

Data Path : Z:\HPCHEM1\BNA M\DATA\BM071516\
 Data File : BM006449.D
 Acq On : 15 Jul 2016 19:17
 Operator : UM/SJ
 Sample : H3992-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BDJ81

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\SOM02.2-EPA-BM071416.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	2.934	54	59	65	rBV	58729	77110	2.22%	0.230%
2	4.181	266	271	279	rVB	27111	40633	1.17%	0.121%
3	4.763	365	370	379	rBV	40746	60237	1.73%	0.180%
4	6.798	709	716	734	rVB	460358	763589	21.97%	2.282%
5	6.957	736	743	756	rVV	328855	516379	14.86%	1.543%
6	7.151	770	776	788	rBV	456462	734050	21.12%	2.193%
7	7.616	849	855	865	rBV	448206	746371	21.47%	2.230%
8	8.328	970	976	992	rVB	516877	884215	25.44%	2.642%
9	8.769	1045	1051	1061	rBV	409352	707675	20.36%	2.115%
10	9.486	1167	1173	1186	rVB	334286	588377	16.93%	1.758%
11	10.028	1257	1265	1279	rBV	488623	892852	25.69%	2.668%
12	10.398	1321	1328	1341	rVB	577047	1034859	29.77%	3.092%
13	10.539	1346	1352	1367	rBV	502426	894648	25.74%	2.673%
14	13.680	1880	1886	1891	rBV	951273	1423574	40.96%	4.254%
15	13.727	1891	1894	1912	rVB	446776	677975	19.51%	2.026%
16	13.963	1927	1934	1946	rBV	1011131	1624202	46.73%	4.853%
17	14.268	1979	1986	1994	rBV2	831705	1380898	39.73%	4.126%
18	14.486	2016	2023	2040	rBV	387537	698967	20.11%	2.089%
19	15.133	2129	2133	2141	rVB	29401	37210	1.07%	0.111%
20	15.268	2149	2156	2164	rBV	1366757	2101908	60.47%	6.281%
21	15.398	2173	2178	2192	rBV2	393960	597934	17.20%	1.787%
22	15.992	2275	2279	2287	rBV2	30468	60901	1.75%	0.182%
23	16.786	2408	2414	2420	rBV2	30727	50761	1.46%	0.152%
24	17.021	2447	2454	2459	rBV	1202653	1807099	51.99%	5.400%
25	17.056	2459	2460	2465	rVV	68276	78704	2.26%	0.235%
26	17.121	2465	2471	2484	rVB	1509668	2402926	69.13%	7.180%
27	19.086	2801	2805	2810	rVB	111657	156632	4.51%	0.468%
28	19.421	2855	2862	2875	rBV	2098812	3119427	89.75%	9.322%
29	20.939	3116	3120	3127	rVV7	53923	110027	3.17%	0.329%
30	21.215	3161	3167	3172	rBV	1931284	2609556	75.08%	7.798%
31	22.250	3340	3343	3346	rBV4	31613	39575	1.14%	0.118%
32	22.750	3424	3428	3434	rBV2	74783	155061	4.46%	0.463%
33	23.274	3510	3517	3533	rVB2	1750992	3475849	100.00%	10.387%
34	23.415	3534	3541	3556	rVB	1301760	2583723	74.33%	7.721%

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35	23.921	3622	3627	3635	rVB2	80673	156035	4.49%	0.466%
36	24.638	3745	3749	3760	rVB3	79926	174917	5.03%	0.523%

