

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
 Data File : BP021097.D
 Acq On : 23 Jul 2024 13:38
 Operator : MA/JU
 Sample : P3238-18
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4CG8

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-BP071024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021097.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.193	32	35	42	rVB	26820	37335	0.86%	0.077%
2	3.487	81	85	91	rBV	26288	47433	1.09%	0.098%
3	3.634	109	110	118	rVV2	13545	22079	0.51%	0.046%
4	3.705	118	122	128	rVV	51910	71440	1.65%	0.148%
5	3.917	155	158	163	rVB	12986	16242	0.37%	0.034%
6	4.328	225	228	232	rBV2	7491	11406	0.26%	0.024%
7	4.464	247	251	259	rVB3	13998	24685	0.57%	0.051%
8	4.817	306	311	320	rVB	48246	77384	1.78%	0.160%
9	4.928	327	330	340	rVB6	8174	17565	0.41%	0.036%
10	5.699	452	461	466	rBV6	4980	12963	0.30%	0.027%
11	5.793	474	477	483	rVB5	5659	10656	0.25%	0.022%
12	6.764	636	642	650	rBV6	12543	28725	0.66%	0.059%
13	6.881	655	662	675	rVV	396916	777081	17.92%	1.609%
14	7.016	679	685	698	rVB	365192	582948	13.44%	1.207%
15	7.216	713	719	727	rBV	466666	780431	18.00%	1.615%
16	7.305	731	734	742	rVV2	11312	26883	0.62%	0.056%
17	7.405	745	751	762	rVB	92137	156853	3.62%	0.325%
18	7.669	789	796	805	rBV	667145	1117300	25.76%	2.313%
19	7.758	806	811	826	rVB	75622	150302	3.47%	0.311%
20	7.946	836	843	844	rBV6	5546	10279	0.24%	0.021%
21	8.181	877	883	890	rVV2	86864	147912	3.41%	0.306%
22	8.399	911	920	927	rBV	487549	914006	21.08%	1.892%
23	8.458	927	930	950	rVB4	44496	133473	3.08%	0.276%
24	8.828	986	993	1003	rBV	448137	750527	17.31%	1.554%
25	8.969	1013	1017	1021	rVB6	10079	16025	0.37%	0.033%
26	9.169	1048	1051	1060	rBV2	7707	16381	0.38%	0.034%
27	9.375	1080	1086	1094	rBV2	34449	63398	1.46%	0.131%
