

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080720\
 Data File : BG046193.D
 Acq On : 7 Aug 2020 14:54
 Operator : CG/JU
 Sample : L3594-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-2

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-BG073020.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	3.155	7	11	27	rVB	481881	748534	9.63%	2.162%
2	5.535	409	416	429	rBV	1652816	2622024	33.74%	7.572%
3	7.121	677	686	697	rBV	1589495	2730649	35.14%	7.886%
4	7.949	820	827	835	rVB	332321	559166	7.20%	1.615%
5	9.101	1013	1023	1033	rBV	1024318	1854458	23.86%	5.355%
6	10.746	1295	1303	1310	rBV	450079	780717	10.05%	2.255%
7	13.202	1714	1721	1728	rBV	2873374	4468972	57.51%	12.906%
8	14.030	1855	1862	1872	rBV	359564	542492	6.98%	1.567%
9	14.577	1948	1955	1965	rBV	740144	1111867	14.31%	3.211%
10	16.063	2200	2208	2229	rBV	2646747	4334658	55.78%	12.518%
11	17.315	2415	2421	2433	rBV	944537	1455668	18.73%	4.204%
12	18.989	2701	2706	2710	rBV	292584	363622	4.68%	1.050%
13	19.395	2770	2775	2780	rBV	63496	84920	1.09%	0.245%
14	19.935	2860	2867	2873	rBV	5916980	7771035	100.00%	22.442%
15	20.840	3017	3021	3027	rBV	164160	200085	2.57%	0.578%
16	21.234	3083	3088	3096	rBV2	495634	699523	9.00%	2.020%
17	21.304	3096	3100	3105	rVB	75019	111601	1.44%	0.322%
18	21.563	3138	3144	3155	rVB	1141614	1794764	23.10%	5.183%
19	21.780	3177	3181	3187	rVB2	51298	79121	1.02%	0.228%
20	22.380	3278	3283	3291	rVB4	91965	185832	2.39%	0.537%
21	23.043	3392	3396	3402	rBV	80592	129401	1.67%	0.374%
22	23.825	3524	3529	3535	rBV2	91231	166503	2.14%	0.481%
23	24.671	3663	3673	3681	rBV2	673755	1832193	23.58%	5.291%

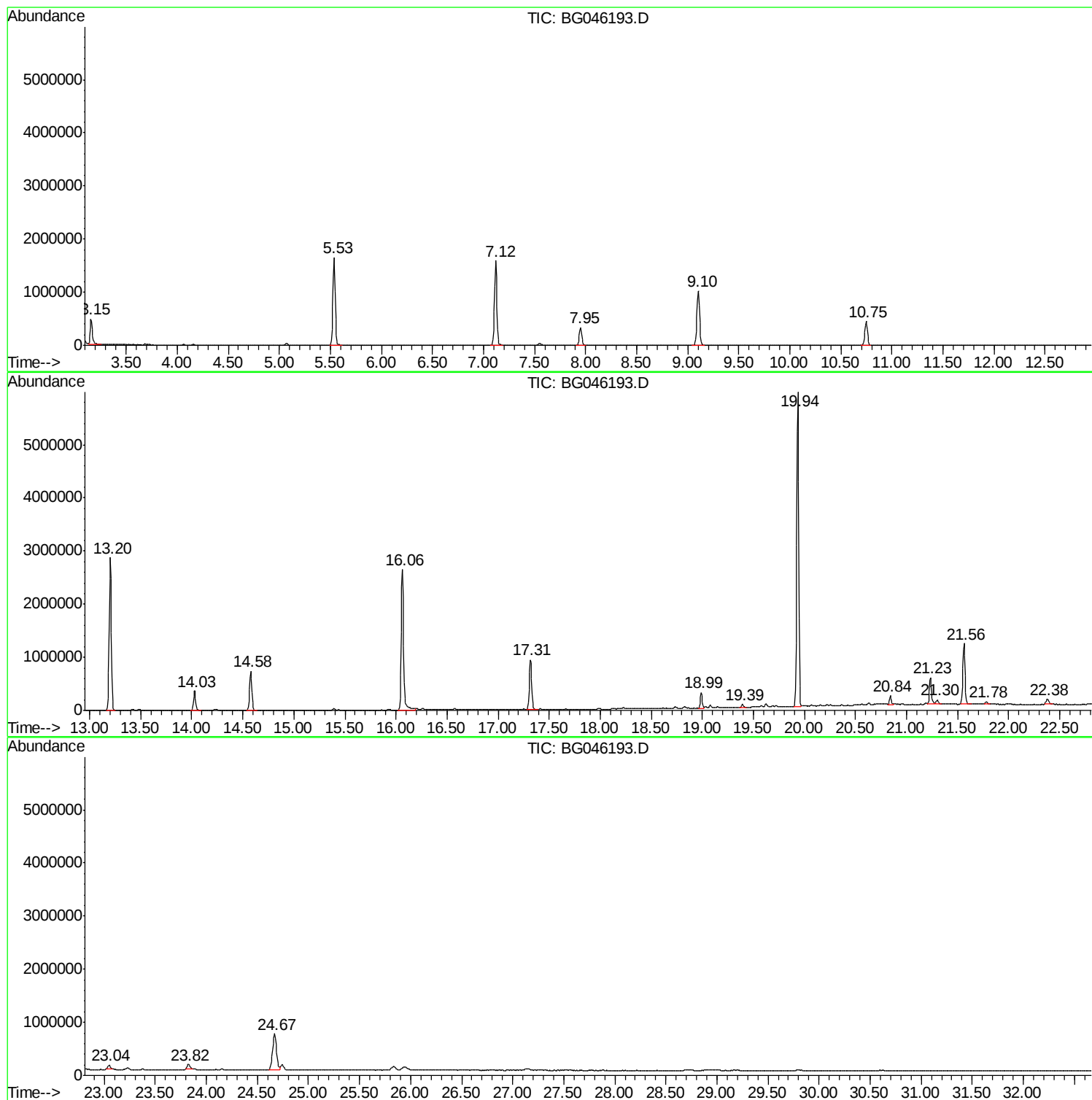
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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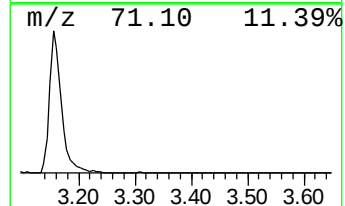
