

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081822\
 Data File : BP011485.D
 Acq On : 18 Aug 2022 16:56
 Operator : CG/JU
 Sample : N4221-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0KN3

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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_P\Methods\SFAM-EPA-BP081622.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011485.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.034	3	8	16	rBV	10175263	13895841	100.00%	34.323%
2	3.099	16	19	31	rVB	794785	955205	6.87%	2.359%
3	3.493	84	86	97	rBV	31205	47666	0.34%	0.118%
4	3.587	99	102	107	rVB	17294	21499	0.15%	0.053%
5	3.693	115	120	125	rVB	16218	23946	0.17%	0.059%
6	3.781	131	135	141	rVB2	12948	17130	0.12%	0.042%
7	3.905	151	156	162	rBV	22201	32723	0.24%	0.081%
8	4.299	216	223	236	rBV	5559676	7648837	55.04%	18.893%
9	5.022	341	346	353	rVB	37151	49014	0.35%	0.121%
10	6.746	633	639	648	rBV	78333	124705	0.90%	0.308%
11	6.904	660	666	680	rBV	292186	454912	3.27%	1.124%
12	7.093	689	698	710	rBV	308965	490644	3.53%	1.212%
13	7.557	770	777	785	rBV	320264	482747	3.47%	1.192%
14	7.634	785	790	794	rBV4	8726	14768	0.11%	0.036%
15	8.275	893	899	910	rBV	214238	354130	2.55%	0.875%
16	8.722	966	975	985	rBV	368134	598692	4.31%	1.479%
17	9.440	1089	1097	1108	rBV	301625	485727	3.50%	1.200%
18	9.975	1180	1188	1199	rBV	385141	644030	4.63%	1.591%
19	10.345	1244	1251	1262	rBV	402113	647729	4.66%	1.600%
20	10.498	1270	1277	1292	rBV	209639	375823	2.70%	0.928%
21	11.169	1385	1391	1399	rBV8	7300	14906	0.11%	0.037%
22	11.951	1519	1524	1533	rVB4	8066	14612	0.11%	0.036%
23	13.657	1805	1814	1827	rBV	719919	981717	7.06%	2.425%
24	13.869	1845	1850	1854	rBV	108940	185041	1.33%	0.457%
25	13.922	1854	1859	1871	rVB	727554	1103956	7.94%	2.727%
26	14.228	1905	1911	1919	rBV	523556	758535	5.46%	1.874%
27	14.469	1946	1952	1963	rBV	39120	87018	0.63%	0.215%
28	14.998	2037	2042	2048	rBV	16067	21867	0.16%	0.054%
29	15.080	2052	2056	2063	rVB2	12338	17201	0.12%	0.042%
30	15.233	2075	2082	2090	rBV	981671	1334082	9.60%	3.295%
31	15.369	2096	2105	2116	rBV	372680	500651	3.60%	1.237%
32	17.004	2377	2383	2392	rVB	604139	848723	6.11%	2.096%
33	17.104	2392	2400	2411	rBV	1125430	1552505	11.17%	3.835%
34	17.704	2497	2502	2506	rBV	20487	26442	0.19%	0.065%
35	17.921	2535	2539	2543	rBV3	28666	41274	0.30%	0.102%
36	18.939	2708	2712	2716	rVB4	13873	18113	0.13%	0.045%

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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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_P\Methods\SFAM-EPA-BP081622.M
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37	19.010	2721	2724	2728	rVB	14490	18309	0.13%	0.045%
38	19.168	2748	2751	2756	rBV6	11513	14658	0.11%	0.036%
39	19.363	2779	2784	2793	rBV	1312394	1707326	12.29%	4.217%
40	20.868	3036	3040	3046	rBV5	42655	83274	0.60%	0.206%
41	21.133	3080	3085	3091	rBV	780874	948320	6.82%	2.342%
42	23.280	3443	3450	3457	rBV	1031939	1824941	13.13%	4.508%
43	23.427	3469	3475	3485	rVB	550888	1016701	7.32%	2.511%