28	9.540	1105	1114	1126	rBV	371788	666120	15.36%	1.379%
29	9.705	1138	1142	1146	rBV6	5483	10343	0.24%	0.021%
30	9.787	1150	1156	1169	rBV5	19701	70551	1.63%	0.146%
31	10.087	1198	1207	1223	rBV	505552	1028073	23.71%	2.128%
32	10.434	1259	1266	1274	rBV	864023	1580627	36.45%	3.272%
33	10.605	1287	1295	1308	rVV	163345	365859	8.44%	0.757%
34	10.987	1353	1360	1366	rBV5	20958	37822	0.87%	0.078%
35	11.193	1388	1395	1404	rBV	89588	193258	4.46%	0.400%
36	11.275	1405	1409	1415	rVB8	5042	10110	0.23%	0.021%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
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37	11.493	1440	1446	1450	rBV4	14352	31228	0.72%	0.065%
38	11.534	1450	1453	1460	rVB4	14927	24621	0.57%	0.051%
39	12.040	1532	1539	1545	rVB4	11250	22653	0.52%	0.047%
40	12.104	1545	1550	1554	rBV4	7278	14368	0.33%	0.030%
41	12.893	1677	1684	1690	rVB10	8508	19775	0.46%	0.041%
42	13.246	1740	1744	1755	rVB6	11208	27710	0.64%	0.057%
43	13.504	1776	1788	1798	rBV	44398	144662	3.34%	0.299%
44	13.616	1801	1807	1812	rVB9	8606	20969	0.48%	0.043%
45	13.716	1812	1824	1836	rBV	910684	1479417	34.11%	3.062%
46	13.998	1861	1872	1888	rBV	1093366	1858144	42.85%	3.846%
47	14.198	1902	1906	1913	rBV9	9579	20982	0.48%	0.043%
48	14.310	1918	1925	1932	rVV	1569450	2174041	50.13%	4.500%
49	14.375	1932	1936	1942	rVB	33206	53716	1.24%	0.111%
50	14.534	1961	1963	1971	rVB	600402	900043	20.75%	1.863%
51	14.610	1971	1976	1991	rBV	243798	550201	12.69%	1.139%
52	14.722	1991	1995	2005	rVB7	36550	87236	2.01%	0.181%
53	15.104	2054	2060	2065	rBV5	12441	25700	0.59%	0.053%
54	15.328	2092	2098	2105	rBV	1449616	2233672	51.51%	4.624%
55	15.387	2105	2108	2118	rVB2	44348	75380	1.74%	0.156%
56	15.481	2118	2124	2136	rBV	406322	689626	15.90%	1.428%
