

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF052920\
 Data File : BF120450.D
 Acq On : 29 May 2020 20:17
 Operator : MA/SJ
 Sample : L2790-11
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3-S

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-BF052820.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	1.922	3	6	10	rVB	651733	782279	9.33%	1.568%
2	1.958	10	12	32	rVV	1729332	2722492	32.46%	5.458%
3	2.099	32	36	55	rVB	1316887	1757136	20.95%	3.523%
4	5.451	601	606	609	rBV	3844235	3504109	41.78%	7.025%
5	6.463	773	778	781	rBV	3838489	3730123	44.47%	7.479%
6	6.834	838	841	845	rBV	1715623	1560096	18.60%	3.128%
7	6.992	864	868	871	rBV	3823708	3245674	38.69%	6.507%
8	7.398	932	937	940	rBV	2824603	2474459	29.50%	4.961%
9	8.116	1055	1059	1063	rBV	2291968	1971447	23.50%	3.953%
10	9.192	1238	1242	1246	rBV	5011166	4696799	55.99%	9.417%
11	9.586	1305	1309	1312	rBV	708941	518432	6.18%	1.039%
12	9.875	1354	1358	1361	rVB	2402387	2073782	24.72%	4.158%
13	10.657	1487	1491	1494	rBV	2894947	2418219	28.83%	4.848%
14	11.357	1606	1610	1613	rBV	2193217	1754955	20.92%	3.519%
15	11.880	1696	1699	1704	rBV	165663	161081	1.92%	0.323%
16	12.939	1875	1879	1883	rBV	3844346	3512081	41.87%	7.041%
17	13.486	1968	1972	1975	rBV	2715567	2242288	26.73%	4.496%
18	13.845	2027	2033	2043	rBV3	441139	794818	9.48%	1.594%
19	13.986	2052	2057	2061	rBV2	7652208	8387898	100.00%	16.817%
20	15.445	2300	2305	2309	rBV	1210212	1411253	16.82%	2.829%
21	15.980	2392	2396	2403	rVB5	75807	157832	1.88%	0.316%

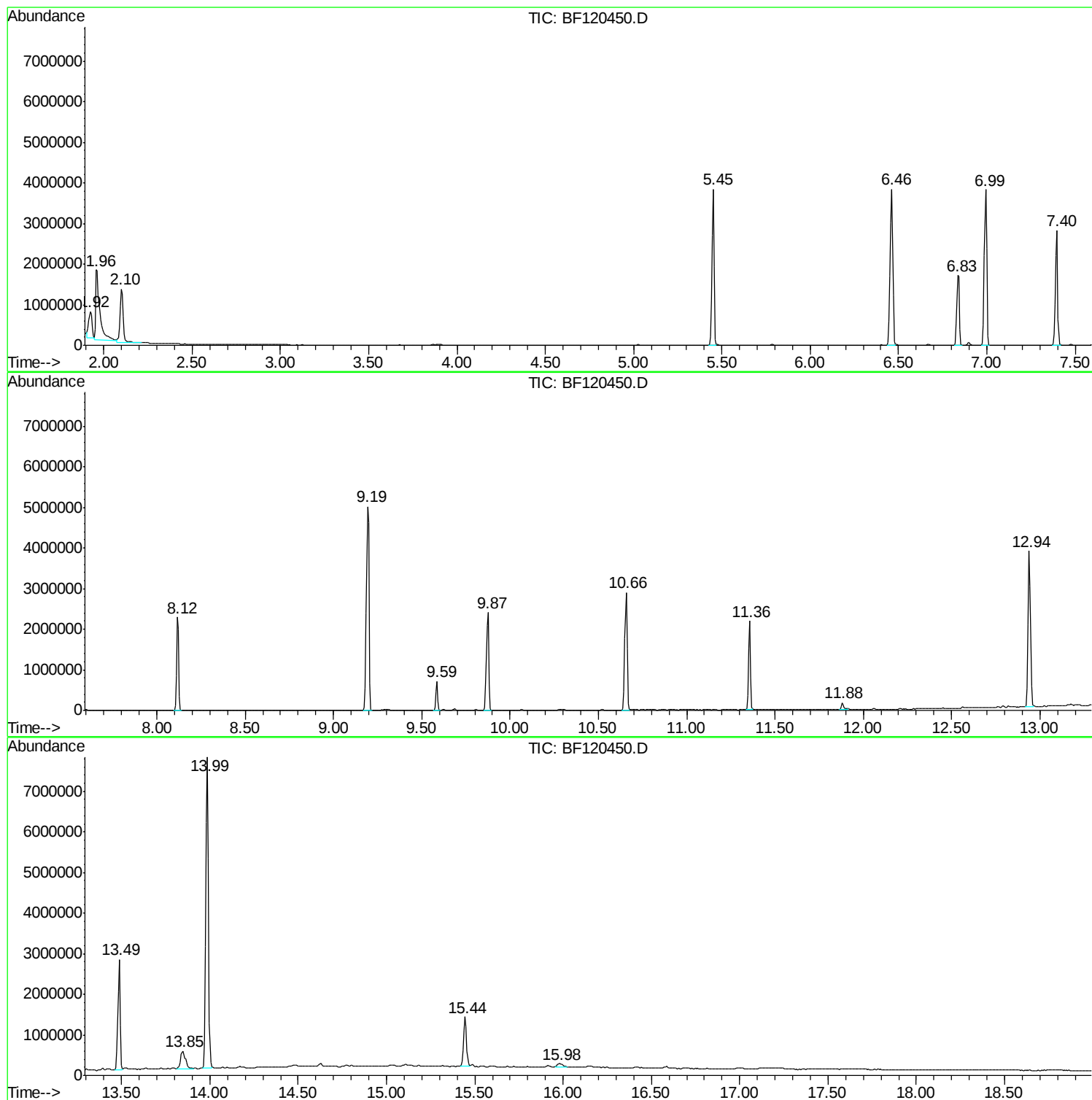
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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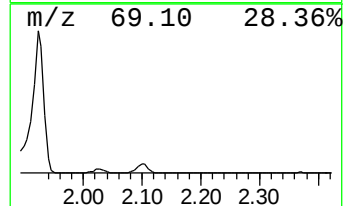
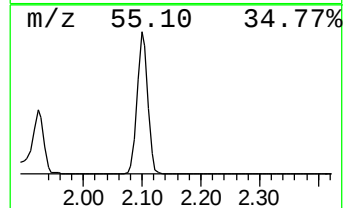
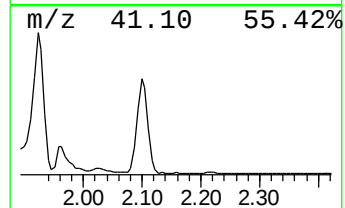
