

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070125\
 Data File : BF142952.D
 Acq On : 01 Jul 2025 16:27
 Operator : RC/JU
 Sample : Q2449-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBR200032

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-BF061925.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF142952.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.081	482	491	499	rBV	79810	113448	8.45%	1.277%
2	5.487	555	560	580	rVB	664555	884463	65.91%	9.959%
3	6.498	727	732	752	rBV	703157	924757	68.92%	10.412%
4	6.869	790	795	804	rBV	307708	375820	28.01%	4.232%
5	7.434	885	891	896	rBV	480706	598157	44.58%	6.735%
6	8.157	1009	1014	1019	rBV	397842	489830	36.50%	5.515%
7	9.234	1191	1197	1202	rBV	1051655	1341875	100.00%	15.109%
8	9.622	1259	1263	1266	rBV2	18091	26349	1.96%	0.297%
9	9.775	1286	1289	1292	rBV4	20436	31013	2.31%	0.349%
10	9.822	1293	1297	1300	rVB2	35862	51422	3.83%	0.579%
11	9.916	1307	1313	1317	rBV	431742	586870	43.74%	6.608%
12	10.316	1379	1381	1385	rBV2	35051	43608	3.25%	0.491%
13	10.628	1431	1434	1437	rBV	59712	63652	4.74%	0.717%
14	10.704	1443	1447	1452	rBV	537072	701255	52.26%	7.896%
15	11.404	1562	1566	1570	rBV	420877	525422	39.16%	5.916%
16	12.992	1831	1836	1843	rVB	803586	1065456	79.40%	11.997%
17	13.892	1985	1989	2007	rVB2	52317	146460	10.91%	1.649%
18	14.045	2011	2015	2026	rVV	263522	368408	27.45%	4.148%
19	14.139	2028	2031	2039	rVB	24725	35177	2.62%	0.396%
20	14.533	2094	2098	2108	rVB	53900	106469	7.93%	1.199%
21	15.539	2263	2269	2277	rVB	248194	377679	28.15%	4.253%
22	16.086	2354	2362	2365	rBV7	8156	23692	1.77%	0.267%