57	15.687	2156	2159	2163	rBV	10523	18315	0.42%	0.038%
58	15.810	2175	2180	2182	rBV6	11440	16933	0.39%	0.035%
59	15.892	2190	2194	2198	rBV6	6726	11623	0.27%	0.024%
60	16.057	2218	2222	2224	rBV4	16184	22632	0.52%	0.047%
61	16.334	2264	2269	2273	rBV6	7553	12784	0.29%	0.026%
62	16.416	2275	2283	2286	rBV7	15832	36053	0.83%	0.075%
63	16.522	2297	2301	2307	rBV5	35416	63833	1.47%	0.132%
64	16.787	2341	2346	2351	rBV	25855	41038	0.95%	0.085%
65	16.845	2351	2356	2359	rBV2	62942	95484	2.20%	0.198%
66	16.939	2368	2372	2378	rVB2	41129	77021	1.78%	0.159%
67	17.028	2385	2387	2393	rVB7	21254	34215	0.79%	0.071%
68	17.128	2397	2404	2408	rBV	1795625	2692655	62.09%	5.574%
69	17.175	2408	2412	2416	rVB	539785	784067	18.08%	1.623%
70	17.234	2416	2422	2426	rBV	1526574	2191167	50.53%	4.536%
71	17.310	2432	2435	2439	rVB6	12898	16605	0.38%	0.034%
72	17.539	2469	2474	2475	rBV4	35973	50079	1.15%	0.104%
73	17.569	2476	2479	2485	rVV	51245	99381	2.29%	0.206%
74	17.628	2487	2489	2492	rVV4	19178	20902	0.48%	0.043%
75	17.663	2492	2495	2499	rVV4	16333	27049	0.62%	0.056%
76	17.739	2503	2508	2514	rVB3	37499	67426	1.55%	0.140%

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77	17.816	2518	2521	2525	rVB2	22688	30909	0.71%	0.064%
78	17.934	2537	2541	2547	rBV9	24452	53135	1.23%	0.110%
79	18.051	2554	2561	2565	rBV2	549114	864400	19.93%	1.789%
80	18.169	2578	2581	2586	rVB	32253	47691	1.10%	0.099%
81	18.233	2586	2592	2596	rBV3	145533	258112	5.95%	0.534%
82	18.269	2596	2598	2601	rVV3	18651	22095	0.51%	0.046%
83	18.339	2607	2610	2611	rBV3	37011	37107	0.86%	0.077%
84	18.616	2653	2657	2664	rVB2	71765	119479	2.76%	0.247%
85	18.933	2707	2711	2714	rBV	40494	59985	1.38%	0.124%
86	19.022	2723	2726	2728	rBV2	43465	59470	1.37%	0.123%
