

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
 Data File : BP020362.D
 Acq On : 13 May 2024 20:47
 Operator : MA/JU
 Sample : P2369-11MEDL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43N5MEDL

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-BP050124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020362.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.646	92	95	100	rBV3	12785	23807	1.81%	0.258%
2	3.911	135	140	142	rBV5	2896	3986	0.30%	0.043%
3	5.881	470	475	482	rBV5	4772	12929	0.98%	0.140%
4	6.211	526	531	534	rBV7	3813	7747	0.59%	0.084%
5	6.269	535	541	545	rVV	16011	23763	1.80%	0.257%
6	6.358	552	556	561	rVB5	4962	8927	0.68%	0.097%
7	6.422	561	567	573	rVB4	7671	13646	1.04%	0.148%
8	6.469	573	575	582	rBV6	4647	9191	0.70%	0.100%
9	6.687	608	612	616	rBV6	5163	9072	0.69%	0.098%
10	6.734	616	620	628	rVV6	18692	39401	2.99%	0.427%
11	6.846	634	639	645	rBV7	15960	43049	3.27%	0.466%
12	6.981	656	662	668	rBV2	23367	47706	3.62%	0.517%
13	7.087	675	680	685	rVB5	12787	21395	1.62%	0.232%
14	7.163	685	693	701	rBV4	26963	75388	5.72%	0.816%
15	7.234	701	705	709	rBV4	6322	10239	0.78%	0.111%
16	7.299	712	716	722	rBV3	11138	20532	1.56%	0.222%
17	7.381	726	730	740	rVB7	14435	37809	2.87%	0.409%
18	7.505	742	751	754	rBV7	22124	55146	4.18%	0.597%
19	7.611	764	769	781	rVB	166580	327041	24.82%	3.541%
20	7.705	781	785	786	rBV4	7046	9916	0.75%	0.107%
21	7.763	790	795	798	rVB4	28032	45479	3.45%	0.492%
22	7.805	799	802	805	rBV3	9303	11342	0.86%	0.123%
23	7.846	807	809	814	rVB5	10377	14637	1.11%	0.159%
24	7.893	814	817	820	rBV4	5953	7719	0.59%	0.084%
25	7.922	820	822	824	rBV2	6665	6914	0.52%	0.075%
26	8.016	834	838	844	rVB8	12176	23238	1.76%	0.252%
27	8.081	844	849	850	rBV5	8078	11768	0.89%	0.127%
28	8.128	850	857	861	rVB4	25655	47059	3.57%	0.510%
29	8.234	871	875	879	rBV5	8161	14097	1.07%	0.153%
30	8.363	891	897	899	rBV6	17591	28594	2.17%	0.310%
31	8.405	899	904	919	rVB9	46796	167388	12.70%	1.813%
32	8.552	924	929	935	rBV5	10367	19284	1.46%	0.209%
33	8.669	940	949	956	rVB5	29506	75092	5.70%	0.813%
34	8.775	962	967	968	rBV5	11832	16874	1.28%	0.183%
35	8.875	980	984	987	rBV2	20576	32971	2.50%	0.357%
36	8.910	987	990	991	rBV3	14111	12275	0.93%	0.133%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

