

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
 Data File : BP008527.D
 Acq On : 23 Dec 2021 16:34
 Operator : CG/JU
 Sample : M5171-09ME
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGLA5ME

Integration Parameters: LSCINT.P

Integrator: RTE

Smoother: 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: BP008527.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.099	183	189	202	rVB	6516586	9385938	6.59%	0.633%
2	5.422	407	414	422	rVV	12266320	19867447	13.94%	1.341%
3	5.558	431	437	455	rVB	25745914	61583367	43.22%	4.157%
4	5.928	493	500	510	rVB2	14445999	28292623	19.86%	1.910%
5	6.405	574	581	589	rBV	6397901	10303300	7.23%	0.695%
6	6.893	657	664	667	rBV2	4172631	8176944	5.74%	0.552%
7	7.005	673	683	688	rBV2	10426479	20887090	14.66%	1.410%
8	7.063	688	693	701	rVB2	6318581	14152563	9.93%	0.955%
9	7.140	701	706	714	rVB	7060803	10999489	7.72%	0.742%
10	7.305	729	734	737	rBV	4510270	6808868	4.78%	0.460%
11	7.558	770	777	785	rBV2	24814499	49252044	34.56%	3.324%
12	7.910	831	837	842	rVB	6527838	10256740	7.20%	0.692%
13	8.028	848	857	867	rVB2	5327893	12125503	8.51%	0.818%
14	8.210	882	888	893	rVB2	2221013	4804423	3.37%	0.324%
15	8.281	893	900	907	rVB3	3174524	7148729	5.02%	0.482%
16	8.463	925	931	937	rVB3	6320425	11916778	8.36%	0.804%
17	8.557	943	947	950	rVV2	5112860	7316323	5.13%	0.494%
18	8.593	950	953	962	rVB2	4933352	10063817	7.06%	0.679%
19	8.693	962	970	978	rBV2	3556027	9460230	6.64%	0.639%
20	9.022	1018	1026	1036	rBV3	5008783	12201263	8.56%	0.824%
21	9.163	1044	1050	1060	rVB	18082501	31480976	22.09%	2.125%
22	9.281	1060	1070	1076	rBV7	2673638	7978906	5.60%	0.539%
23	9.399	1086	1090	1098	rVB5	1692898	4440383	3.12%	0.300%
24	9.604	1119	1125	1130	rVB2	2679884	4778800	3.35%	0.323%
25	10.116	1200	1212	1217	rBV9	4624540	17459356	12.25%	1.178%
26	10.322	1242	1247	1257	rVB5	1966244	6086926	4.27%	0.411%
27	10.710	1306	1313	1317	rVV2	13524268	23141663	16.24%	1.562%
28	10.781	1317	1325	1331	rVV2	37501313	92755378	65.10%	6.260%
29	10.863	1331	1339	1342	rVV2	3072774	8278232	5.81%	0.559%
30	10.893	1342	1344	1349	rVV	2971684	4347413	3.05%	0.293%
31	11.651	1465	1473	1479	rBV3	1801869	4178513	2.93%	0.282%
32	11.763	1485	1492	1499	rBV2	4067675	8415129	5.91%	0.568%
33	12.157	1552	1559	1567	rBV	11704824	18859426	13.24%	1.273%
34	12.375	1588	1596	1602	rVB	19256720	33749847	23.69%	2.278%
35	12.587	1626	1632	1637	rBV	8916220	14245727	10.00%	0.962%
36	12.969	1691	1697	1705	rVB2	4836469	8641670	6.06%	0.583%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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