87	19.263	2759	2767	2773	rBV	1321609	2000491	46.13%	4.141%
88	19.392	2785	2789	2798	rVB4	223983	363270	8.38%	0.752%
89	19.616	2821	2827	2830	rBV	2348525	3438784	79.30%	7.118%
90	19.639	2830	2831	2841	rVV	996170	1038818	23.95%	2.150%
91	19.716	2842	2844	2849	rVB2	61723	78557	1.81%	0.163%
92	20.163	2917	2920	2925	rVB2	180402	311202	7.18%	0.644%
93	20.269	2935	2938	2942	rBV	87489	114638	2.64%	0.237%
94	21.222	3095	3100	3110	rVB3	771946	1273453	29.36%	2.636%
95	21.475	3140	3143	3147	rVB	273506	341529	7.88%	0.707%
96	21.575	3152	3160	3164	rVV3	2227039	4336689	100.00%	8.977%
97	21.622	3165	3168	3178	rVB	538938	875381	20.19%	1.812%
98	23.851	3542	3547	3554	rBV	399870	922604	21.27%	1.910%
99	24.674	3681	3687	3694	rVB	777698	1742838	40.19%	3.608%
100	24.910	3719	3727	3736	rBV	1145192	3101113	71.51%	6.419%

Sum of corrected areas: 48309631

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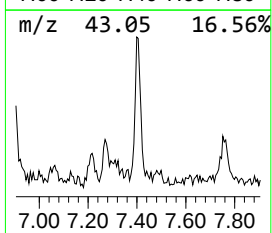
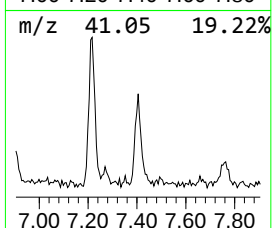
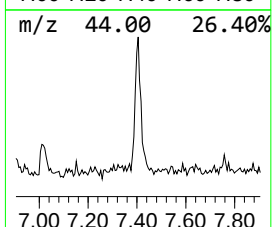
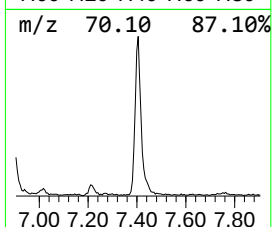
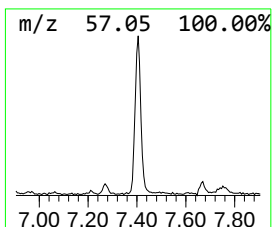
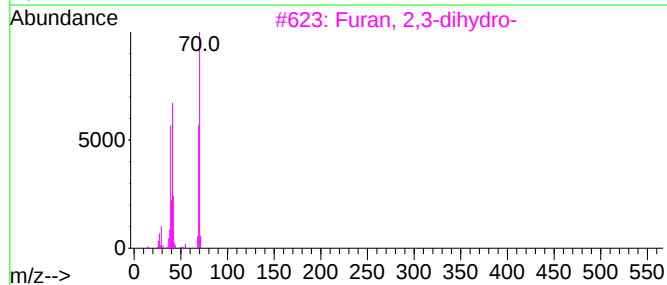
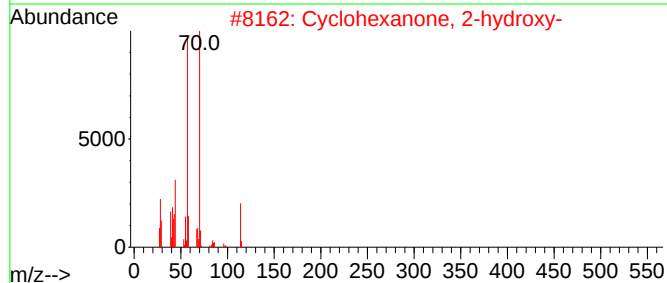
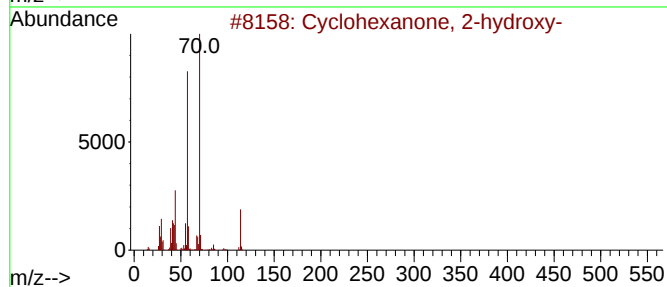
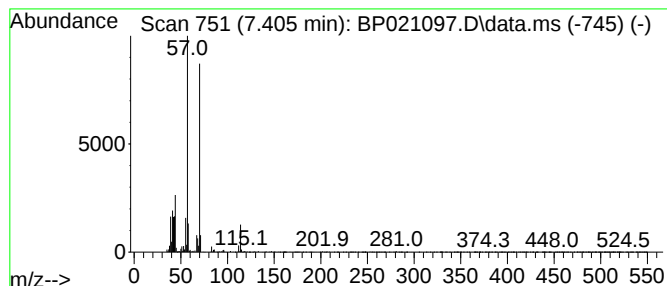
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Cyclohexanone, 2-hydroxy- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.405	2.81 ng/ul	156853	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 2-hydroxy-	114	C6H10O2	000533-60-8	80
2			Cyclohexanone, 2-hydroxy-	114	C6H10O2	000533-60-8	53
3			Furan, 2,3-dihydro-	70	C4H6O	001191-99-7	37
4			(S)-(+)-2-(Methoxymethyl)pyrroli...	115	C6H13NO	063126-47-6	32
5			Proline	115	C5H9NO2	000147-85-3	32



