

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070119\
 Data File : BG041561.D
 Acq On : 1 Jul 2019 16:25
 Operator : HP/JU
 Sample : K3555-25
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GWMH-13

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_G\METHODS\8270-BG062619.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	4.562	246	251	260	rVB2	55171	83186	2.90%	0.702%
2	5.567	417	422	433	rBV	248853	429721	15.00%	3.627%
3	7.077	675	679	685	rVB2	35361	49617	1.73%	0.419%
4	7.189	692	698	714	rVB	176461	327912	11.45%	2.768%
5	7.559	754	761	774	rVB	449596	762733	26.62%	6.438%
6	8.029	835	841	850	rBV	116171	206050	7.19%	1.739%
7	8.352	889	896	908	rBV	412912	717377	25.04%	6.055%
8	9.210	1035	1042	1051	rBV	312912	593518	20.72%	5.010%
9	10.855	1315	1322	1330	rBV	130044	245451	8.57%	2.072%
10	13.293	1730	1737	1748	rBV	944940	1441008	50.30%	12.163%
11	14.674	1966	1972	1989	rVB2	236232	397868	13.89%	3.358%
12	16.166	2220	2226	2243	rBV	756986	1226096	42.79%	10.349%
13	17.424	2433	2440	2452	rBV2	358283	555143	19.38%	4.686%
14	18.282	2578	2586	2596	rBV6	20138	38547	1.35%	0.325%
15	20.027	2877	2883	2893	rBV	2248807	2865111	100.00%	24.183%
16	20.291	2924	2928	2933	rBV5	27087	30981	1.08%	0.261%
17	20.785	3009	3012	3017	rVB3	28220	32510	1.13%	0.274%
18	21.290	3094	3098	3104	rBV2	39522	60397	2.11%	0.510%
19	21.695	3161	3167	3180	rVB	436012	741701	25.89%	6.260%
20	21.836	3186	3191	3195	rVB	44444	62885	2.19%	0.531%
21	22.447	3289	3295	3303	rVB	40013	66905	2.34%	0.565%
22	23.141	3407	3413	3420	rBV3	45220	91909	3.21%	0.776%
23	23.945	3544	3550	3557	rVB2	31502	67181	2.34%	0.567%
24	24.897	3706	3712	3717	rBV3	23473	54765	1.91%	0.462%
25	24.980	3717	3726	3739	rVB2	232525	699030	24.40%	5.900%

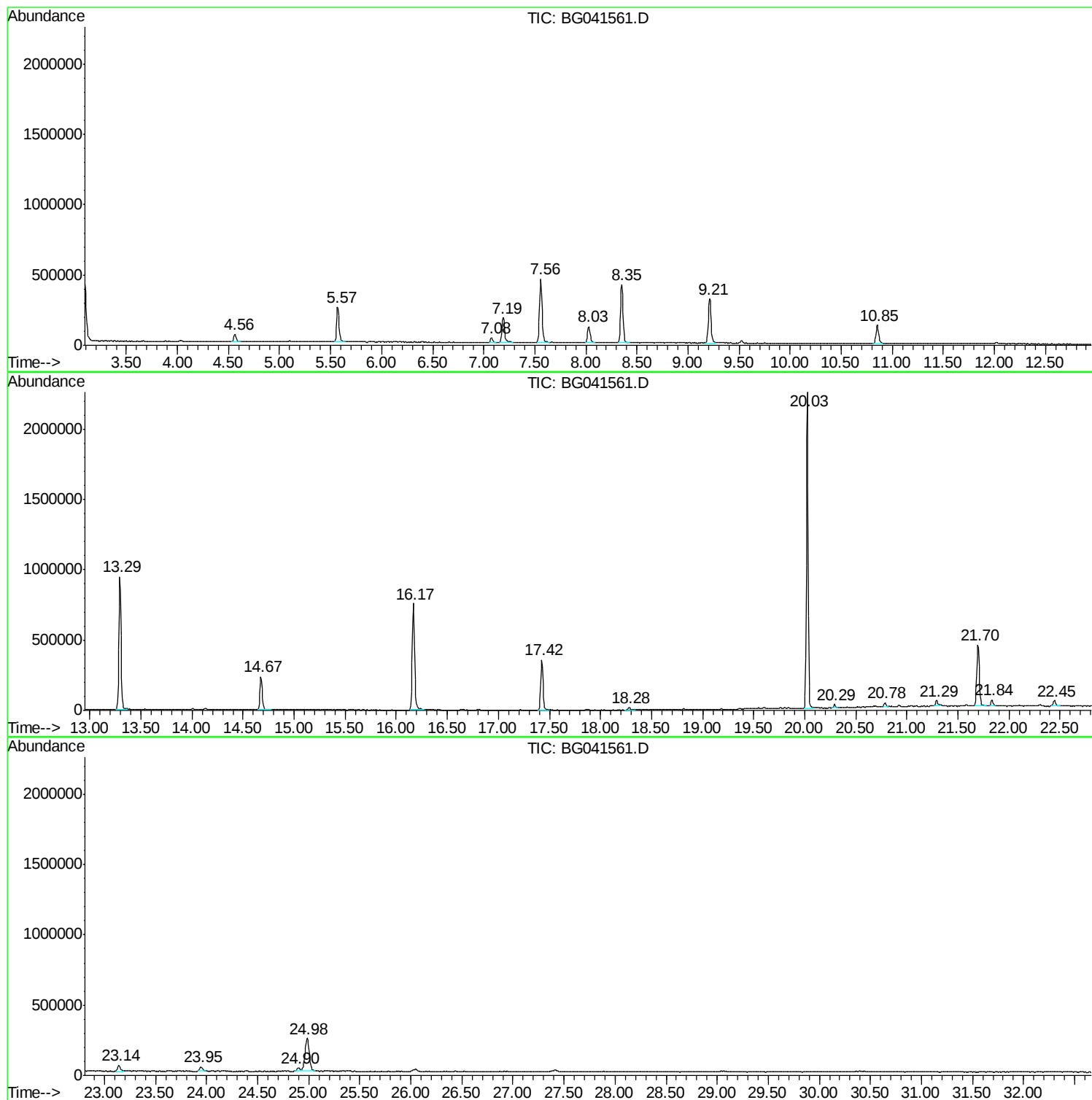
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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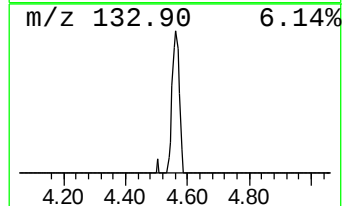