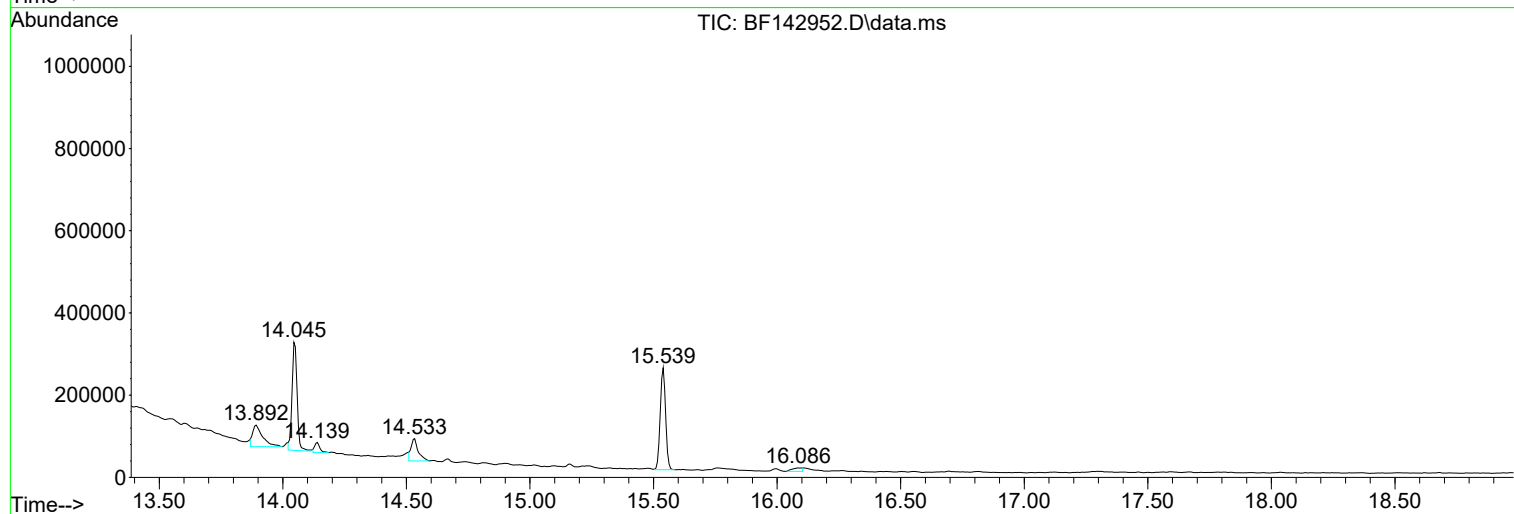
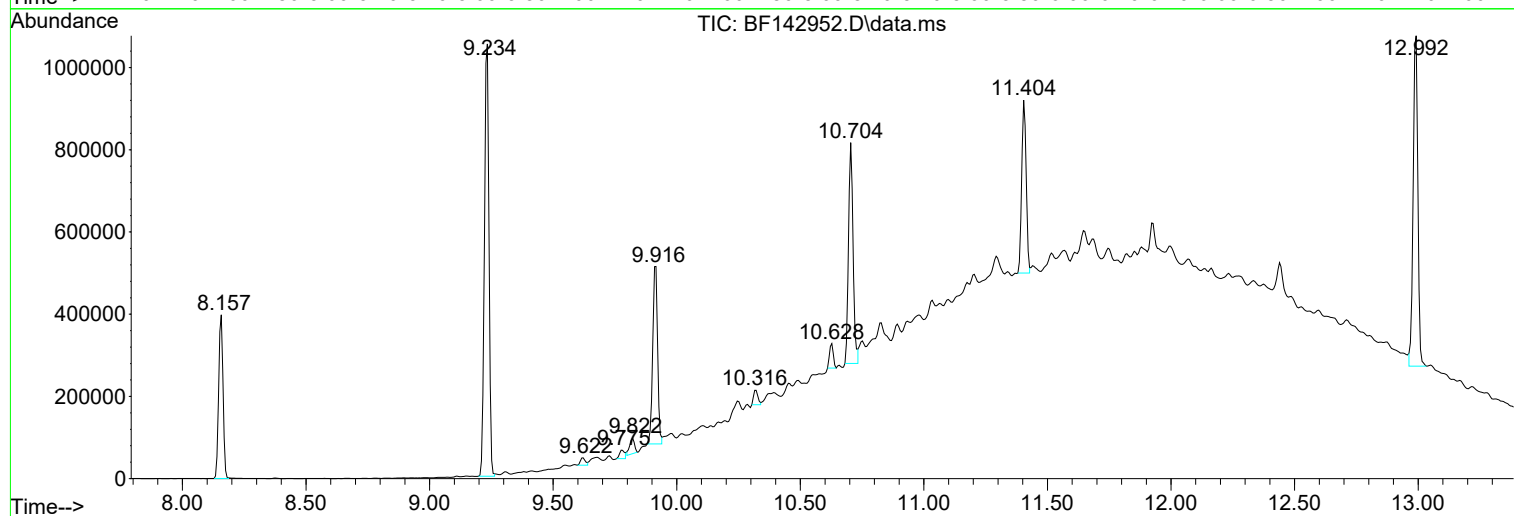
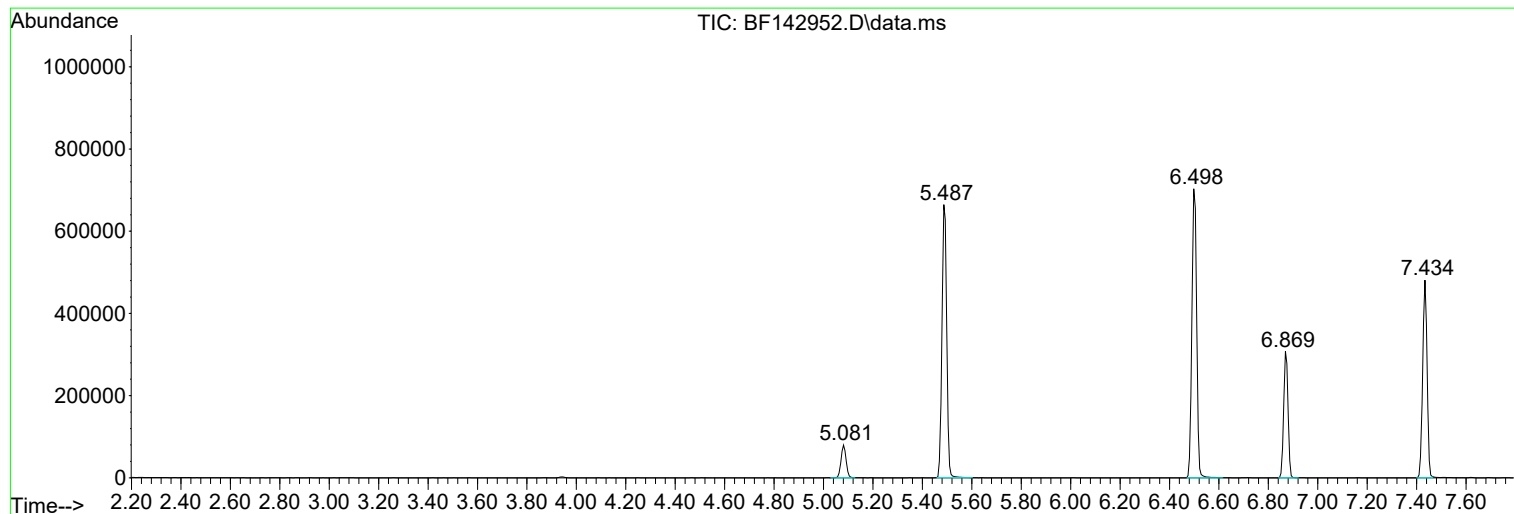
Sum of corrected areas: 8881282

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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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Misc :
ALS Vial : 13 Sample Multiplier: 1

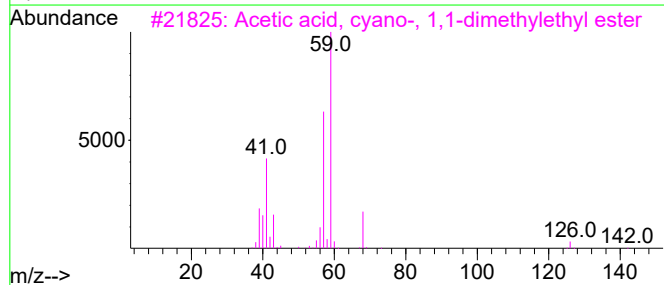
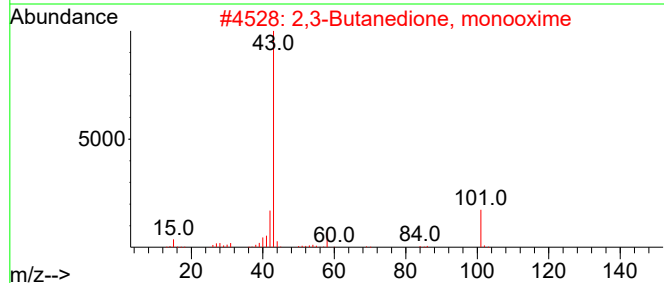
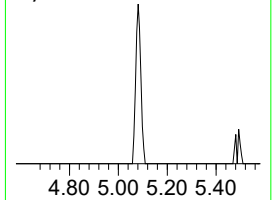
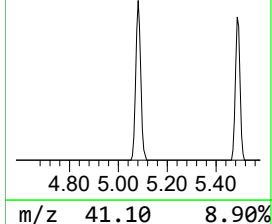
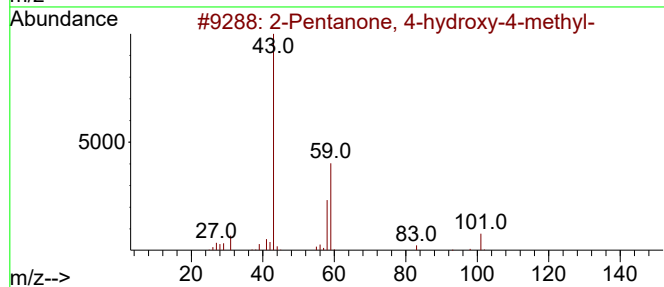
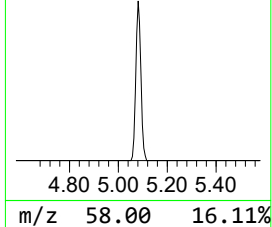
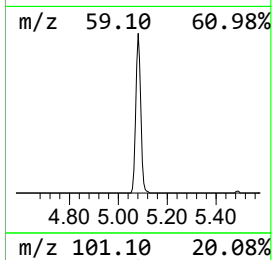
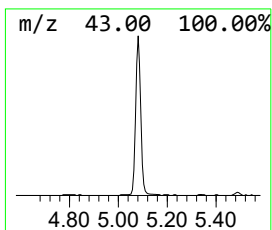
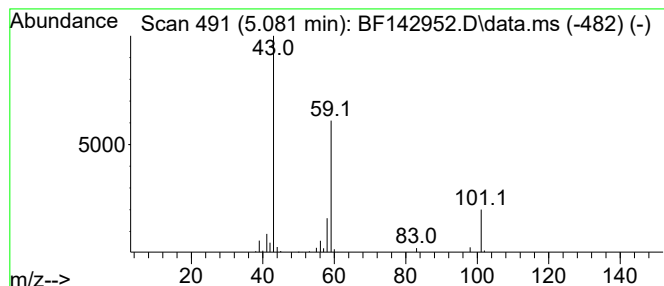
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD			R.T.
