

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032515\
 Data File : BF077957.D
 Acq On : 25 Mar 2015 17:59
 Operator : TP/IZ
 Sample : G1612-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-2

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:\HPCHEM1\BNA_F\METHODS\8270-BF031115.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	1.355	38	41	44	rVB	156949	221128	13.00%	2.251%
2	1.515	52	55	58	rBV	190220	216299	12.71%	2.202%
3	1.618	62	64	70	rVB2	7695	19306	1.13%	0.197%
4	1.744	74	75	81	rBV	69225	100938	5.93%	1.028%
5	1.950	91	93	100	rVB	39188	57837	3.40%	0.589%
6	4.830	344	345	352	rVB	31183	44053	2.59%	0.448%
7	5.379	390	393	399	rVB	328769	415062	24.39%	4.225%
8	6.670	504	506	509	rBV	282565	261091	15.34%	2.658%
9	6.807	515	518	520	rBV	851025	817016	48.02%	8.317%
10	7.093	540	543	546	rBV	252509	221481	13.02%	2.255%
11	7.276	557	559	562	rBV	735306	752613	44.23%	7.662%
12	7.790	602	604	606	rBV	703196	620595	36.47%	6.318%
13	8.670	679	681	684	rBV	354860	305680	17.96%	3.112%
14	10.019	797	799	801	rBV	1535540	1338892	78.69%	13.630%
15	10.773	861	865	869	rBV2	12577	22195	1.30%	0.226%
16	10.842	869	871	873	rVB	361675	360205	21.17%	3.667%
17	11.813	954	956	959	rBV2	671588	801174	47.08%	8.156%
18	12.671	1029	1031	1034	rBV	307843	388195	22.81%	3.952%
19	14.671	1204	1206	1209	rBV	1276901	1701580	100.00%	17.323%
20	15.905	1311	1314	1321	rBV2	21211	41322	2.43%	0.421%
21	16.008	1321	1323	1326	rBV	382803	457457	26.88%	4.657%
22	17.917	1485	1490	1493	rBV	320770	509315	29.93%	5.185%
23	18.431	1532	1535	1538	rBV	27152	54122	3.18%	0.551%
24	18.980	1580	1583	1586	rBV2	23154	51720	3.04%	0.527%