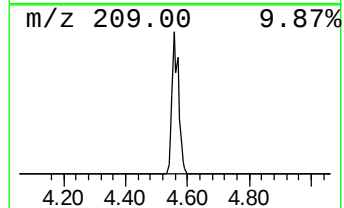
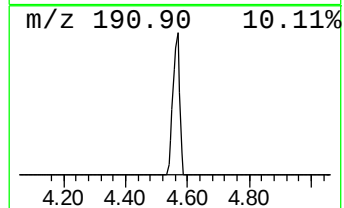
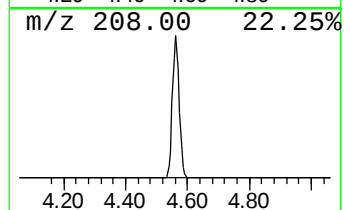
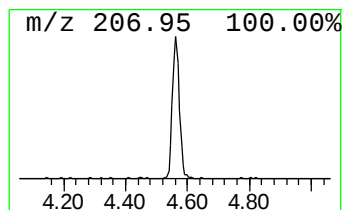
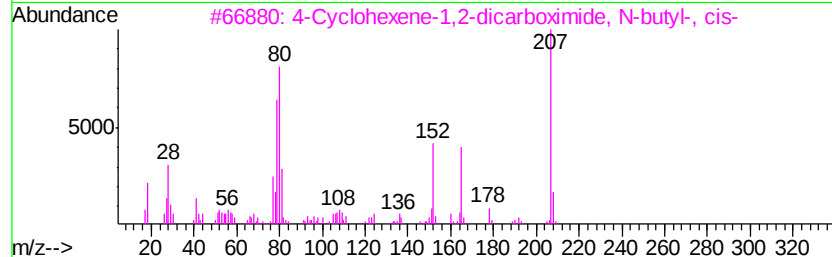
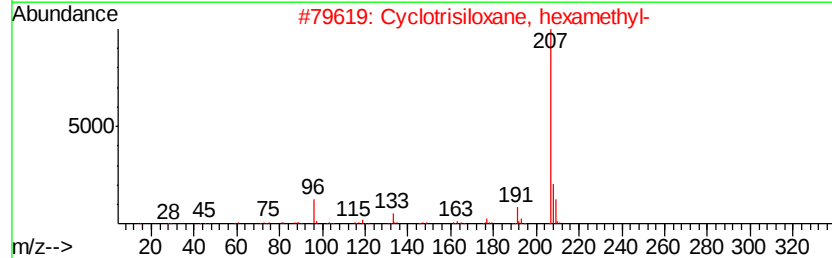
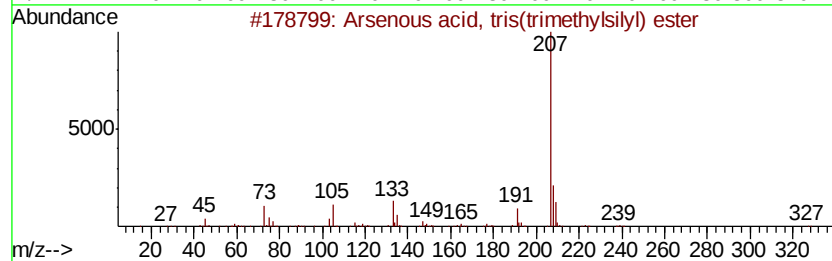
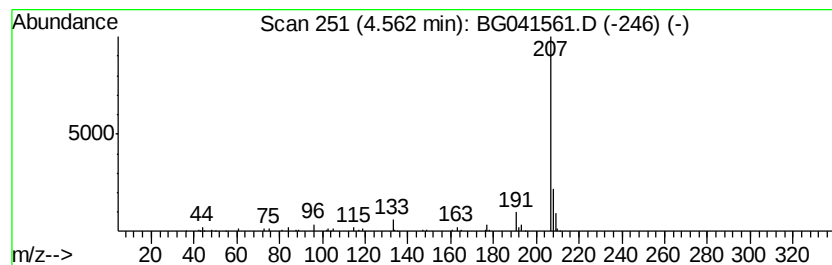
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG062619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Arsenous acid, tris(trimeth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.56	8.07 ng	83186	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Arsenous acid, tris(trimethylsil...	342	C9H27AsO3Si3	055429-29-3	78
2		Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	72
3		4-Cyclohexene-1,2-dicarboximide,...	207	C12H17N02	028916-00-9	42
4		Benzo[h]quinoline, 2,4-dimethyl-	207	C15H13N	000605-67-4	39
5		1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	39



