

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090222\
 Data File : BF130116.D
 Acq On : 02 Sep 2022 17:50
 Operator : CG\JU
 Sample : N4466-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAP2-T-5-(15-35)

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-BF083122.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF130116.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.899	439	444	456	rVB	692986	847703	16.82%	2.409%
2	5.310	508	514	517	rBV	3866677	4324137	85.80%	12.287%
3	5.716	579	583	588	rVB	73723	65189	1.29%	0.185%
4	6.334	682	688	691	rBV	3862739	4288871	85.10%	12.187%
5	6.704	747	751	754	rBV	1372056	1186722	23.55%	3.372%
6	7.269	841	847	850	rBV	2648477	2765565	54.88%	7.858%
7	7.904	951	955	964	rBV2	55463	167677	3.33%	0.476%
8	7.981	964	968	972	rVB	1892096	1617188	32.09%	4.595%
9	9.057	1146	1151	1154	rBV	5103371	5039572	100.00%	14.320%
10	9.733	1261	1266	1269	rBV	2108280	1843872	36.59%	5.239%
11	10.516	1395	1399	1402	rBV	3443590	3245090	64.39%	9.221%
12	11.210	1513	1517	1521	rBV	2132351	1798184	35.68%	5.110%
13	11.280	1527	1529	1532	rVB	73827	59932	1.19%	0.170%
14	11.745	1604	1608	1612	rBV	100736	95452	1.89%	0.271%
15	12.798	1783	1787	1791	rBV	4649751	4640405	92.08%	13.186%
16	13.380	1882	1886	1888	rBV	63967	54059	1.07%	0.154%
17	13.704	1938	1941	1945	rBV	90069	75715	1.50%	0.215%
18	13.839	1960	1964	1968	rBV	1513410	1328857	26.37%	3.776%
19	13.945	1977	1982	1990	rVB	144328	183960	3.65%	0.523%
20	14.021	1992	1995	1999	rVB	104348	86534	1.72%	0.246%
21	14.327	2043	2047	2050	rBV	88951	77714	1.54%	0.221%
22	14.621	2094	2097	2102	rVB	83906	73544	1.46%	0.209%
23	14.939	2147	2151	2156	rBV	57800	66621	1.32%	0.189%
24	15.239	2197	2202	2207	rBV	889177	1060317	21.04%	3.013%
25	15.286	2207	2210	2217	rVB	60272	78953	1.57%	0.224%
26	15.692	2275	2279	2284	rBV	48828	66722	1.32%	0.190%
27	16.157	2354	2358	2363	rVB	34043	54307	1.08%	0.154%

