

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112817\
 Data File : BF100963.D
 Acq On : 29 Nov 2017 5:18
 Operator : SJ/JU
 Sample : I6603-05
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 112217-JETT-F-D-B

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-BF110817.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	5.075	505	508	515	rBV	103061	108411	5.06%	0.924%
2	5.475	572	576	588	rBV	644165	583703	27.23%	4.976%
3	6.481	744	747	756	rBV	477991	429923	20.05%	3.665%
4	6.628	768	772	775	rBV	1262746	1055860	49.25%	9.001%
5	6.857	808	811	815	rBV	458735	366322	17.09%	3.123%
6	7.016	834	838	841	rBV	1717678	1510489	70.46%	12.877%
7	7.428	902	908	911	rBV	1134693	1163411	54.27%	9.918%
8	8.139	1025	1029	1032	rBV	575017	452927	21.13%	3.861%
9	9.216	1207	1212	1215	rBV	2331603	2143821	100.00%	18.276%
10	9.898	1322	1328	1334	rBV	488612	475692	22.19%	4.055%
11	10.686	1458	1462	1473	rBV	718136	699814	32.64%	5.966%
12	11.380	1576	1580	1584	rBV	480201	411659	19.20%	3.509%
13	12.968	1845	1850	1853	rBV	1821158	1640746	76.53%	13.988%
14	13.857	1998	2001	2009	rBV2	26380	42008	1.96%	0.358%
15	14.021	2026	2029	2036	rBV	390887	348538	16.26%	2.971%
16	15.492	2275	2279	2287	rBV	229842	296680	13.84%	2.529%

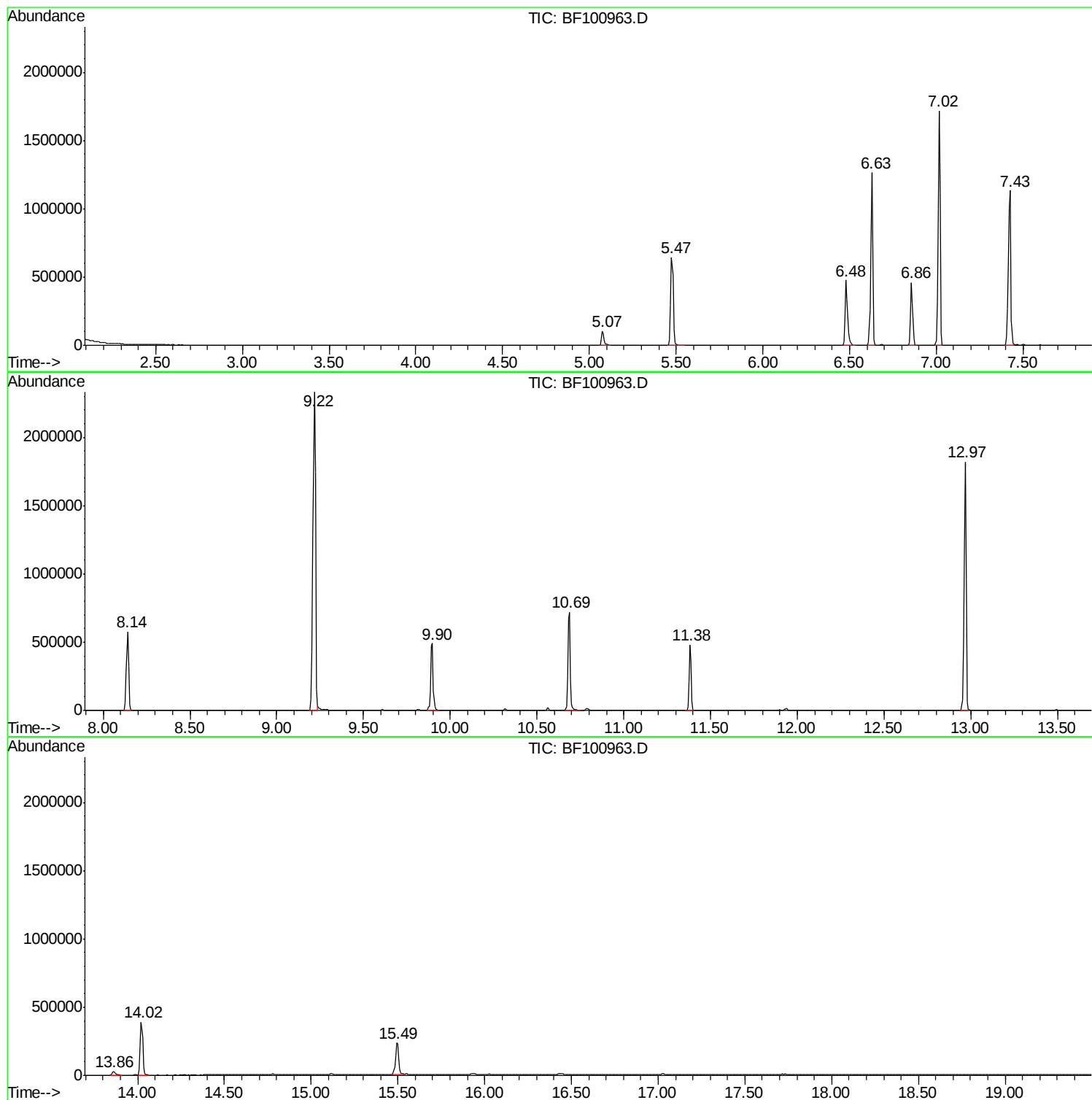
Sum of corrected areas: 11730004

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.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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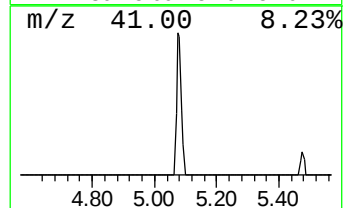