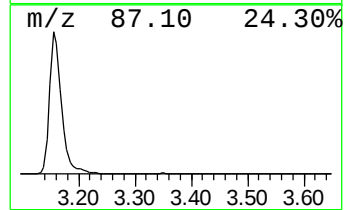
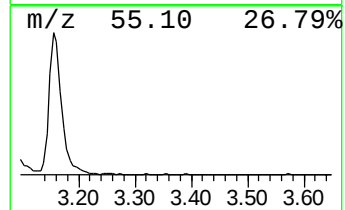
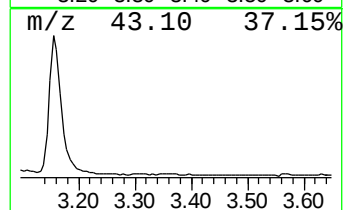
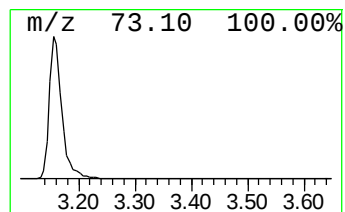
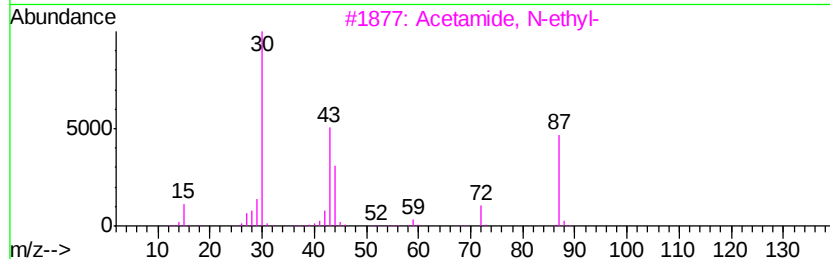
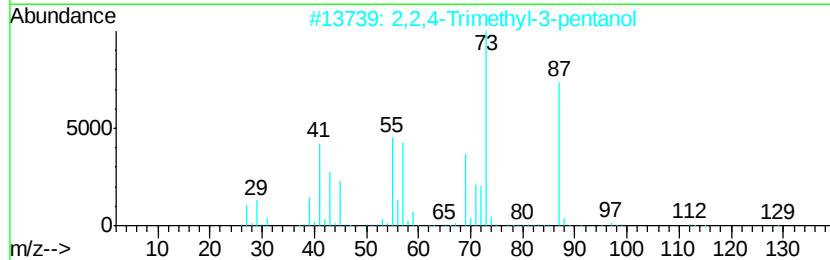
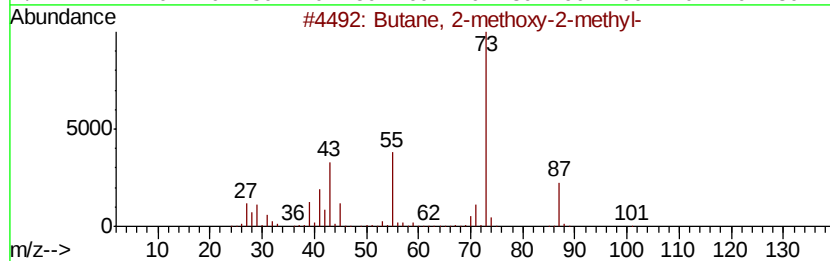
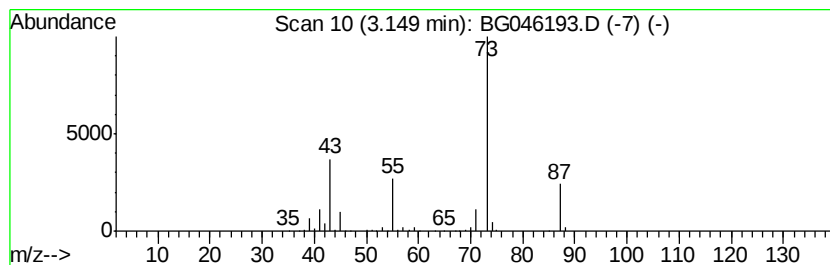
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	26.77 ng	748534	1,4-Dichlorobenzene-d4	7.95

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		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23



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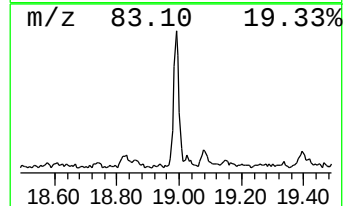
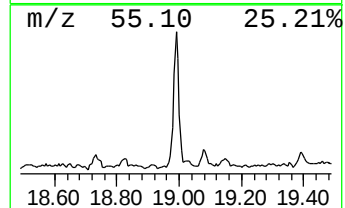
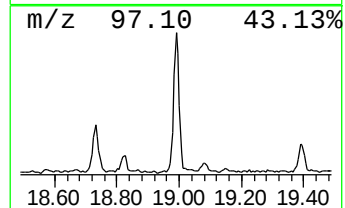
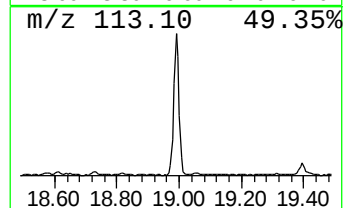
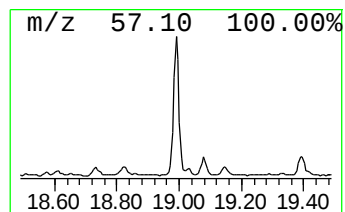
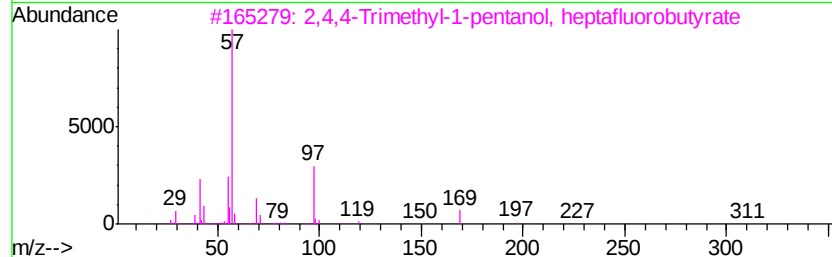
