

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091019\
 Data File : BP000195.D
 Acq On : 11 Sep 2019 11:52
 Operator : HP/JU
 Sample : PB122977BL
 Misc : GCMS CONFIRMATION
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK77

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % 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_P\METHODS\SOM-EPA-BP090519MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.117	3	5	11	rVB	48258	46574	1.54%	0.279%
2	2.193	13	18	27	rVB2	96401	132359	4.38%	0.792%
3	2.334	38	42	55	rVV	218237	392431	13.00%	2.348%
4	2.428	55	58	65	rVB	16153	22924	0.76%	0.137%
5	2.575	77	83	91	rVB	223685	303648	10.06%	1.817%
6	3.122	171	176	184	rVB2	13469	22894	0.76%	0.137%
7	4.005	322	326	332	rVB	13960	17514	0.58%	0.105%
8	4.381	386	390	394	rBV	3904	4159	0.14%	0.025%
9	4.522	411	414	417	rBV3	3056	3672	0.12%	0.022%
10	4.881	472	475	482	rBV8	2006	3729	0.12%	0.022%
11	5.110	511	514	517	rBV	4480	4754	0.16%	0.028%
12	6.893	813	817	820	rBV	2168744	1693343	56.08%	10.131%
13	8.175	1031	1035	1039	rBV	2832834	2273993	75.31%	13.605%
14	8.228	1042	1044	1052	rVB3	4193	5057	0.17%	0.030%
15	9.945	1332	1336	1340	rBV	3447015	2799625	92.71%	16.750%
16	10.469	1422	1425	1428	rVB	7232	6166	0.20%	0.037%
17	11.328	1568	1571	1575	rVB	11852	9817	0.33%	0.059%
18	11.451	1588	1592	1596	rBV	3546618	2958391	97.97%	17.700%
19	11.504	1599	1601	1611	rVB	16285	17727	0.59%	0.106%
20	14.151	2046	2051	2060	rBV	2946505	3019649	100.00%	18.066%
21	15.704	2308	2315	2323	rBV	1878941	2976006	98.55%	17.805%