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Stop Thrs : 0

Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Title : SVOA CALIBRATION

37	8.993	1000	1004	1005	rBV4	10066	11386	0.86%	0.123%
38	9.028	1007	1010	1016	rVB6	12945	25061	1.90%	0.271%
39	9.087	1016	1020	1023	rBV6	6983	12113	0.92%	0.131%
40	9.281	1051	1053	1059	rVB4	22886	33630	2.55%	0.364%
41	9.440	1075	1080	1087	rVB6	16496	31827	2.42%	0.345%
42	9.540	1087	1097	1102	rBV6	37366	89888	6.82%	0.973%
43	9.587	1102	1105	1109	rVV4	15554	24972	1.90%	0.270%
44	9.640	1112	1114	1117	rVB4	4784	5066	0.38%	0.055%
45	9.705	1123	1125	1130	rVB6	4790	6302	0.48%	0.068%
46	9.946	1161	1166	1172	rBV10	7179	15846	1.20%	0.172%
47	10.093	1183	1191	1193	rBV8	10042	22946	1.74%	0.248%
48	10.240	1212	1216	1221	rBV8	5247	7730	0.59%	0.084%
49	10.381	1233	1240	1253	rVB	215350	444322	33.72%	4.811%
50	10.522	1261	1264	1270	rVB2	9606	14653	1.11%	0.159%
51	10.746	1298	1302	1308	rVB6	6478	12673	0.96%	0.137%
52	10.840	1316	1318	1325	rVB7	3750	5737	0.44%	0.062%
53	11.393	1407	1412	1418	rBV8	4770	9671	0.73%	0.105%
54	11.651	1449	1456	1461	rBV8	3210	5846	0.44%	0.063%
55	11.804	1478	1482	1484	rBV3	3368	3982	0.30%	0.043%
56	11.851	1486	1490	1496	rVB8	4184	7008	0.53%	0.076%
57	12.251	1553	1558	1561	rVV6	3190	6694	0.51%	0.072%
58	12.775	1644	1647	1656	rVB10	2863	6511	0.49%	0.071%
59	13.081	1693	1699	1702	rBV7	4031	7373	0.56%	0.080%
60	13.428	1754	1758	1763	rVB6	6009	9481	0.72%	0.103%
61	13.498	1766	1770	1773	rBV6	4214	6773	0.51%	0.073%
62	13.581	1781	1784	1787	rBV4	3858	4146	0.31%	0.045%
63	13.710	1800	1806	1817	rBV	55505	130540	9.91%	1.414%
64	13.893	1833	1837	1839	rBV3	10821	13492	1.02%	0.146%
65	13.922	1839	1842	1845	rVV4	8050	9103	0.69%	0.099%
66	13.969	1845	1850	1857	rVV	71312	132210	10.03%	1.432%
67	14.034	1859	1861	1864	rVB3	7760	7638	0.58%	0.083%
68	14.069	1864	1867	1870	rBV4	4150	5559	0.42%	0.060%
69	14.216	1889	1892	1896	rBV4	6777	8596	0.65%	0.093%
70	14.281	1896	1903	1910	rBV	337212	607521	46.10%	6.579%
71	14.340	1910	1913	1918	rVB2	24510	36023	2.73%	0.390%
72	14.434	1925	1929	1933	rVB7	4447	7295	0.55%	0.079%
73	14.510	1933	1942	1955	rBV	133494	334707	25.40%	3.624%
74	14.598	1955	1957	1961	rVB5	3313	4650	0.35%	0.050%