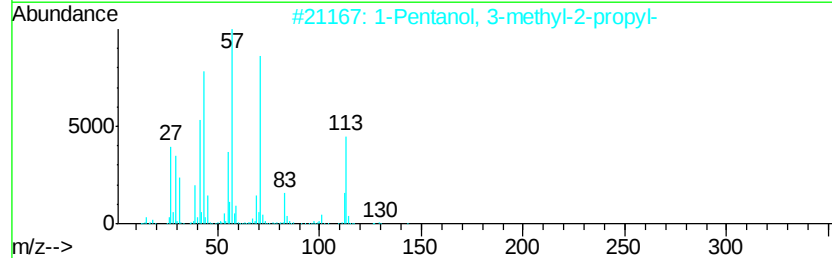
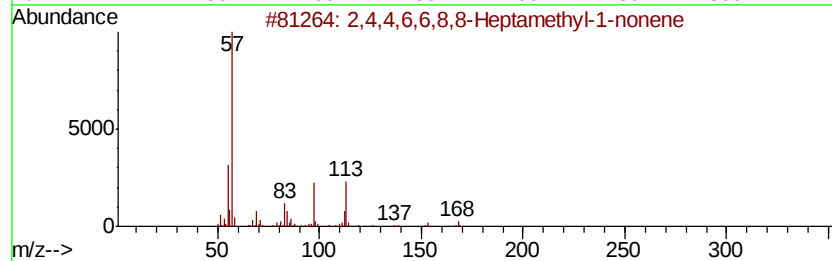
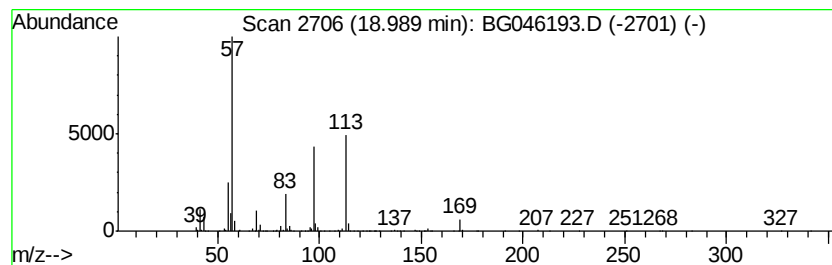
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 Peak Number 2 unknown18.99 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.99	5.00 ng	363622	Phenanthrene-d10	17.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	43
2		1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	33
3		2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1000365-01-4	32
4		2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	27
5		Octane, 2-bromo-	192	C8H17Br	000557-35-7	25



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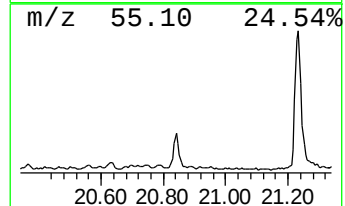
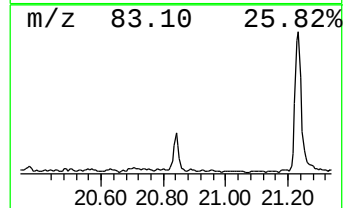
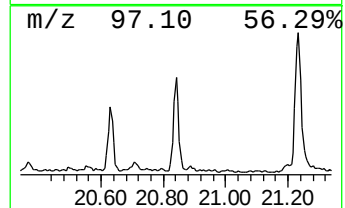
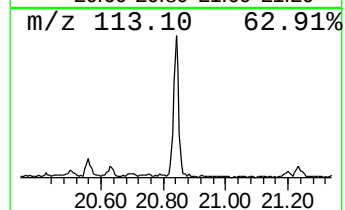
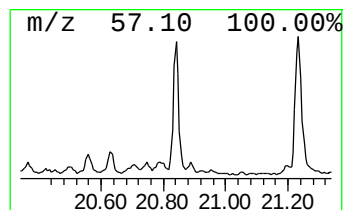
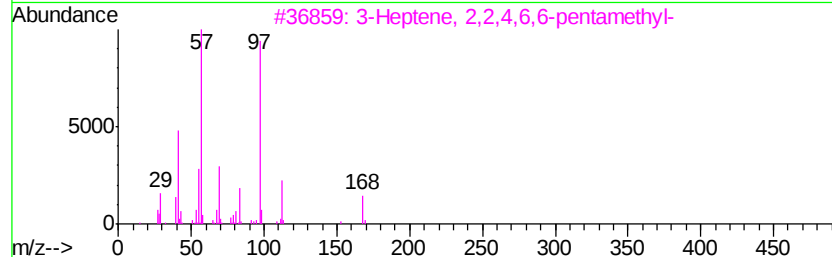
