

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
 Data File : BP020802.D
 Acq On : 06 Jul 2024 01:13
 Operator : MA/JU
 Sample : P2964-04
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4B51

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-BP062524.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020802.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.052	9	11	22	rVB	104521	152917	5.19%	0.408%
2	3.269	45	48	58	rVB	24780	37838	1.28%	0.101%
3	3.540	90	94	106	rVB	64765	105027	3.56%	0.280%
4	3.687	114	119	129	rBV2	110475	174424	5.92%	0.466%
5	3.775	129	134	141	rVV	119447	168070	5.70%	0.449%
6	3.846	143	146	152	rVB	14933	21012	0.71%	0.056%
7	3.987	167	170	179	rVB2	12699	21182	0.72%	0.057%
8	4.146	191	197	204	rVB8	9709	17860	0.61%	0.048%
9	4.358	229	233	237	rVB4	6058	9060	0.31%	0.024%
10	4.487	251	255	258	rBV2	9428	14961	0.51%	0.040%
11	4.534	258	263	267	rVV	44346	65358	2.22%	0.175%
12	4.581	267	271	282	rVB	33634	56776	1.93%	0.152%
13	4.669	282	286	294	rBV5	11602	24364	0.83%	0.065%
14	4.793	299	307	315	rBV5	8548	23204	0.79%	0.062%
15	4.881	316	322	329	rBV	139236	212335	7.20%	0.567%
16	4.987	336	340	347	rVB2	15381	27598	0.94%	0.074%
17	5.846	482	486	494	rVB	14000	23042	0.78%	0.062%
18	6.010	509	514	524	rVB	5361	12243	0.42%	0.033%
19	6.134	532	535	541	rVB8	4669	9066	0.31%	0.024%
20	6.287	558	561	566	rVB3	12438	16493	0.56%	0.044%
21	6.375	570	576	580	rBV7	8949	16347	0.55%	0.044%
22	6.610	611	616	625	rVB7	9117	23242	0.79%	0.062%
23	6.716	631	634	638	rVV2	13211	15904	0.54%	0.042%
24	6.763	639	642	648	rVB7	5368	9771	0.33%	0.026%
25	6.916	655	668	679	rVV	382626	657323	22.30%	1.755%
26	7.069	686	694	705	rBV	313075	527822	17.91%	1.410%
27	7.210	715	718	721	rBV4	6525	8790	0.30%	0.023%
