

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
 Data File : BM014415.D
 Acq On : 28 Feb 2018 23:08
 Operator : SJ/JU
 Sample : J1679-08
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BE614

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.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	6.722	629	635	648	rBV	94376	227343	5.13%	0.614%
2	6.863	654	659	672	rBV	349743	612354	13.83%	1.655%
3	7.057	685	692	703	rBV	412516	770751	17.41%	2.083%
4	7.516	764	770	778	rBV	344749	579685	13.09%	1.567%
5	7.592	778	783	794	rVB5	34096	80704	1.82%	0.218%
6	8.239	887	893	906	rBV2	324904	638007	14.41%	1.724%
7	8.675	961	967	980	rVB	601696	1055233	23.83%	2.852%
8	9.386	1081	1088	1101	rBV	539566	1010667	22.82%	2.731%
9	9.928	1173	1180	1198	rBV	479434	1223088	27.62%	3.305%
10	10.286	1234	1241	1253	rVB	547680	931840	21.04%	2.518%
11	13.598	1798	1804	1817	rVB	1691425	2357599	53.24%	6.371%
12	13.863	1842	1849	1862	rBV	1836181	2715929	61.34%	7.340%
13	14.174	1895	1902	1911	rBV2	954225	1434808	32.40%	3.877%
14	14.474	1945	1953	1964	rBV6	49941	155649	3.52%	0.421%
15	15.180	2066	2073	2084	rBV	2265705	3260441	73.63%	8.811%
16	15.339	2094	2100	2110	rBV2	700137	1133399	25.60%	3.063%
17	15.421	2110	2114	2124	rVB6	35688	65572	1.48%	0.177%
18	16.933	2363	2371	2382	rBV2	1229571	1779177	40.18%	4.808%
19	17.033	2382	2388	2401	rBV	2440648	3690611	83.35%	9.974%
20	17.604	2480	2485	2491	rBV2	222950	292984	6.62%	0.792%
21	18.980	2715	2719	2725	rBV2	31526	46534	1.05%	0.126%
22	19.345	2775	2781	2794	rBV	3244723	4378540	98.88%	11.833%
23	21.150	3083	3088	3095	rBV	1585019	2085540	47.10%	5.636%
24	23.191	3428	3435	3444	rVB	2492745	4427994	100.00%	11.966%
25	23.327	3451	3458	3468	rVB	1062536	2049714	46.29%	5.539%