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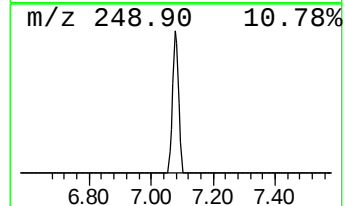
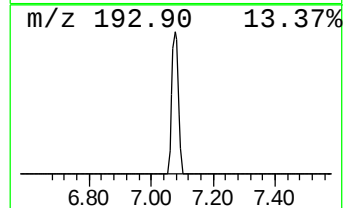
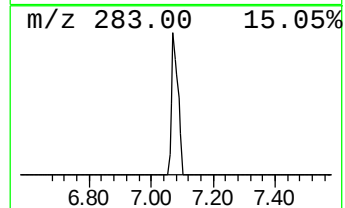
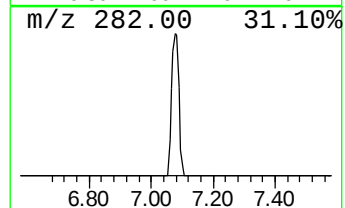
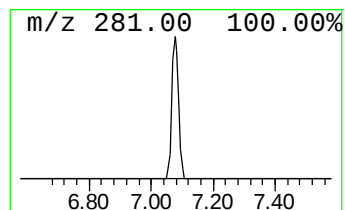
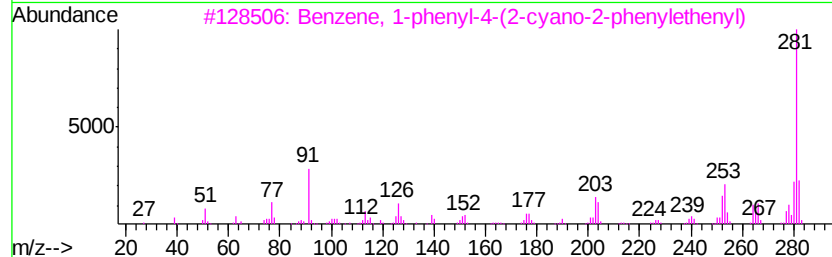
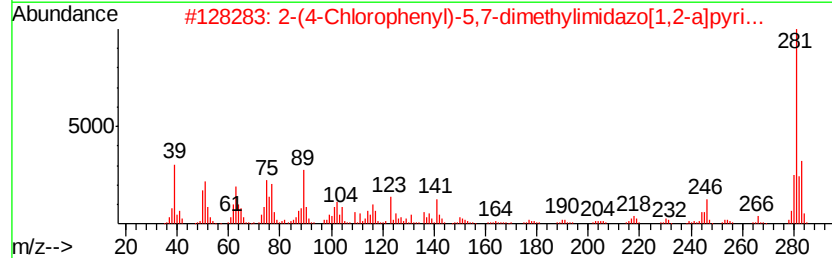
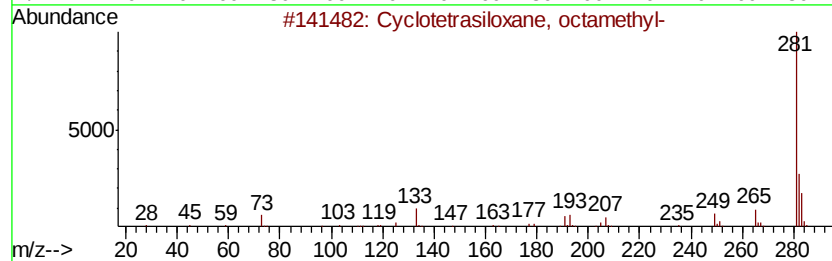
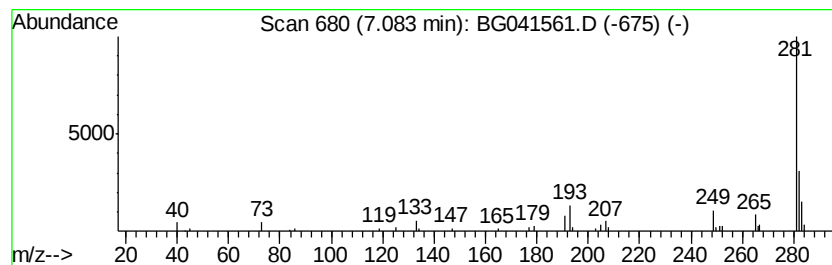
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Peak Number 2 Cyclotetrasiloxane, octamet... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.08	4.82 ng	49617	1,4-Dichlorobenzene-d4	8.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	91
2			2-(4-Chlorophenyl)-5,7-dimethyl-1H-imidazo[1,2-a]pyridine	281	C16H12ClN3	1000305-50-9	50
3			Benzene, 1-phenyl-4-(2-cyano-2-phenylethenyl)-	281	C21H15N	027869-56-3	50
4			6-Chloro-3-ethyl-2-methyl-4-phenyl-1,2,4-triazole-3-thiol, 4-allyl-	281	C18H16ClN	022609-09-2	50
5			4H-1,2,4-Triazole-3-thiol, 4-allyl-	281	C16H15N3S	031803-13-1	47