28	7.269	722	728	736	rVV	465926	694264	23.56%	1.854%
29	7.340	736	740	748	rVB2	21312	45683	1.55%	0.122%
30	7.534	769	773	780	rVB8	4056	8246	0.28%	0.022%
31	7.622	787	788	794	rVV6	4266	7159	0.24%	0.019%
32	7.722	794	805	812	rVV	695885	1164308	39.51%	3.109%
33	7.787	814	816	821	rVB4	8292	12456	0.42%	0.033%
34	7.993	848	851	856	rVV3	9909	17577	0.60%	0.047%
35	8.357	906	913	918	rBV7	5809	9353	0.32%	0.025%
36	8.428	918	925	932	rBV	298249	508383	17.25%	1.358%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
 Title : SVOA CALIBRATION

37	8.540	940	944	950	rVV	20638	37113	1.26%	0.099%
38	8.804	984	989	994	rVB9	5346	7591	0.26%	0.020%
39	8.875	994	1001	1013	rBV	381964	671564	22.79%	1.793%
40	8.999	1018	1022	1026	rVV6	8941	18574	0.63%	0.050%
41	9.034	1026	1028	1037	rVB9	6994	11817	0.40%	0.032%
42	9.246	1060	1064	1072	rVB4	7566	11624	0.39%	0.031%
43	9.428	1091	1095	1106	rVB	41637	65485	2.22%	0.175%
44	9.587	1114	1122	1129	rBV	319312	566072	19.21%	1.512%
45	9.822	1159	1162	1169	rVB9	5197	10271	0.35%	0.027%
46	10.128	1204	1214	1236	rBV	488585	891286	30.24%	2.380%
47	10.498	1266	1277	1289	rBV	989416	1660788	56.35%	4.435%
48	10.634	1293	1300	1317	rVV	195141	396939	13.47%	1.060%
49	11.028	1361	1367	1377	rBV	12894	33033	1.12%	0.088%
50	11.240	1395	1403	1410	rBV	99234	191639	6.50%	0.512%
51	11.910	1510	1517	1522	rBV	4074	7985	0.27%	0.021%
52	12.081	1541	1546	1551	rVV3	9699	17668	0.60%	0.047%
53	12.310	1578	1585	1586	rBV7	4154	7784	0.26%	0.021%
54	12.334	1586	1589	1592	rVV3	4945	8280	0.28%	0.022%
55	13.157	1725	1729	1737	rVB7	6070	13376	0.45%	0.036%
56	13.528	1785	1792	1799	rBV9	10411	29226	0.99%	0.078%
57	13.675	1811	1817	1821	rBV9	5295	9915	0.34%	0.026%
58	13.763	1823	1832	1838	rBV	930183	1306184	44.32%	3.488%