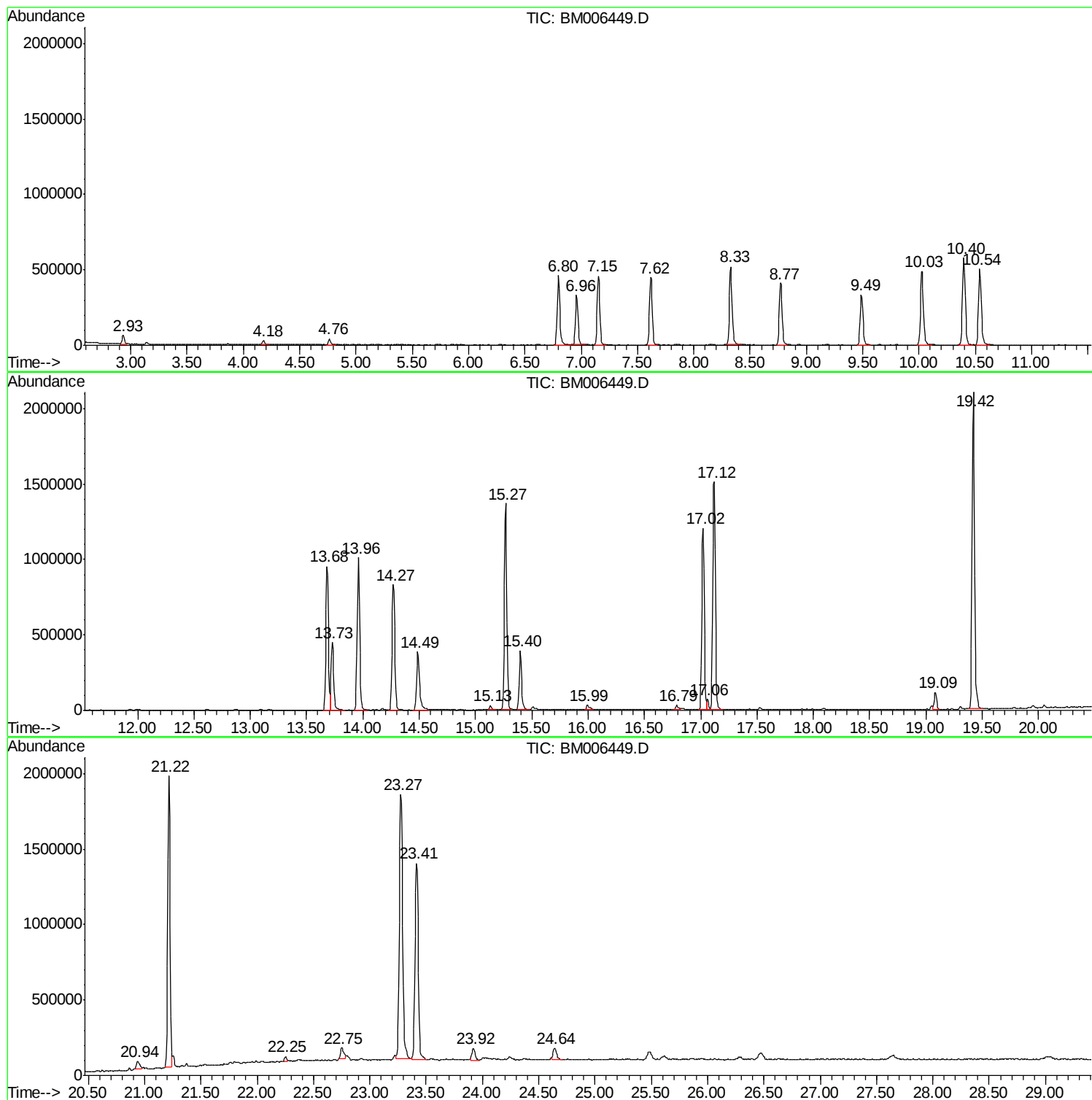
Sum of corrected areas: 33464856

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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM071416.M
Quant Title : SVOA CALIBRATION