75	14.681	1967	1971	1973	rBV5	5072	6969	0.53%	0.075%
76	14.722	1973	1978	1980	rVV5	3017	4908	0.37%	0.053%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Stop Thrs : 0

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Title : SVOA CALIBRATION

77	14.869	1999	2003	2008	rVB6	7167	9538	0.72%	0.103%
78	14.945	2011	2016	2020	rBV4	14075	25640	1.95%	0.278%
79	15.134	2039	2048	2054	rVB7	9459	19679	1.49%	0.213%
80	15.316	2072	2079	2092	rBV	96467	193408	14.68%	2.094%
81	15.498	2102	2110	2118	rBV10	13916	41139	3.12%	0.445%
82	15.787	2152	2159	2162	rBV9	3182	6146	0.47%	0.067%
83	15.916	2177	2181	2184	rBV5	4173	6042	0.46%	0.065%
84	16.275	2237	2242	2247	rBV9	6195	11830	0.90%	0.128%
85	16.763	2318	2325	2339	rBV	318271	690745	52.42%	7.480%
86	17.116	2378	2385	2399	rBV	454079	840268	63.77%	9.099%
87	17.228	2399	2404	2421	rVB2	114851	231538	17.57%	2.507%
88	17.545	2454	2458	2461	rBV6	7201	11289	0.86%	0.122%
89	17.657	2472	2477	2479	rBV6	3001	4995	0.38%	0.054%
90	18.081	2544	2549	2557	rBV8	17196	31297	2.38%	0.339%
91	19.045	2709	2713	2718	rBV5	13206	21108	1.60%	0.229%
92	19.622	2807	2811	2816	rBV	154549	268720	20.39%	2.910%
93	19.663	2816	2818	2823	rVB	43464	49104	3.73%	0.532%
94	20.228	2910	2914	2920	rBV	72551	95481	7.25%	1.034%
95	20.757	2999	3004	3008	rBV	110042	147543	11.20%	1.598%
96	21.292	3090	3095	3102	rVB2	98412	162834	12.36%	1.763%
97	21.622	3144	3151	3167	rVB2	572112	1154082	87.58%	12.497%
98	21.869	3189	3193	3201	rVB2	71404	118404	8.99%	1.282%
99	24.727	3674	3679	3692	rVB2	100361	299791	22.75%	3.246%
100	24.951	3709	3717	3731	rBV	397576	1317731	100.00%	14.269%

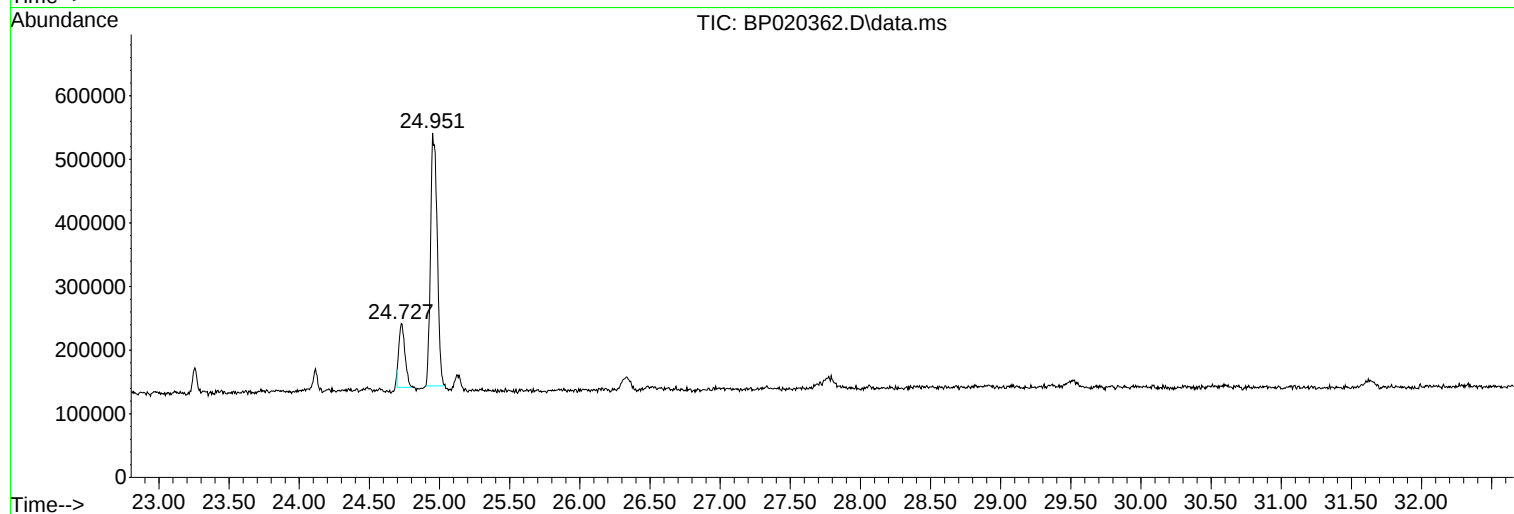
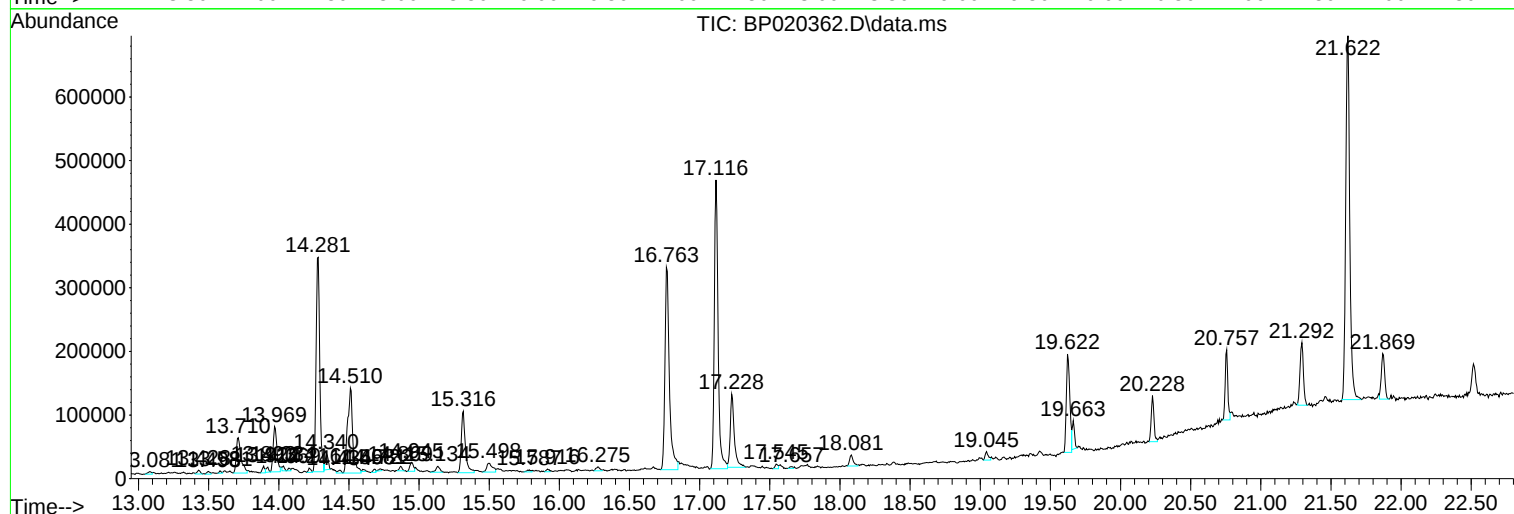
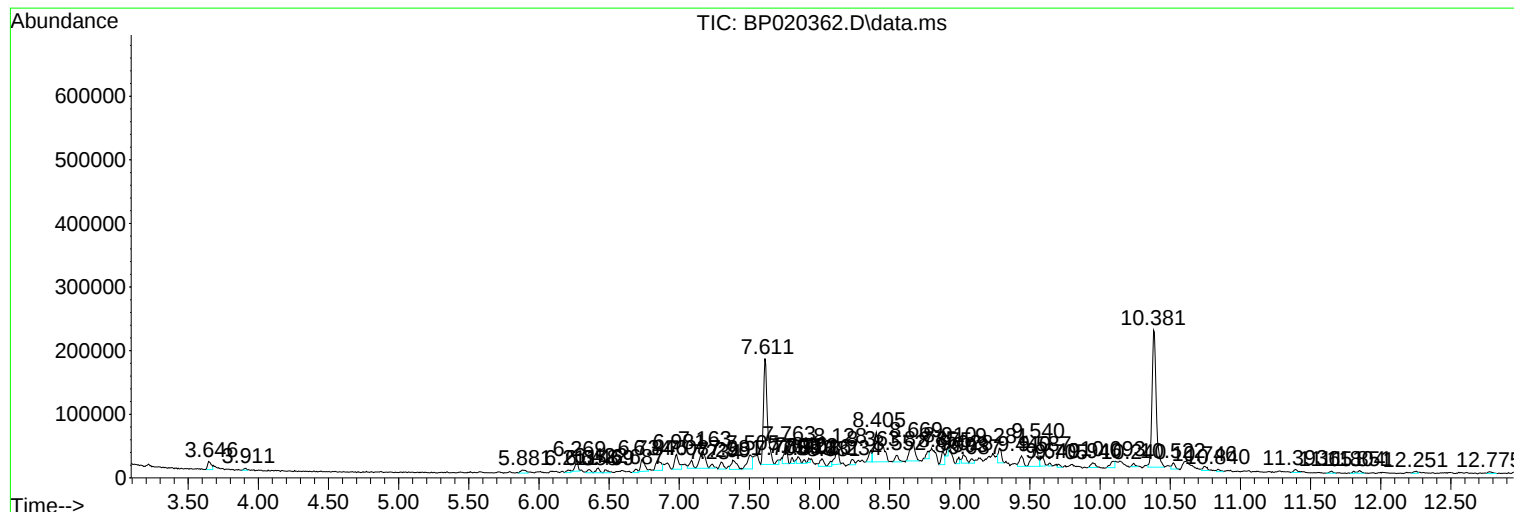
Sum of corrected areas: 9234621

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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
A43N5MEDL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
BNA_P
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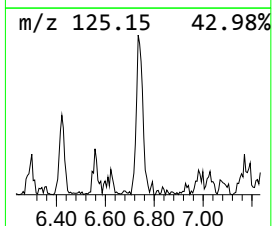
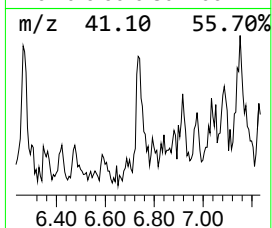
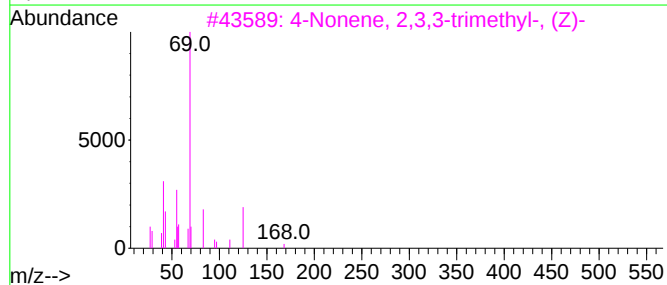
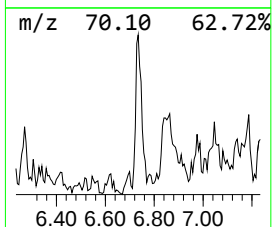
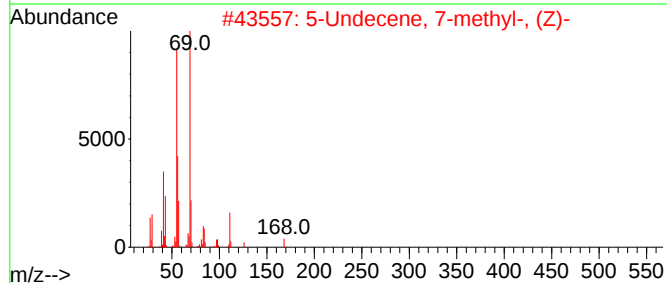