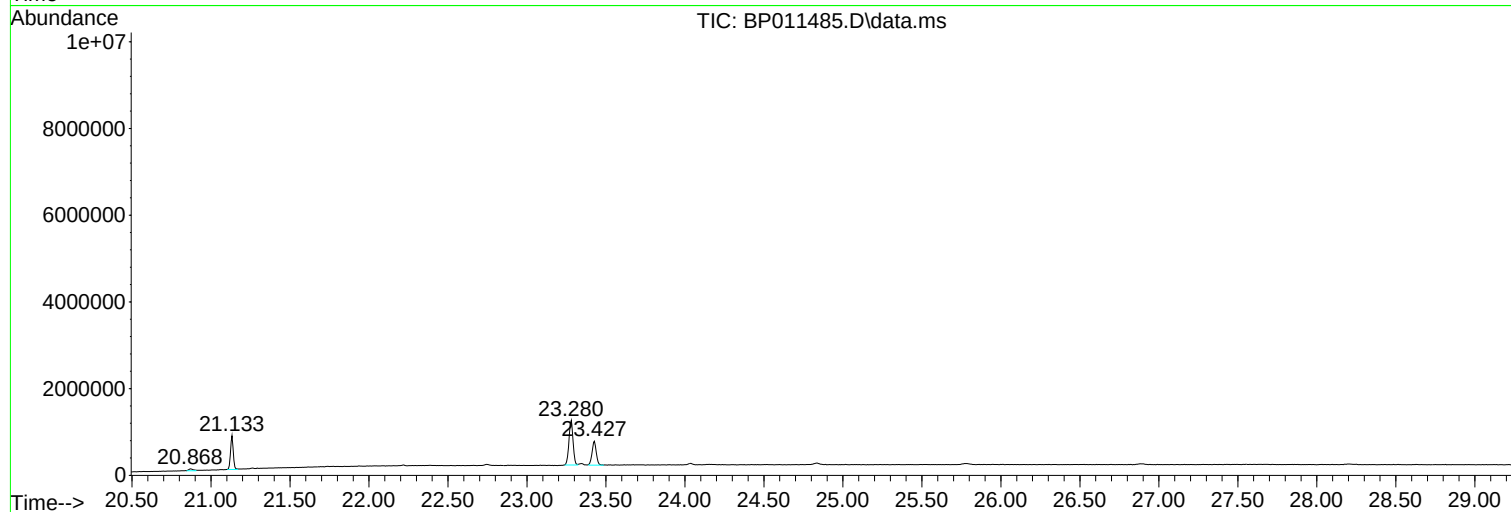
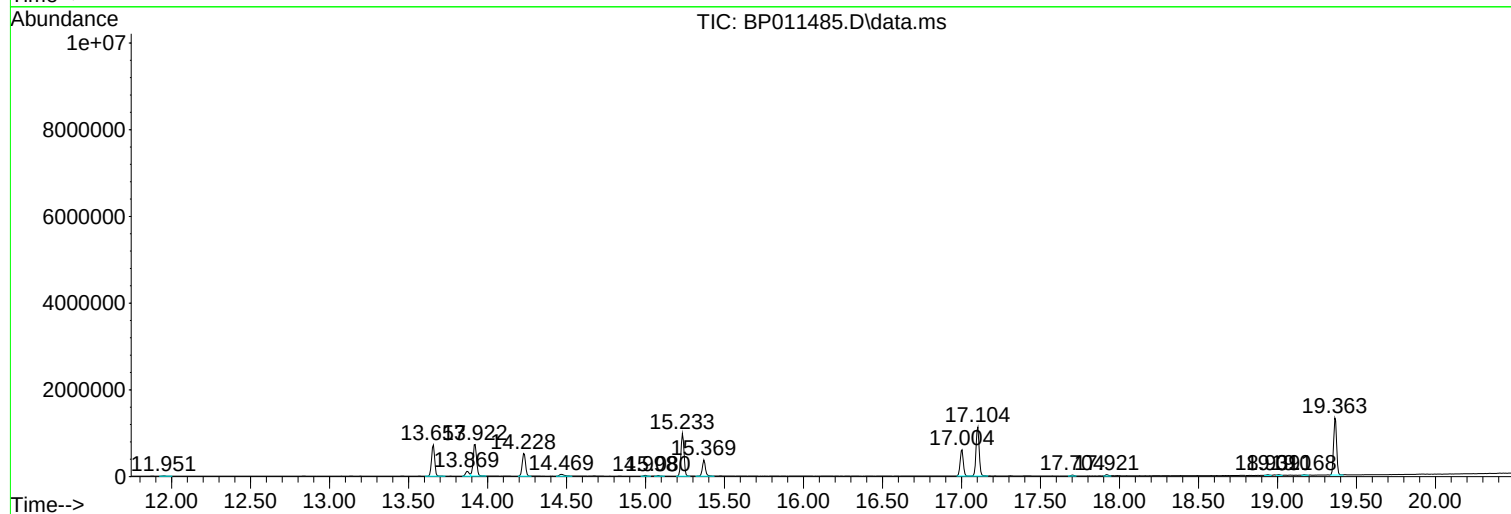