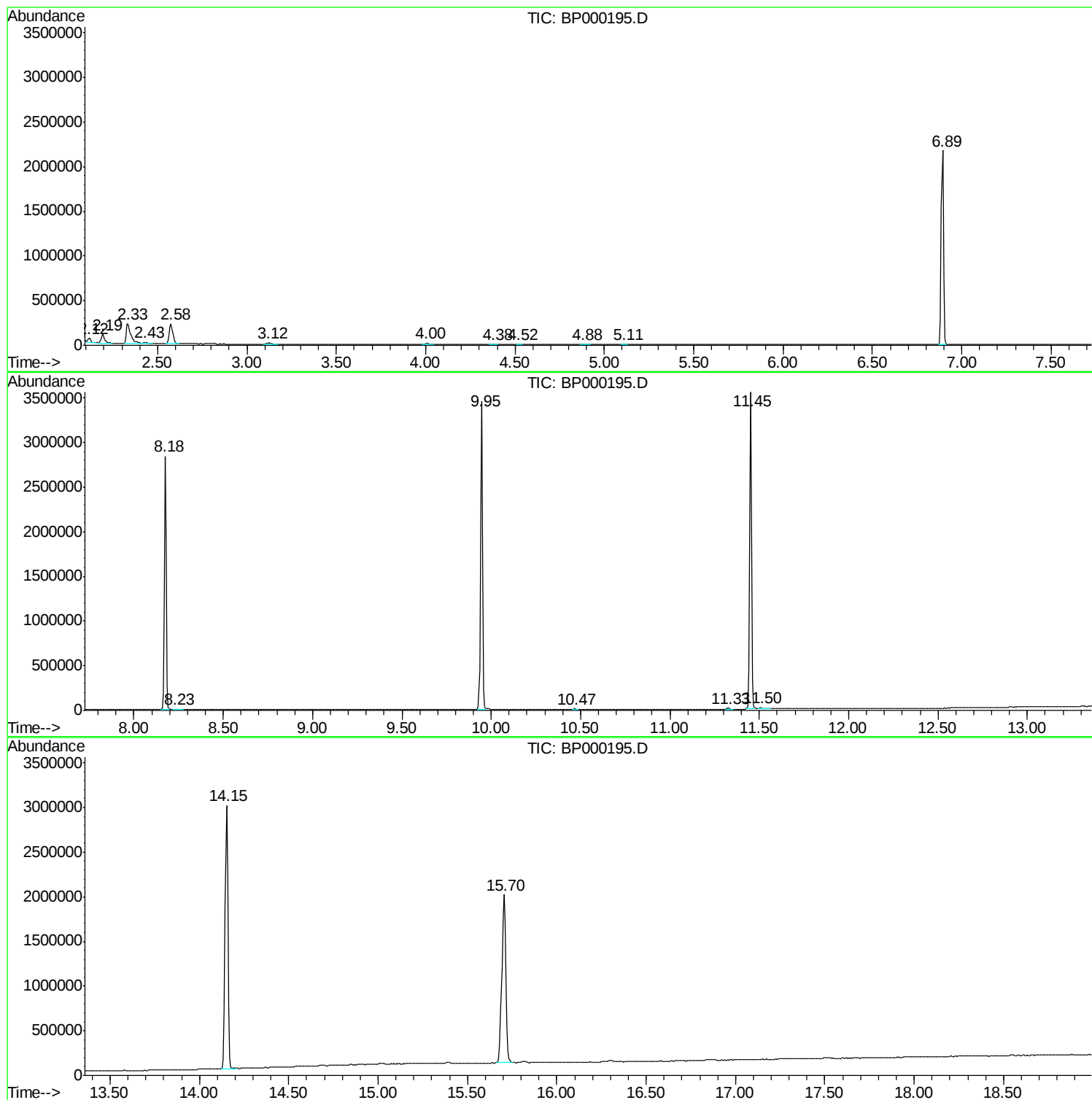
Sum of corrected areas: 16714432

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP090519MA.M
Quant Title : SVOA CALIBRATION