59	14.004	1868	1873	1874	rBV5	12072	16222	0.55%	0.043%
60	14.051	1874	1881	1894	rVB	1106600	1593003	54.05%	4.254%
61	14.169	1899	1901	1908	rVB8	7219	12752	0.43%	0.034%
62	14.245	1908	1914	1921	rVB5	8707	13663	0.46%	0.036%
63	14.357	1925	1933	1940	rVV	1500690	2215694	75.18%	5.917%
64	14.416	1940	1943	1949	rVB	15679	23736	0.81%	0.063%
65	14.492	1951	1956	1959	rBV	16144	23336	0.79%	0.062%
66	14.592	1964	1973	1991	rBV3	244995	560155	19.01%	1.496%
67	14.734	1993	1997	2003	rBV4	16642	35206	1.19%	0.094%
68	15.169	2065	2071	2077	rBV8	14283	38275	1.30%	0.102%
69	15.216	2077	2079	2082	rVB4	7208	7251	0.25%	0.019%
70	15.363	2098	2104	2111	rBV	1331154	1965947	66.71%	5.250%
71	15.422	2111	2114	2122	rVB	34202	52887	1.79%	0.141%
72	15.498	2122	2127	2138	rBV	294120	464732	15.77%	1.241%
73	15.628	2144	2149	2157	rVB	71054	110028	3.73%	0.294%
74	16.098	2225	2229	2231	rBV3	13817	18893	0.64%	0.050%
75	16.257	2254	2256	2258	rBV3	6684	7584	0.26%	0.020%
76	16.375	2272	2276	2282	rVB	36839	54248	1.84%	0.145%

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77	16.686	2324	2329	2338	rBV5	69999	135636	4.60%	0.362%
78	16.980	2376	2379	2384	rVB	16519	20581	0.70%	0.055%
79	17.039	2386	2389	2391	rBV3	29112	29230	0.99%	0.078%
80	17.069	2391	2394	2398	rVB	34255	43438	1.47%	0.116%
81	17.163	2404	2410	2414	rBV	1778094	2599033	88.19%	6.941%
82	17.204	2414	2417	2422	rVB	354783	497975	16.90%	1.330%
83	17.263	2422	2427	2438	rVB2	1273385	2020754	68.57%	5.396%
84	17.345	2438	2441	2446	rBV	56546	89786	3.05%	0.240%
85	17.627	2487	2489	2492	rBV4	17463	21477	0.73%	0.057%
86	17.780	2510	2515	2521	rBV3	52671	107576	3.65%	0.287%
87	18.104	2562	2570	2579	rBV2	322450	754680	25.61%	2.015%
88	18.269	2594	2598	2603	rBV2	91902	132191	4.49%	0.353%
89	18.386	2614	2618	2620	rBV3	45396	66446	2.25%	0.177%