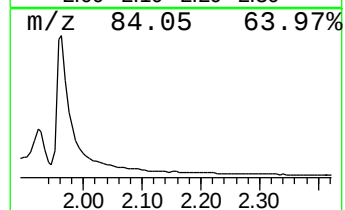
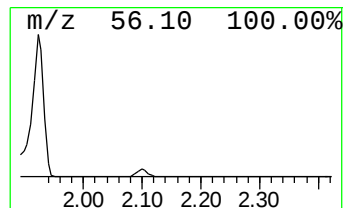
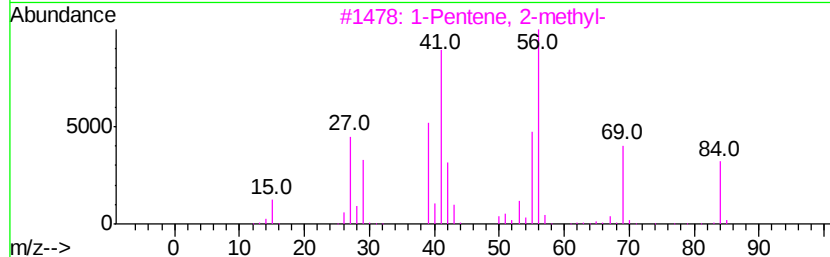
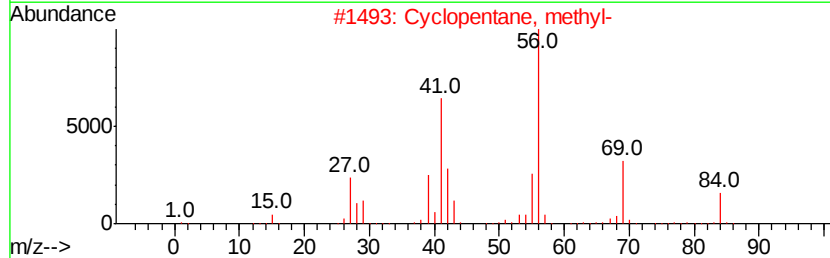
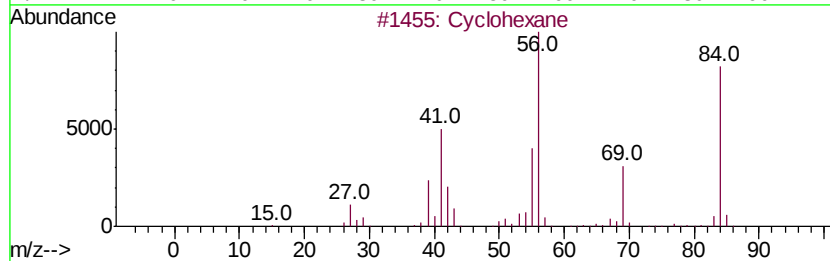
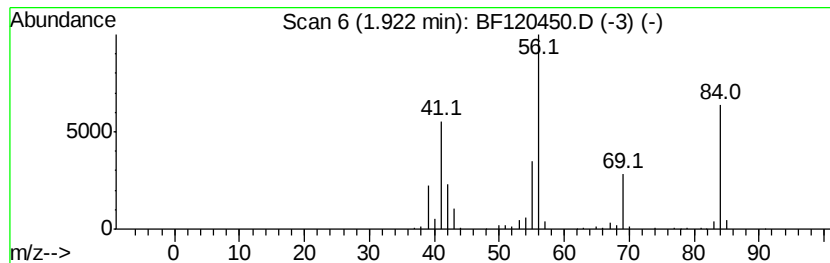
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclohexane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.92	10.03 ng	782279	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane	84	C6H12	000110-82-7	91
2		Cyclopentane, methyl-	84	C6H12	000096-37-7	59
3		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	42
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	40
5		2H-Pyran-2,6(3H)-dione, dihydro-...	142	C7H10O3	004160-82-1	37



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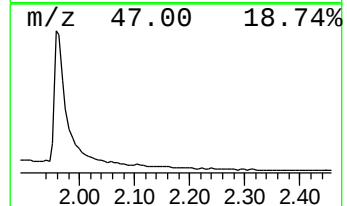
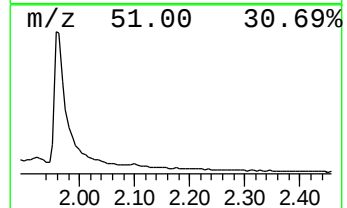
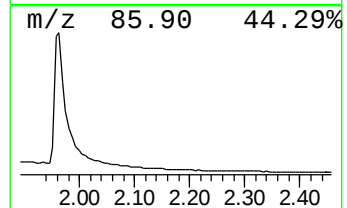