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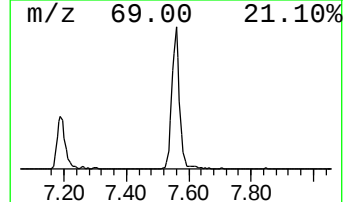
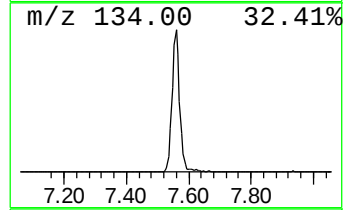
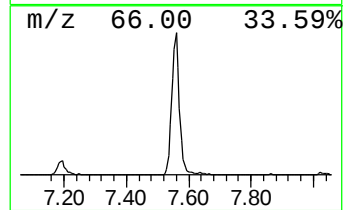
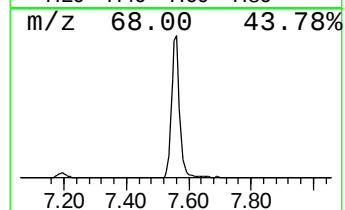
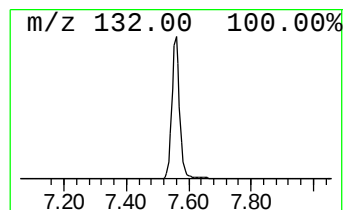
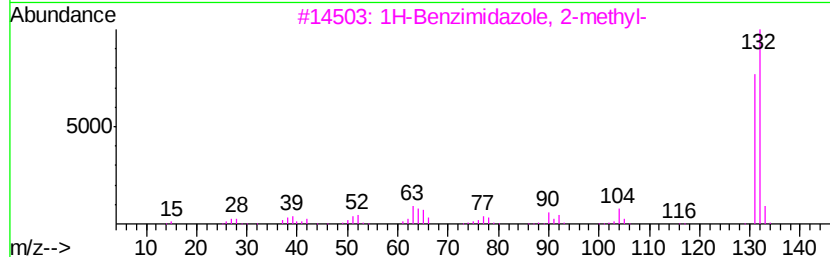
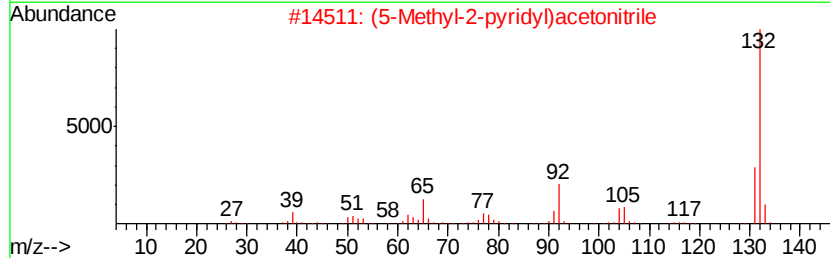
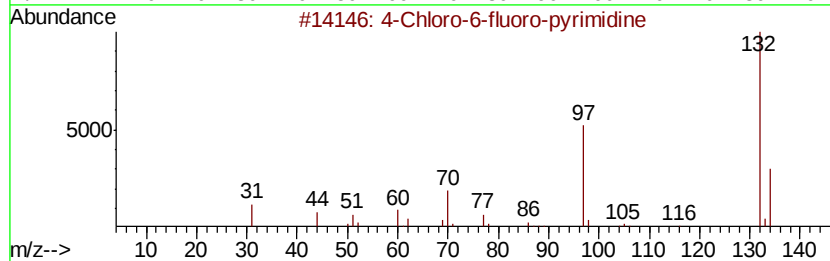
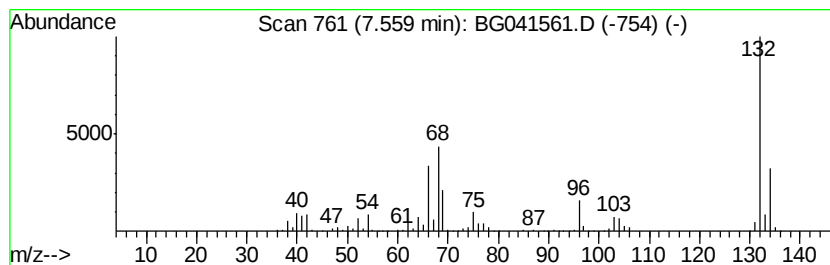
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Peak Number 3 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	74.03 ng	762733	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	22
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Aminoindole	132	C8H8N2	005192-03-0	22
5		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	22



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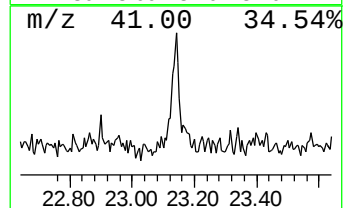
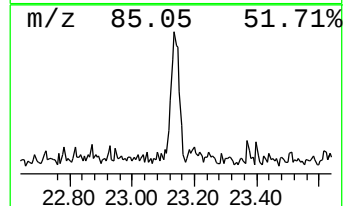
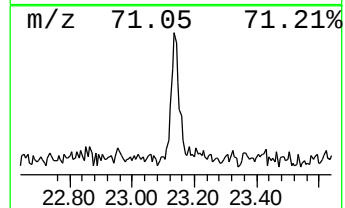
