

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122721\
 Data File : BP008557.D
 Acq On : 27 Dec 2021 15:56
 Operator : CG/JU
 Sample : M5201-15
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P080-TW001-A-01

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-BP122321.M
 Title : SVOA CALIBRATION

Signal : TIC: BP008557.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.017	3	5	15	rVB	1226738	1642571	12.41%	1.377%
2	3.093	15	18	35	rVB2	296725	591065	4.47%	0.495%
3	3.352	58	62	75	rVB	134297	230250	1.74%	0.193%
4	3.764	129	132	141	rBV	117849	180426	1.36%	0.151%
5	4.099	186	189	195	rVB2	29846	40831	0.31%	0.034%
6	7.058	686	692	702	rVB	504078	815251	6.16%	0.683%
7	7.228	714	721	740	rBV	1746208	2931408	22.15%	2.457%
8	7.434	748	756	770	rVV	1752879	2922347	22.08%	2.450%
9	7.540	772	774	782	rVB7	10599	18065	0.14%	0.015%
10	7.905	828	836	845	rBV	1646642	2736916	20.68%	2.294%
11	8.152	871	878	882	rBV9	7175	15552	0.12%	0.013%
12	8.605	948	955	971	rBV	1449266	2329666	17.60%	1.953%
13	9.058	1021	1032	1043	rBV	2057408	3393173	25.64%	2.844%
14	9.528	1107	1112	1117	rVB3	15051	23303	0.18%	0.020%
15	9.781	1147	1155	1163	rBV	1363842	2217786	16.76%	1.859%
16	10.316	1238	1246	1259	rBV	2456229	4224297	31.92%	3.541%
17	10.705	1304	1312	1321	rBV	2267851	3887453	29.37%	3.259%
18	10.834	1328	1334	1341	rVB2	27414	46974	0.35%	0.039%
19	12.140	1551	1556	1562	rVB9	9268	16892	0.13%	0.014%
20	12.287	1576	1581	1588	rVB	51129	89070	0.67%	0.075%
21	13.504	1784	1788	1792	rVB2	9499	13618	0.10%	0.011%
22	13.940	1854	1862	1870	rBV	5145519	7186225	54.30%	6.024%
23	14.222	1900	1910	1925	rBV	5952526	8691958	65.68%	7.286%
24	14.534	1955	1963	1971	rBV	3433498	5328879	40.27%	4.467%
25	14.710	1986	1993	1999	rBV	333119	511015	3.86%	0.428%
26	15.351	2096	2102	2108	rBV4	13460	23400	0.18%	0.020%
27	15.522	2124	2131	2141	rBV	7745817	11472500	86.69%	9.617%
28	15.634	2143	2150	2168	rVV	1275813	1845244	13.94%	1.547%