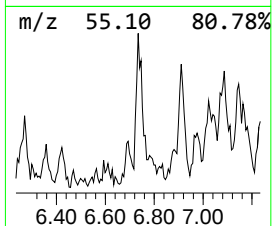
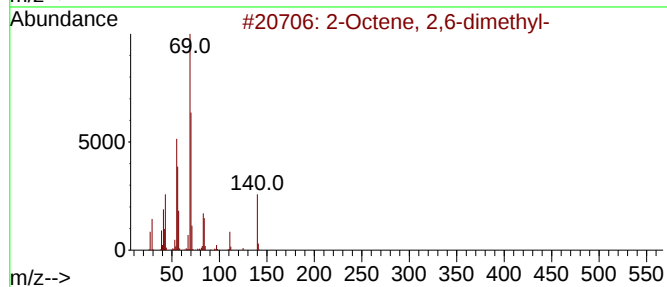
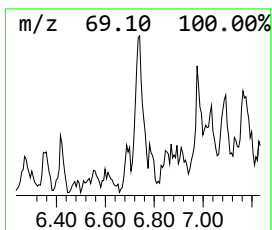
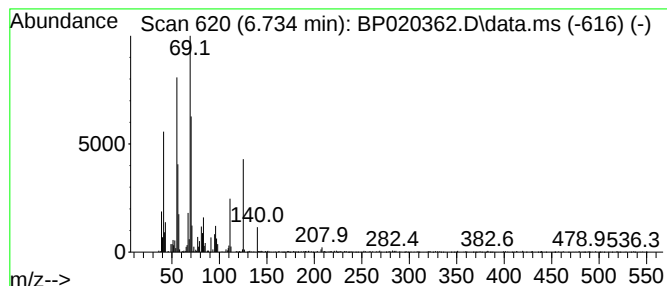
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.734	2.41 ng/ul	39401	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Octene, 2,6-dimethyl-			140	C10H20	004057-42-5	50
2	5-Undecene, 7-methyl-, (Z)-			168	C12H24	074630-62-9	43
3	4-Nonene, 2,3,3-trimethyl-, (Z)-			168	C12H24	063830-68-2	43
4	3-Ethyl-3-octene			140	C10H20	019781-31-8	35
5	Cyclooctanemethanol			142	C9H18O	003637-63-6	35



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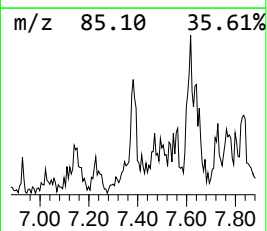
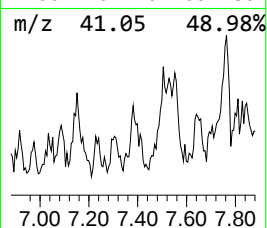
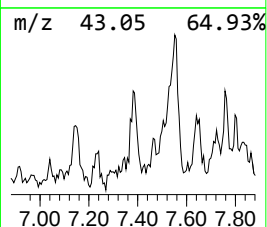
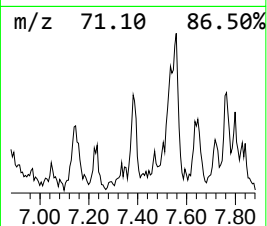
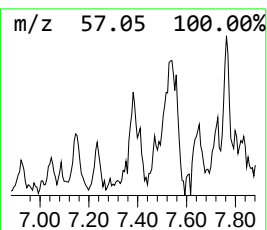
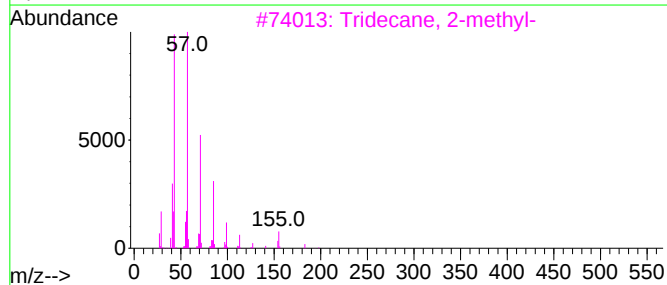
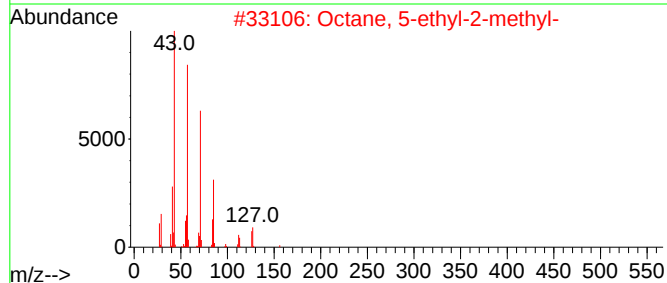
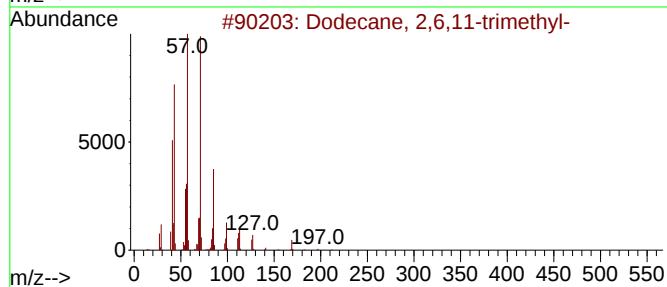
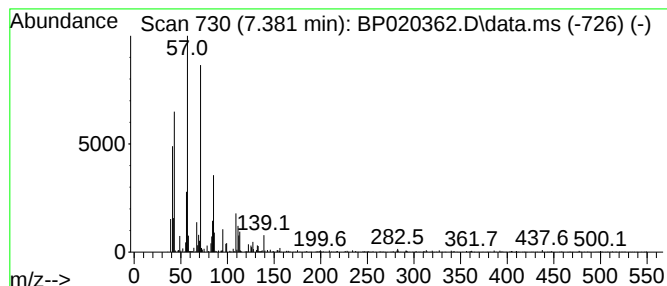
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
7.381	2.31 ng/ul	37809	1,4-Dichlorobenzene-d4		7.611	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Dodecane, 2,6,11-trimethyl-		212	C15H32	031295-56-4	64
2	Octane, 5-ethyl-2-methyl-		156	C11H24	062016-18-6	64
3	Tridecane, 2-methyl-		198	C14H30	001560-96-9	58
4	Octane, 2-methyl-		128	C9H20	003221-61-2	58
5	Triacontane, 1-bromo-		500	C30H61Br	004209-22-7	53



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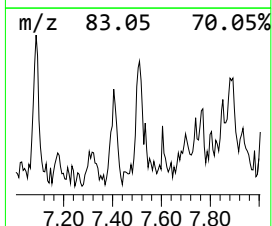
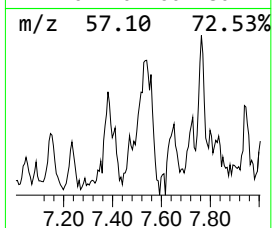
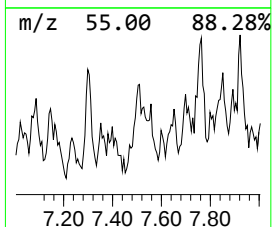
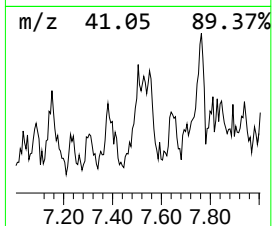
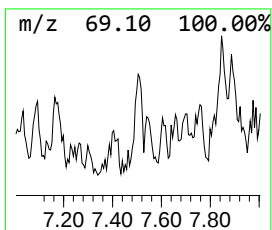
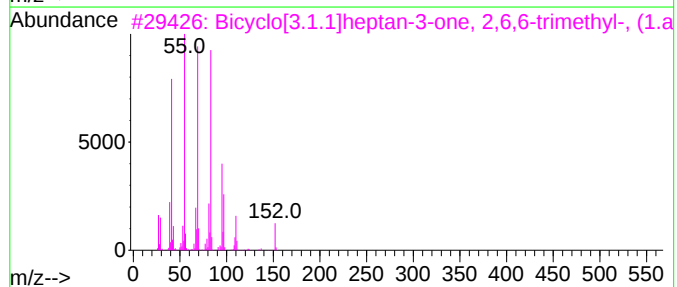
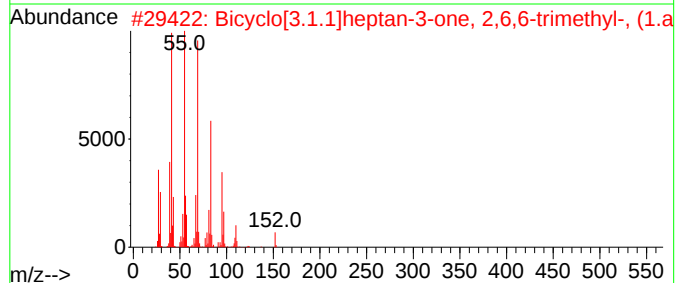
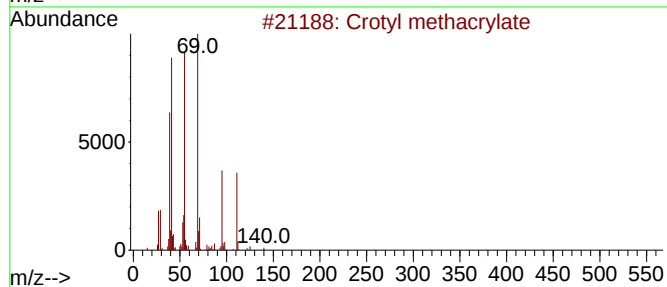
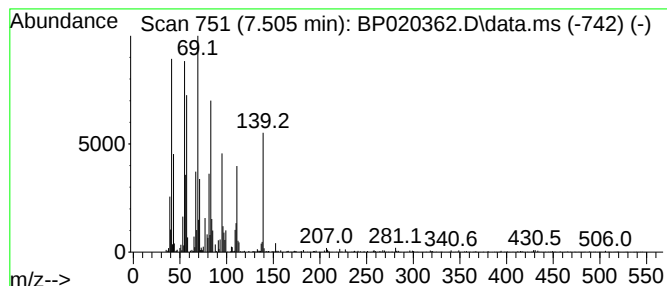
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Crotyl methacrylate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.505	3.37 ng/ul	55146	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Crotyl methacrylate	140	C8H12O2	007376-45-6	55
2			Bicyclo[3.1.1]heptan-3-one, 2,6,...	152	C10H16O	015358-88-0	49
3			Bicyclo[3.1.1]heptan-3-one, 2,6,...	152	C10H16O	000547-60-4	46