90	18.874	2694	2701	2704	rBV3	206342	384566	13.05%	1.027%
91	19.298	2768	2773	2779	rVB	847694	1164994	39.53%	3.111%
92	19.433	2792	2796	2805	rVB2	164898	275397	9.34%	0.735%
93	19.592	2820	2823	2827	rBV5	103960	169628	5.76%	0.453%
94	19.645	2827	2832	2836	rBV	1622689	2446930	83.03%	6.534%
95	21.286	3109	3111	3118	rVB	476780	677870	23.00%	1.810%
96	21.616	3161	3167	3172	rBV3	1666049	2947154	100.00%	7.870%
97	21.663	3173	3175	3180	rVB	273730	349831	11.87%	0.934%
98	22.498	3315	3317	3323	rBV3	218159	316465	10.74%	0.845%
99	24.727	3691	3696	3704	rVB3	536350	1322400	44.87%	3.531%
100	24.968	3729	3737	3746	rBV	1036336	2765536	93.84%	7.385%

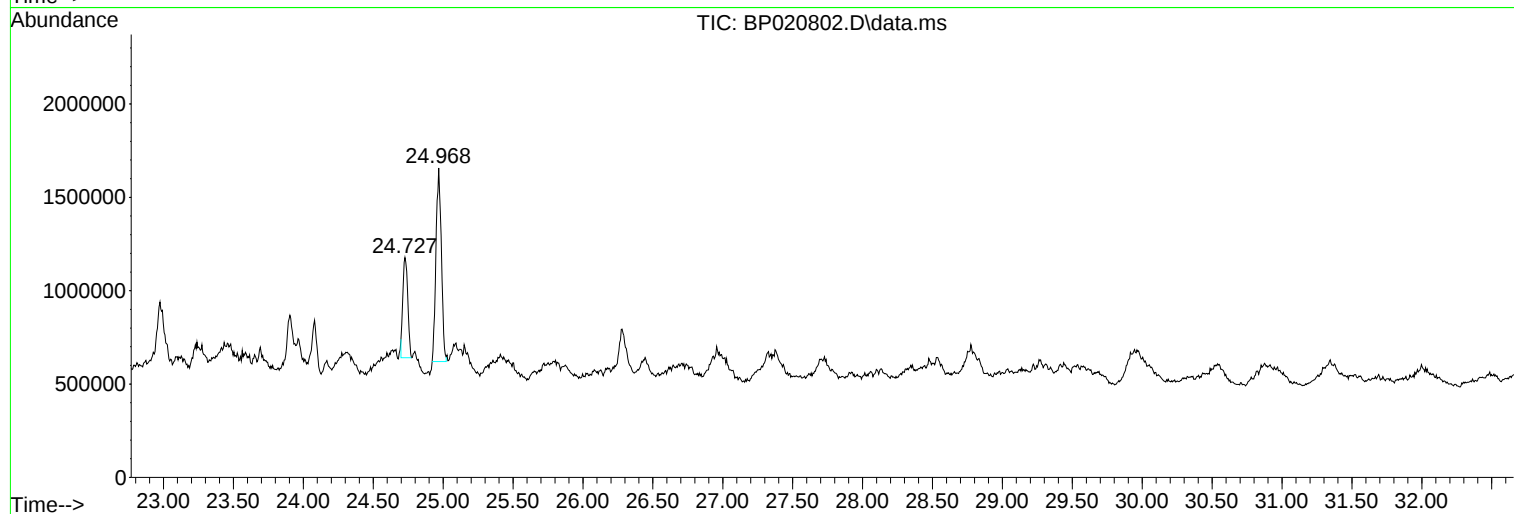
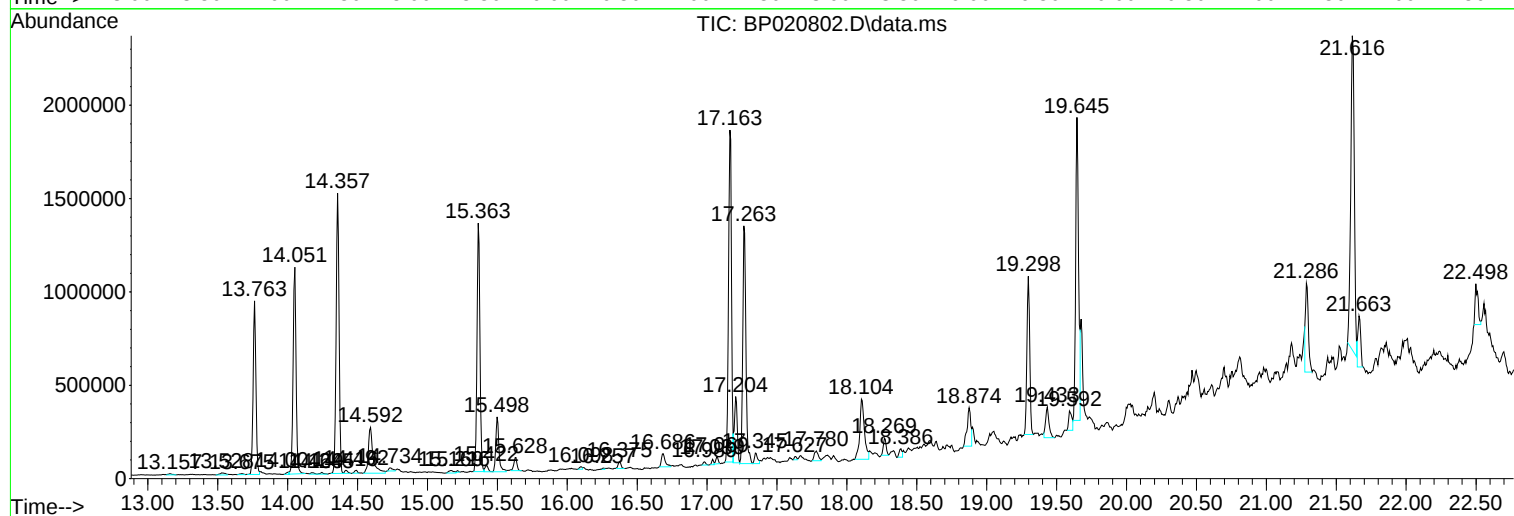
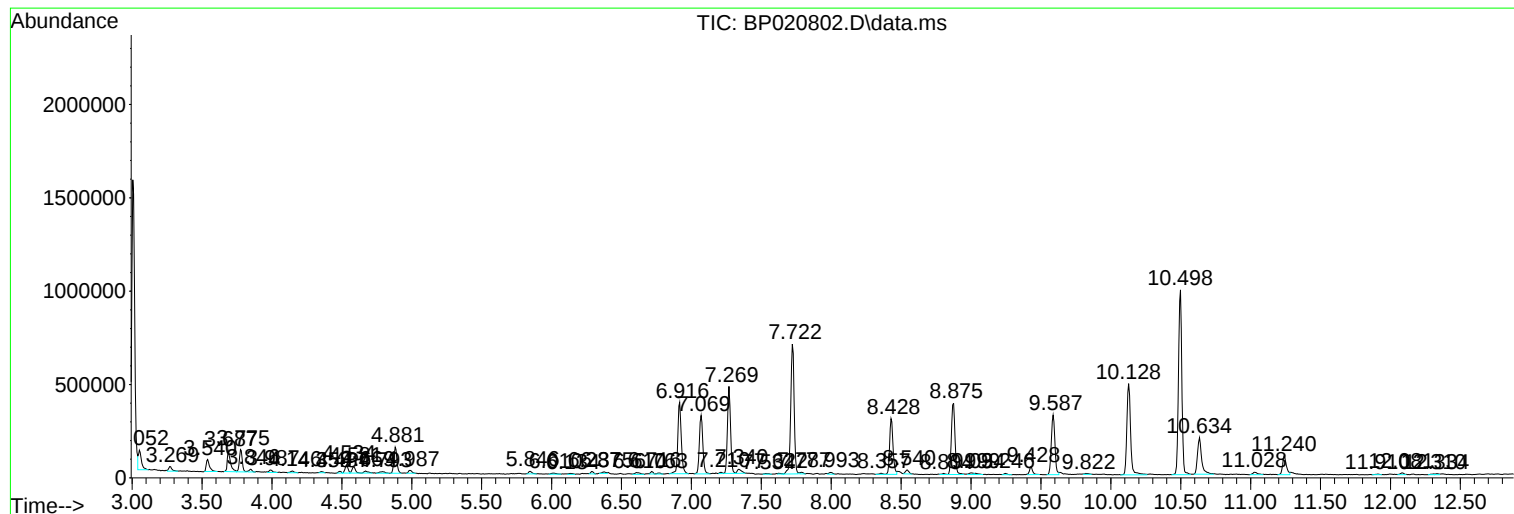
Sum of corrected areas: 37446858

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Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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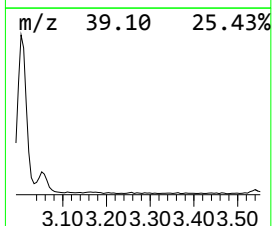
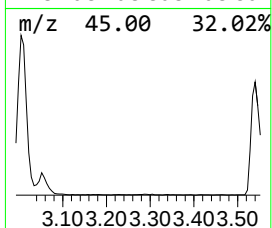
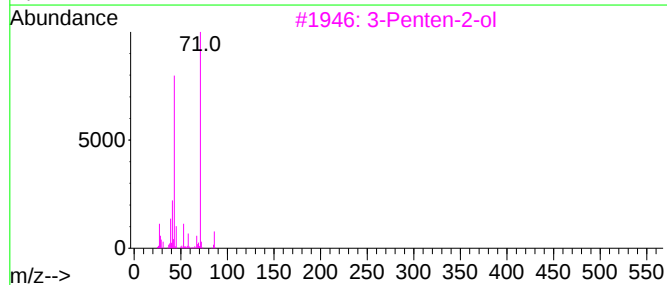
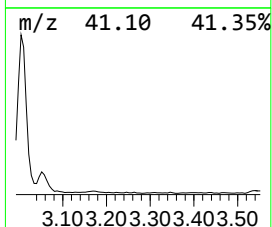
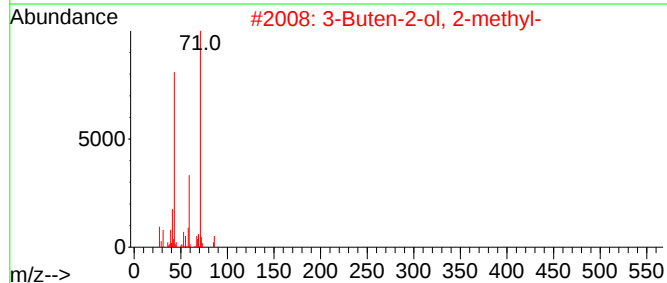
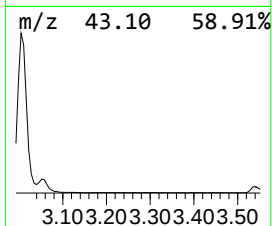
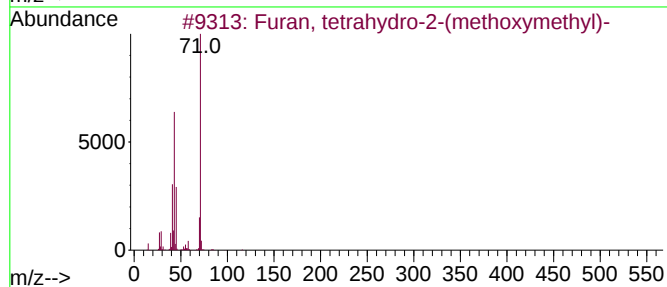
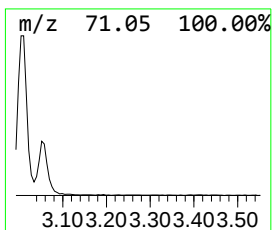
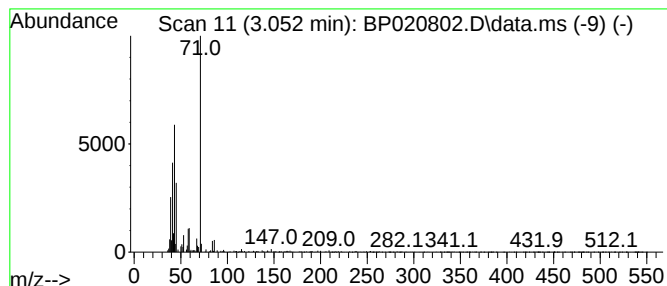
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Furan, tetrahydro-2-(methox... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.052	2.63 ng/ul	152917	1,4-Dichlorobenzene-d4	7.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Furan, tetrahydro-2-(methoxymeth...	116	C6H12O2	019354-27-9	59
2			3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	58
3			3-Penten-2-ol	86	C5H10O	001569-50-2	58