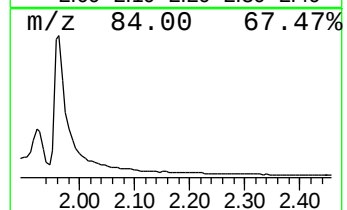
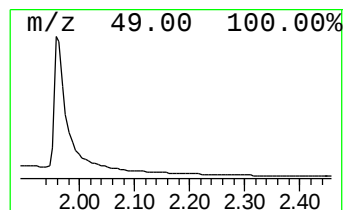
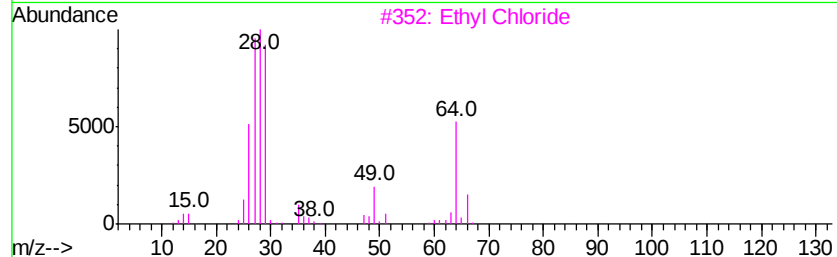
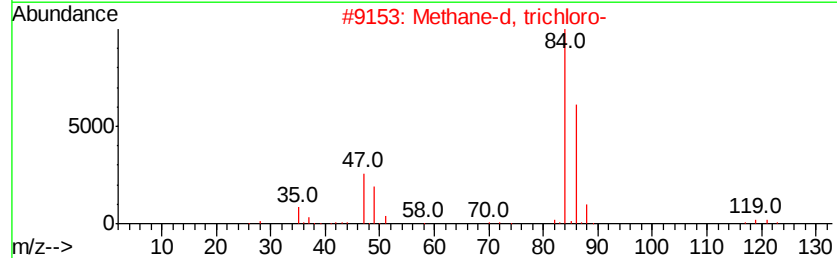
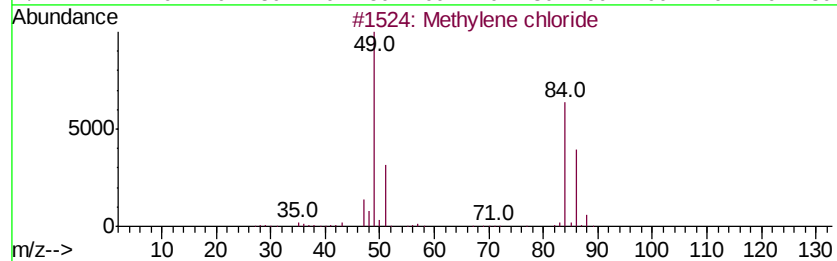
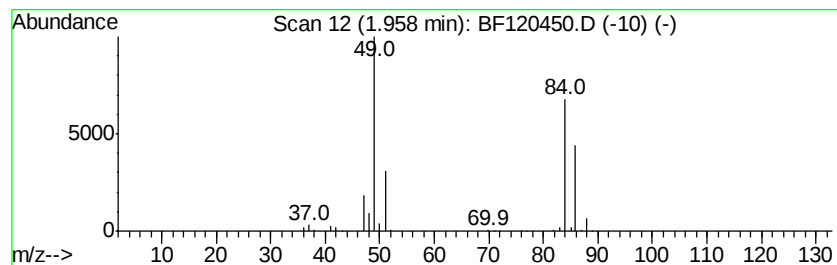
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Methylene chloride Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.96	34.90 ng	2722490	1,4-Dichlorobenzene-d4	6.84

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methylene chloride	84	CH2Cl2	000075-09-2	95
2		Methane-d, trichloro-	119	CDCl3	000865-49-6	10
3		Ethyl Chloride	64	C2H5Cl	000075-00-3	4
4		Acetic acid, chloro-, ethyl ester	122	C4H7ClO2	000105-39-5	2
5		2-Chloroethanol	80	C2H5ClO	000107-07-3	1



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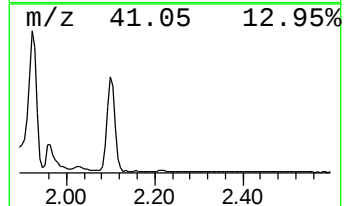
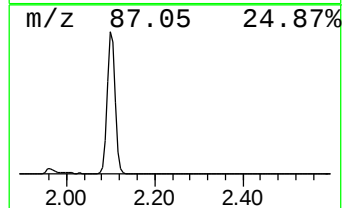