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



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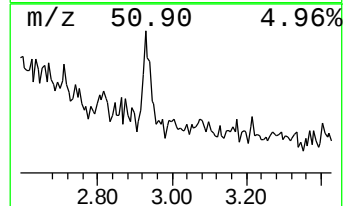
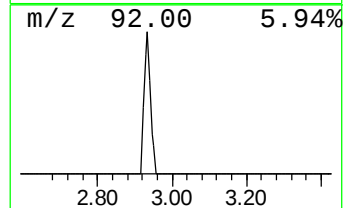
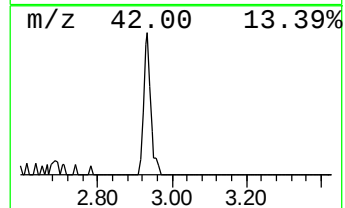
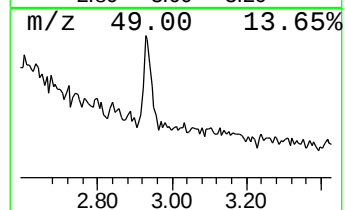
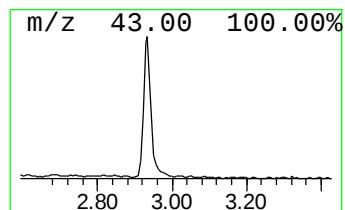
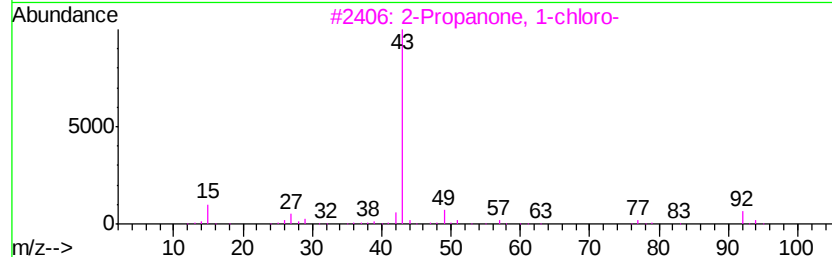
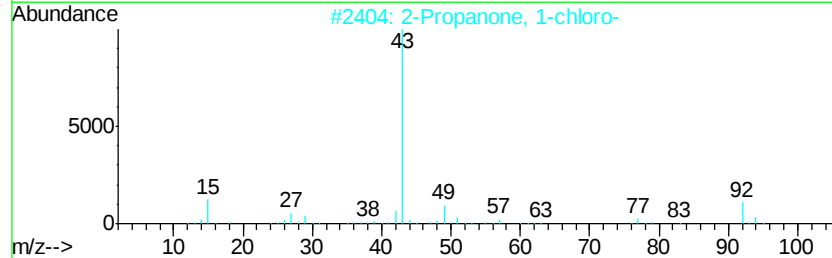
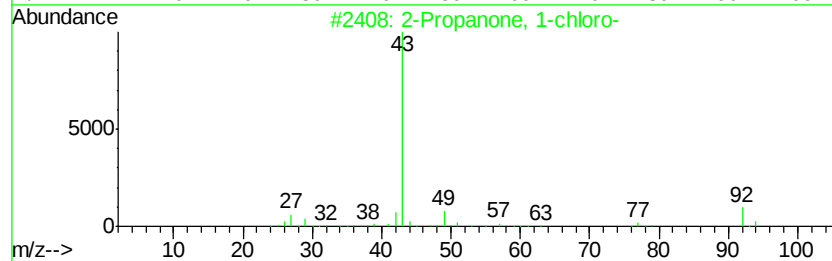
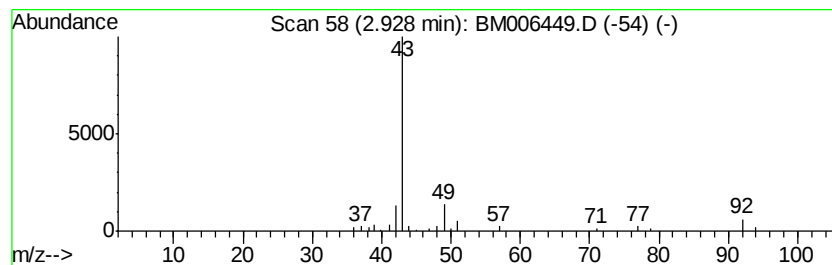
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 1 2-Propanone, 1-chloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.93	2.07 ng/ul	77110	1,4-Dichlorobenzene-d4	7.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanone, 1-chloro-	92	C3H5ClO	000078-95-5	78
2			2-Propanone, 1-chloro-	92	C3H5ClO	000078-95-5	72
3			2-Propanone, 1-chloro-	92	C3H5ClO	000078-95-5	64
4			Chloroacetamidine	92	C2H5ClN2	020846-52-0	59
5			2-Propanone, 1-chloro-	92	C3H5ClO	000078-95-5	58



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Propanone, 1-ch...	2.93	2.1	ng/ul	77110	1	7.62	746371	20.0