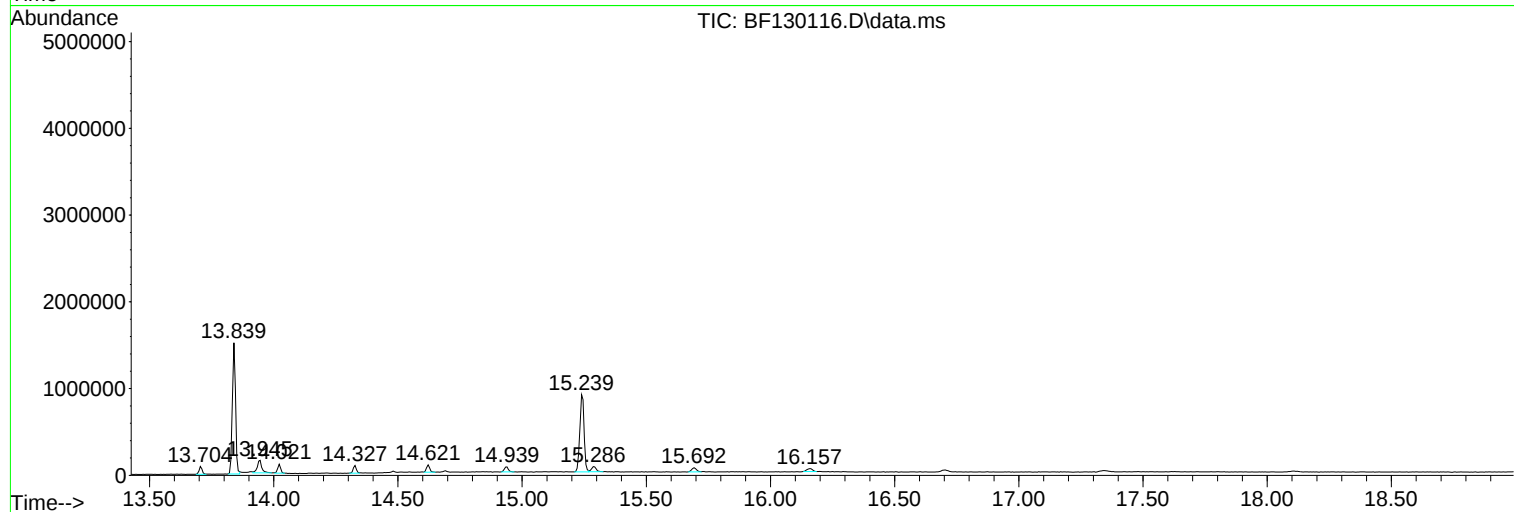
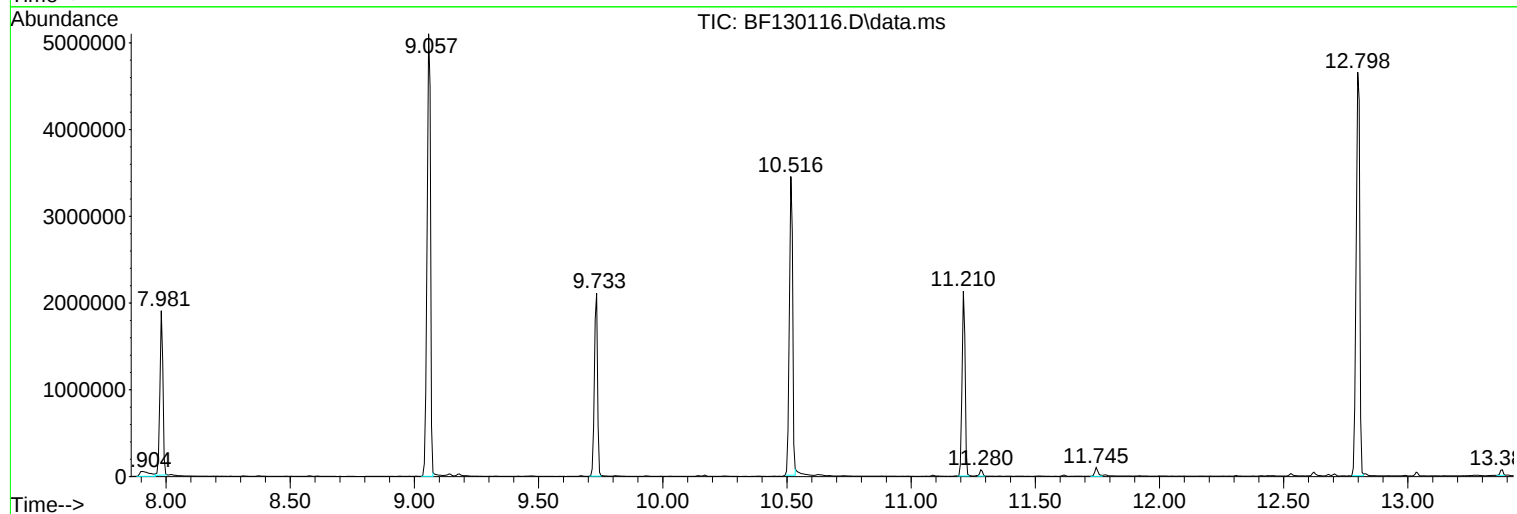
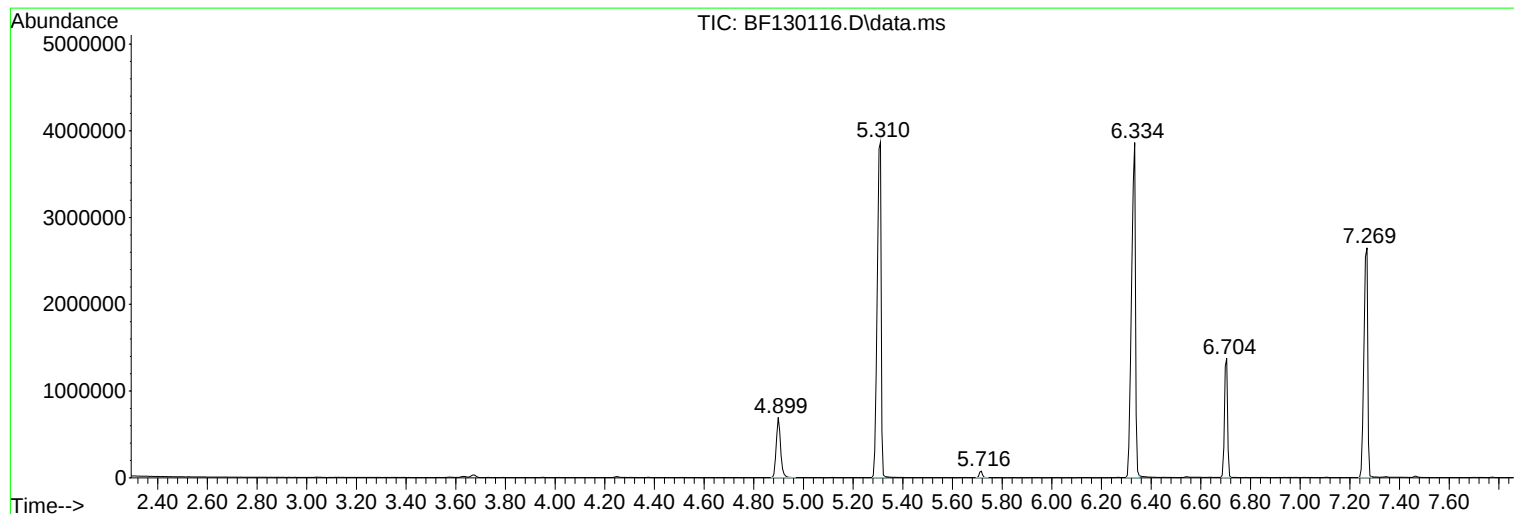
Sum of corrected areas: 35192862