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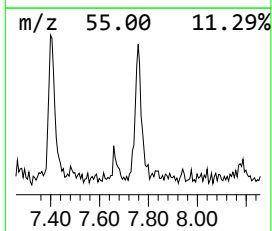
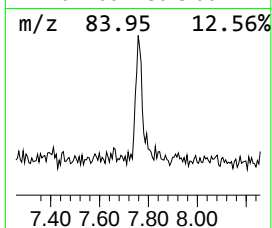
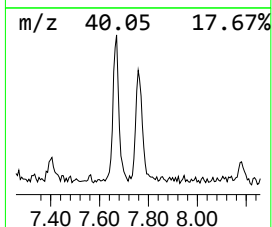
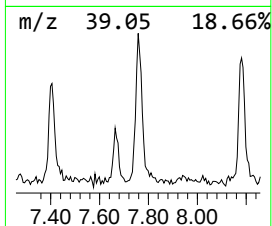
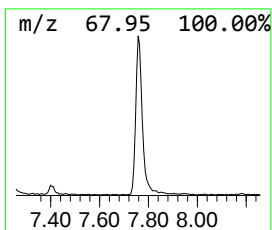
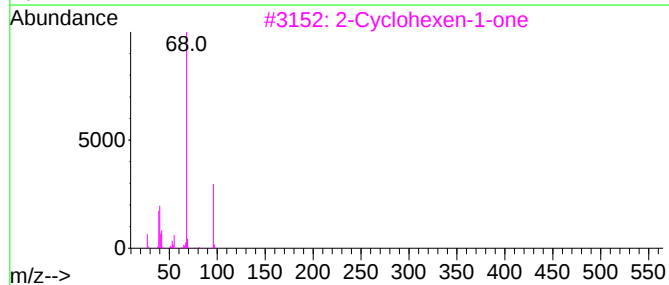
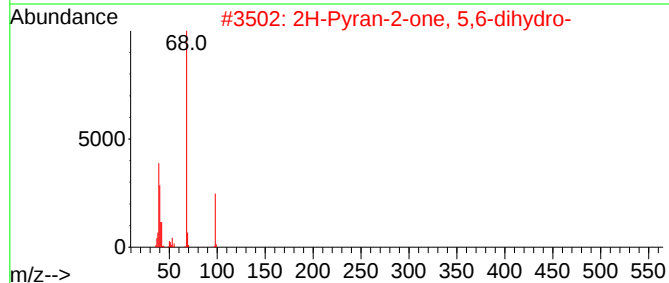
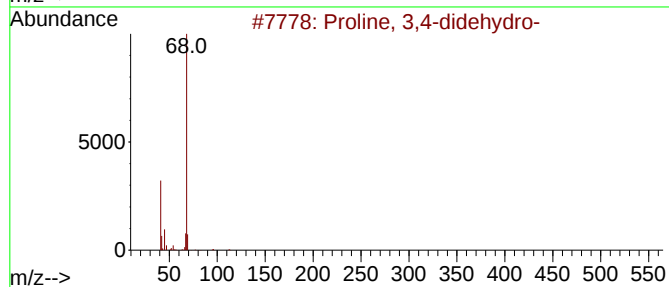
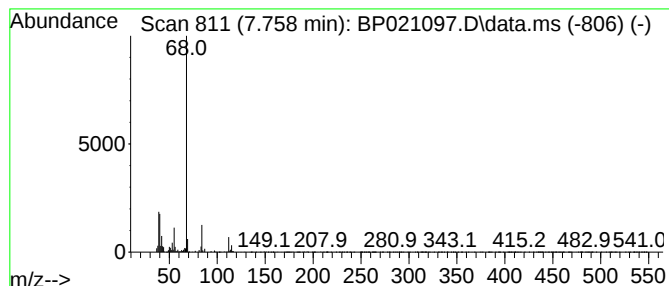
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Proline, 3,4-didehydro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.758	2.69 ng/ul	150302	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Proline, 3,4-didehydro-	113	C5H7NO2	003395-35-5	64
2			2H-Pyran-2-one, 5,6-dihydro-	98	C5H6O2	003393-45-1	64
3			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	59
4			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	45
5			Dihydro-3-methylene-5-methyl-2-f...	112	C6H8O2	062873-16-9	42



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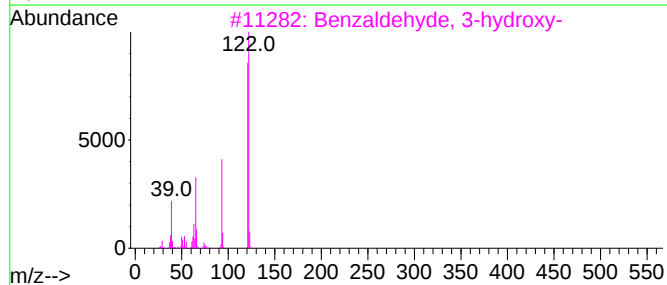
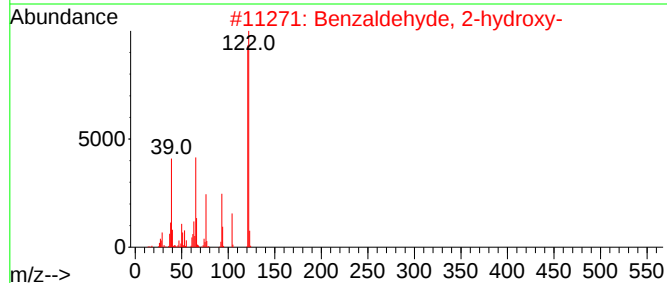
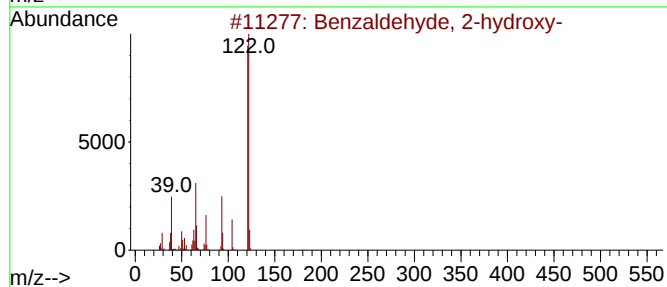
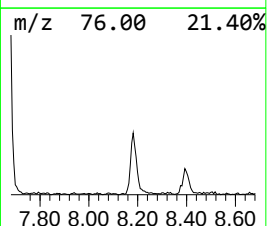
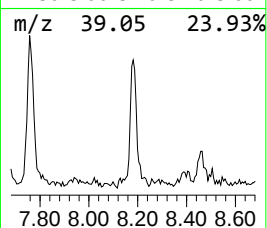
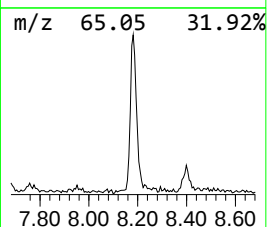
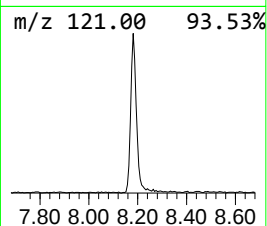
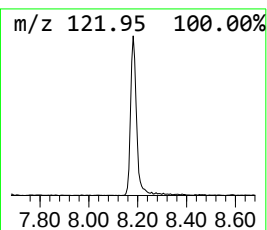
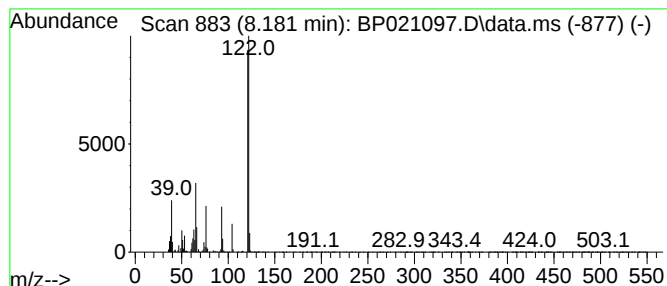
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzaldehyde, 2-hydroxy- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.181	2.65 ng/ul	147912	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 2-hydroxy-	122	C7H6O2	000090-02-8	96
2			Benzaldehyde, 2-hydroxy-	122	C7H6O2	000090-02-8	95
3			Benzaldehyde, 3-hydroxy-	122	C7H6O2	000100-83-4	95
4			Benzaldehyde, 3-hydroxy-	122	C7H6O2	000100-83-4	93