25	19.631	1638	1640	1650	rVB5	16743	43650	2.57%	0.444%

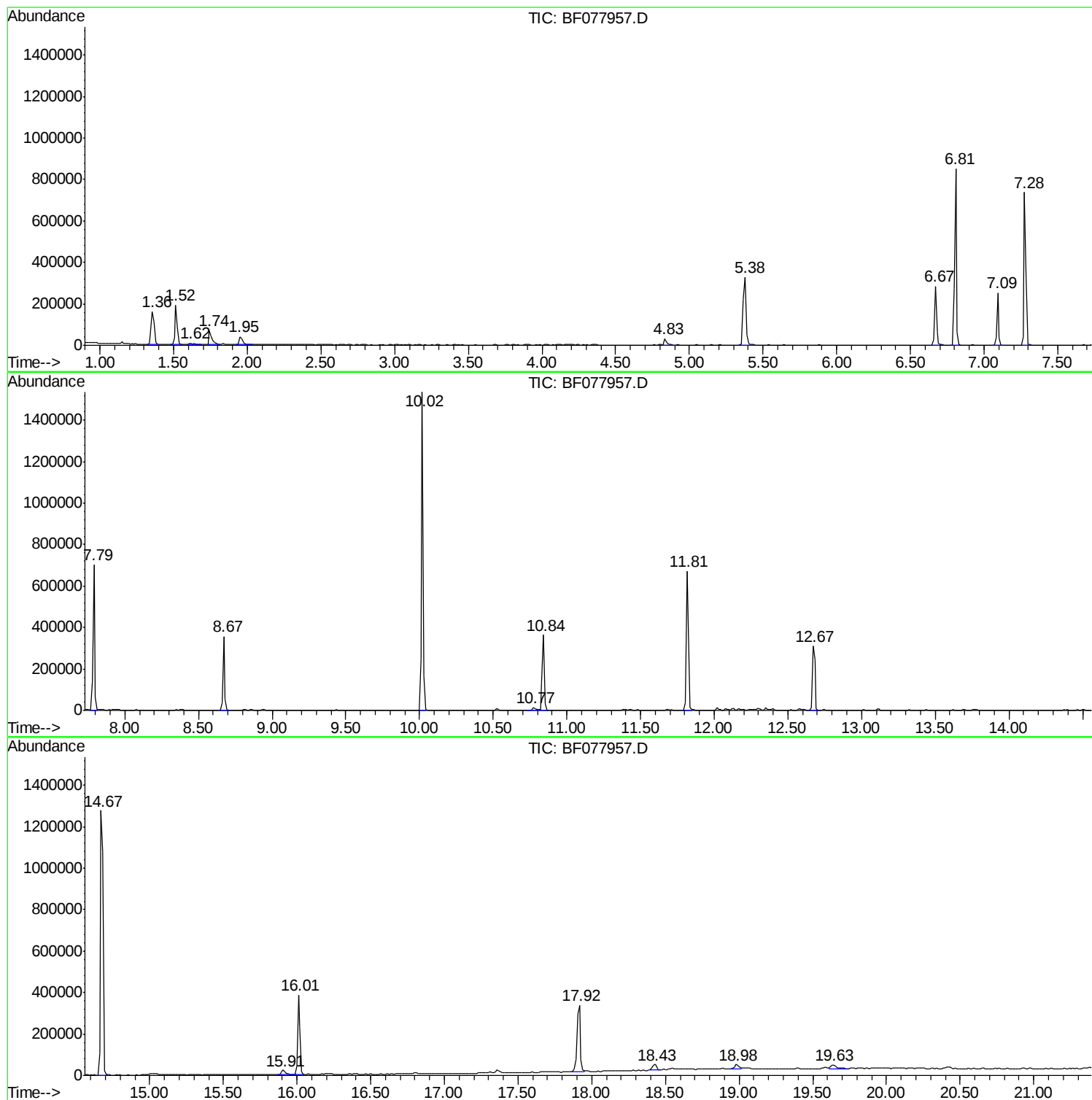
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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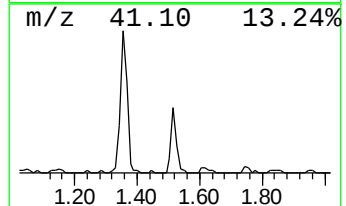
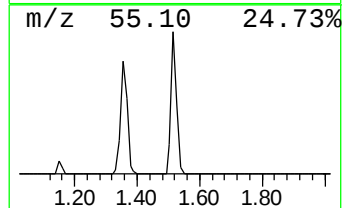
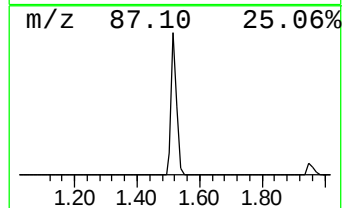
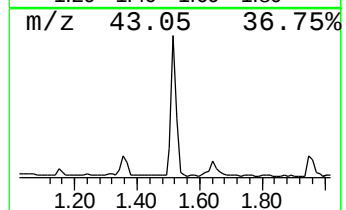
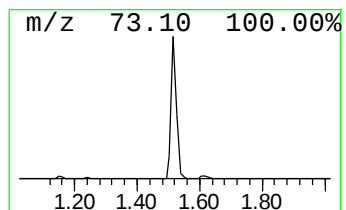
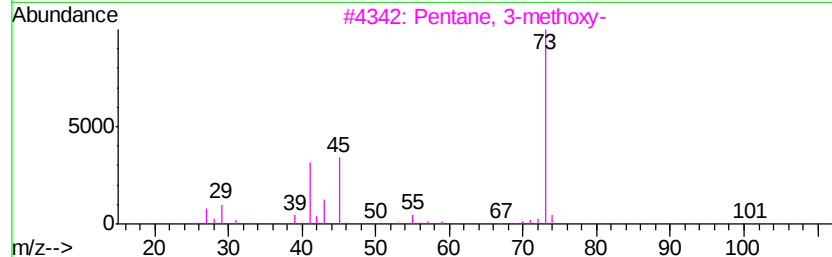
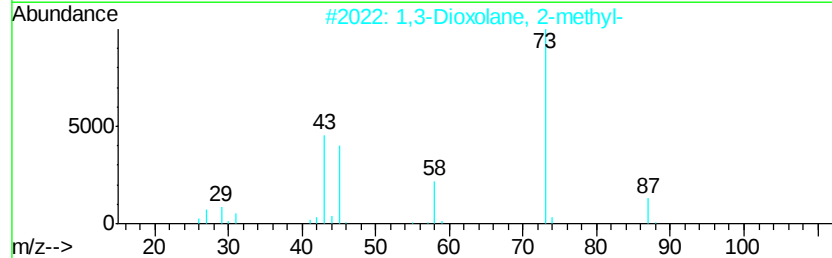
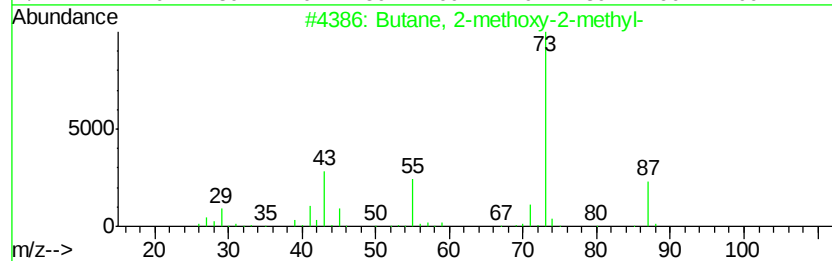
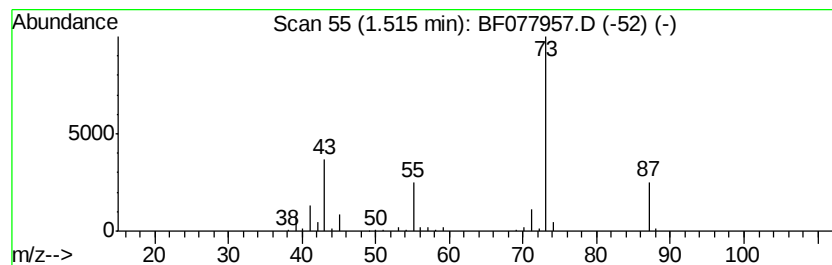
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.52	19.53 ng	216299	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	10



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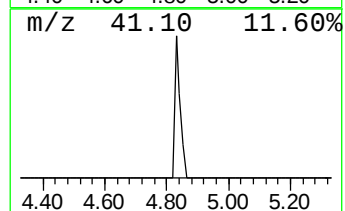
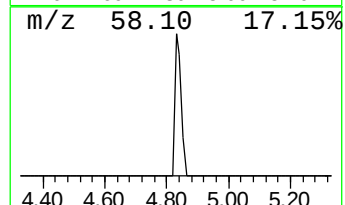