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



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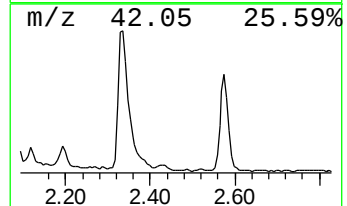
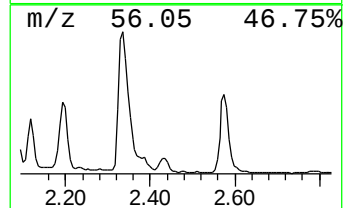
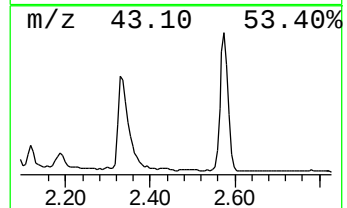
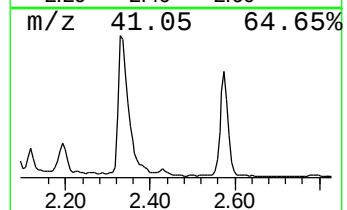
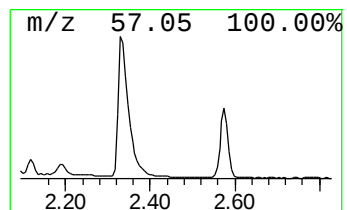
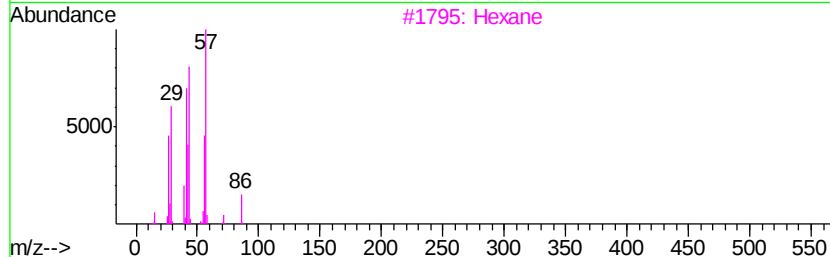
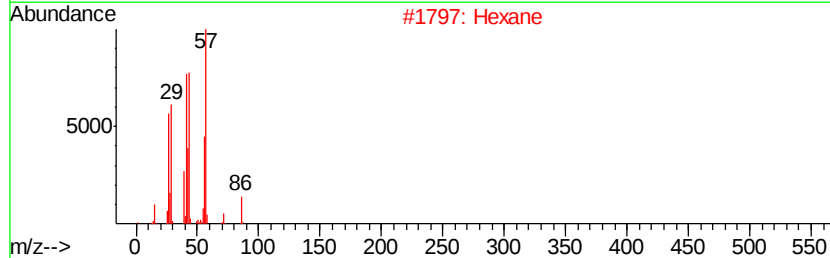
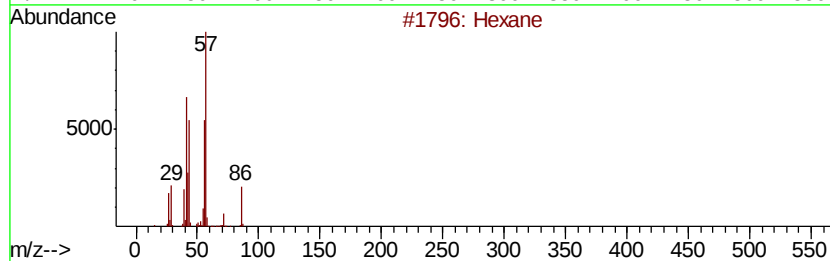
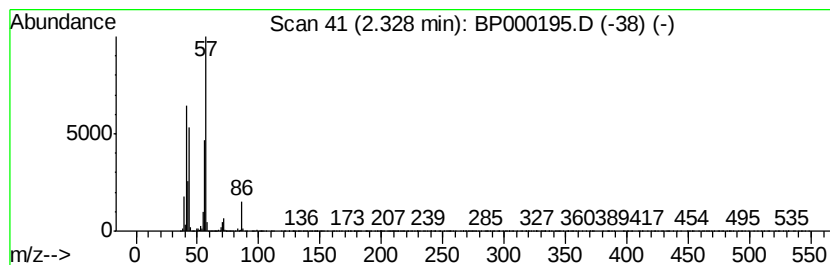
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP090519MA.M
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Hexane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.33	4.63 ng/ul	392431	1,4-Dichlorobenzene-d4	6.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane			86	C6H14	000110-54-3	94
2	Hexane			86	C6H14	000110-54-3	58
3	Hexane			86	C6H14	000110-54-3	53
4	Pentane, 2,2,3,4-tetramethyl-			128	C9H20	001186-53-4	50
5	Furan, tetrahydro-3-methyl-			86	C5H10O	013423-15-9	46