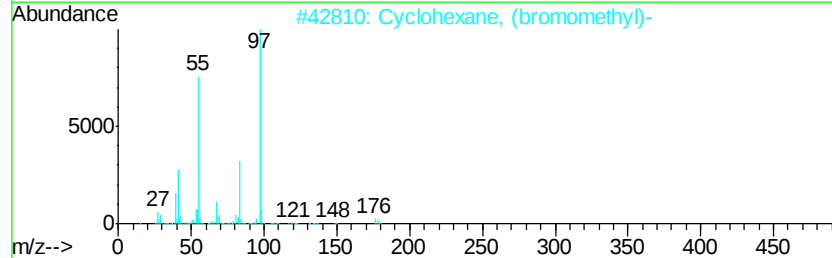
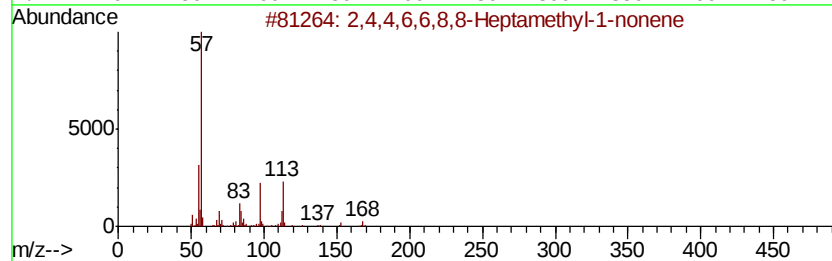
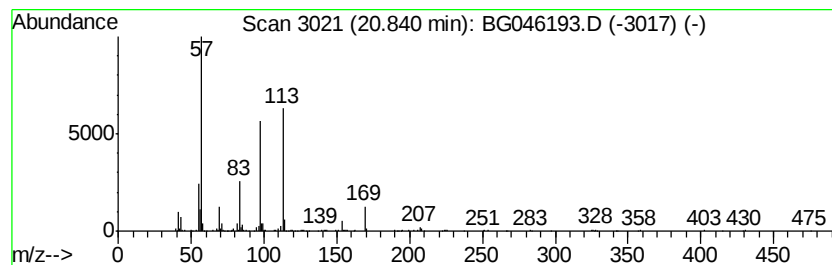
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Peak Number 3 2,4,4,6,6,8,8-Heptamethyl-1... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.84	2.23 ng	200085	Chrysene-d12	21.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	56
2		Cyclohexane, (bromomethyl)-	176	C7H13Br	002550-36-9	27
3		3-Heptene, 2,2,4,6,6-pentamethyl-	168	C12H24	000123-48-8	25
4		Octane, 2-bromo-	192	C8H17Br	000557-35-7	25
5		Sulfurous acid, cyclohexylmethyl...	234	C11H22O3S	1000309-21-3	25



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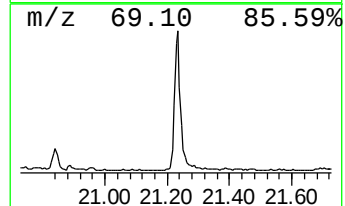
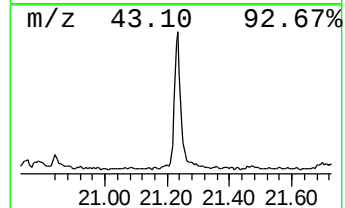
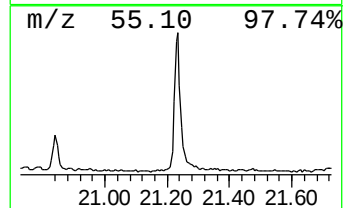
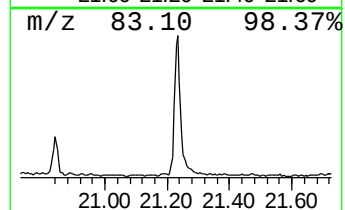
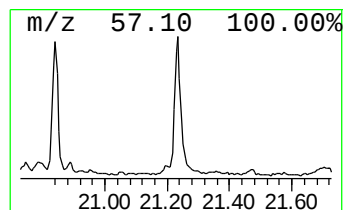
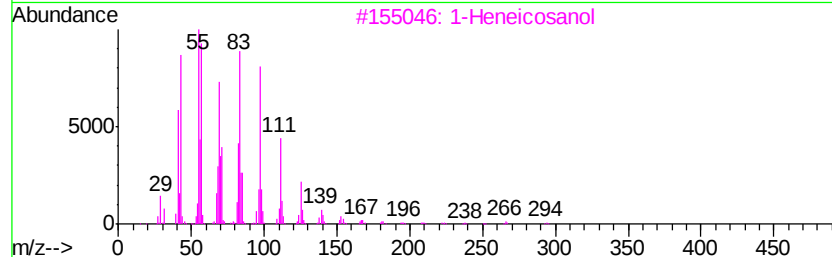