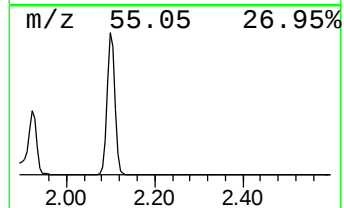
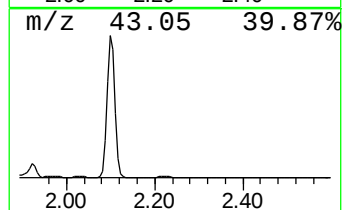
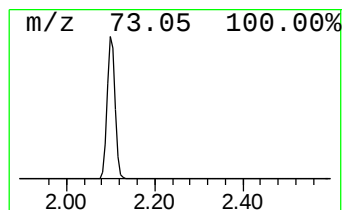
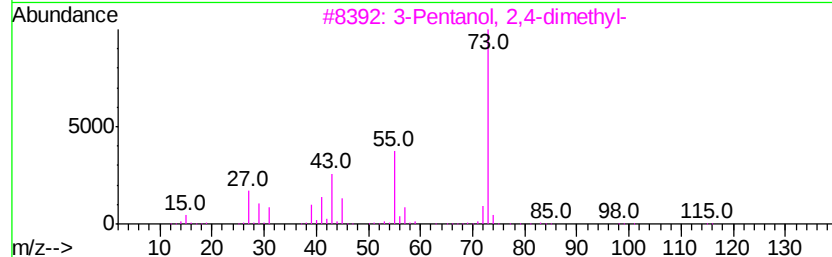
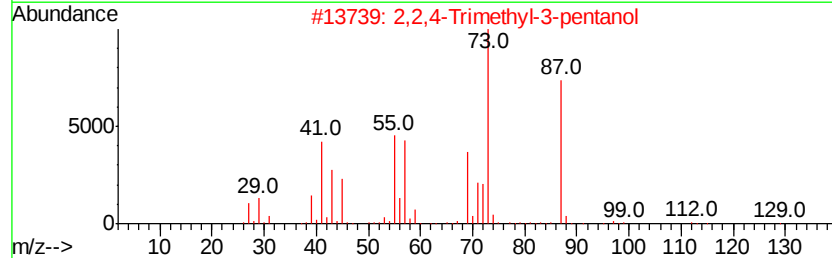
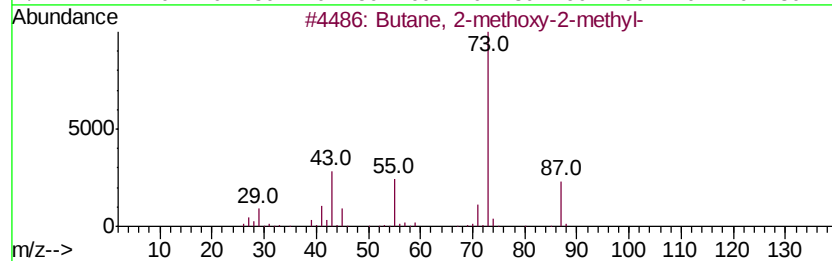
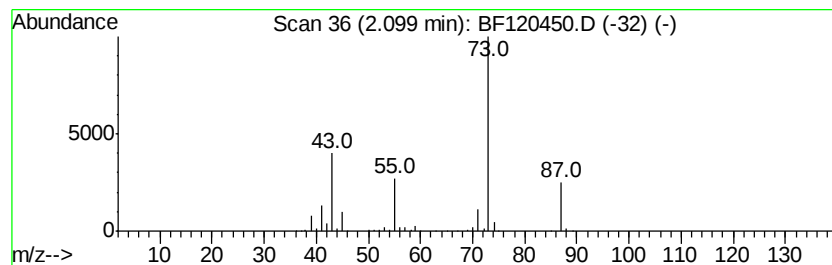
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Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.10	22.53 ng	1757140	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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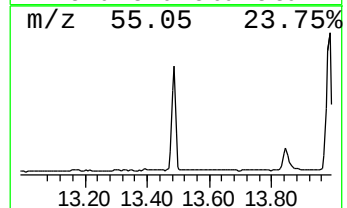
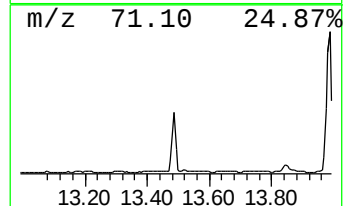
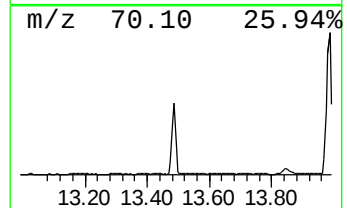
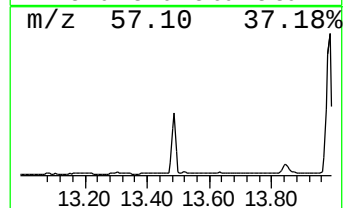
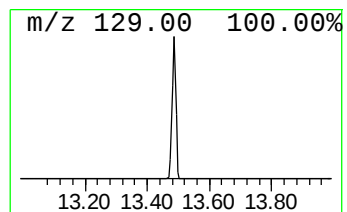
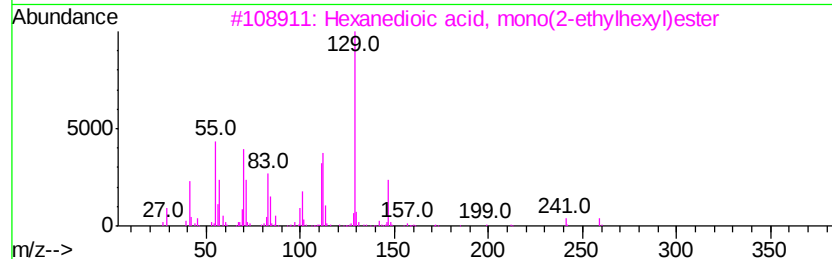
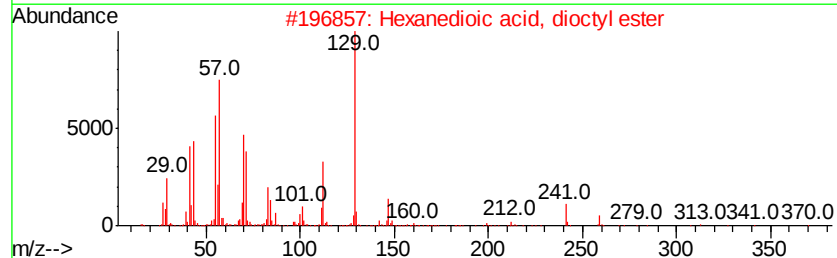
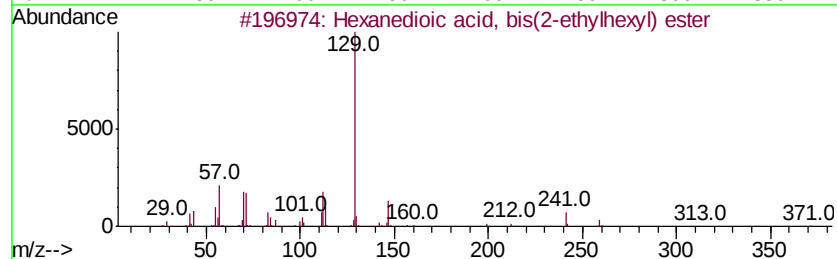
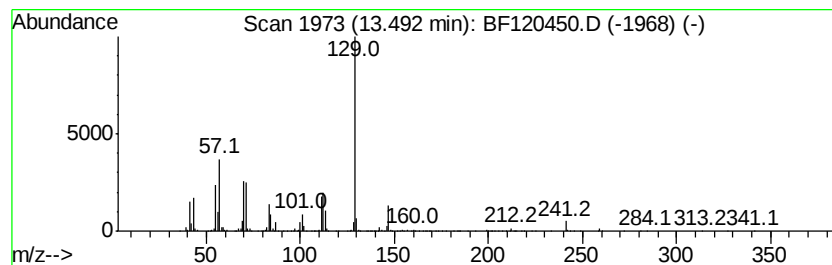