37	13.104	1715	1720	1724	rVB	3554126	4905868	3.44%	0.331%
38	13.393	1761	1769	1775	rVB2	17595000	30165810	21.17%	2.036%
39	13.704	1816	1822	1823	rBV	4694842	6518171	4.57%	0.440%
40	13.869	1838	1850	1854	rBV2	10524172	20722605	14.54%	1.399%
41	13.916	1854	1858	1861	rVV	4485227	6401558	4.49%	0.432%
42	14.045	1872	1880	1883	rBV2	4314639	6837256	4.80%	0.461%
43	14.092	1883	1888	1892	rVV2	2683335	4687946	3.29%	0.316%
44	14.457	1945	1950	1955	rBV	15356442	21431625	15.04%	1.447%
45	14.534	1955	1963	1968	rBV3	4519873	9678979	6.79%	0.653%
46	14.669	1982	1986	1991	rVB	7761379	10241108	7.19%	0.691%
47	14.745	1991	1999	2010	rVB3	2256315	6322685	4.44%	0.427%
48	14.939	2022	2032	2036	rBV2	4696399	10545103	7.40%	0.712%
49	15.016	2041	2045	2049	rVV	3456469	5152455	3.62%	0.348%
50	15.075	2051	2055	2062	rVB3	3213290	5153071	3.62%	0.348%
51	15.145	2063	2067	2074	rBV2	2540145	5383780	3.78%	0.363%
52	15.357	2091	2103	2107	rBV5	3259491	8571509	6.02%	0.579%
53	15.410	2107	2112	2116	rBV	11105490	14172811	9.95%	0.957%
54	15.816	2174	2181	2185	rBV5	4034920	8909352	6.25%	0.601%
55	16.275	2254	2259	2261	rBV2	9948364	14271094	10.02%	0.963%
56	16.298	2261	2263	2268	rVB	5035006	6325913	4.44%	0.427%
57	16.386	2273	2278	2284	rBV	3506873	5685884	3.99%	0.384%
58	16.457	2285	2290	2297	rVB2	2922406	4382486	3.08%	0.296%
59	17.069	2389	2394	2398	rBV	7316682	9535768	6.69%	0.644%
60	17.204	2414	2417	2425	rVB2	3396905	5004220	3.51%	0.338%
61	17.286	2426	2431	2434	rVV	3451459	5359784	3.76%	0.362%
62	17.328	2434	2438	2442	rVV2	2886122	4758703	3.34%	0.321%
63	17.381	2442	2447	2456	rVB2	3600754	6856553	4.81%	0.463%
64	17.804	2515	2519	2524	rVB	6072483	7470921	5.24%	0.504%
65	18.198	2581	2586	2590	rVV	12824307	17892944	12.56%	1.208%
66	18.245	2590	2594	2599	rVB	5748664	7552892	5.30%	0.510%
67	18.475	2629	2633	2638	rVB	4552585	5101156	3.58%	0.344%
68	19.033	2721	2728	2732	rBV4	2593659	4781458	3.36%	0.323%
69	19.080	2732	2736	2741	rVB2	5521683	7212289	5.06%	0.487%
70	19.327	2766	2778	2787	rBV4	20736692	60032831	42.13%	4.052%
71	19.427	2792	2795	2803	rVV	6784046	9917365	6.96%	0.669%
72	19.498	2803	2807	2814	rVB2	5410525	7346160	5.16%	0.496%
73	19.616	2822	2827	2828	rBV	3867169	4891566	3.43%	0.330%
74	19.639	2828	2831	2835	rVB2	6256679	8039370	5.64%	0.543%
75	19.916	2874	2878	2882	rBV	4265679	5652390	3.97%	0.382%
76	20.157	2916	2919	2924	rVB	5685553	6787004	4.76%	0.458%

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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
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77	20.539	2976	2984	2988	rBV4	51140701	142492072	100.00%	9.617%
78	20.645	2998	3002	3006	rBV	5269808	5748980	4.03%	0.388%
79	20.857	3033	3038	3045	rVB	9768285	12096801	8.49%	0.816%
80	21.098	3074	3079	3084	rVB	14387063	18501297	12.98%	1.249%
81	21.298	3107	3113	3120	rVV2	47161734	98412370	69.07%	6.642%
82	21.369	3122	3125	3130	rVB	5736647	7448986	5.23%	0.503%
83	21.539	3147	3154	3169	rVB3	8442763	14220339	9.98%	0.960%
84	21.857	3193	3208	3211	rBV2	6313804	21319618	14.96%	1.439%
85	21.916	3211	3218	3222	rVV3	12207249	22551797	15.83%	1.522%
86	22.016	3230	3235	3240	rBV2	9694187	14945328	10.49%	1.009%
87	22.163	3256	3260	3263	rBV	4712754	6994540	4.91%	0.472%
88	22.233	3268	3272	3279	rVB	7340678	12841004	9.01%	0.867%
89	22.521	3317	3321	3328	rVB	4954498	7900109	5.54%	0.533%
90	22.592	3329	3333	3338	rVB	3782334	5463491	3.83%	0.369%
91	22.857	3373	3378	3383	rBV2	3200270	6207633	4.36%	0.419%
92	23.098	3415	3419	3425	rVB	3823964	6033743	4.23%	0.407%
93	23.351	3458	3462	3471	rVB	6199231	14546567	10.21%	0.982%
94	23.539	3489	3494	3505	rVB4	3468579	9122498	6.40%	0.616%
95	23.745	3524	3529	3534	rVV2	2826019	5133693	3.60%	0.346%
96	23.804	3534	3539	3553	rVB2	3206346	8170610	5.73%	0.551%
97	24.492	3650	3656	3665	rVB	2286682	5200485	3.65%	0.351%
98	24.804	3703	3709	3722	rVB	3399541	7847567	5.51%	0.530%
99	24.927	3724	3730	3738	rVB2	3005564	6928873	4.86%	0.468%
100	25.780	3868	3875	3896	rVB3	1593595	6981483	4.90%	0.471%