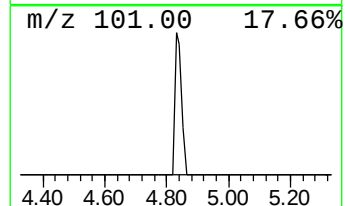
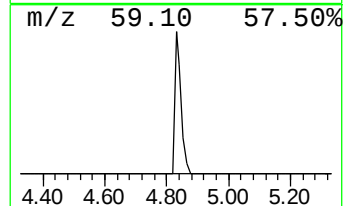
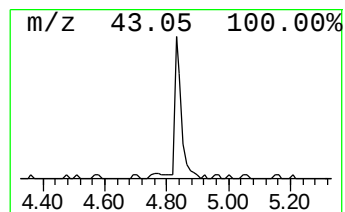
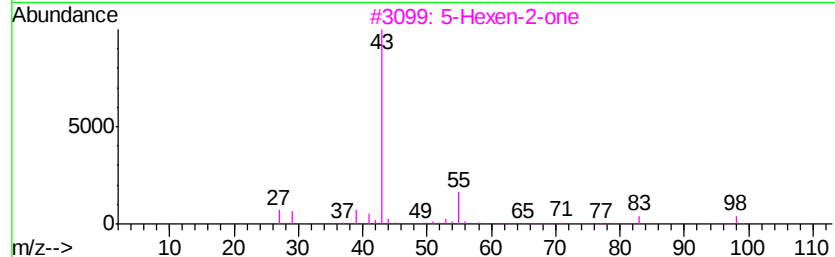
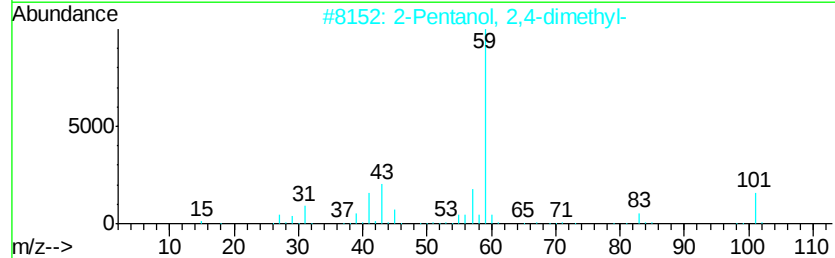
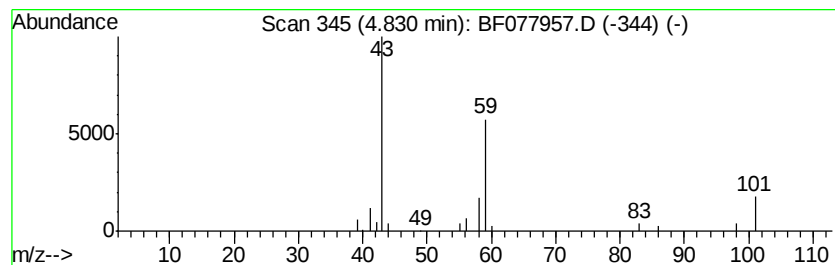
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	3.98 ng	44053	1,4-Dichlorobenzene-d4	7.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	50
3	5-Hexen-2-one	98	C6H10O	000109-49-9	9
4	1-Butanol, 3-methoxy-	104	C5H12O2	002517-43-3	9
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9



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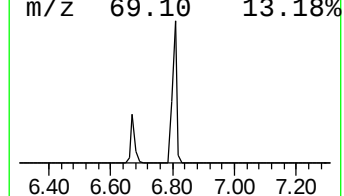
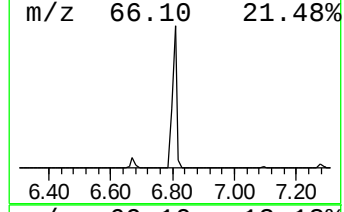
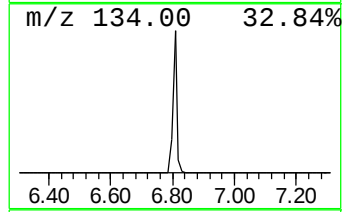
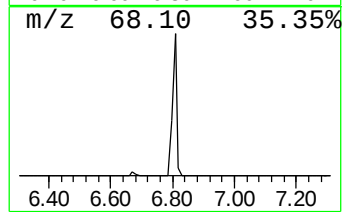
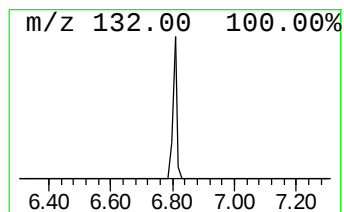