4			Bicyclo[3.1.1]heptan-3-one, 2,6,...	152	C10H16O	015358-88-0	43
5			Cyclooctyl bromide	190	C8H15Br	001556-09-8	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020362.D
Acq On : 13 May 2024 20:47
Operator : MA/JU
Sample : P2369-11MEDL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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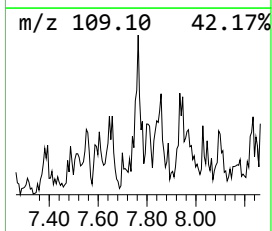
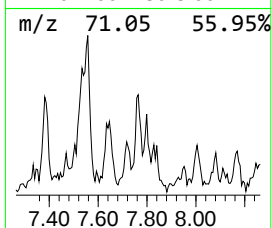
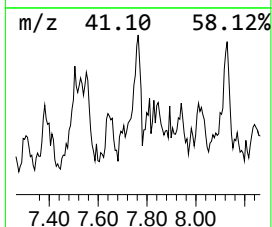
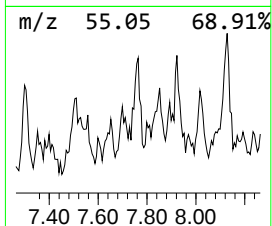
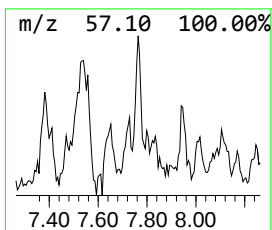
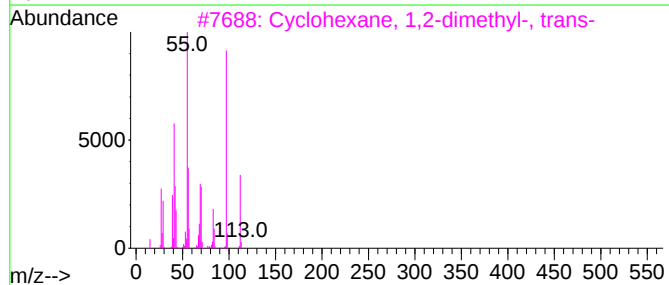
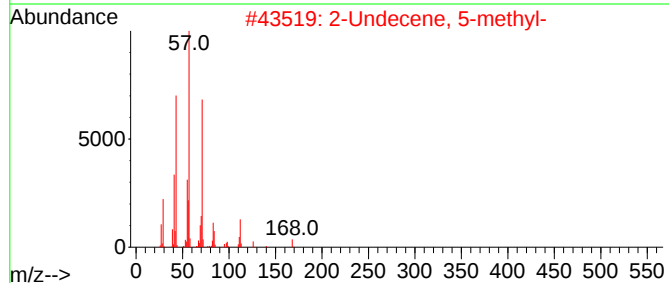
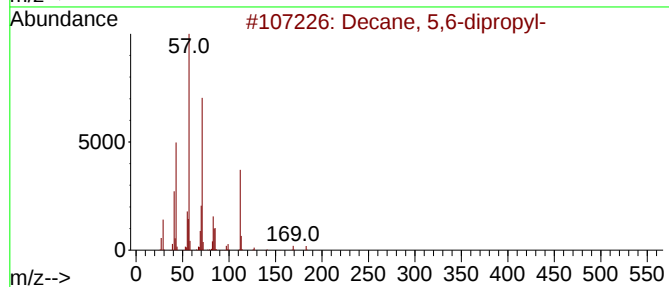
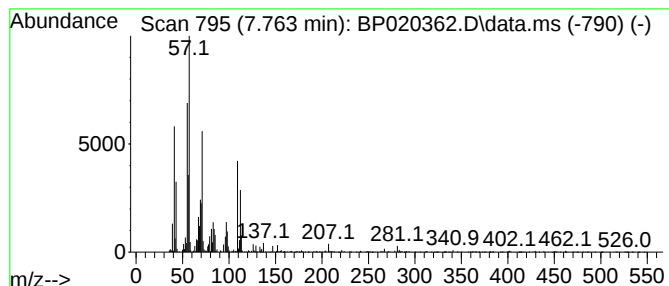
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.763	2.78 ng/ul	45479	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	43
2			2-Undecene, 5-methyl-	168	C12H24	056851-34-4	38
3			Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	38
4			Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	38
5			Oxalic acid, bis(2-ethylhexyl) e...	314	C18H34O4	1000309-39-0	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020362.D
Acq On : 13 May 2024 20:47
Operator : MA/JU
Sample : P2369-11MEDL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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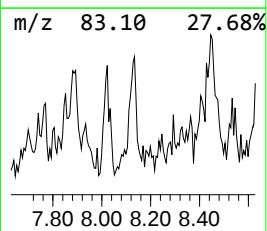
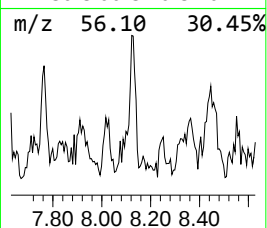
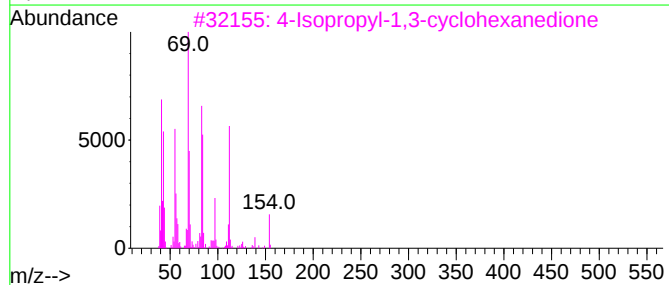
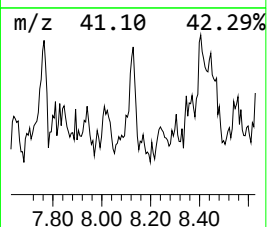
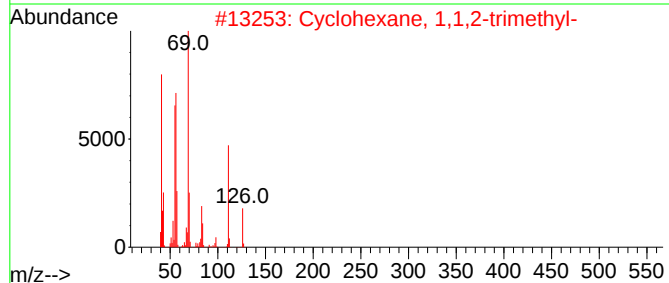
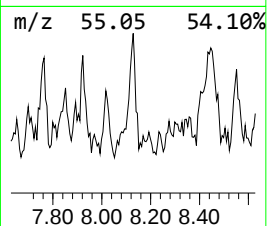
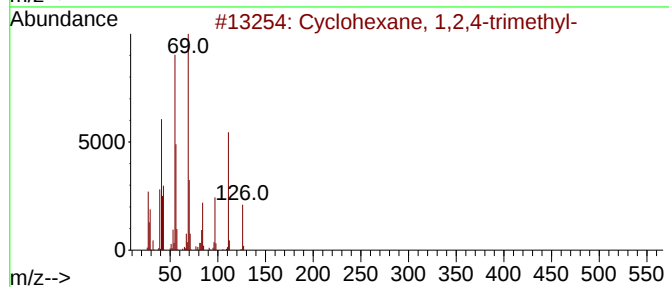
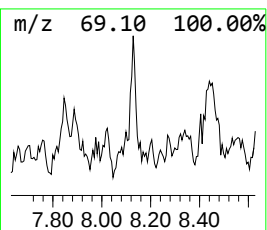
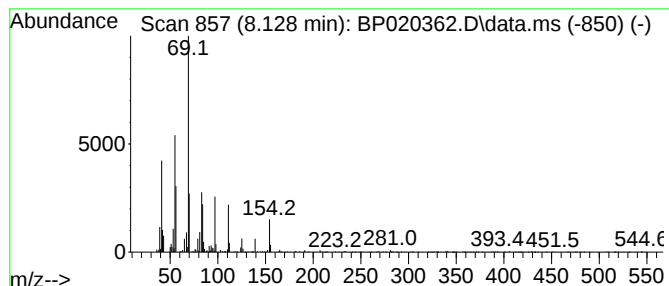
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Cyclic8.128 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.128	2.88 ng/ul	47059	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	59
2			Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	58
3			4-Isopropyl-1,3-cyclohexanedione	154	C9H14O2	062831-62-3	47
4			1,1,4-Trimethylcyclohexane	126	C9H18	007094-27-1	47
5			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020362.D
Acq On : 13 May 2024 20:47
Operator : MA/JU