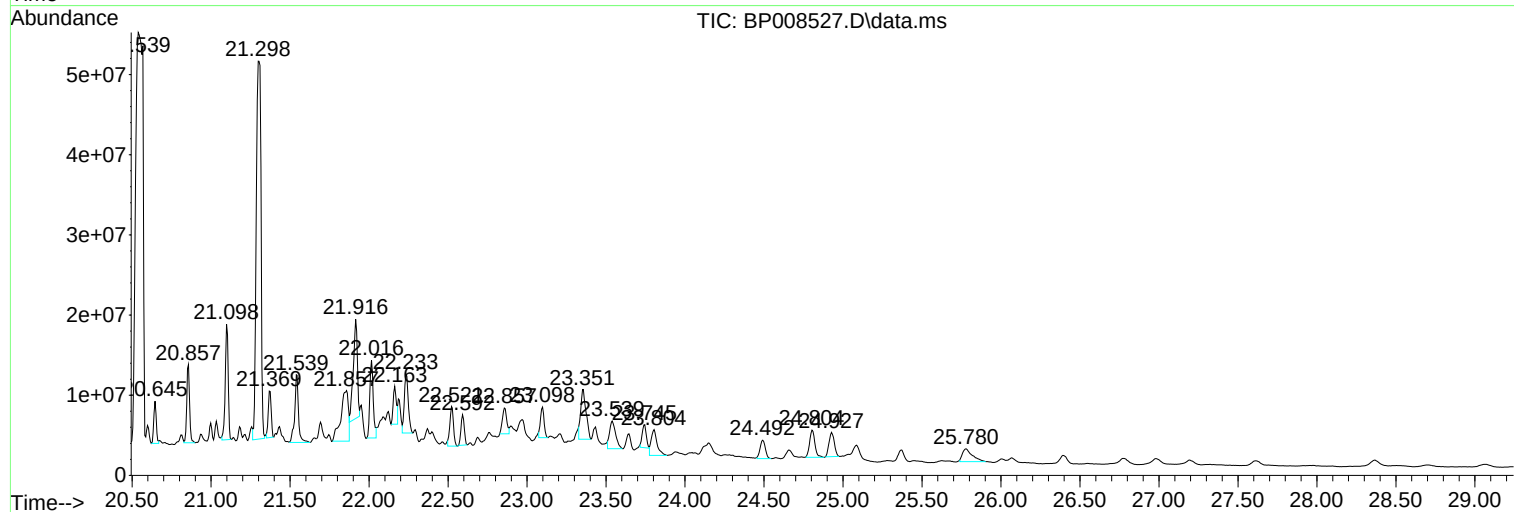
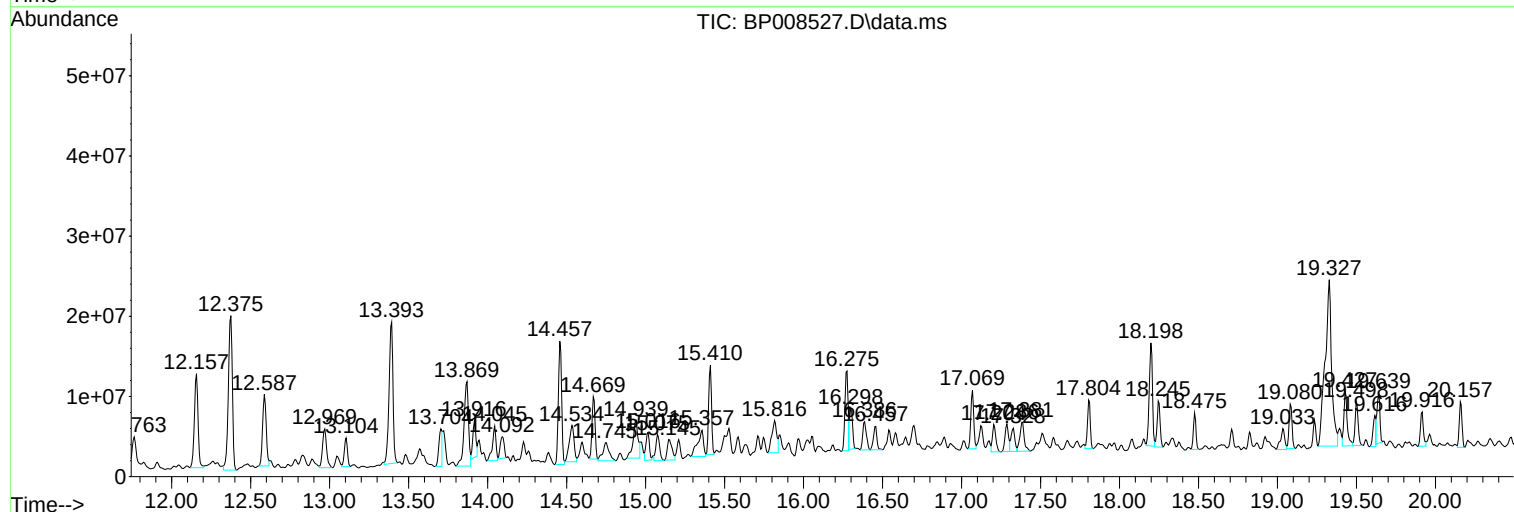
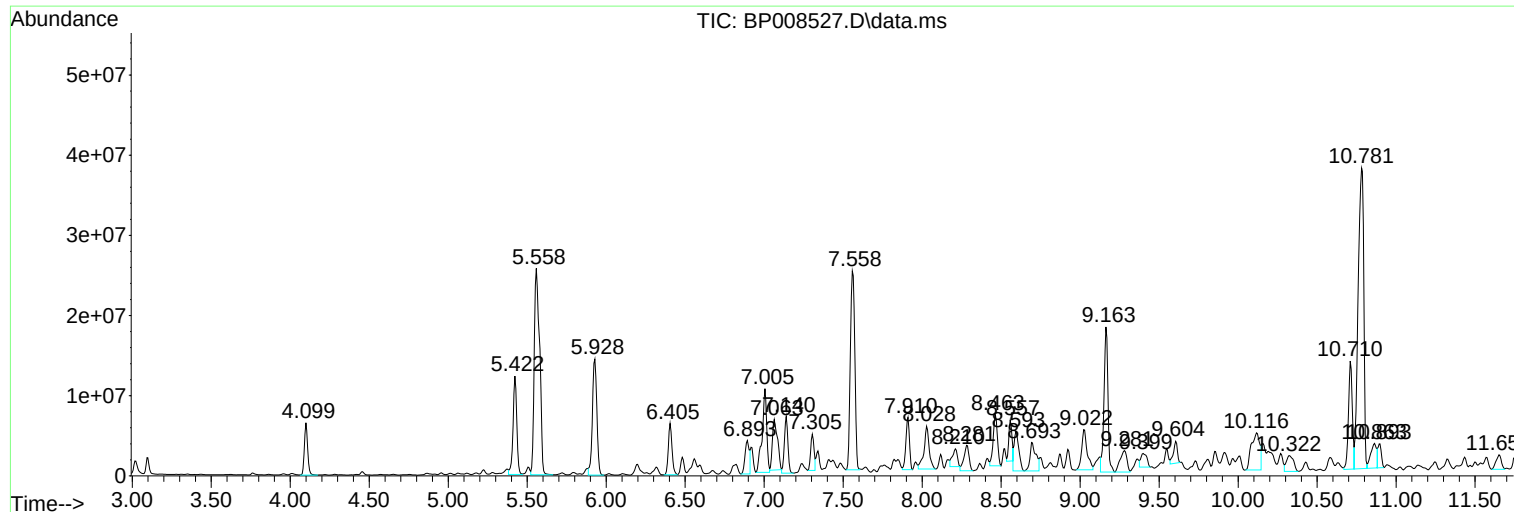
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Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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