5			Benzaldehyde, 4-hydroxy-	122	C7H6O2	000123-08-0	91



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ClientSampleId :
A4CG8

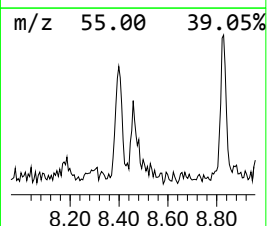
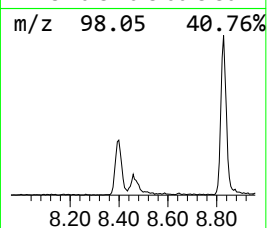
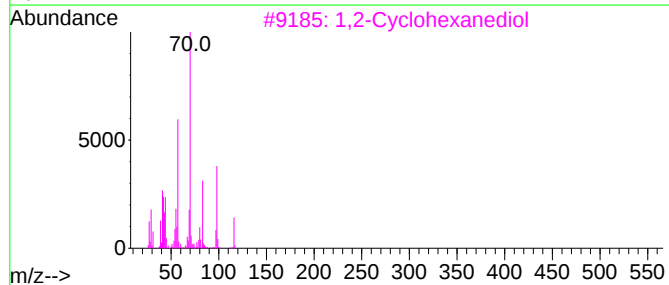
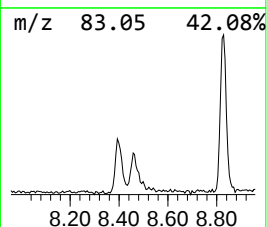
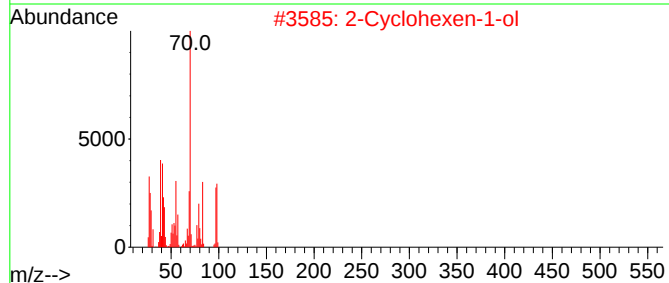
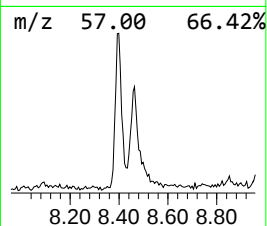
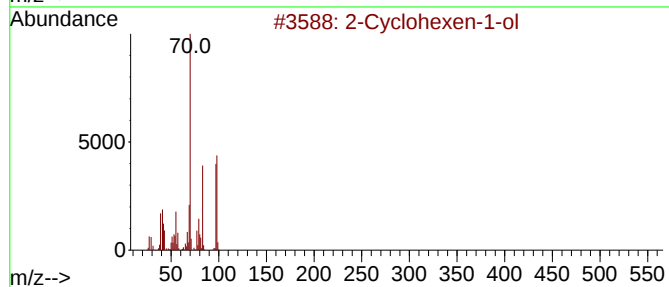
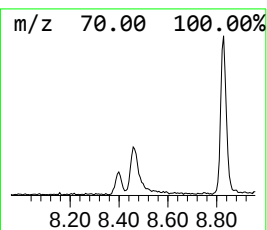
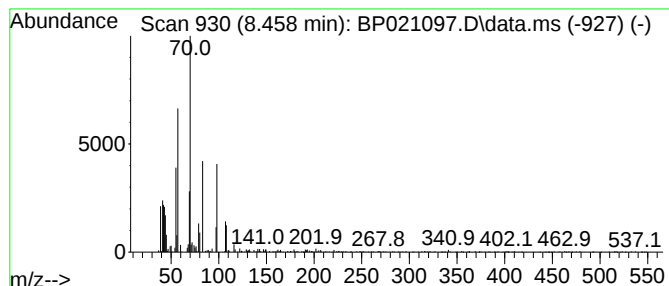
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Cyclohexen-1-ol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.458	2.39 ng/ul	133473	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	72
2	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	64
3	1,2-Cyclohexanediol			116	C6H12O2	000931-17-9	53
4	1,2-Cyclohexanediol, trans-			116	C6H12O2	001460-57-7	50
5	1,7-Diazabicyclo[2.2.0]heptane			98	C5H10N2	000279-42-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021097.D
Acq On : 23 Jul 2024 13:38
Operator : MA/JU
Sample : P3238-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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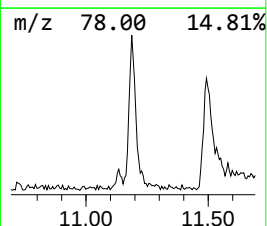
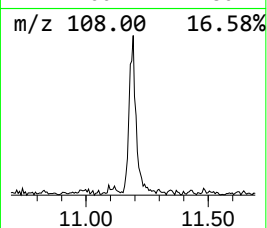
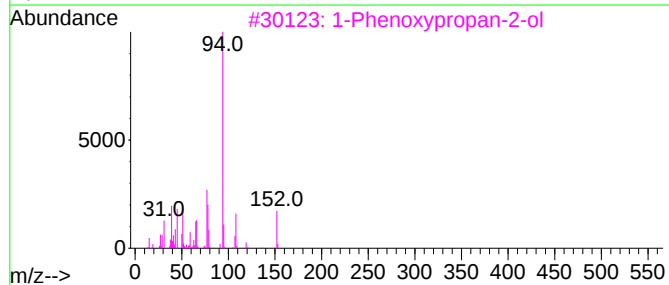
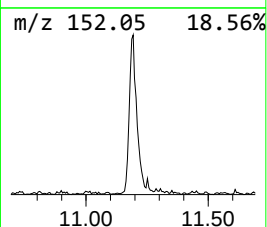
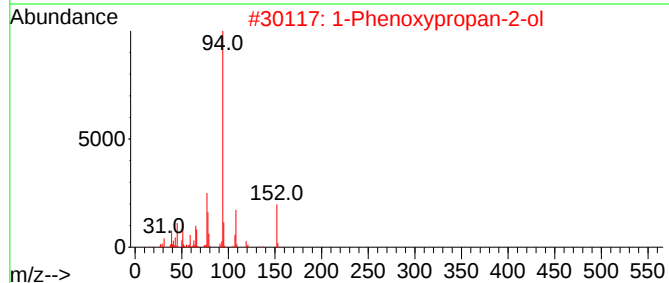
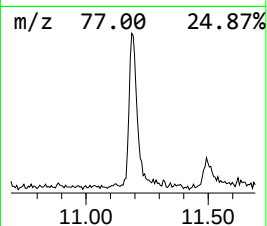
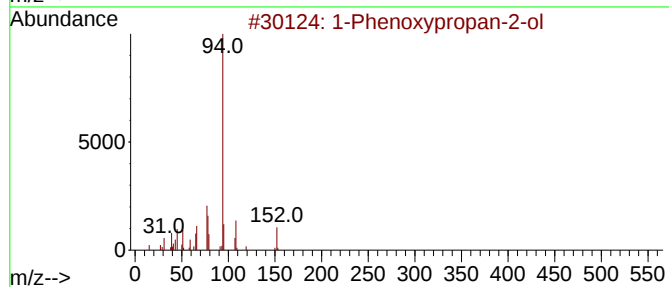
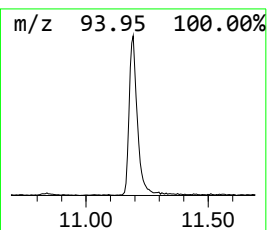
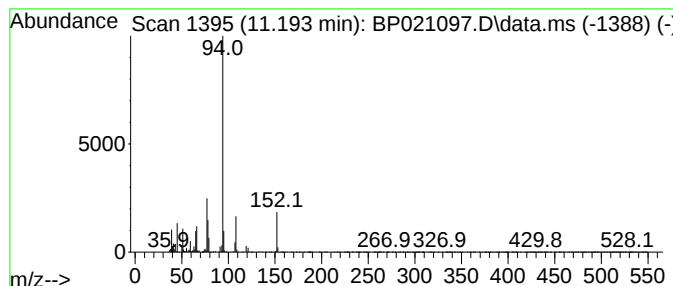
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1-Phenoxypropan-2-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.193	2.45 ng/ul	193258	Naphthalene-d8	10.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	97