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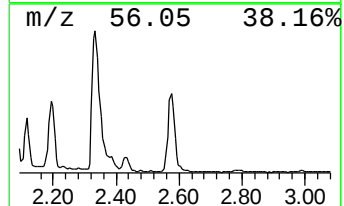
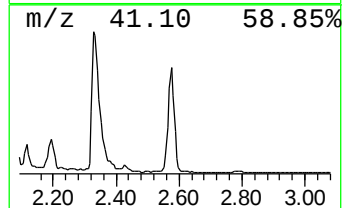
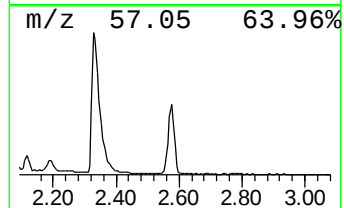
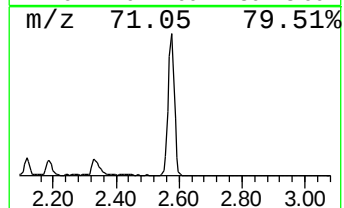
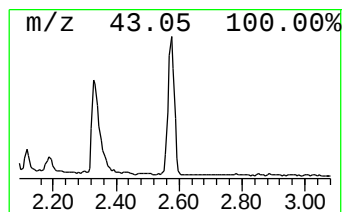
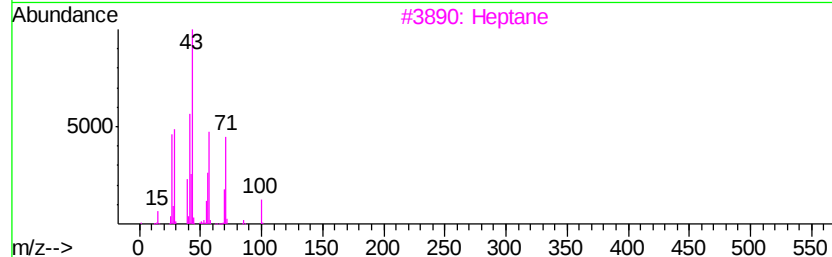
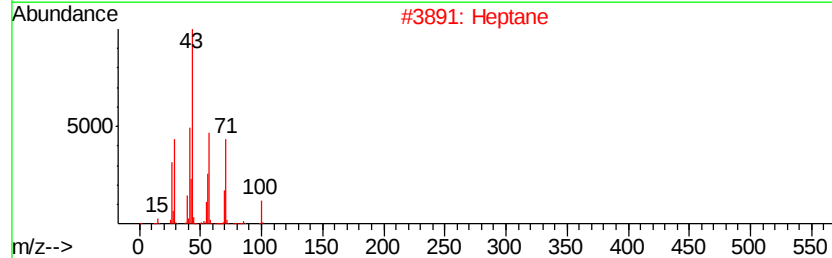
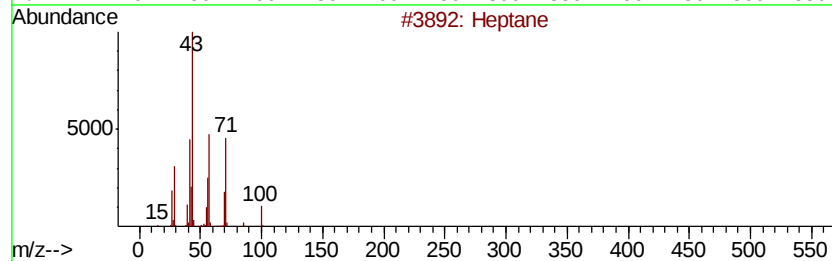
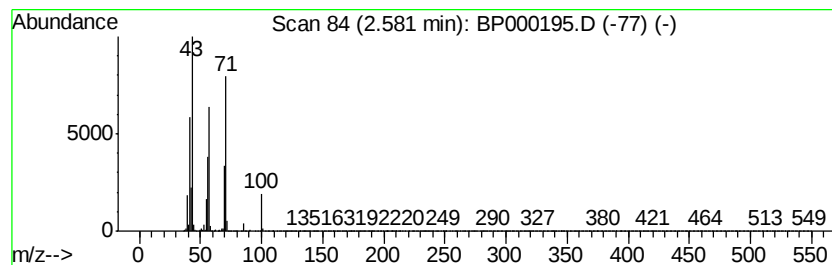
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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.58	3.59 ng/ul	303648	1,4-Dichlorobenzene-d4	6.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane			100	C7H16	000142-82-5	94
2	Heptane			100	C7H16	000142-82-5	91
3	Heptane			100	C7H16	000142-82-5	91
4	Heptane			100	C7H16	000142-82-5	80
5	Butane, 2,2-dimethyl-			86	C6H14	000075-83-2	53



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Hexane	2.33	4.6	ng/ul	392431	1	6.89	1693340	20.0
(DEL) Alkane: Str...	2.58	3.6	ng/ul	303648	1	6.89	1693340	20.0