29	15.763	2168	2172	2179	rVB	37901	57750	0.44%	0.048%
30	16.310	2259	2265	2272	rBV	11239	20483	0.15%	0.017%
31	16.957	2372	2375	2382	rVB6	9648	13410	0.10%	0.011%
32	17.063	2388	2393	2399	rVB	17093	30908	0.23%	0.026%
33	17.281	2423	2430	2440	rBV	4144496	5751844	43.46%	4.822%
34	17.381	2440	2447	2466	rVV	8443505	12301773	92.95%	10.312%
35	17.575	2475	2480	2488	rBV	185766	230025	1.74%	0.193%
36	17.651	2489	2493	2499	rVB6	8633	15830	0.12%	0.013%

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37	17.957	2541	2545	2549	rBV6	13309	17343	0.13%	0.015%
38	18.169	2577	2581	2586	rBV2	61169	80698	0.61%	0.068%
39	19.186	2750	2754	2760	rBV3	98636	130196	0.98%	0.109%
40	19.257	2761	2766	2773	rVB	96528	139929	1.06%	0.117%
41	19.404	2787	2791	2795	rBV	60824	78062	0.59%	0.065%
42	19.610	2820	2826	2834	rBV	10173321	13234337	100.00%	11.094%
43	21.363	3119	3124	3135	rVB	4140828	5396425	40.78%	4.524%
44	23.633	3503	3510	3522	rBV2	6299499	12534974	94.72%	10.508%