4			Z-1-Methoxy-2-pentene	100	C6H12O	1000279-59-4	56
5			Trichloroacetic acid, 2-tetrahyd...	246	C7H9Cl3O3	004743-27-5	50



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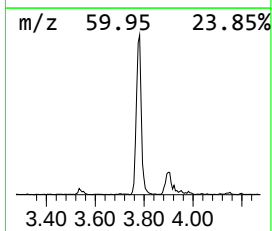
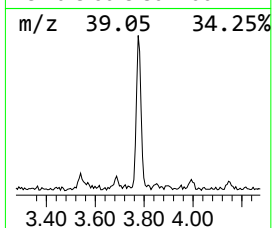
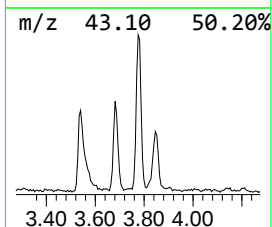
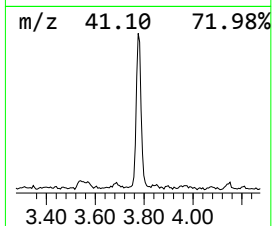
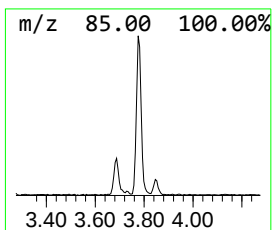
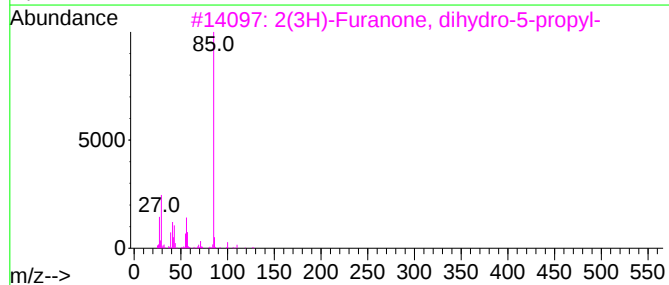
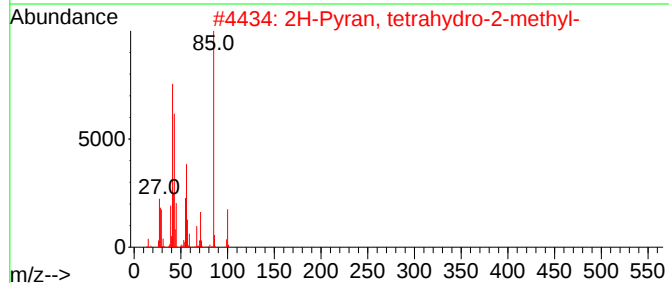
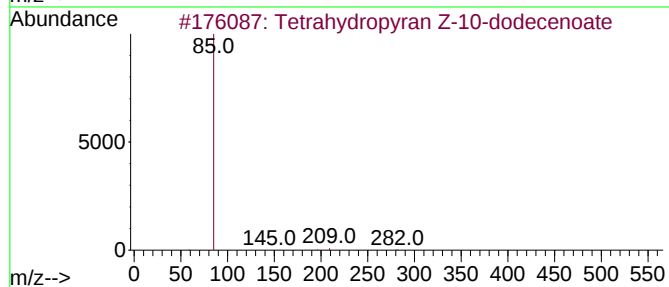
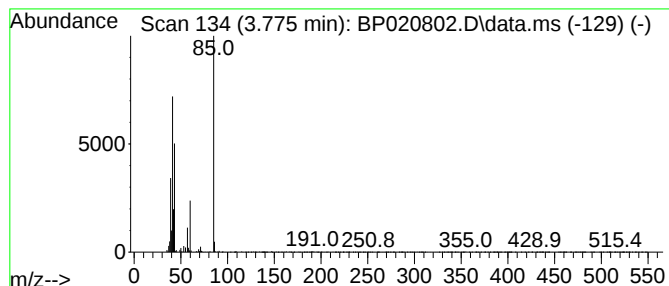
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Tetrahydropyran Z-10-dodece... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
3.775	2.89 ng/ul	168070	1,4-Dichlorobenzene-d4	7.722	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetrahydropyran Z-10-dodecenoate	282	C17H30O3	1000130-96-5	64
2	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	38
3	2(3H)-Furanone, dihydro-5-propyl-	128	C7H12O2	000105-21-5	38
4	2-(Chloromethyl)tetrahydropyran	134	C6H11ClO	018420-41-2	9
5	2-Pyrrolidinone	85	C4H7NO	000616-45-5	9



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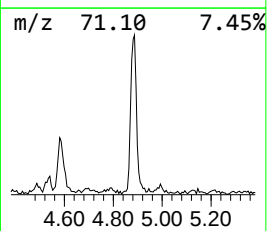
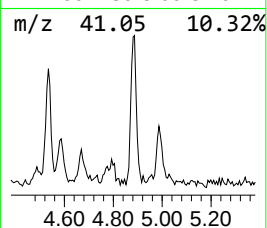
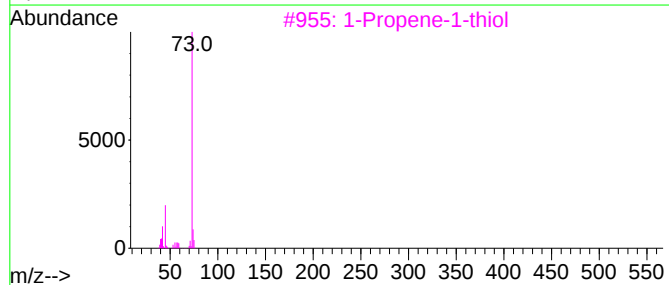
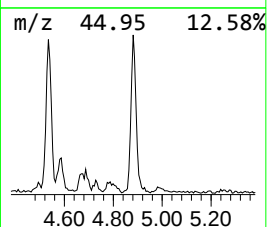
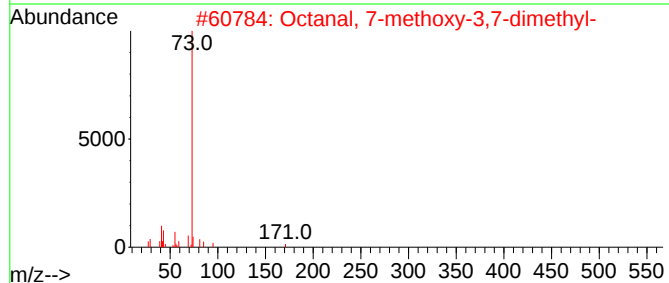
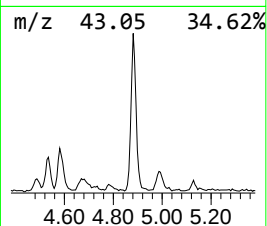
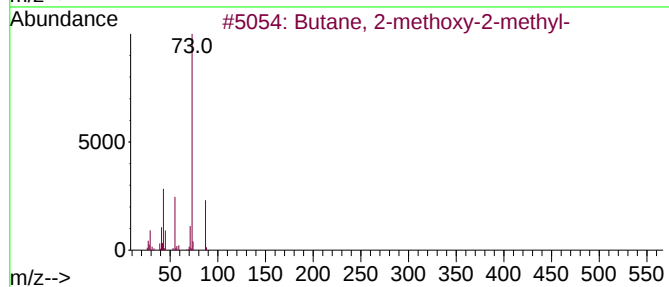
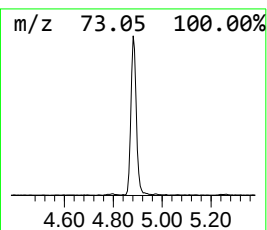
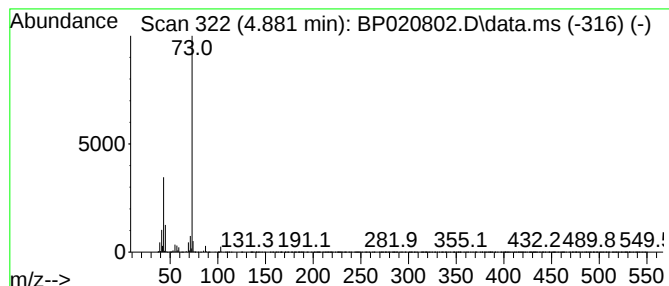
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.881	3.65 ng/ul	212335	1,4-Dichlorobenzene-d4	7.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	33
2			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	28