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Peak Number 5 Hexanedioic acid, bis(2-eth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	5.35 ng	2242290	Chrysene-d12	13.99

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
2		Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	90
3		Hexanedioic acid, mono(2-ethylhe...	258	C14H26O4	004337-65-9	68
4		Diisooctyl adipate	370	C22H42O4	001330-86-5	53
5		Adipic acid, 2-ethylhexyl isohex...	342	C20H38O4	1000324-48-3	53



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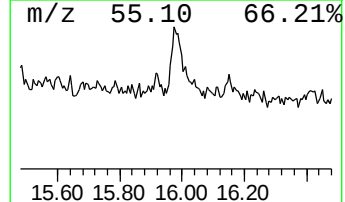
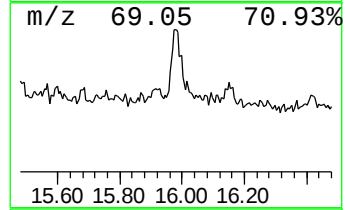
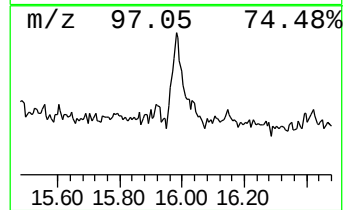
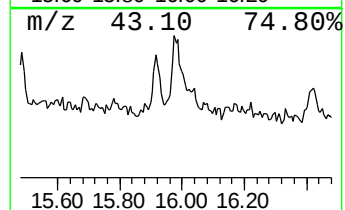
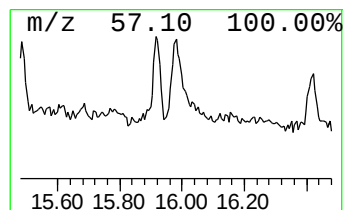
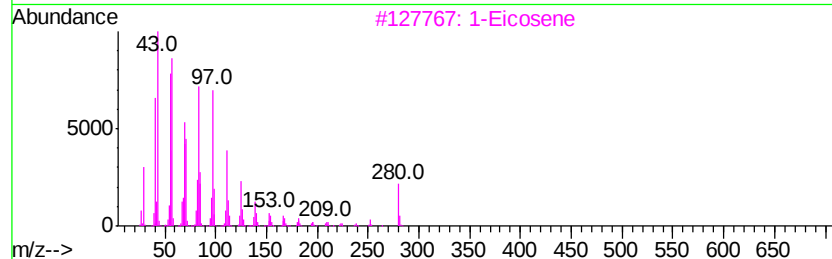
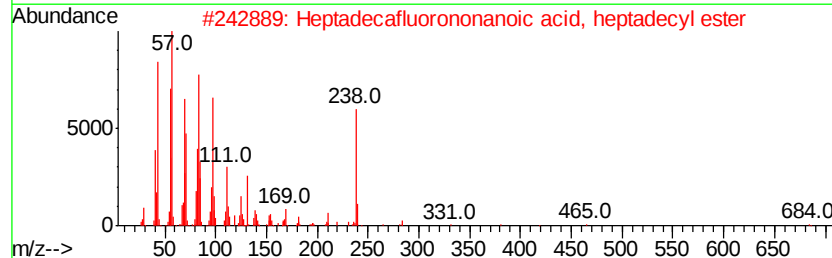
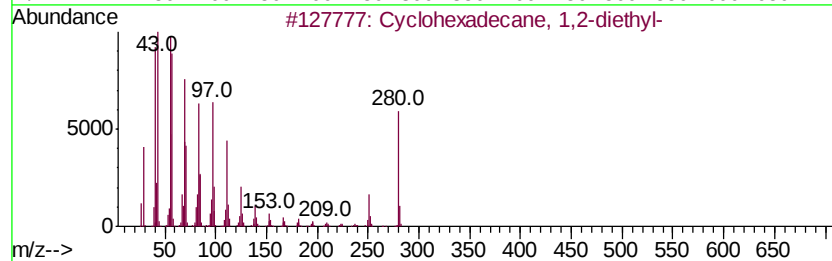