45	23.792	3530	3537	3556	rVB2	2872998	5833016	44.07%	4.890%

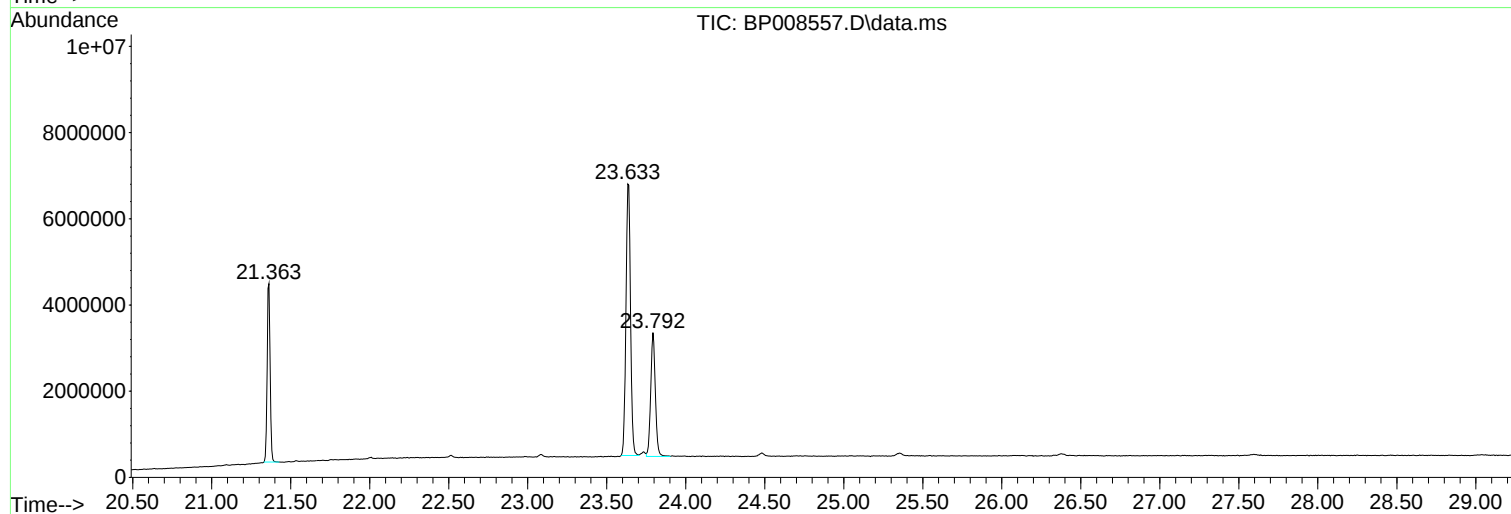
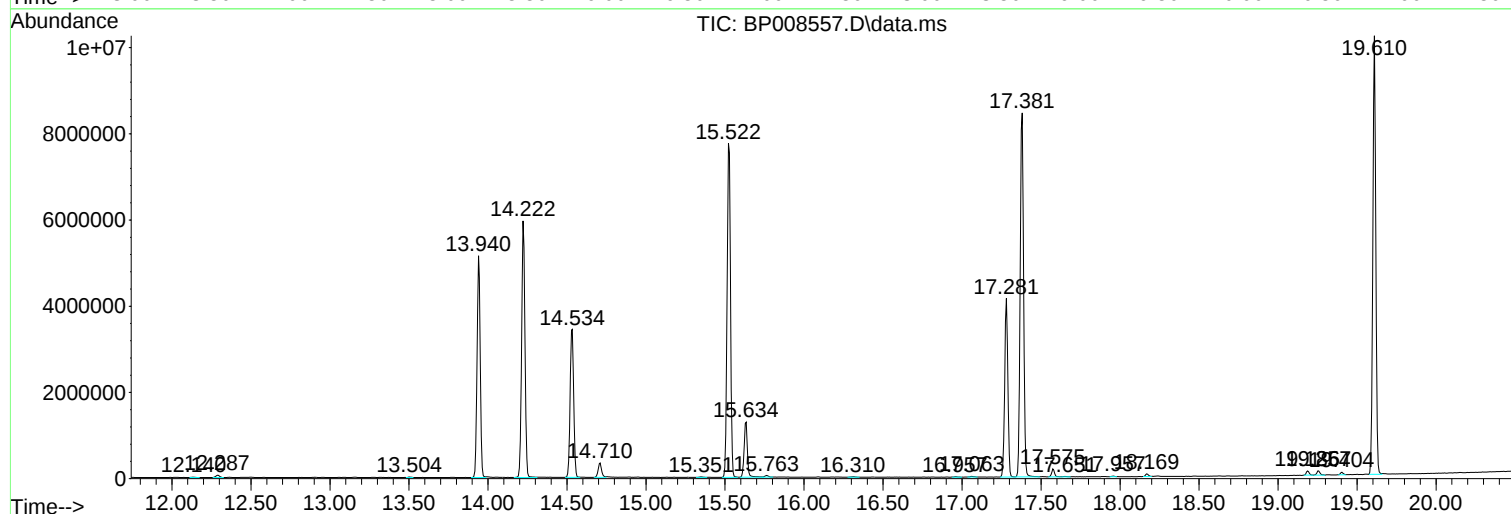
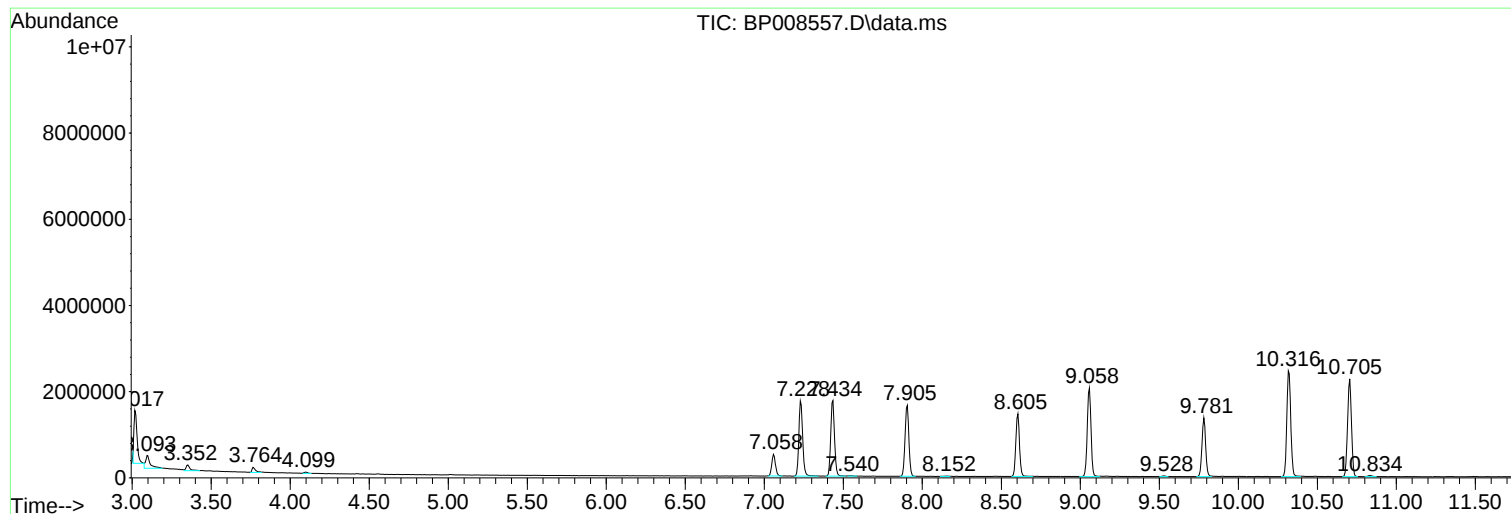
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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



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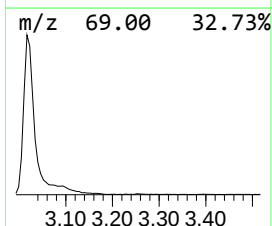
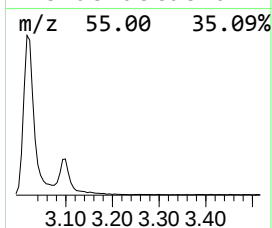
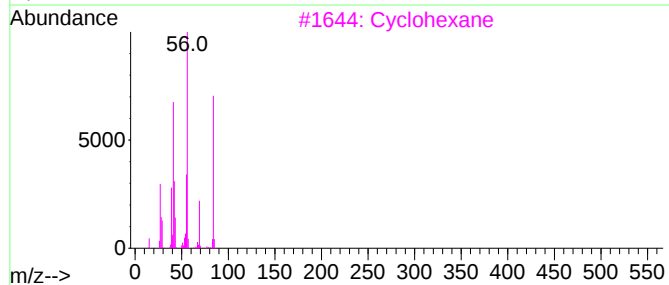
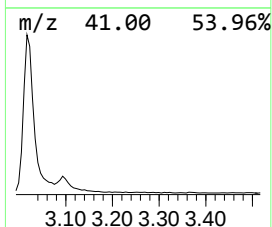
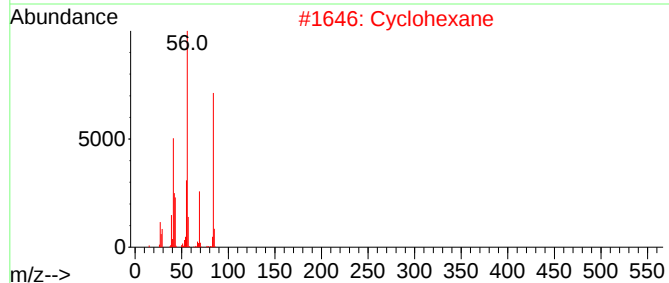
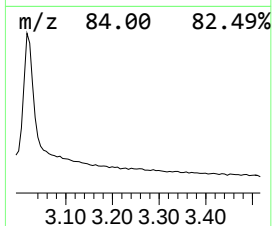
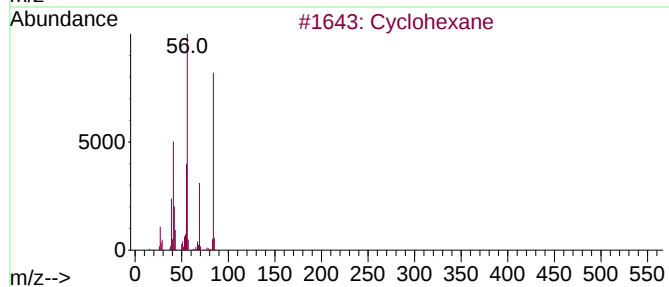
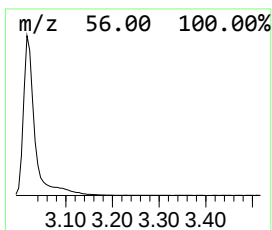
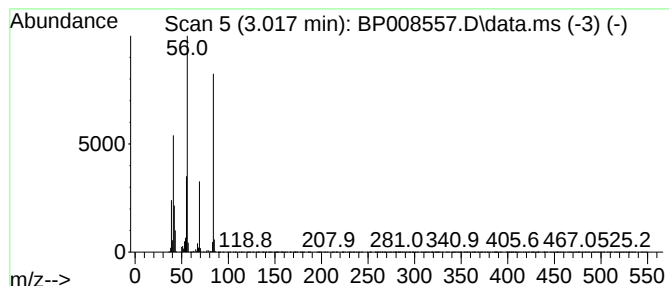
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.017 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.017	12.00 ng/ul	1642570	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane	84	C6H12	000110-82-7	91
2			Cyclohexane	84	C6H12	000110-82-7	91