3			1-Propene-1-thiol	74	C3H6S	000925-89-3	9
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5			Isopropyl 5,11-dihydroxy-3,7,11-...	314	C18H34O4	1000223-34-1	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020802.D
Acq On : 06 Jul 2024 01:13
Operator : MA/JU
Sample : P2964-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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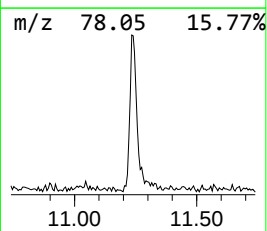
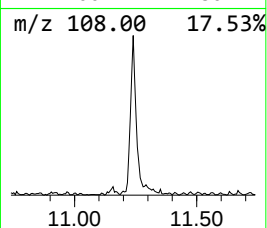
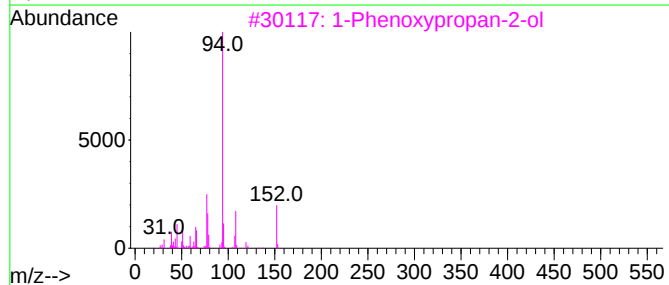
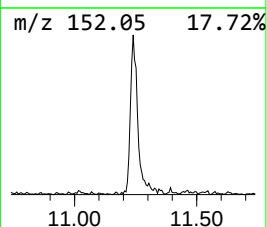
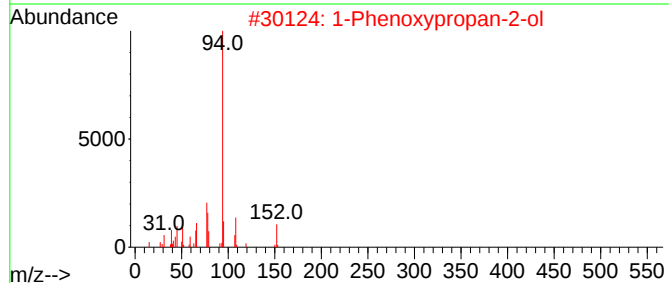
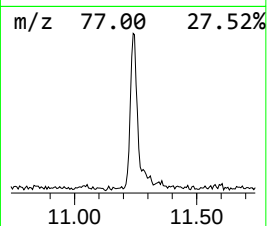
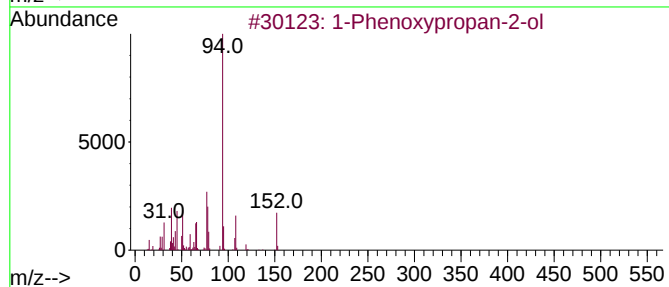
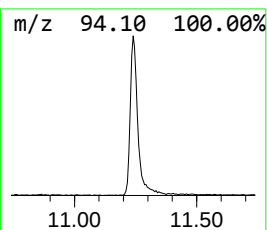
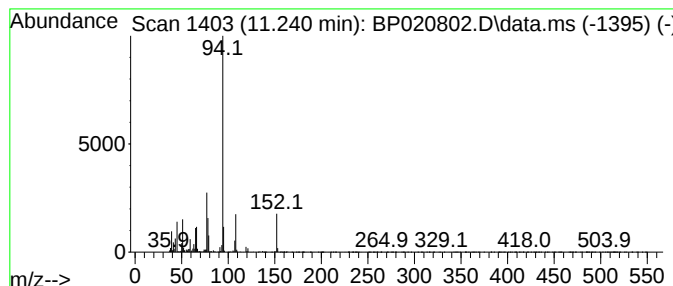
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.240	2.31 ng/ul	191639	Naphthalene-d8	10.499

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020802.D
Acq On : 06 Jul 2024 01:13
Operator : MA/JU
Sample : P2964-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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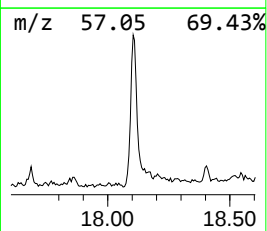
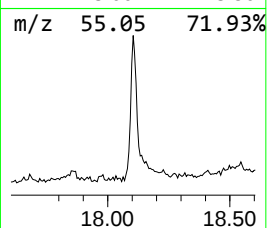
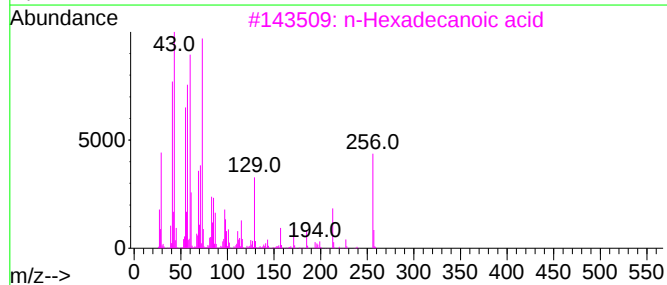
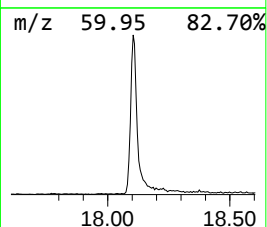
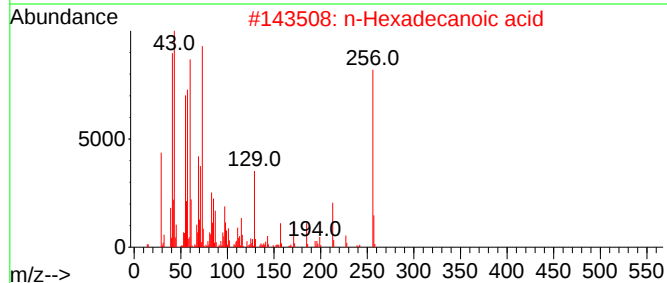
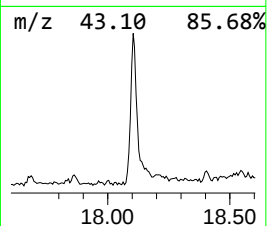
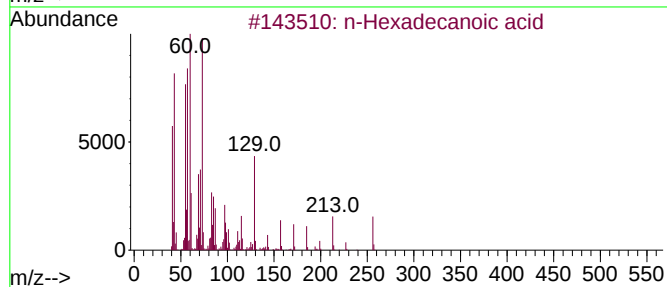
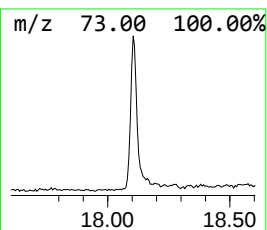
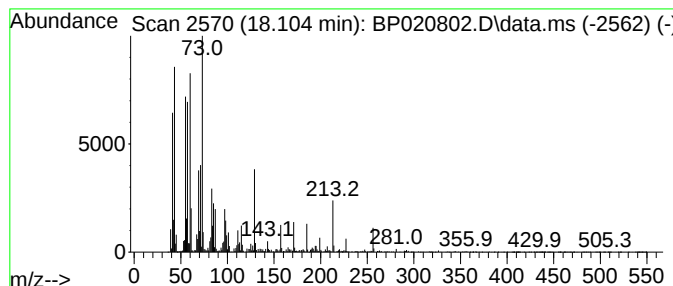
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.104	5.81 ng/ul	754680	Phenanthrene-d10	17.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020802.D
Acq On : 06 Jul 2024 01:13
Operator : MA/JU