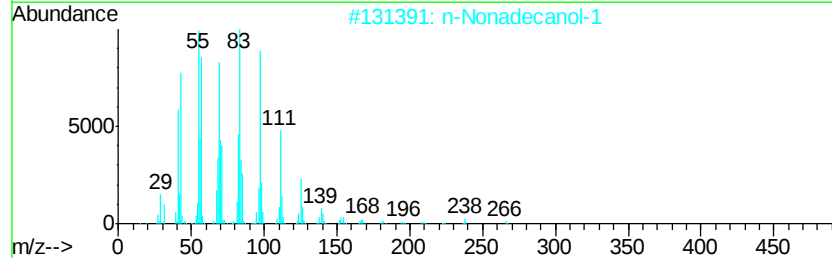
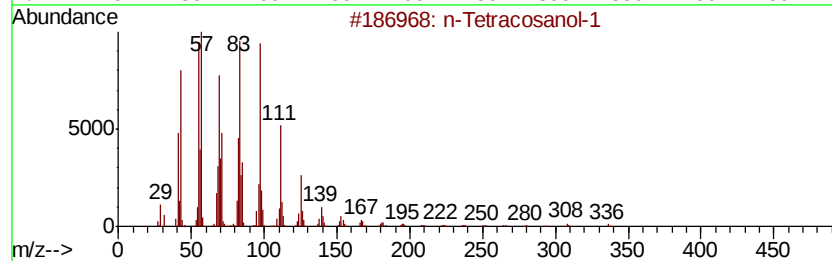
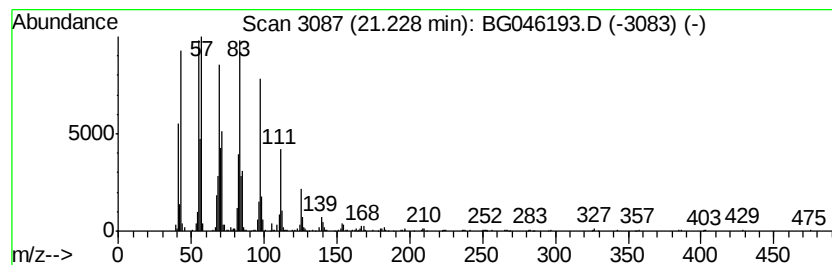
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Peak Number 4 n-Tetracosanol-1 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.23	7.80 ng	699523	Chrysene-d12	21.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Tetracosanol-1	354	C24H50O	000506-51-4	94	
2	n-Nonadecanol-1	284	C19H40O	001454-84-8	94	
3	1-Heneicosanol	312	C21H44O	015594-90-8	94	
4	Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94	
5	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94	



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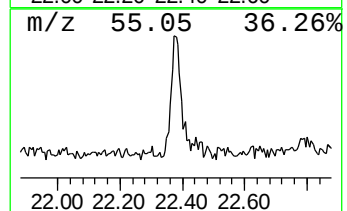
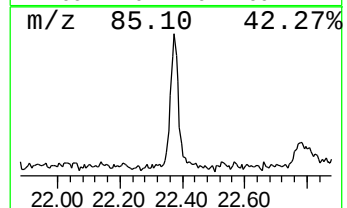
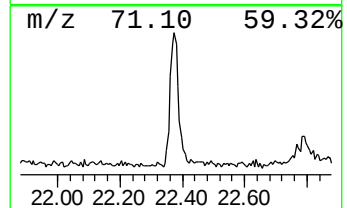
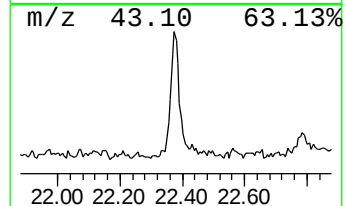
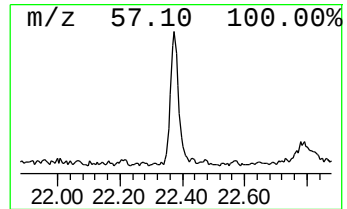