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.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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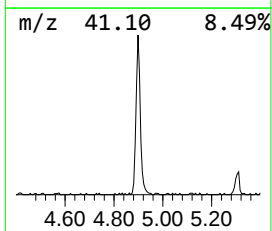
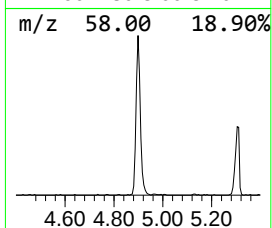
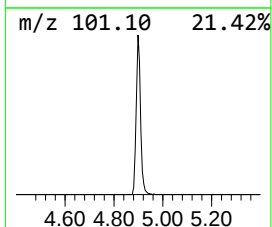
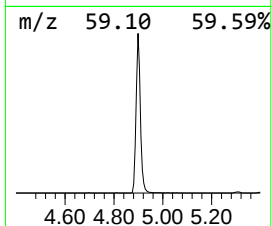
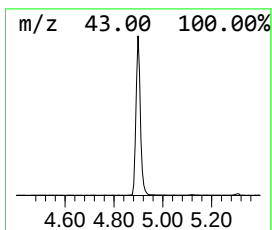
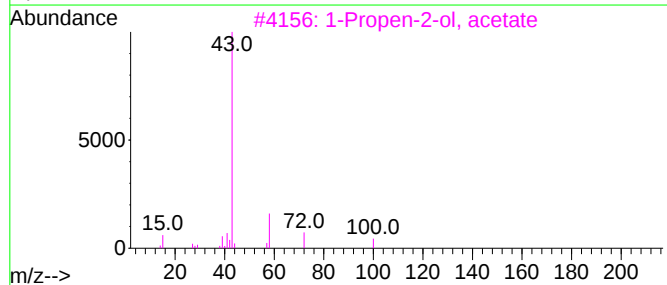
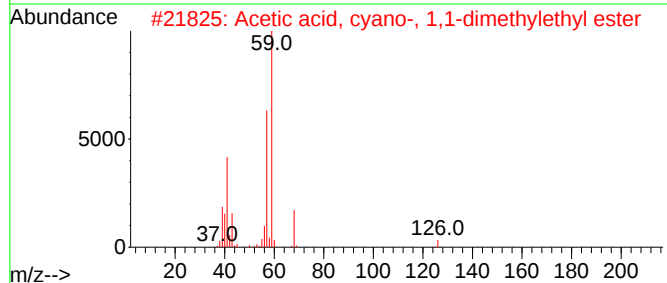
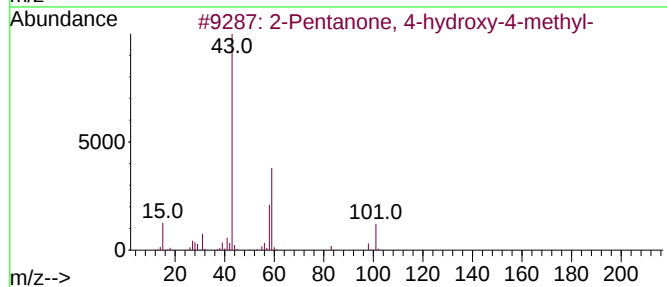
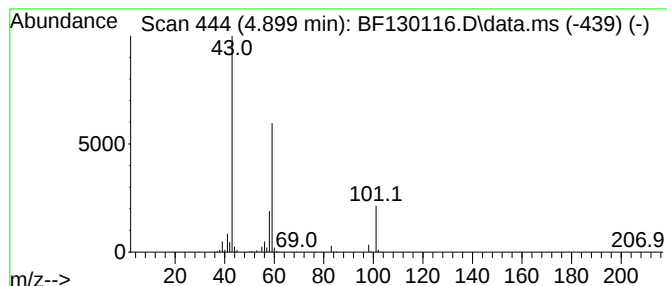
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.899	14.29 ng	847703	1,4-Dichlorobenzene-d4	6.704