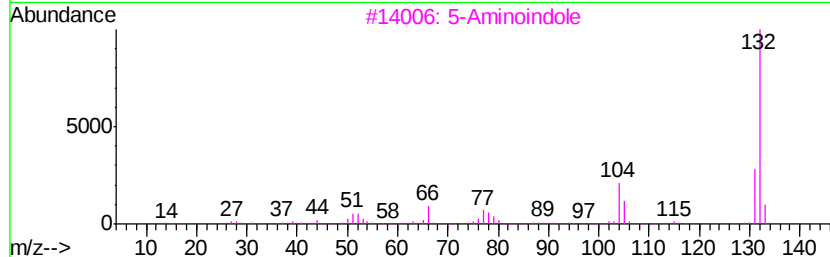
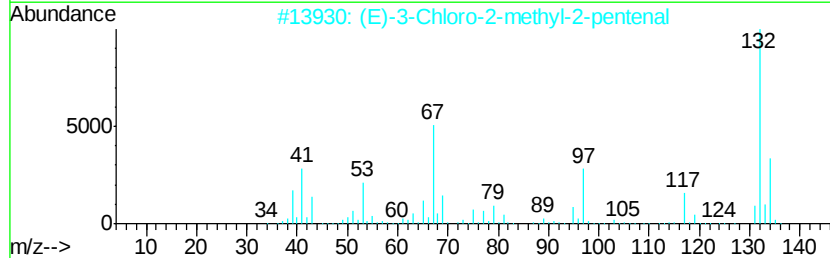
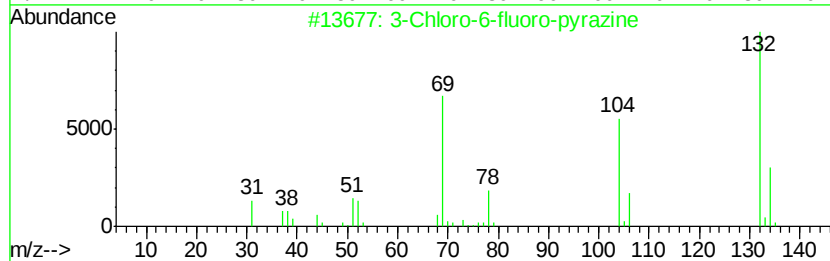
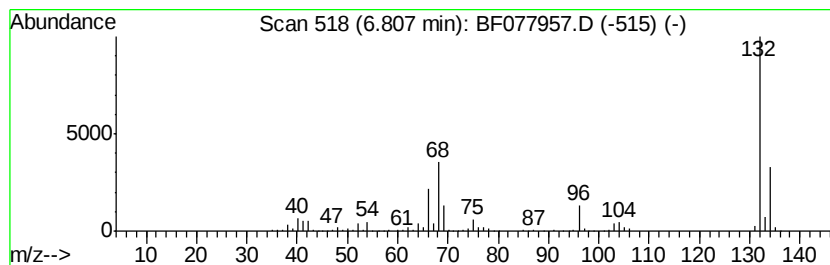
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Peak Number 5 unknown6.81 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.81	73.78 ng	817016	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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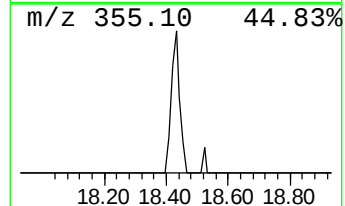
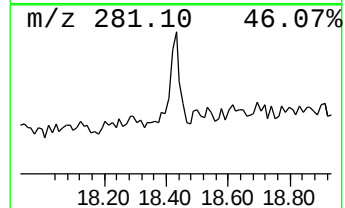
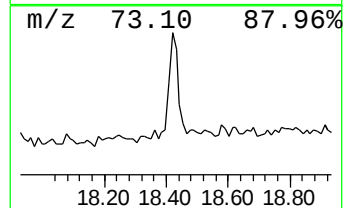
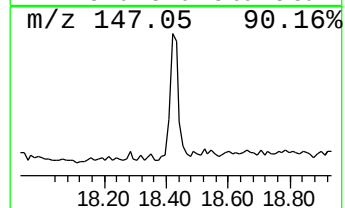
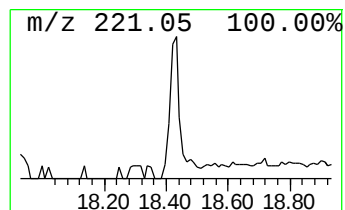
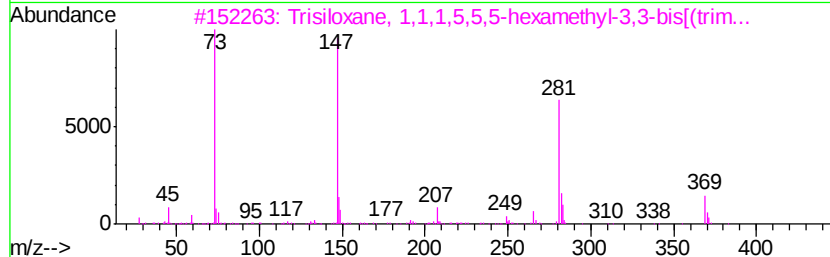
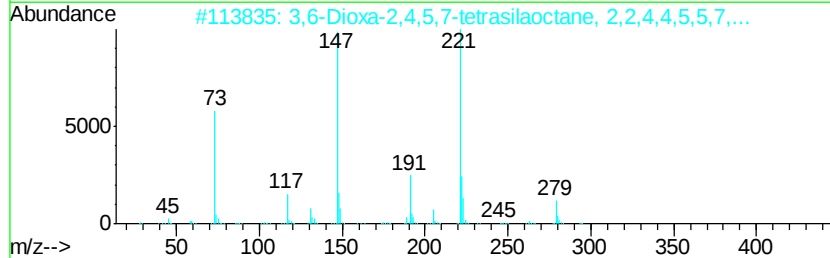
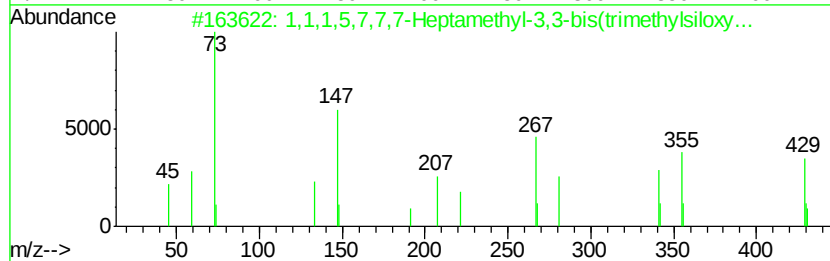
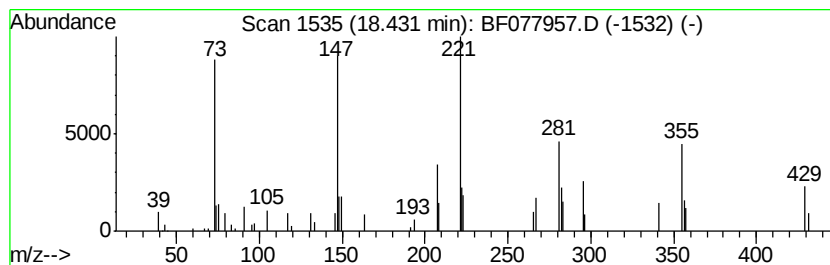