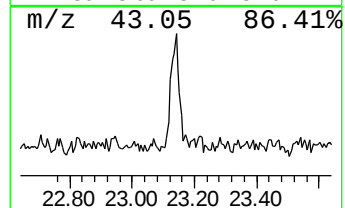
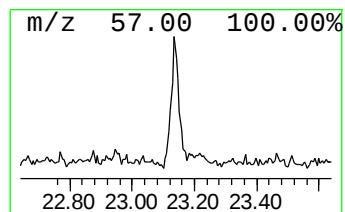
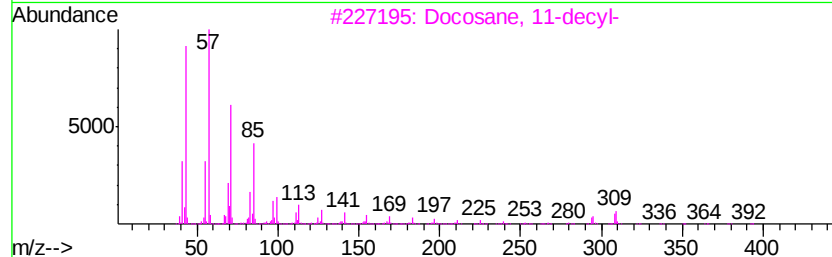
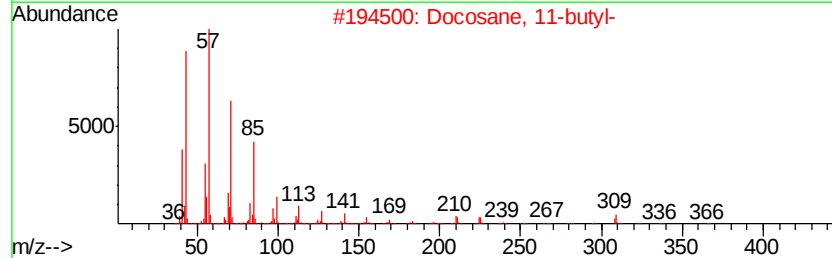
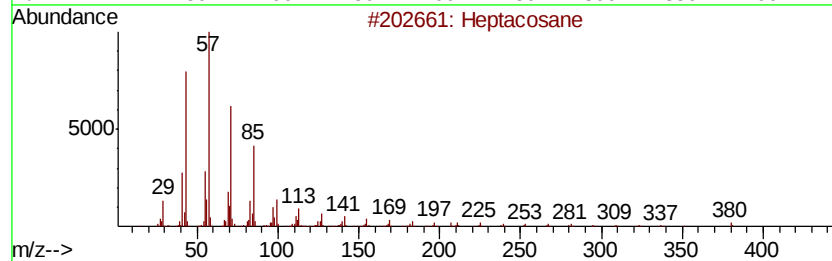
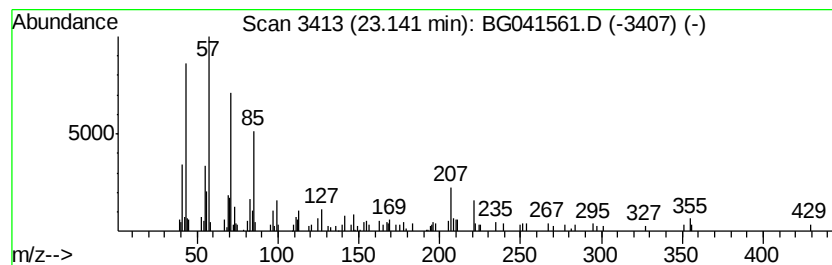
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Peak Number 5 Heptacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.14	2.48 ng	91909	Chrysene-d12	21.70

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	60
2	Docosane, 11-butyl-		366	C26H54	013475-76-8	60
3	Docosane, 11-decyl-		451	C32H66	055401-55-3	60
4	Pentadecane, 8-hexyl-		296	C21H44	013475-75-7	53
5	Hexacosane		366	C26H54	000630-01-3	53



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Arsenous acid, tr...	4.56	8.1 ng		83186	1	8.03	206050	20.0
Cyclotetrasiloxan...	7.08	4.8 ng		49617	1	8.03	206050	20.0
unknown7.56	7.56	74.0 ng		762733	1	8.03	206050	20.0
Heptacosane	23.14	2.5 ng		91909	5	21.70	741701	20.0