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, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Acetone	58	C3H6O	000067-64-1	9



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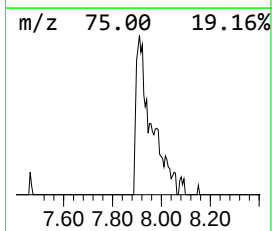
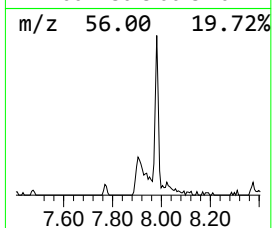
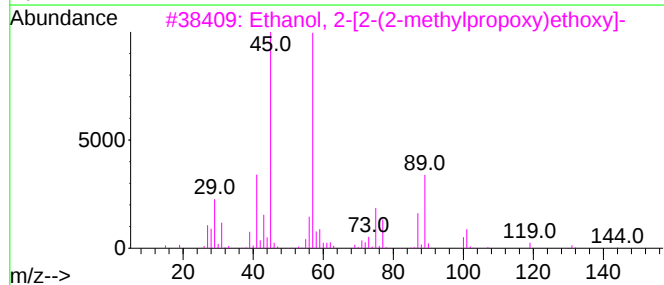
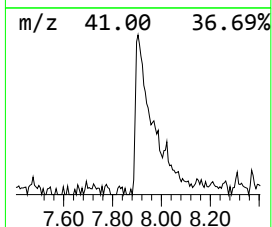
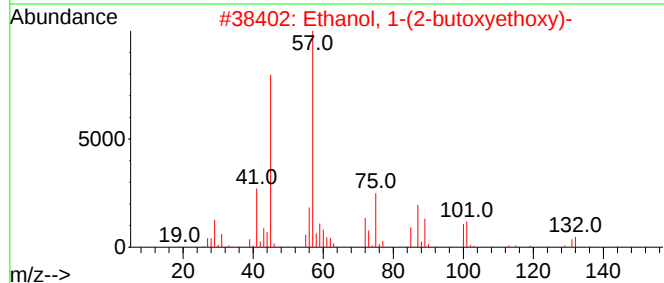
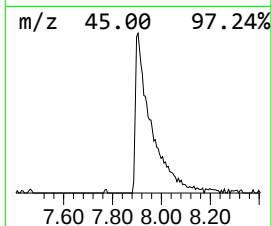
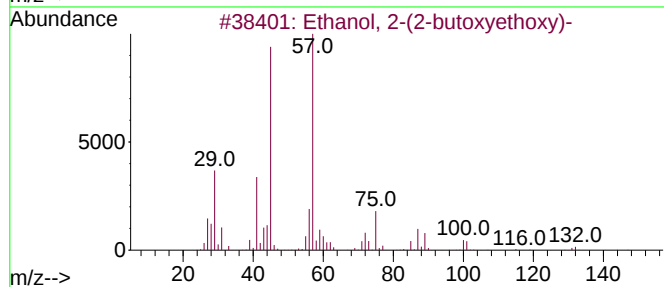
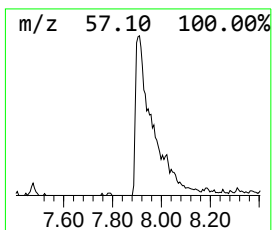
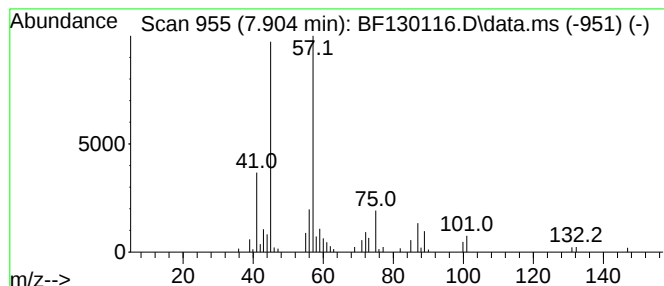
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Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.904	2.07 ng	167677	Naphthalene-d8	7.981

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	78
3			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	72
4			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	59
5			Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	47