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021097.D
Acq On : 23 Jul 2024 13:38
Operator : MA/JU
Sample : P3238-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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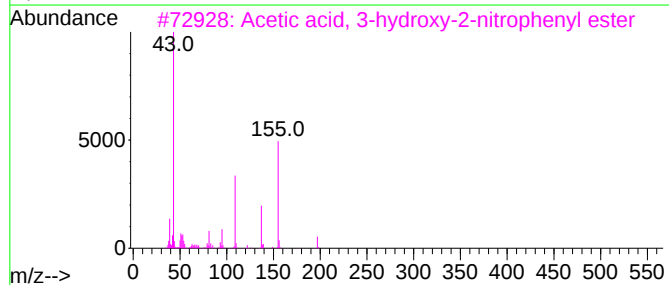
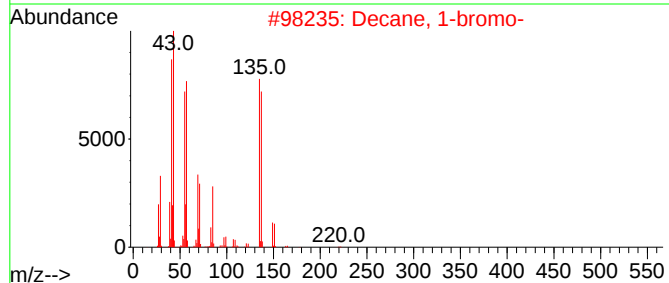
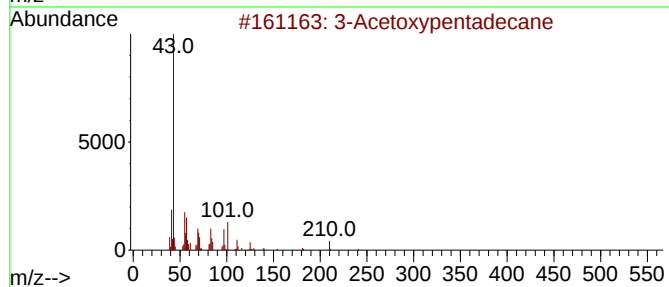
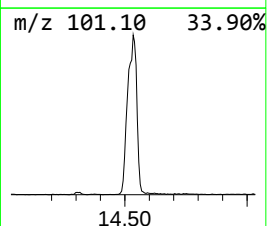
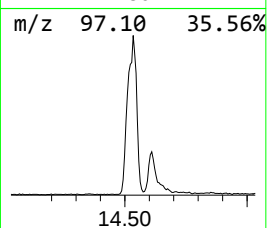
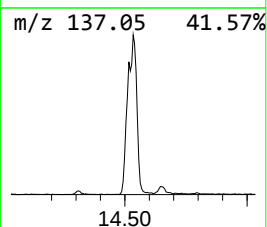
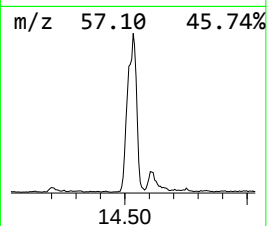
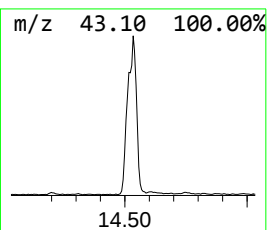
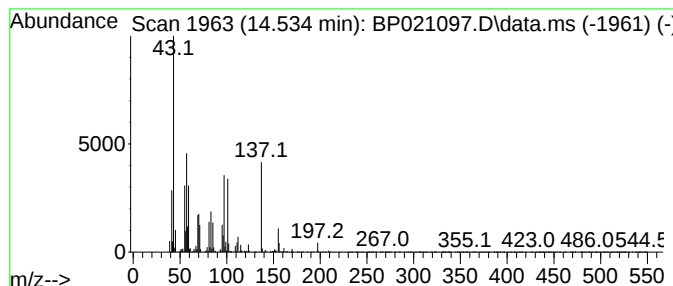
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.534	8.28 ng/ul	900043	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Acetoxyptadecane	270	C17H34O2	1000245-62-1	12
2			Decane, 1-bromo-	220	C10H21Br	000112-29-8	12
3			Acetic acid, 3-hydroxy-2-nitroph...	197	C8H7NO5	1000269-41-1	10
4			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	10
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021097.D
Acq On : 23 Jul 2024 13:38
Operator : MA/JU
Sample : P3238-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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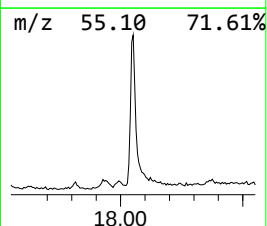
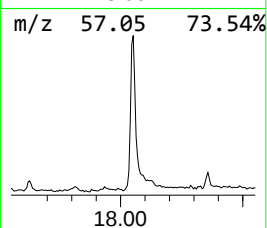
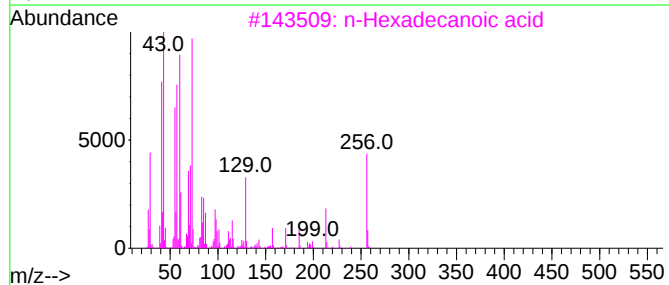
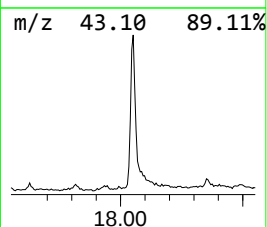
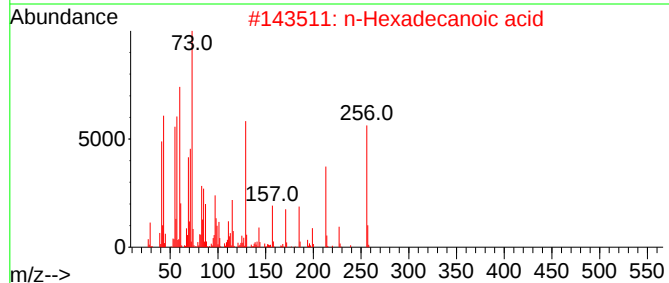
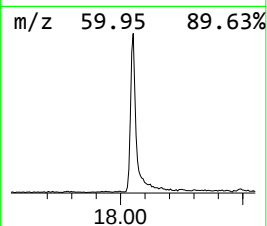
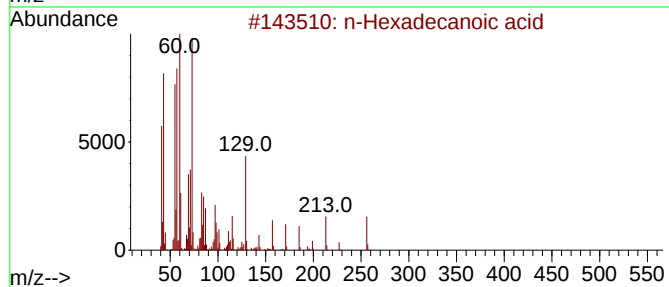
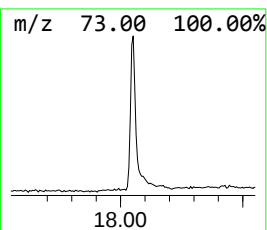
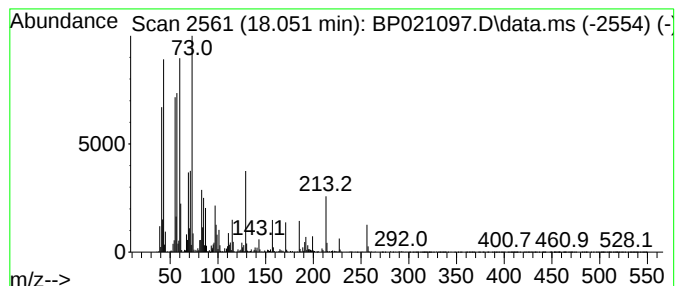
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.051	6.42 ng/ul	864400	Phenanthrene-d10	17.128

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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021097.D