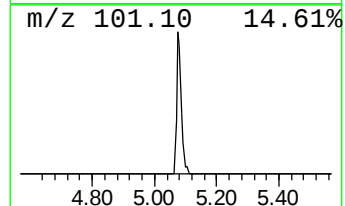
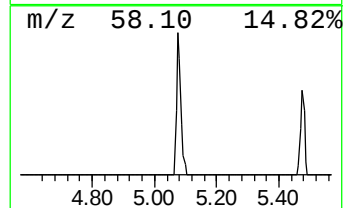
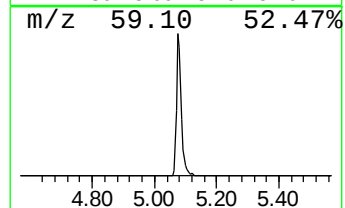
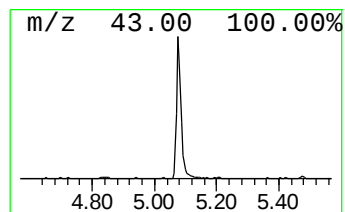
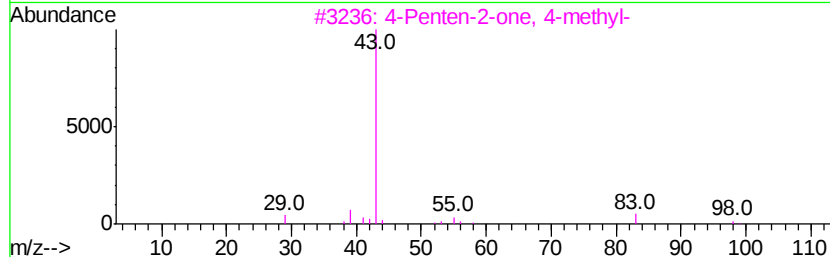
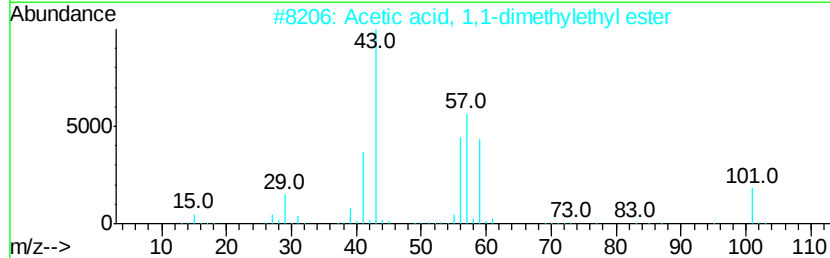
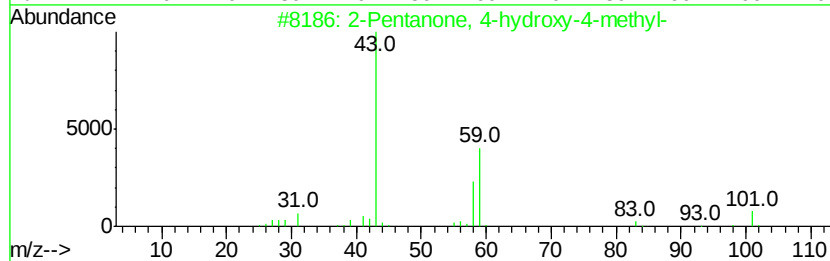
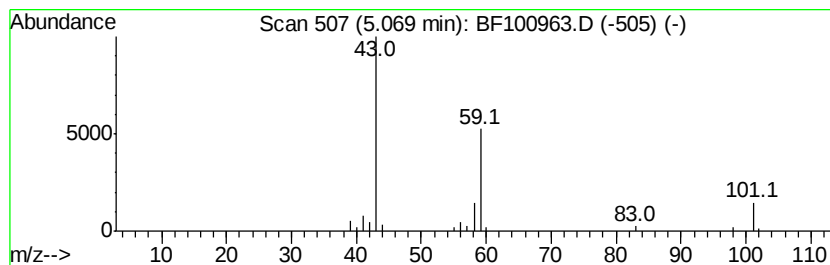
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	5.92 ng	108411	1,4-Dichlorobenzene-d4	6.86

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, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	9



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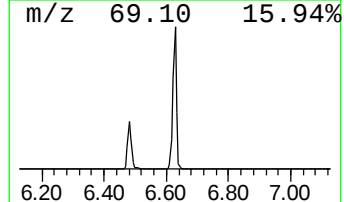
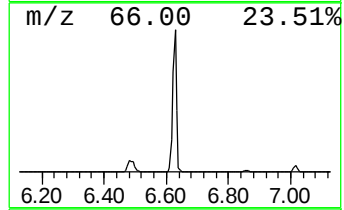
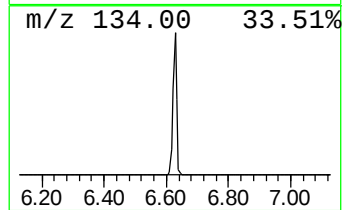
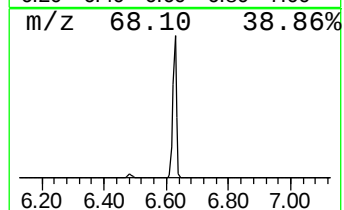