Sample : P2964-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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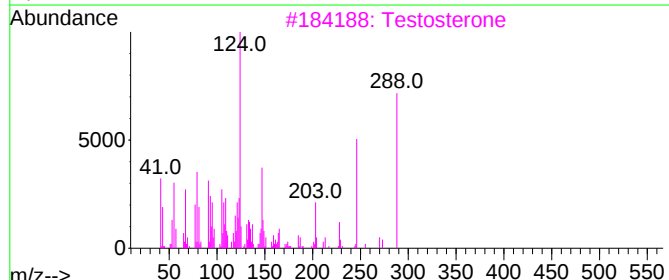
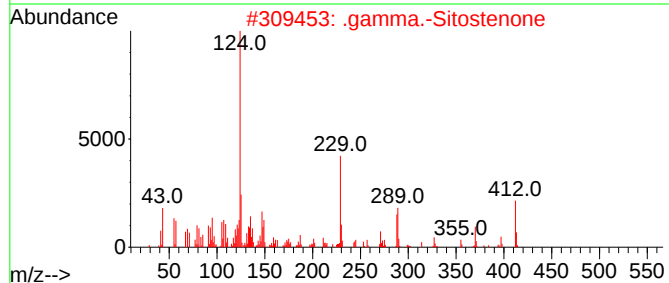
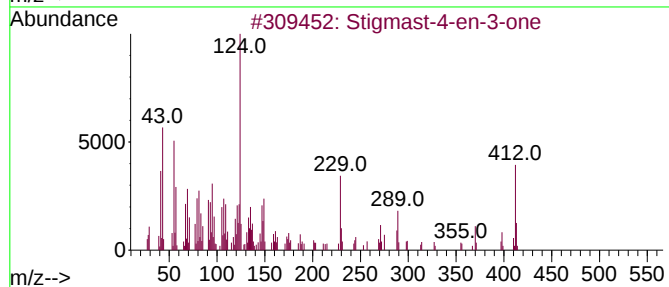
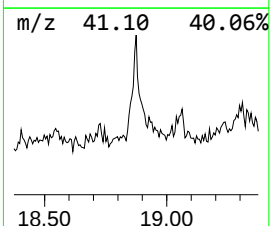
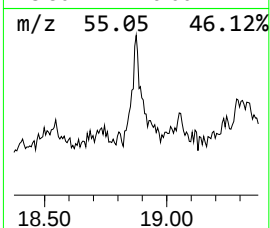
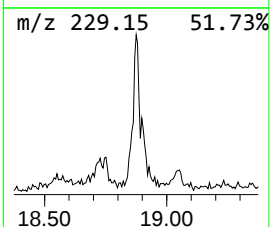
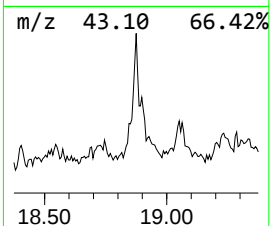
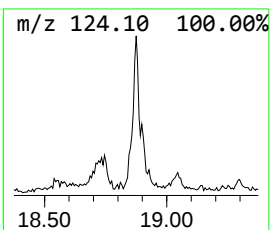
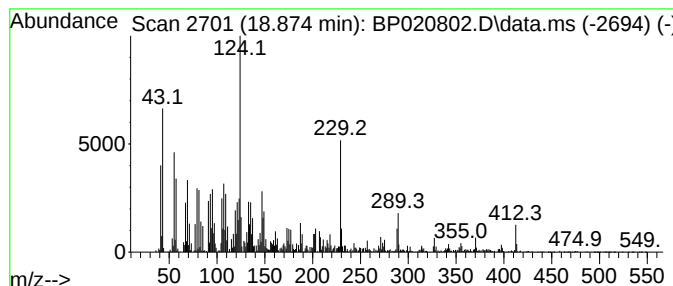
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Stigmast-4-en-3-one Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.875	2.96 ng/ul	384566	Phenanthrene-d10	17.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Stigmast-4-en-3-one	412	C29H48O	001058-61-3	95
2			.gamma.-Sitostenone	412	C29H48O	084924-96-9	90
3			Testosterone	288	C19H28O2	000058-22-0	86
4			Testosterone	288	C19H28O2	000058-22-0	70
5			Androst-4-en-3-one, 17-hydroxy-,...	288	C19H28O2	000604-39-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020802.D
Acq On : 06 Jul 2024 01:13
Operator : MA/JU
Sample : P2964-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

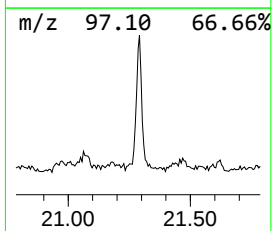
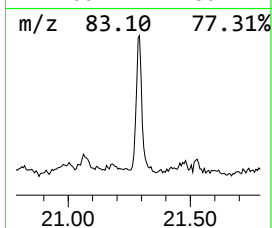
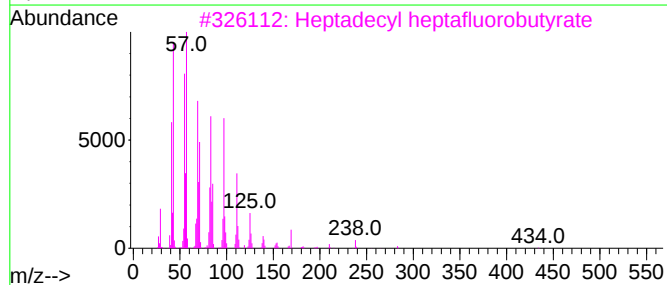
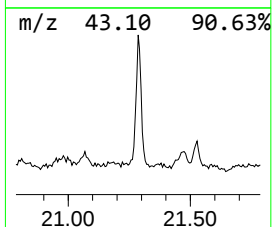
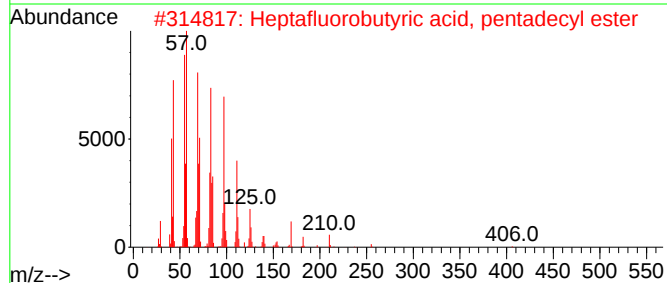
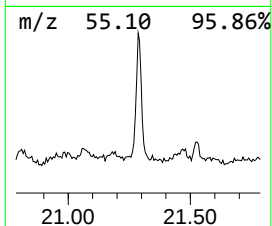
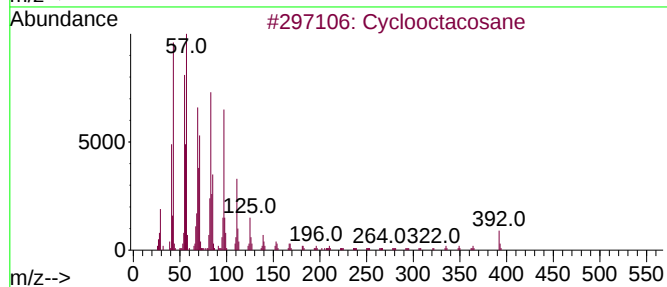
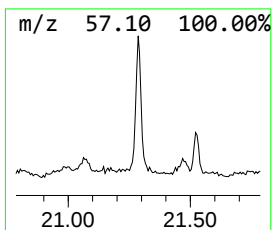
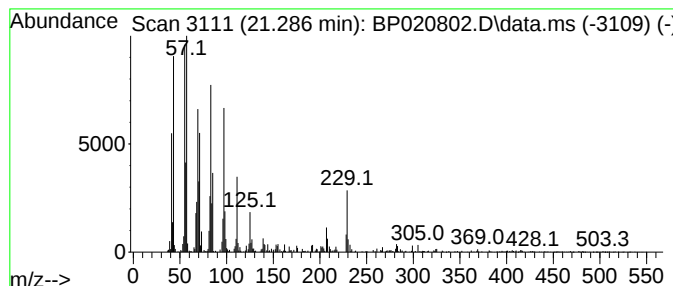
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Cyclic21.286 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.286	4.60 ng/ul	677870	Chrysene-d12	21.616	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclooctacosane	392	C28H56	000297-24-5	94
2	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93