5.081	6.04 ng	113448	1,4-Dichlorobenzene-d4			6.869
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	56
2	2,3-Butanedione, monooxime		101	C4H7NO2	000057-71-6	35
3	Acetic acid, cyano-, 1,1-dimethy...		141	C7H11NO2	001116-98-9	23
4	1-Propen-2-ol, acetate		100	C5H8O2	000108-22-5	10
5	Butane, 1-ethoxy-		102	C6H14O	000628-81-9	9



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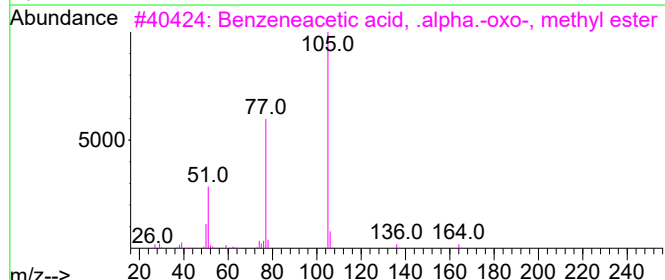
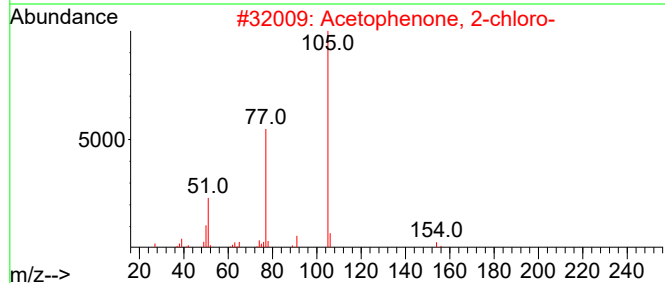
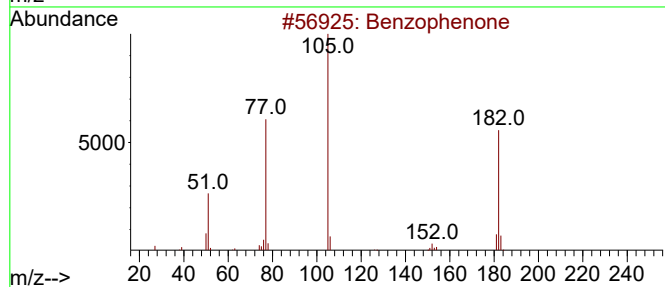
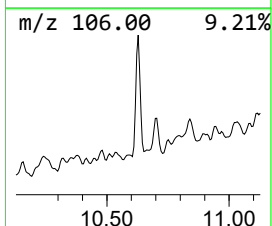
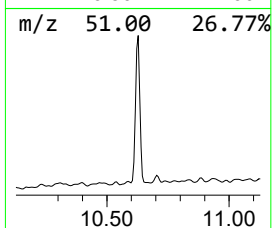
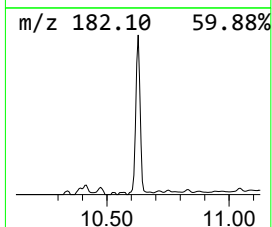
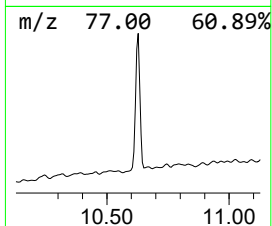
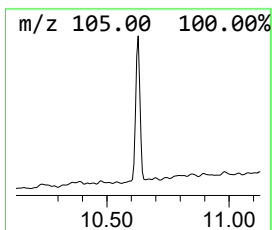
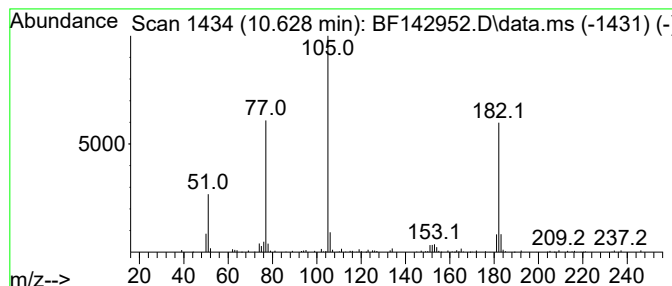