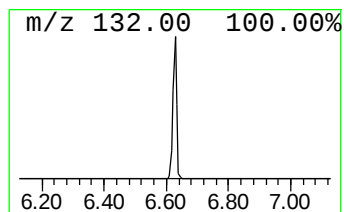
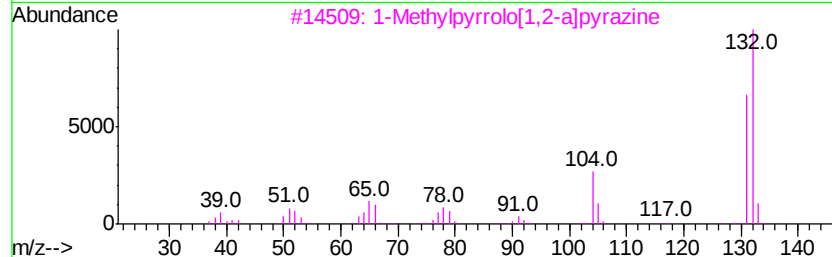
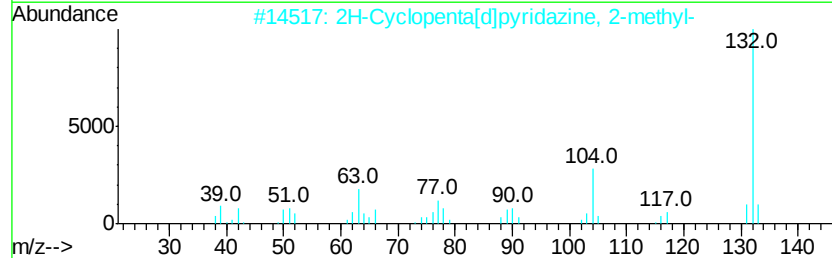
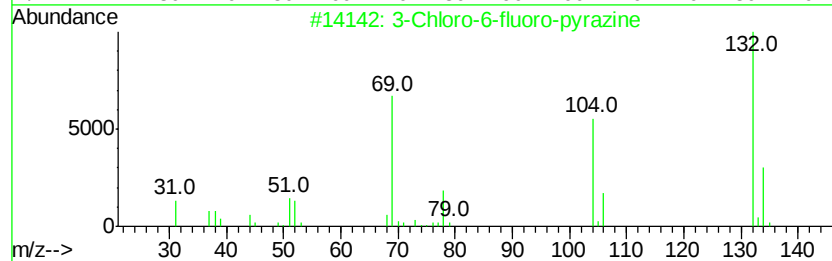
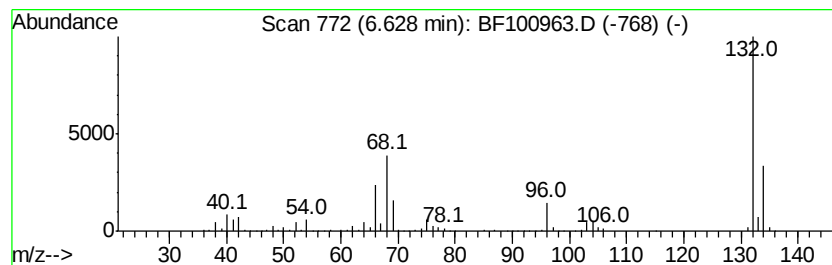
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Peak Number 2 unknown6.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	57.65 ng	1055860	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
3		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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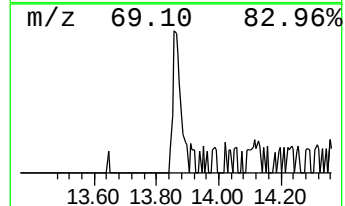
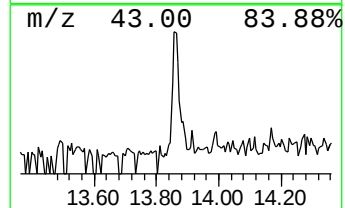
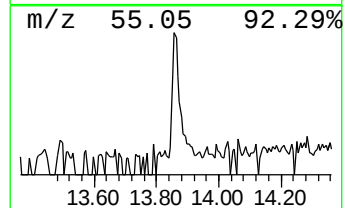
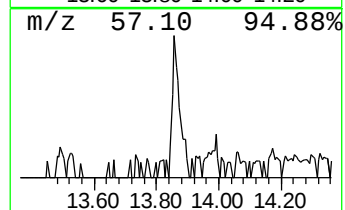
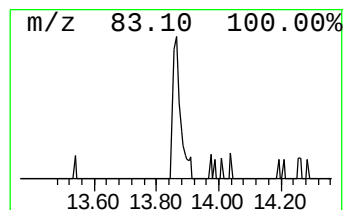
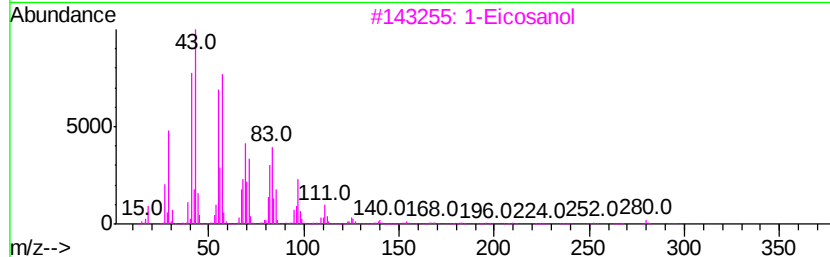