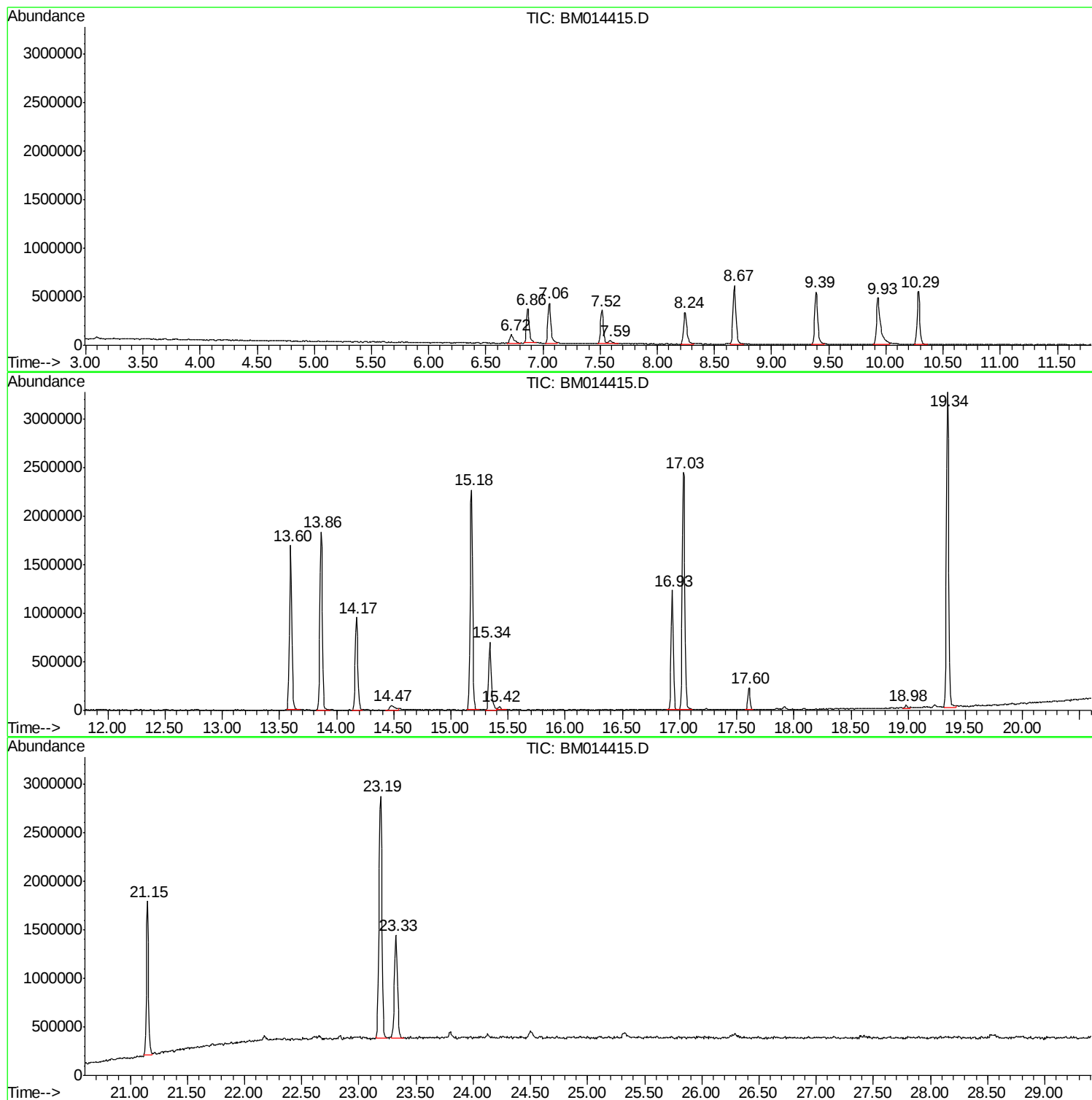
Sum of corrected areas: 37004163

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
Data File : BM014415.D
Acq On : 28 Feb 2018 23:08
Operator : SJ/JU
Sample : J1679-08
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE614

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
Data File : BM014415.D
Acq On : 28 Feb 2018 23:08
Operator : SJ/JU
Sample : J1679-08
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE614