Sample : P2369-11MEDL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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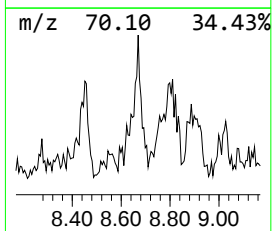
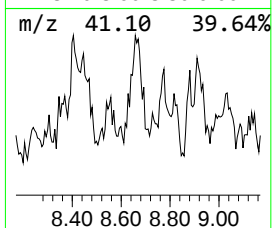
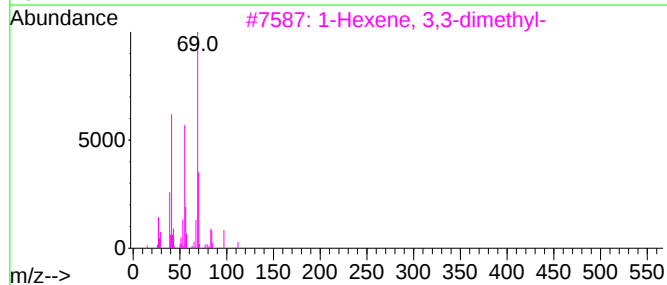
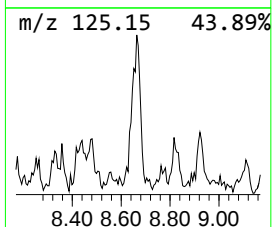
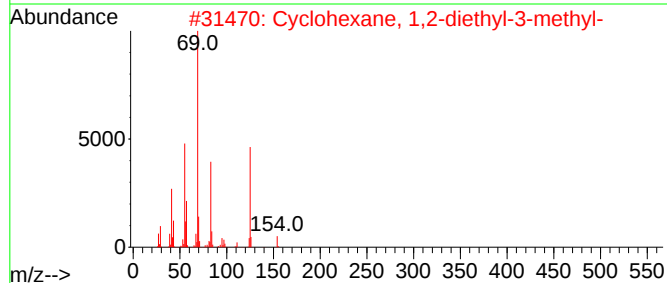
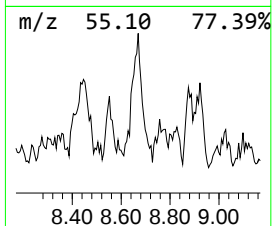
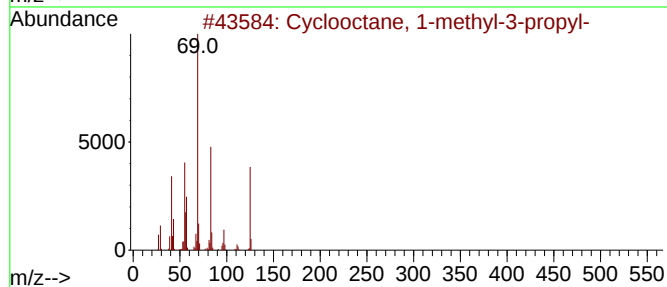
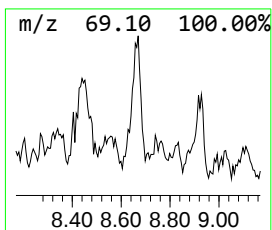
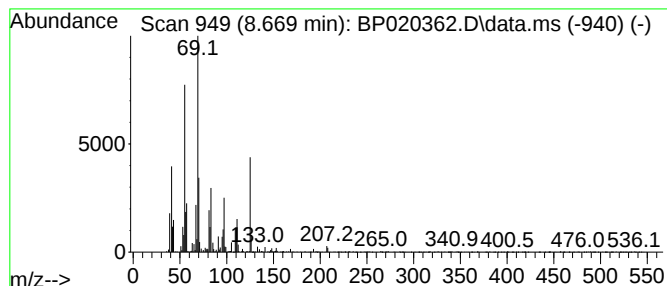
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic8.669 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.669	4.59 ng/ul	75092	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclooctane, 1-methyl-3-propyl-	168	C12H24	255885-37-1	53
2			Cyclohexane, 1,2-diethyl-3-methyl-	154	C11H22	061141-80-8	53
3			1-Hexene, 3,3-dimethyl-	112	C8H16	003404-77-1	43
4			Cyclohexane, 1-ethyl-2-propyl-	154	C11H22	062238-33-9	43
5			Methanone, dicyclopentyl-	110	C7H10O	001121-37-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020362.D
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Operator : MA/JU
Sample : P2369-11MEDL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

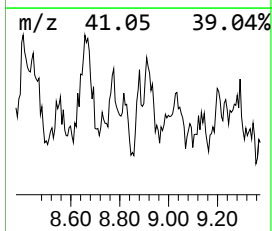
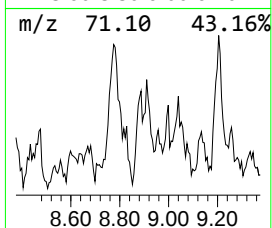
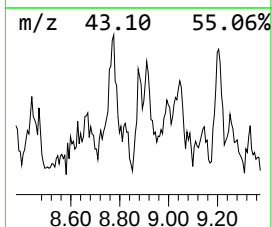
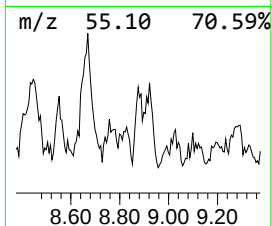
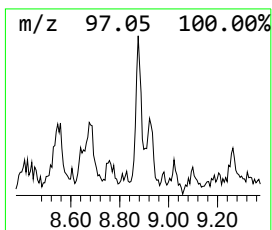
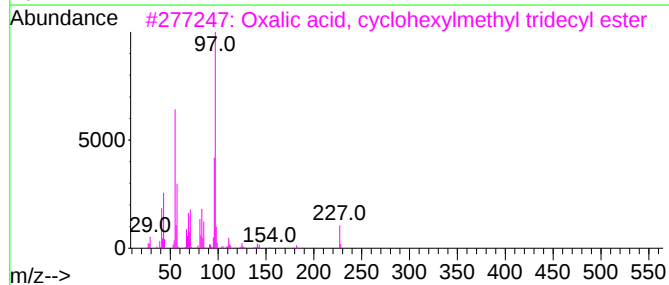
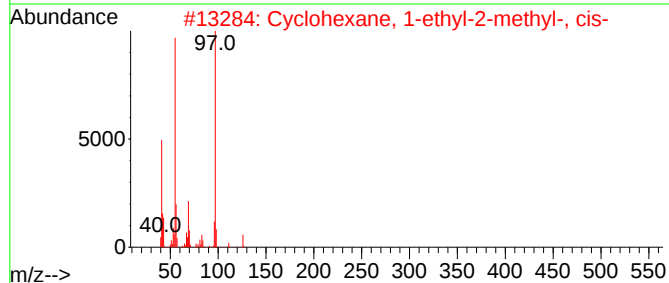
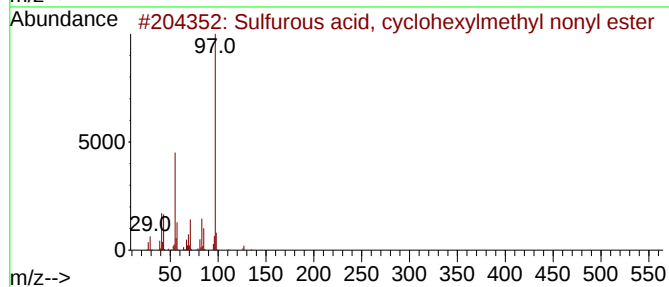
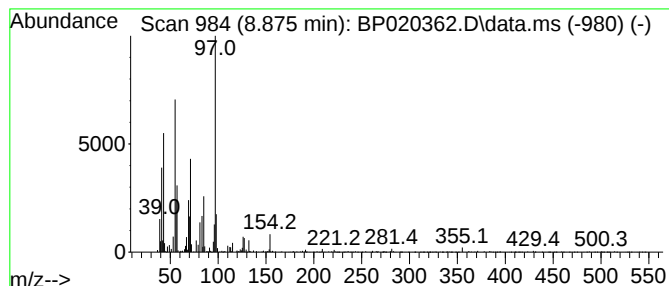
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Sulfurous acid, cyclohexylm... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.875	2.02 ng/ul	32971	1,4-Dichlorobenzene-d4	7.611	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfurous acid, cyclohexylmethyl...	304	C16H32O3S	1000309-21-8	50
2	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	49
3	Oxalic acid, cyclohexylmethyl tr...	368	C22H40O4	1000309-69-1	47
4	Sulfurous acid, cyclohexylmethyl...	360	C20H40O3S	1000309-22-1	47
5	Sulfurous acid, cyclohexylmethyl...	374	C21H42O3S	1000309-22-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020362.D
Acq On : 13 May 2024 20:47
Operator : MA/JU