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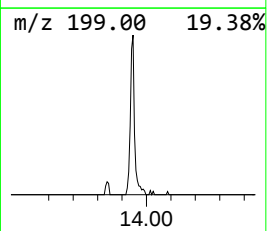
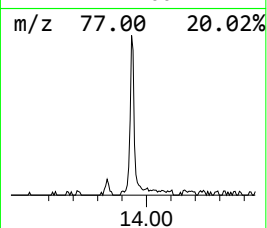
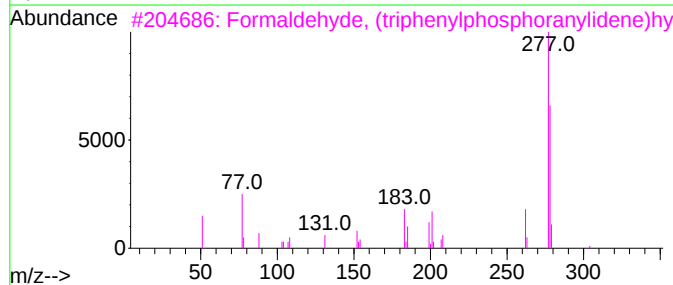
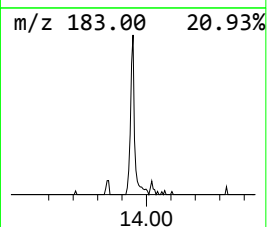
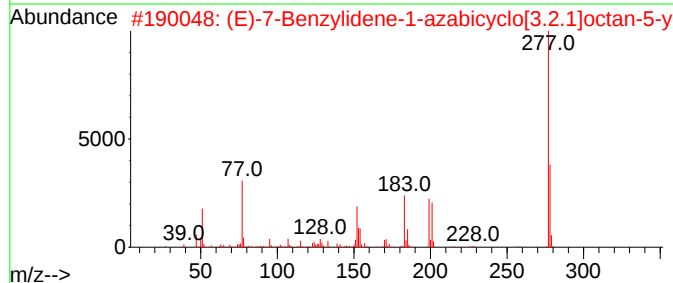
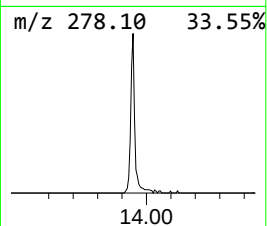
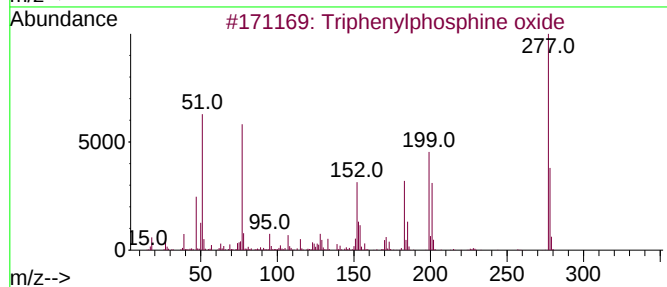
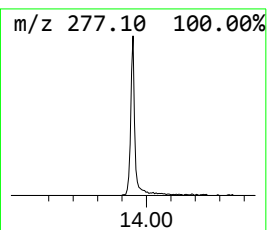
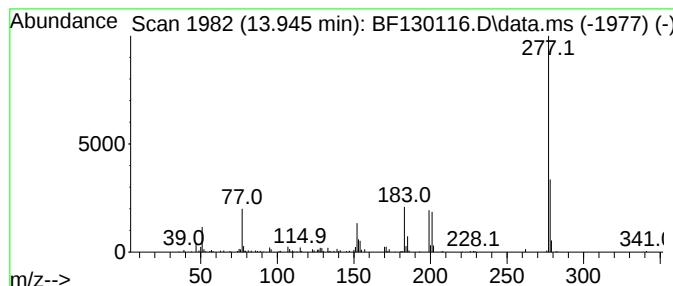
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 Triphenylphosphine oxide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.945	2.77 ng	183960	Chrysene-d12	13.839

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	97
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19N03S	1000483-83-4	91
3			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	42
4			(Carbethoxymethyl)-triphenylphos...	428	C22H22BrO2P	001530-45-6	38
5			5-Methyl-8-phenylfuro[3',2':5,6]...	277	C16H11N3O2	1000460-07-9	28



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.899	14.3	ng	847703	1	6.704	1186720	20.0
Ethanol, 2-(2-b...	7.904	2.1	ng	167677	2	7.981	1617190	20.0
Triphenylphosph...	13.945	2.8	ng	183960	5	13.839	1328860	20.0