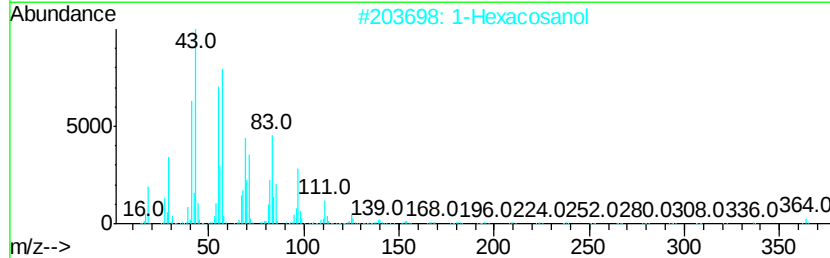
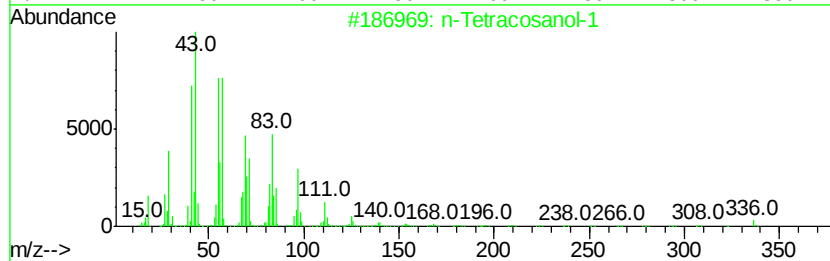
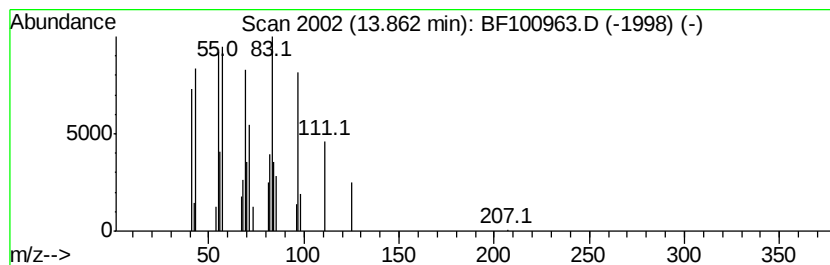
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Peak Number 4 n-Tetracosanol-1 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	2.41 ng	42008	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Tetracosanol-1	354	C24H50O	000506-51-4	72
2		1-Hexacosanol	382	C26H54O	000506-52-5	64
3		1-Eicosanol	298	C20H42O	000629-96-9	64
4		5-Octadecene, (E)-	252	C18H36	007206-21-5	50
5		Octadecane, 1-(ethenyl)-	296	C20H40	000930-02-9	46



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.07	5.9	ng	108411	1	6.86	366322	20.0
unknown6.63	6.63	57.6	ng	1055860	1	6.86	366322	20.0
n-Tetracosanol-1	13.86	2.4	ng	42008	5	14.02	348538	20.0