Sample : P2369-11MEDL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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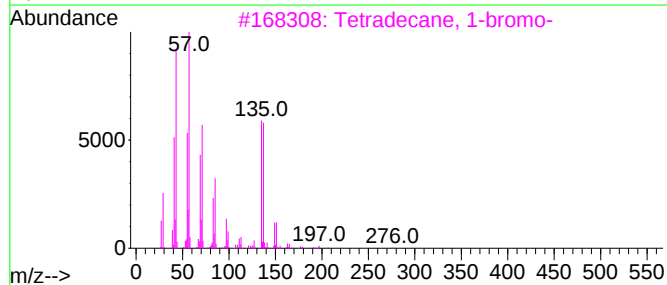
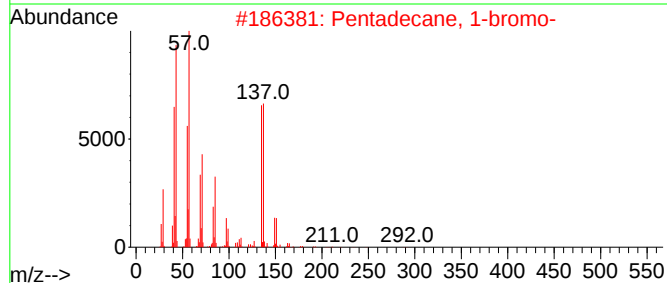
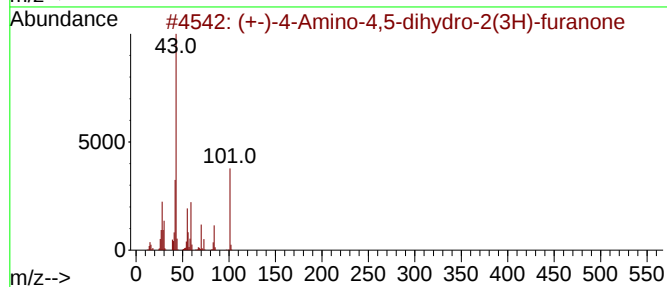
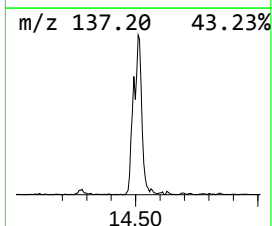
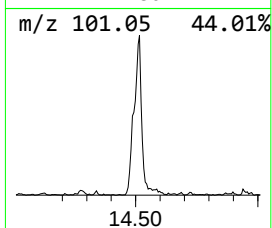
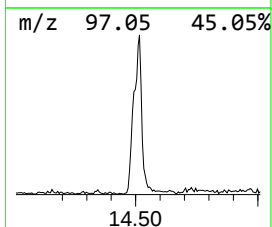
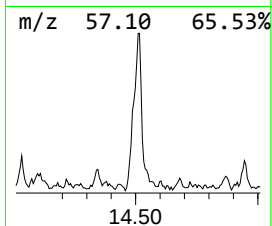
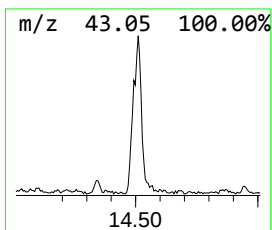
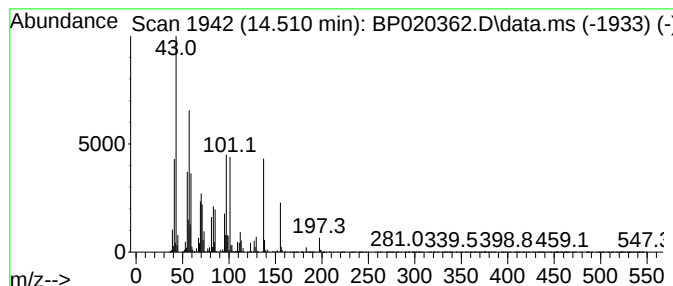
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.510	11.02 ng/ul	334707	Acenaphthene-d10	14.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	20		
2	Pentadecane, 1-bromo-	290	C15H31Br	000629-72-1	20		
3	Tetradecane, 1-bromo-	276	C14H29Br	000112-71-0	12		
4	Heptadecane, 1-bromo-	318	C17H35Br	003508-00-7	10		
5	1,3-Dioxane, 4,4-dimethyl-	116	C6H12O2	000766-15-4	10		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
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Acq On : 13 May 2024 20:47
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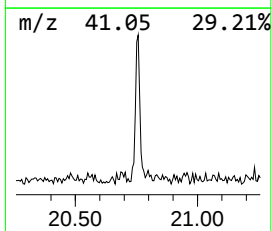
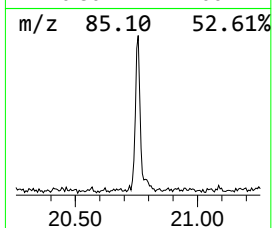
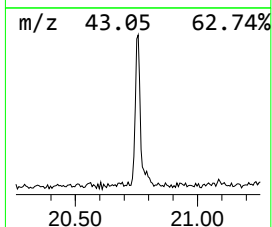
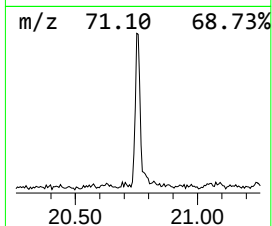
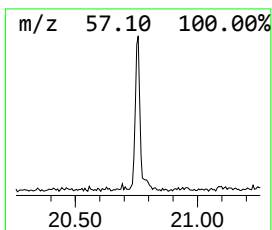
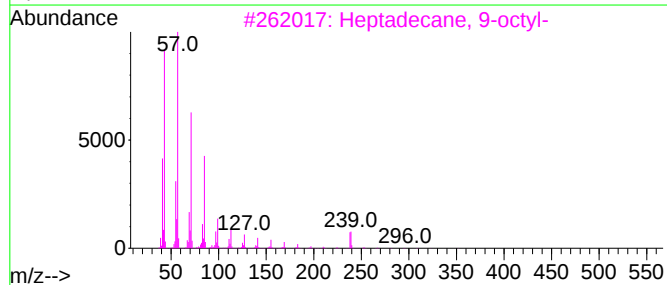
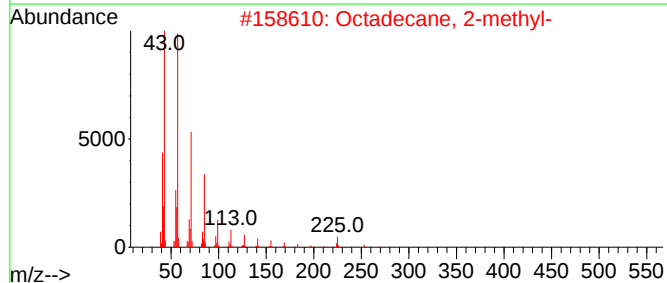
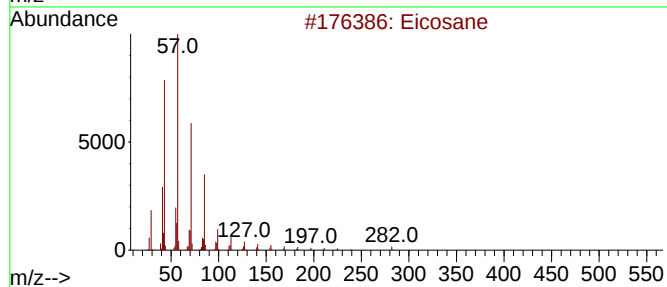
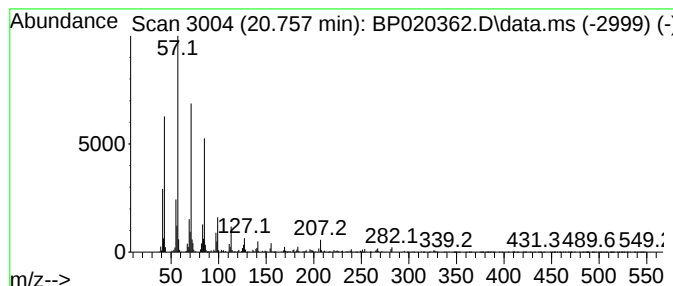
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.757	2.56 ng/ul	147543	Chrysene-d12	21.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	95
2		Octadecane, 2-methyl-	268	C19H40	001560-88-9	89
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
4		Tetracosane, 1-iodo-	464	C24H49I	1000406-32-0	87
5		Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020362.D
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Sample : P2369-11MEDL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

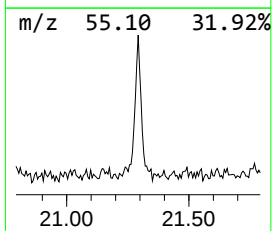
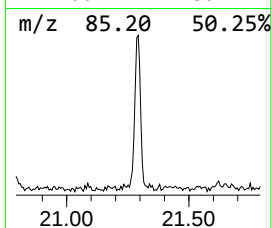
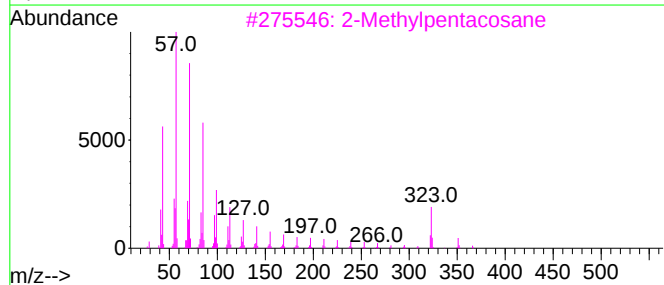
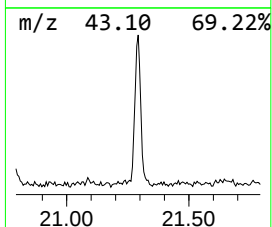
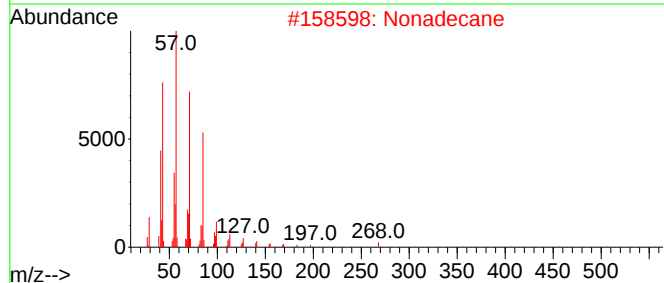
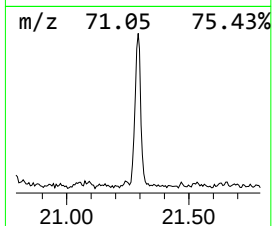
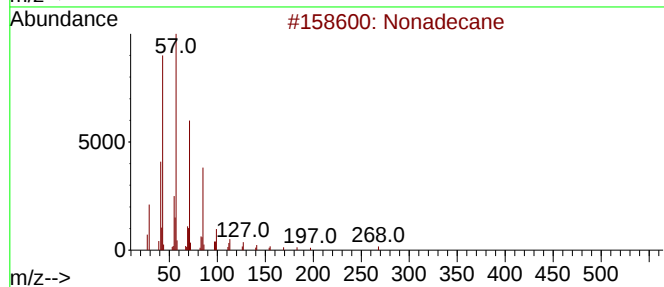
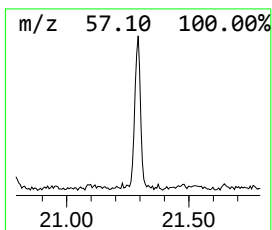