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Peak Number 2 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.628	2.17 ng	63652	Acenaphthene-d10	9.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	95
2			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	53
3			Benzeneacetic acid, .alpha.-oxo-...	164	C9H8O3	015206-55-0	47
4			Ethanedione, diphenyl-	210	C14H10O2	000134-81-6	47
5			Ethanone, 2-hydroxy-1-phenyl-	136	C8H8O2	000582-24-1	47



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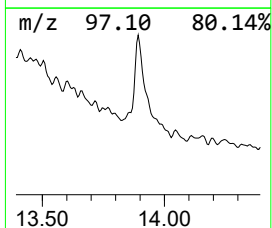
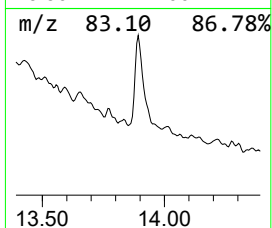
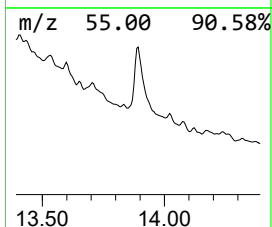
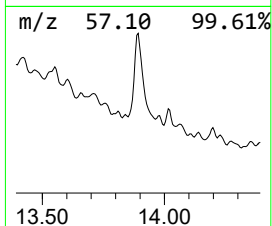
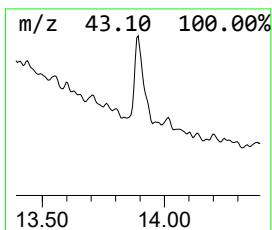
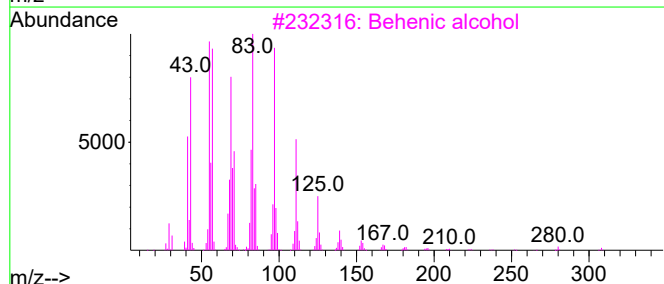
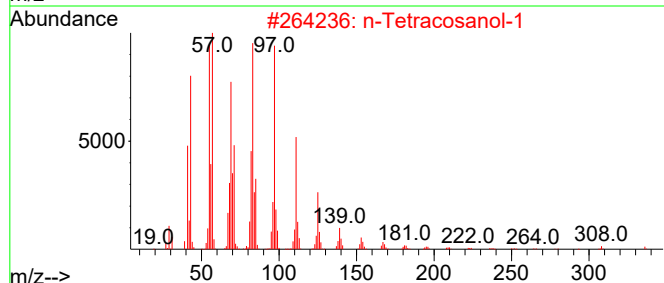
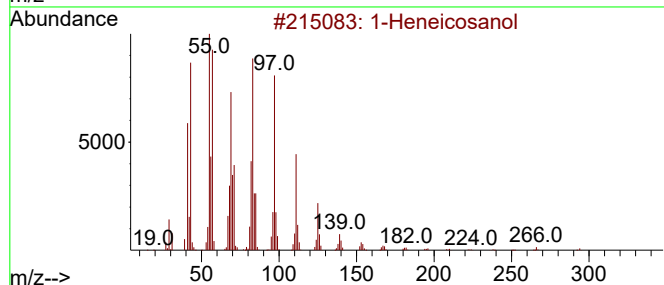