Acq On : 23 Jul 2024 13:38
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Sample : P3238-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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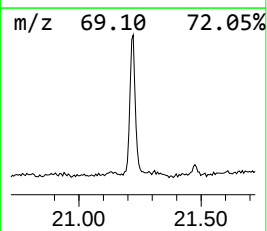
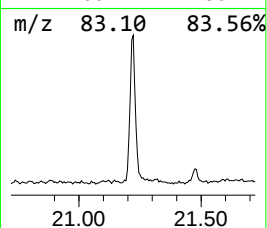
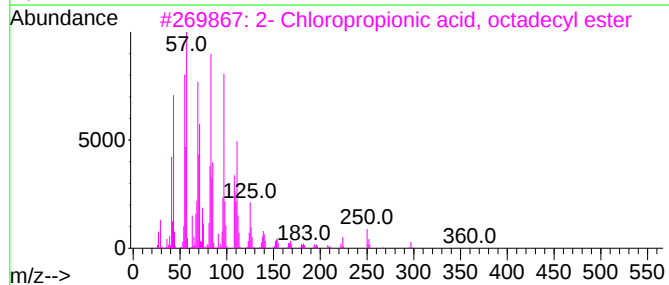
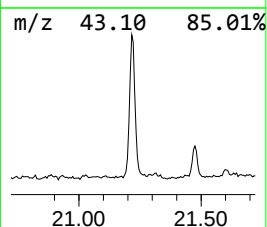
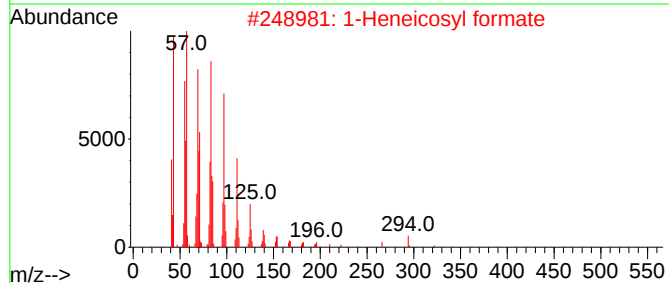
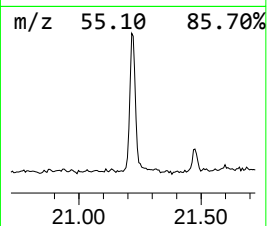
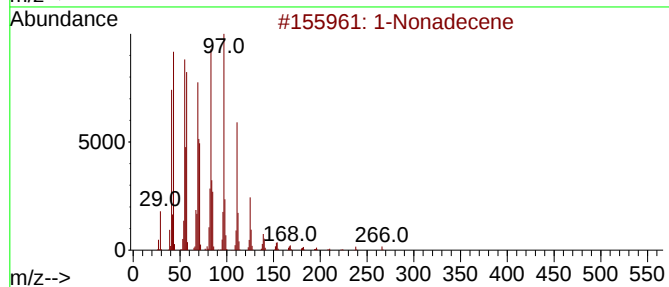
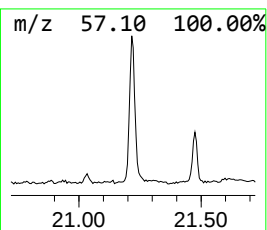
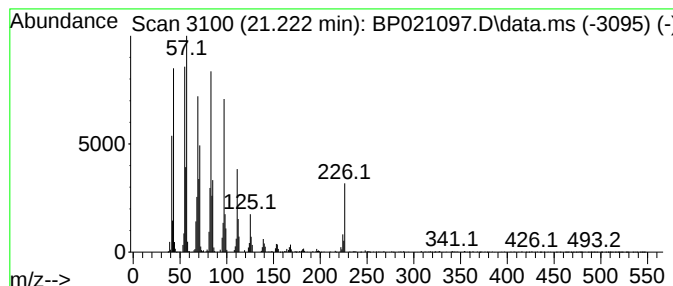
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.222	5.87 ng/ul	1273450	Chrysene-d12	21.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene			266	C19H38	018435-45-5	95
2	1-Heneicosyl formate			340	C22H44O2	077899-03-7	94
3	2- Chloropropionic acid, octadec...			360	C21H41ClO2	088104-31-8	93
4	Cyclohexadecane			224	C16H32	000295-65-8	93
5	Heptafluorobutyric acid, hexadec...			438	C20H33F7O2	006385-15-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021097.D
Acq On : 23 Jul 2024 13:38
Operator : MA/JU
Sample : P3238-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
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
Cyclohexanone, ...	7.405	2.8	ng/ul	156853	1	7.669	1117300	20.0
Proline, 3,4-di...	7.758	2.7	ng/ul	150302	1	7.669	1117300	20.0
Benzaldehyde, 2...	8.181	2.6	ng/ul	147912	1	7.669	1117300	20.0
2-Cyclohexen-1-ol	8.458	2.4	ng/ul	133473	1	7.669	1117300	20.0
1-Phenoxypropan...	11.193	2.5	ng/ul	193258	2	10.434	1580630	20.0
unknown-01	14.534	8.3	ng/ul	900043	3	14.310	2174040	20.0
n-Hexadecanoic ...	18.051	6.4	ng/ul	864400	4	17.128	2692660	20.0
(DEL) Alkane: S...	21.222	5.9	ng/ul	1273450	5	21.575	4336690	20.0