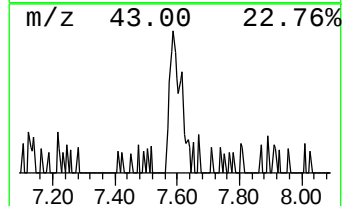
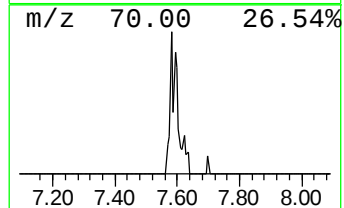
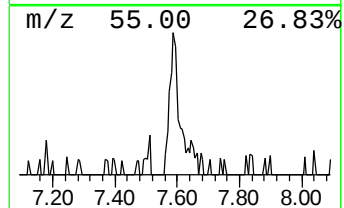
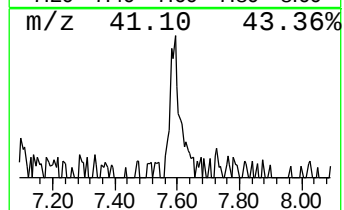
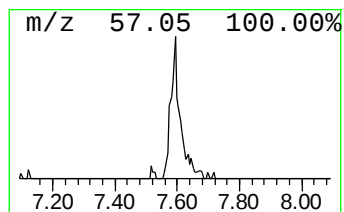
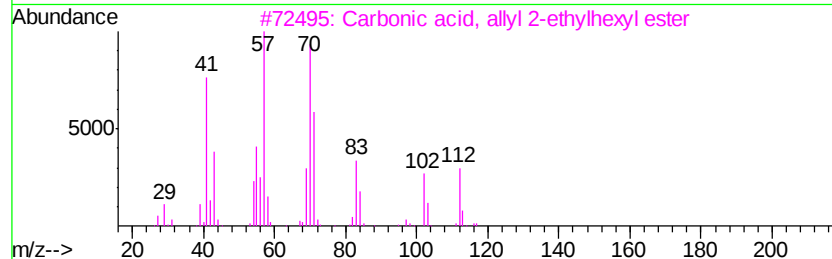
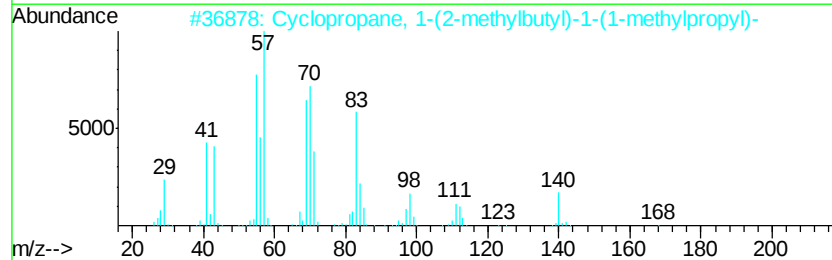
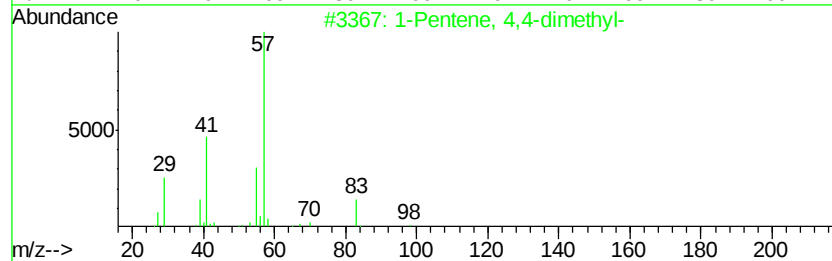
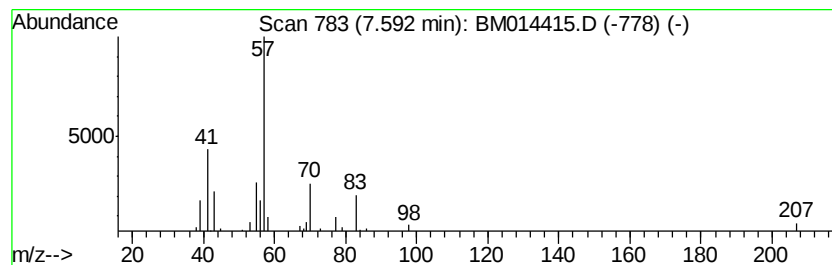
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic7.59 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	2.78 ng/ul	80704	1,4-Dichlorobenzene-d4	7.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentene, 4,4-dimethyl-	98	C7H14	000762-62-9	43
2		Cyclopropane, 1-(2-methylbutyl)-...	168	C12H24	064723-36-0	25
3		Carbonic acid, allyl 2-ethylhexyl...	214	C12H22O3	1000357-80-6	17
4		1-Hexene, 4-methyl-	98	C7H14	003769-23-1	17
5		1-Penten-3-ol	86	C5H10O	000616-25-1	12



Data Path : Z:\HPCHEM1\BNA M\DATA\BM022818\
Data File : BM014415.D
Acq On : 28 Feb 2018 23:08
Operator : SJ/JU
Sample : J1679-08
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE614