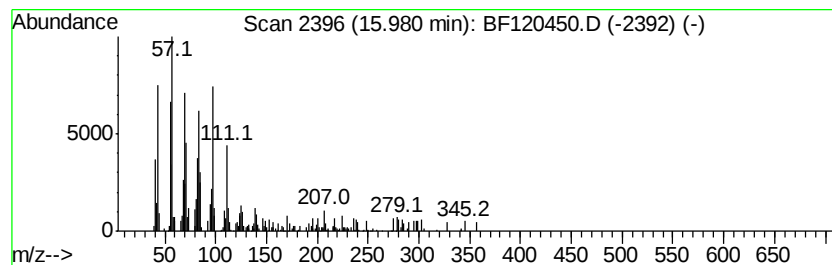
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Cyclohexadecane, 1,2-diethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.98	2.24 ng	157832	Perylene-d12	15.44

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexadecane, 1,2-diethyl-	280	C20H40	1000155-85-3	87
2		Heptadecafluorononanoic acid, he...	702	C26H35F17O2	1000356-03-3	87
3		1-Eicosene	280	C20H40	003452-07-1	87
4		Cycloeicosane	280	C20H40	000296-56-0	86
5		Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	83



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclohexane	1.92	10.0	ng	782279	1	6.84	1560100	20.0
Methylene chloride	1.96	34.9	ng	2722490	1	6.84	1560100	20.0
Butane, 2-methoxy...	2.10	22.5	ng	1757140	1	6.84	1560100	20.0
Hexanedioic acid,...	13.49	5.3	ng	2242290	5	13.99	8387900	20.0
Cyclohexadecane, ...	15.98	2.2	ng	157832	6	15.44	1411250	20.0