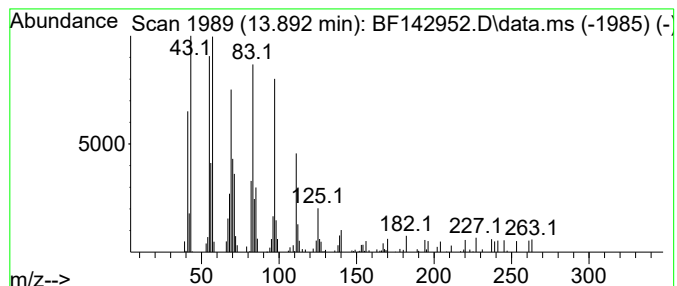
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TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Heneicosanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.892	7.95 ng	146460	Chrysene-d12	14.045	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	n-Tetracosanol-1	354	C24H50O	000506-51-4	95
3	Behenic alcohol	326	C22H46O	000661-19-8	95
4	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
5	1-Tetracosene	336	C24H48	010192-32-2	91



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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.081	6.0	ng	113448	1	6.869	375820	20.0
Benzophenone	10.628	2.2	ng	63652	3	9.916	586870	20.0
1-Heneicosanol	13.892	8.0	ng	146460	5	14.045	368408	20.0
Strychnidin-10-...	14.533	5.8	ng	106469	5	14.045	368408	20.0