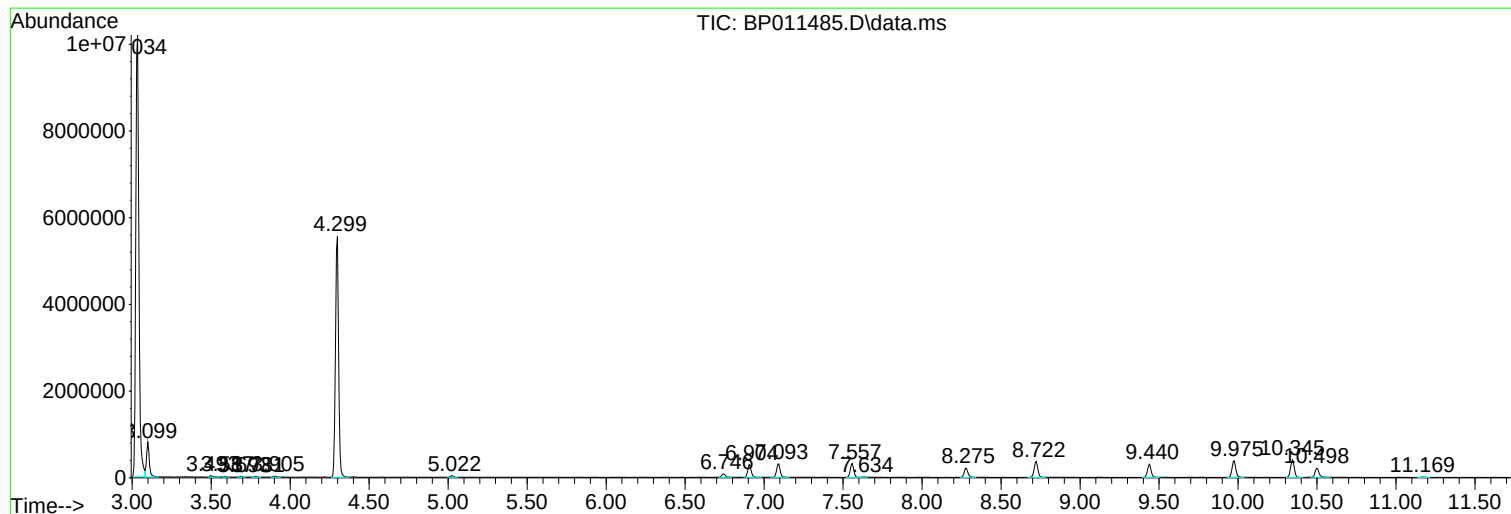
Sum of corrected areas: 40485940