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 7 unknown18.43 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.43	2.13 ng	54122	Perylene-d12	17.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	37
2		3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	32
3		Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	12
4		2,6-Dimethyl-3,4-bis(trimethylsi...	311	C15H29N02Si2	1000079-52-1	12
5		1,2-Benzenedicarboxylic acid, bi...	310	C14H22O4Si2	002078-22-0	10



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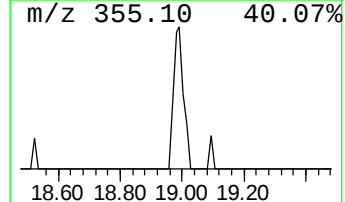
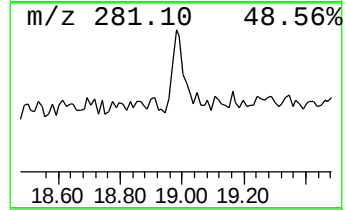
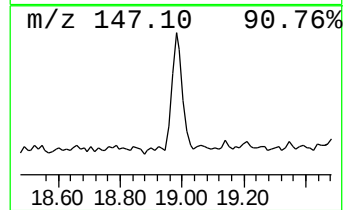
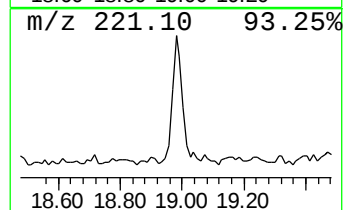
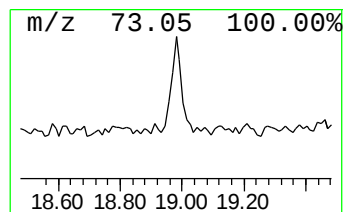
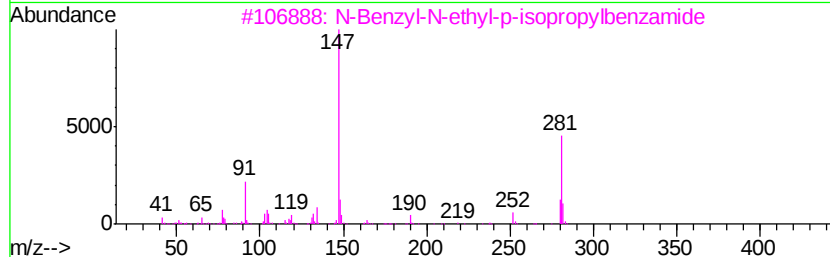
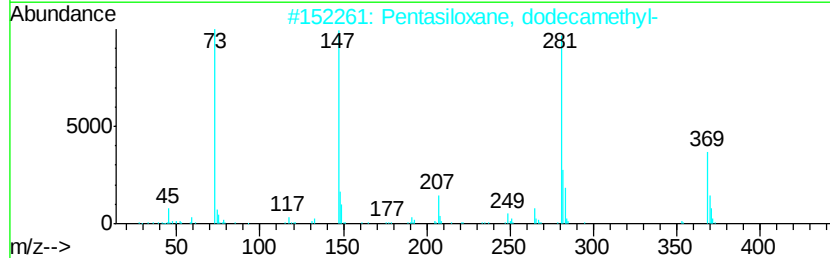
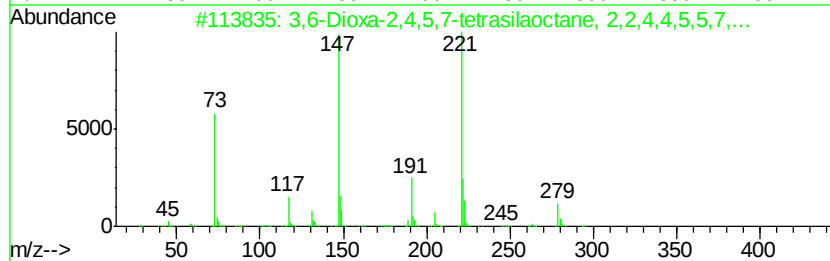