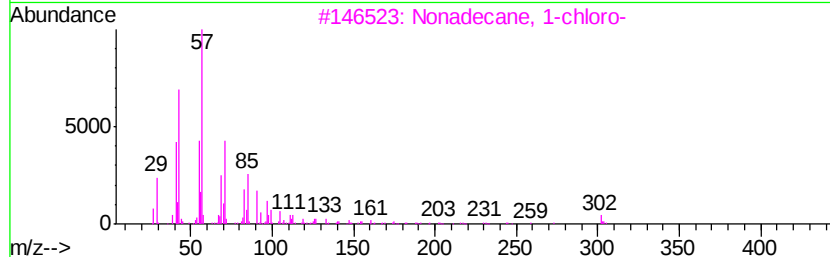
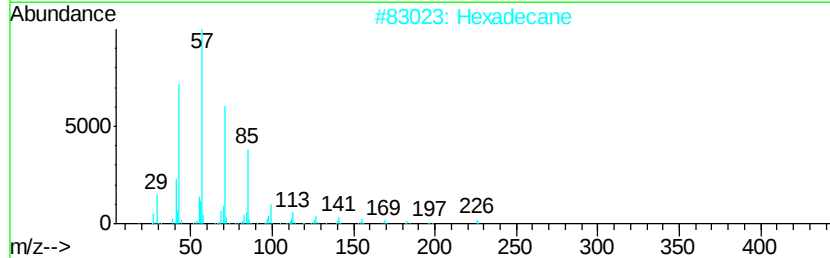
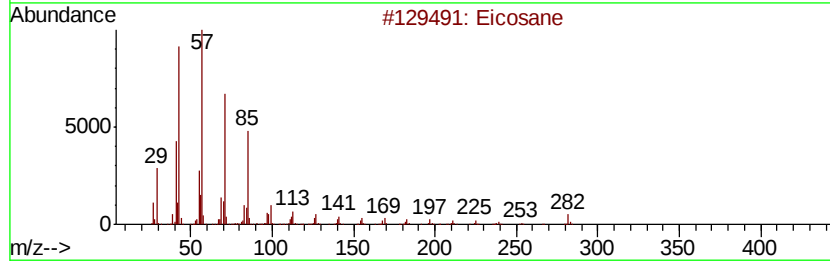
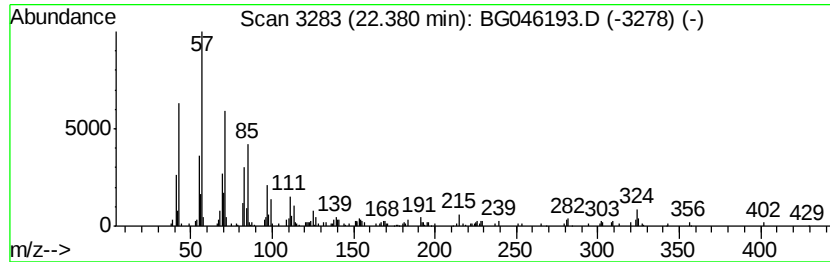
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Peak Number 5 Eicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.38	2.07 ng	185832	Chrysene-d12	21.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	96
2	Hexadecane		226	C16H34	000544-76-3	96
3	Nonadecane, 1-chloro-		302	C19H39Cl	062016-76-6	93
4	Tricosane		324	C23H48	000638-67-5	93
5	Sulfurous acid, octadecyl 2-prop...		376	C21H44O3S	1000309-12-7	76



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.15	26.8	ng	748534	1	7.95	559166	20.0
unknown18.99	18.99	5.0	ng	363622	4	17.31	1455670	20.0
2,4,4,6,6,8,8-Hep...	20.84	2.2	ng	200085	5	21.56	1794760	20.0
n-Tetracosanol-1	21.23	7.8	ng	699523	5	21.56	1794760	20.0
Eicosane	22.38	2.1	ng	185832	5	21.56	1794760	20.0