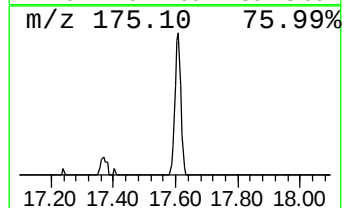
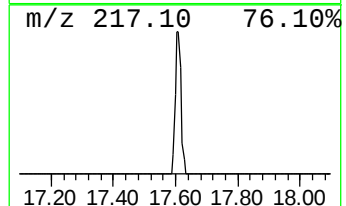
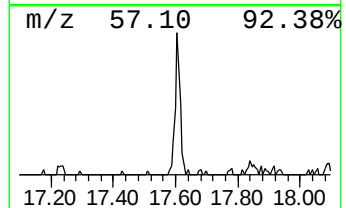
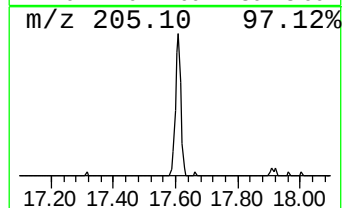
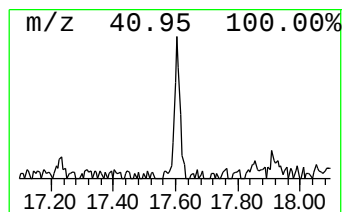
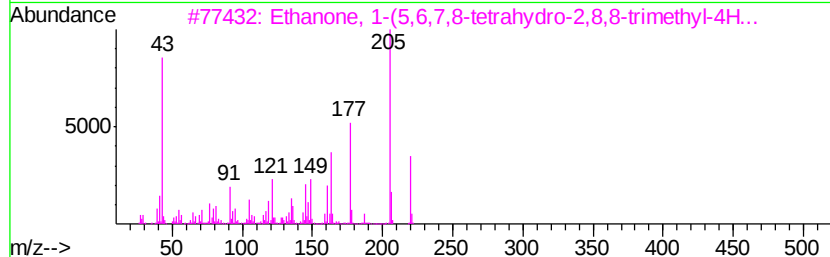
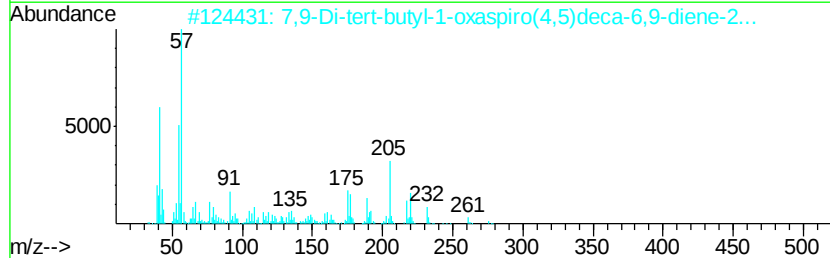
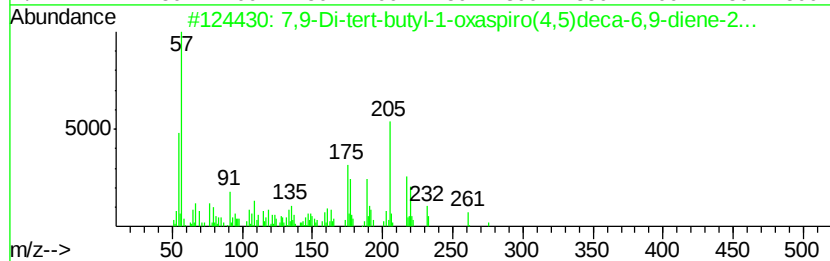
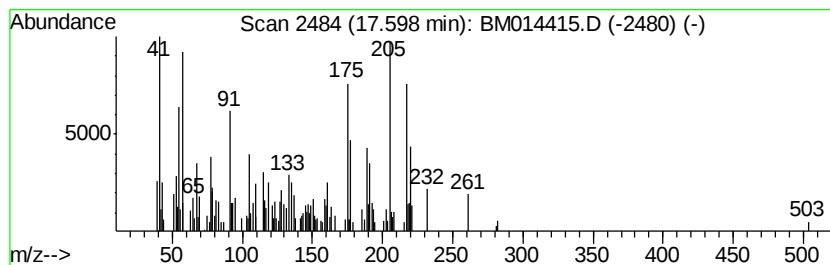
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.60	3.29 ng/ul	292984	Phenanthrene-d10	16.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	93		
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	70		
3	Ethanone, 1-(5,6,7,8-tetrahydro-...	220	C14H20O2	071596-88-8	56		
4	2,5-Cyclohexadien-1-one, 2,6-bis...	232	C16H24O	006738-27-8	42		
5	9-Acetylphenanthrene	220	C16H12O	002039-77-2	38		



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM022818\
Data File : BM014415.D
Acq On : 28 Feb 2018 23:08
Operator : SJ/JU
Sample : J1679-08
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE614

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Cyc...	7.59	2.8	ng/ul	80704	1	7.52	579685	20.0
7,9-Di-tert-butyl...	17.60	3.3	ng/ul	292984	4	16.93	1779180	20.0