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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



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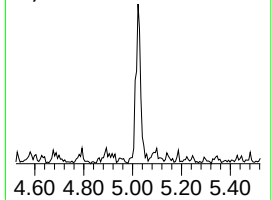
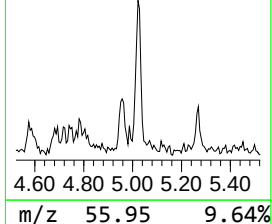
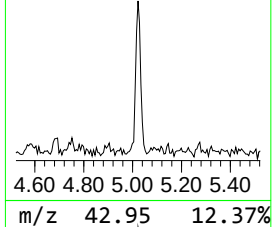
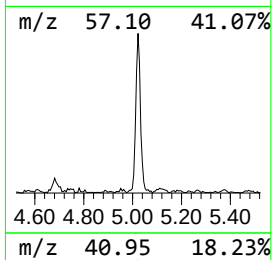
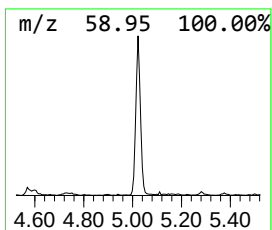
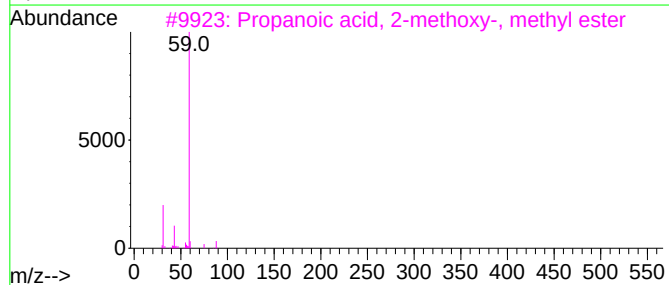
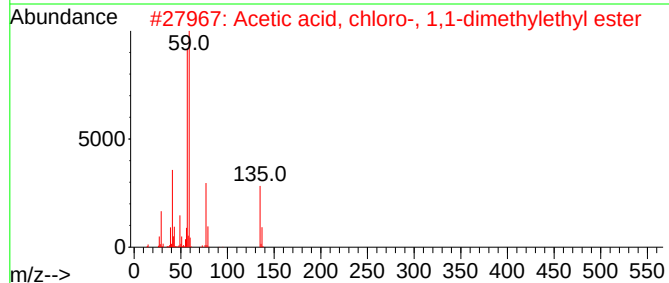
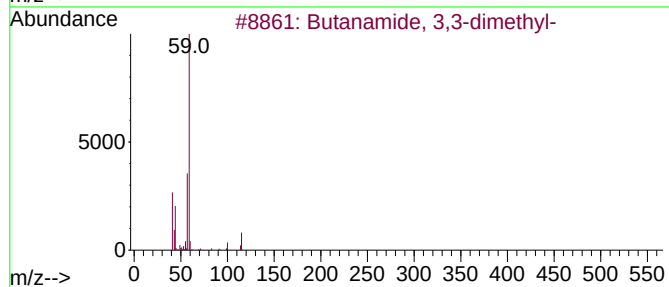
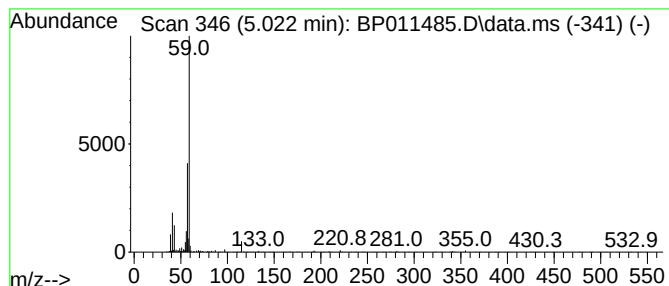
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.022	2.03 ng/ul	49014	1,4-Dichlorobenzene-d4	7.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanamide, 3,3-dimethyl-	115	C6H13NO	000926-04-5	45
2			Acetic acid, chloro-, 1,1-dimeth...	150	C6H11ClO2	000107-59-5	38
3			Propanoic acid, 2-methoxy-, meth...	118	C5H10O3	017639-76-8	36
4			N-Chloroacetyl-dl-erythro-O-meth...	209	C7H12ClNO4	1010214-48-1	36
5			Ether, hexyl t-butyl	158	C10H22O	069775-79-7	23



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	5.022	2.0	ng/u1	49014	1	7.557	482747	20.0