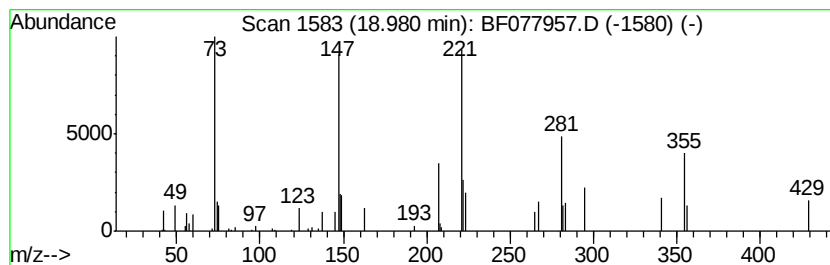
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Peak Number 8 unknown18.98 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.98	2.03 ng	51720	Perylene-d12	17.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	37
2		Pentasiloxane, dodecamethyl-	384	C12H36O4Si5	000141-63-9	27
3		N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	16
4		Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	16
5		Piperidine, 1-(5-trifluoromethyl...	295	C15H16F3N3	1000268-74-7	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	1.52	19.5	ng	216299	1	7.09	221481	20.0
2-Pentanone, 4-hy...	4.83	4.0	ng	44053	1	7.09	221481	20.0
unknown6.81	6.81	73.8	ng	817016	1	7.09	221481	20.0
unknown18.43	18.43	2.1	ng	54122	6	17.92	509315	20.0
unknown18.98	18.98	2.0	ng	51720	6	17.92	509315	20.0