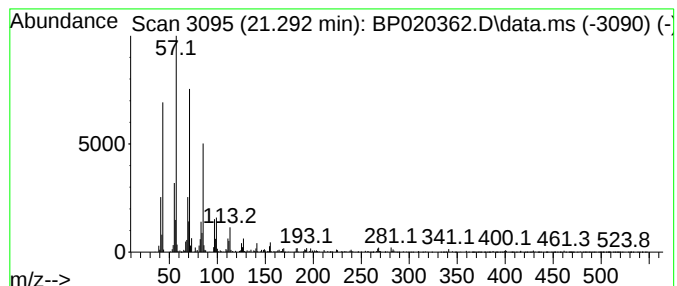
Instrument :
BNA_P
ClientSampleId :
A43N5MEDL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.292	2.82 ng/ul	162834	Chrysene-d12		21.622	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	95
2	Nonadecane		268	C19H40	000629-92-5	95
3	2-Methylpentacosane		366	C26H54	000629-87-8	95
4	Eicosane		282	C20H42	000112-95-8	93
5	Octacosane		394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020362.D
Acq On : 13 May 2024 20:47
Operator : MA/JU
Sample : P2369-11MEDL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N5MEDL

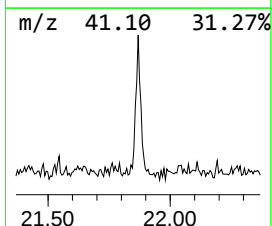
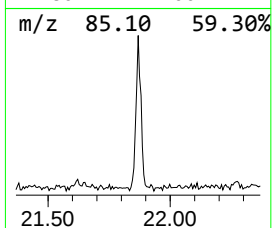
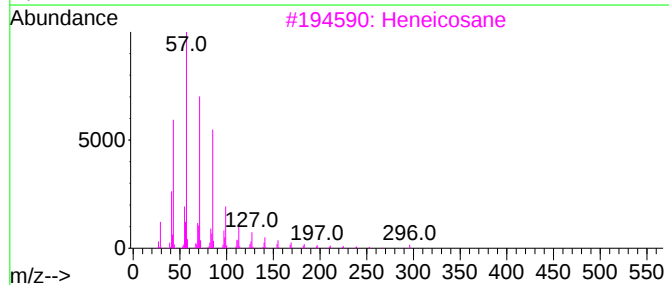
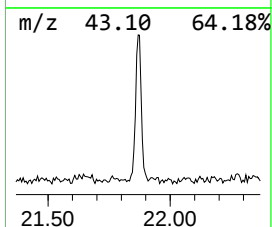
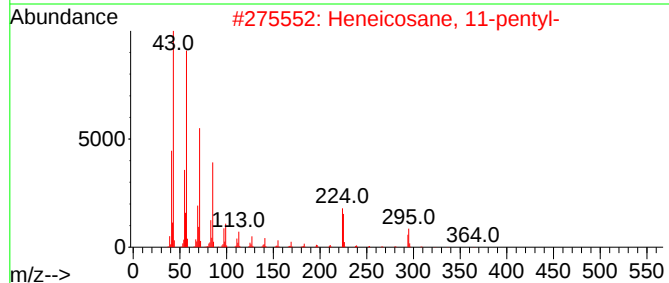
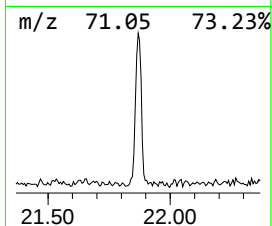
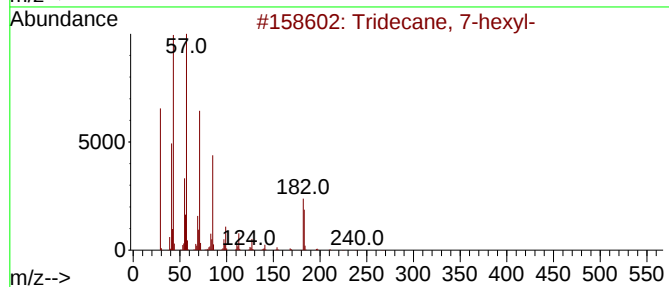
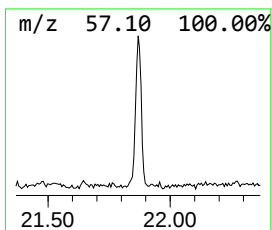
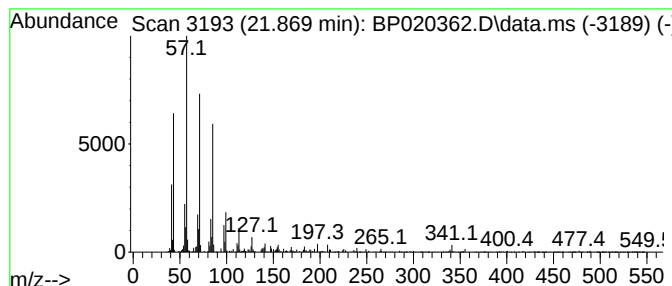
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.869	2.05 ng/ul	118404	Chrysene-d12	21.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	93
2		Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Docosane, 9-butyl-	366	C26H54	055282-14-9	90
5		Docosane, 5-butyl-	366	C26H54	055282-16-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020362.D
Acq On : 13 May 2024 20:47
Operator : MA/JU
Sample : P2369-11MEDL 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N5MEDL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.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
(DEL) Alkane: S...	6.734	2.4	ng/u1	39401	1	7.611	327041	20.0
(DEL) Alkane: S...	7.381	2.3	ng/u1	37809	1	7.611	327041	20.0
Crotyl methacry...	7.505	3.4	ng/u1	55146	1	7.611	327041	20.0
(DEL) Alkane: S...	7.763	2.8	ng/u1	45479	1	7.611	327041	20.0
(DEL) Alkane: C...	8.128	2.9	ng/u1	47059	1	7.611	327041	20.0
(DEL) Alkane: C...	8.669	4.6	ng/u1	75092	1	7.611	327041	20.0
Sulfurous acid,...	8.875	2.0	ng/u1	32971	1	7.611	327041	20.0
unknown-01	14.510	11.0	ng/u1	334707	3	14.281	607521	20.0
(DEL) Alkane: S...	20.757	2.6	ng/u1	147543	5	21.622	1154080	20.0
(DEL) Alkane: S...	21.292	2.8	ng/u1	162834	5	21.622	1154080	20.0
(DEL) Alkane: S...	21.869	2.0	ng/u1	118404	5	21.622	1154080	20.0