3			Cyclohexane	84	C6H12	000110-82-7	91
4			Cyclohexane	84	C6H12	000110-82-7	90
5			Cyclopentane, methyl-	84	C6H12	000096-37-7	83



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Instrument :
BNA_P
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P080-TW001-A-01

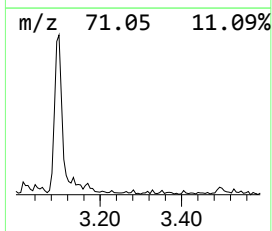
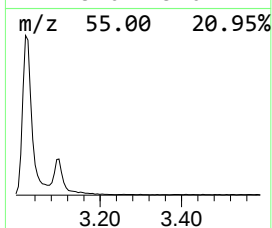
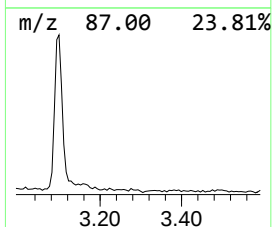
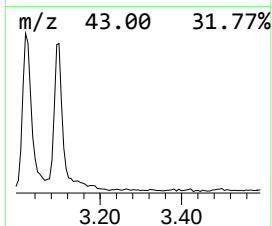
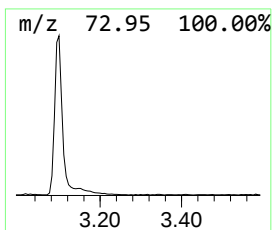
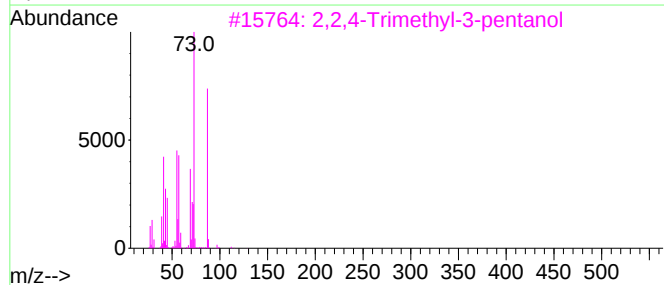
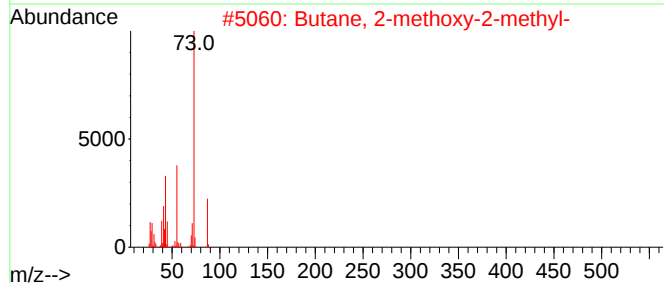
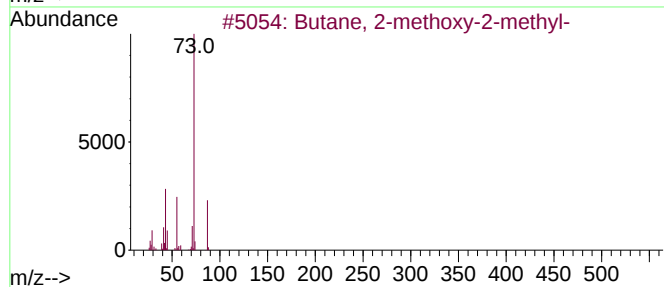
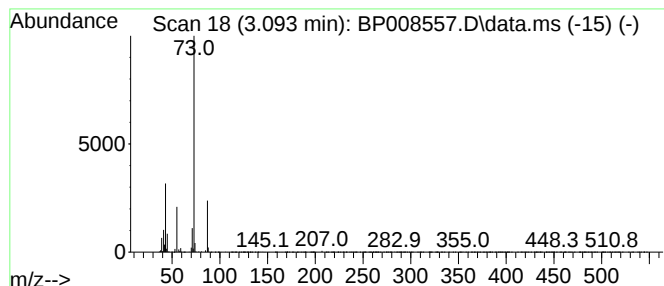
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.093	4.32 ng/ul	591065	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	36
5			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: C...	3.017	12.0	ng/ul	1642570	1	7.905	2736920	20.0
Butane, 2-metho...	3.093	4.3	ng/ul	591065	1	7.905	2736920	20.0