3	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	90
4	9-Tricosene, (Z)-	322	C23H46	027519-02-4	90
5	Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020802.D
Acq On : 06 Jul 2024 01:13
Operator : MA/JU
Sample : P2964-04
Misc :
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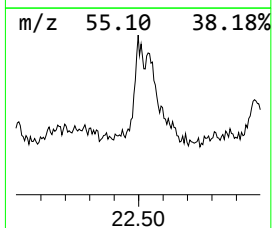
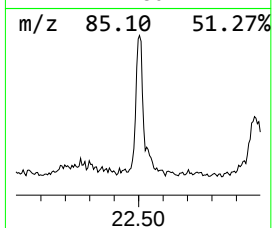
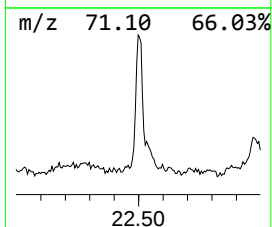
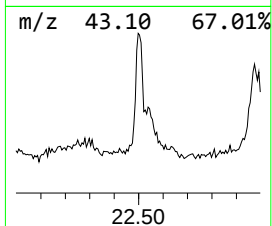
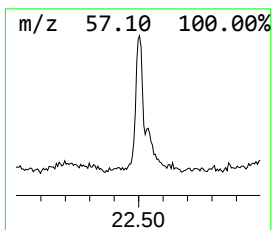
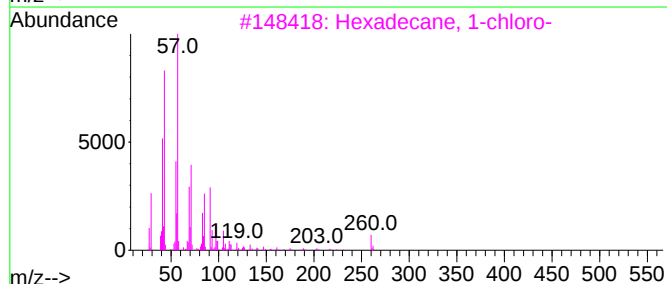
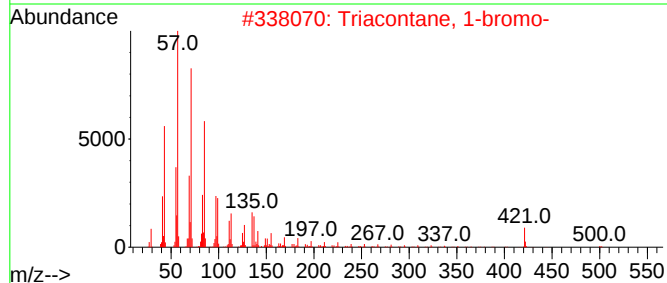
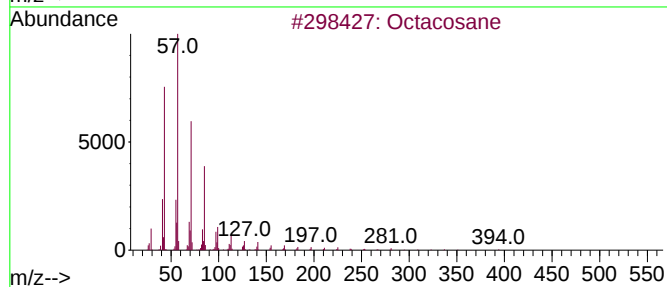
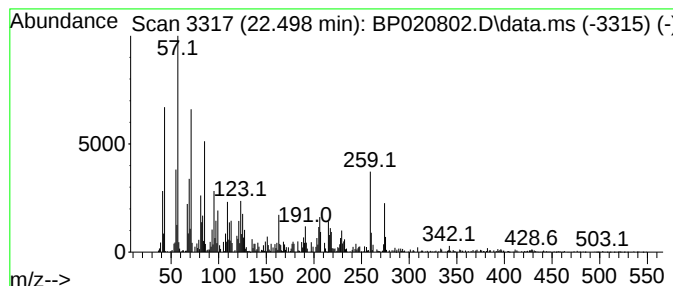
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.498	2.15 ng/ul	316465	Chrysene-d12		21.616	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octacosane		394	C28H58	000630-02-4	93
2	Triacontane, 1-bromo-		500	C30H61Br	004209-22-7	56
3	Hexadecane, 1-chloro-		260	C16H33Cl	004860-03-1	53
4	Docosane, 11-butyl-		366	C26H54	013475-76-8	47
5	2-Methylhexacosane		380	C27H56	001561-02-0	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020802.D
Acq On : 06 Jul 2024 01:13
Operator : MA/JU
Sample : P2964-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Furan, tetrahyd...	3.052	2.6	ng/u1	152917	1	7.722	1164310	20.0
Tetrahydropyran...	3.775	2.9	ng/u1	168070	1	7.722	1164310	20.0
unknown-01	4.881	3.6	ng/u1	212335	1	7.722	1164310	20.0
1-Phenoxypropan...	11.240	2.3	ng/u1	191639	2	10.499	1660790	20.0
n-Hexadecanoic ...	18.104	5.8	ng/u1	754680	4	17.163	2599030	20.0
Stigmast-4-en-3...	18.875	3.0	ng/u1	384566	4	17.163	2599030	20.0
(DEL) Alkane: C...	21.286	4.6	ng/u1	677870	5	21.616	2947150	20.0
(DEL) Alkane: S...	22.498	2.1	ng/u1	316465	5	21.616	2947150	20.0