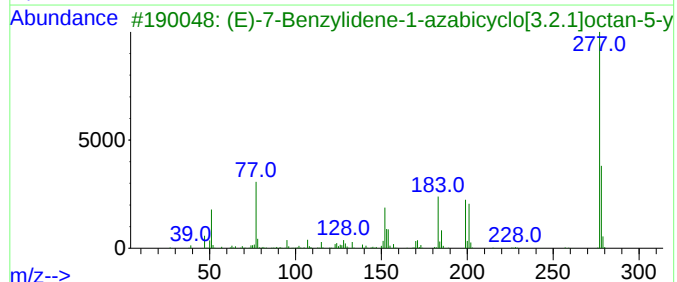
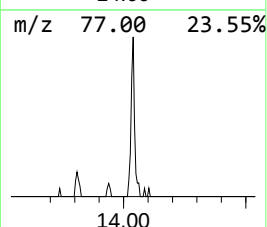
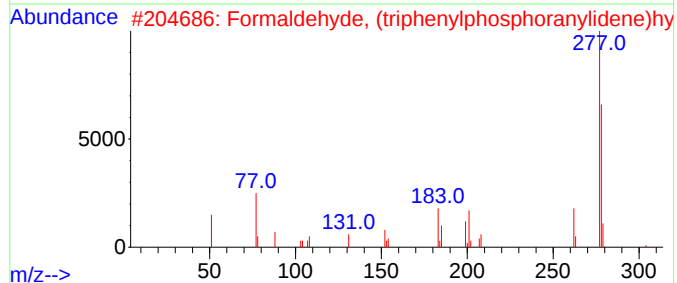
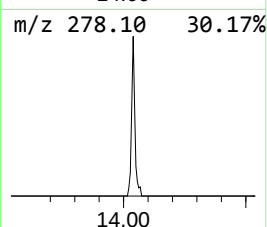
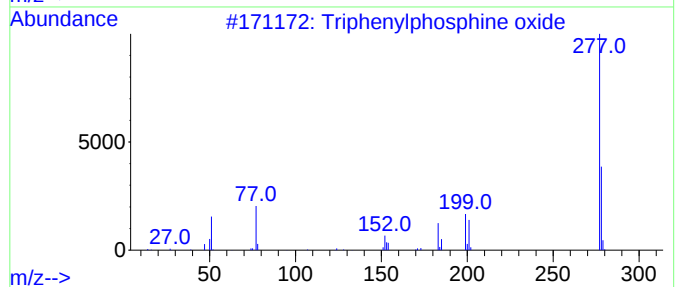
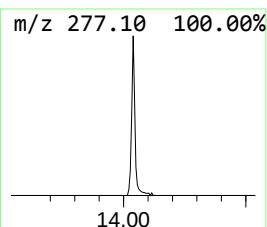
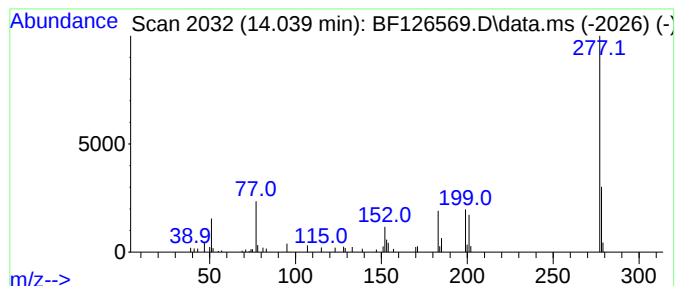
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Peak Number 4 Triphenylphosphine oxide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.039	3.99 ng	80624	Chrysene-d12	13.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	97
2			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	50
3			(E)-7-Benzylidene-1-azabicyclo[3...	293	C15H19N03S	1000483-83-4	43
4			Carbazol-1-one, 1,2,3,4-tetrahyd...	277	C18H15NO2	299962-12-2	40
5			Cobalt,(.eta.<5>-2,4-cyclopentad...	277	C16H15BCo	036682-12-9	23



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.975	2.6	ng	65089	1	6.769	494247	20.0
Ethanol, 2-(2-b...	7.957	7.5	ng	256568	2	8.051	681591	20.0
Trifluoroacetox...	13.792	6.3	ng	128041	5	13.939	404007	20.0
Triphenylphosph...	14.039	4.0	ng	80624	5	13.939	404007	20.0