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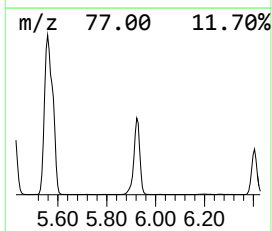
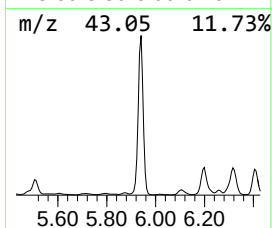
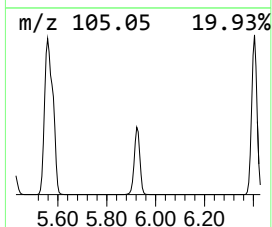
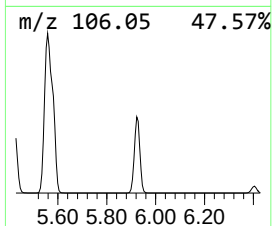
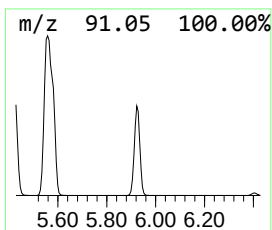
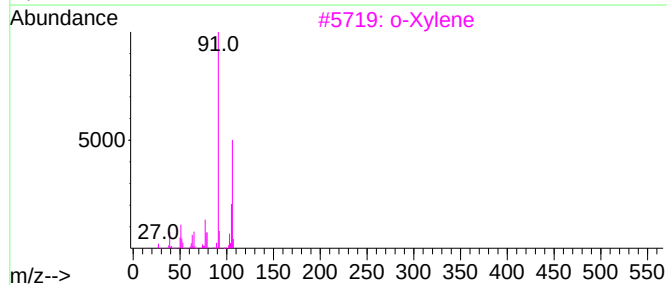
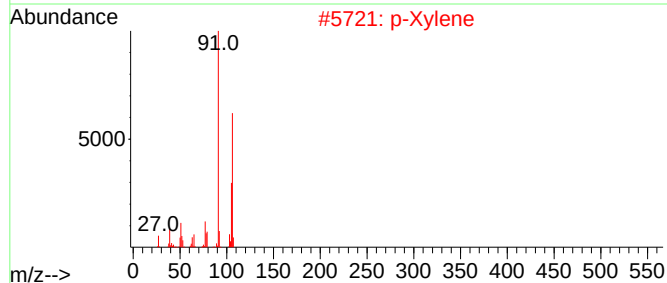
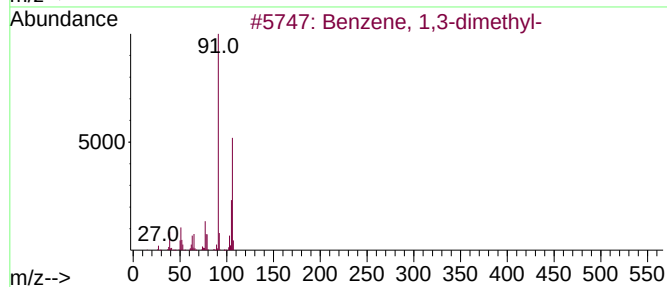
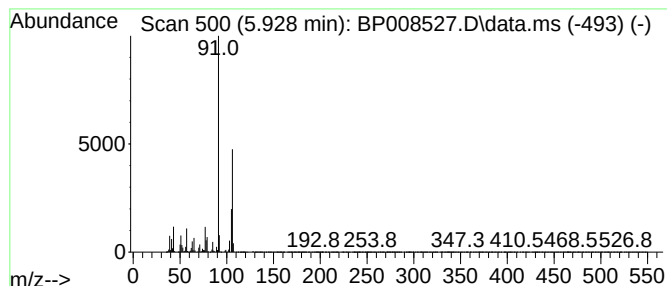
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.928	55.17 ng/ul	28292600	1,4-Dichlorobenzene-d4	7.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94
2			p-Xylene	106	C8H10	000106-42-3	94
3			o-Xylene	106	C8H10	000095-47-6	94
4			p-Xylene	106	C8H10	000106-42-3	93
5			o-Xylene	106	C8H10	000095-47-6	93



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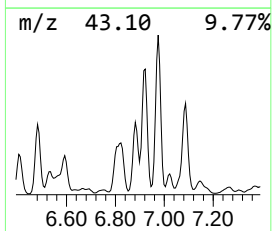
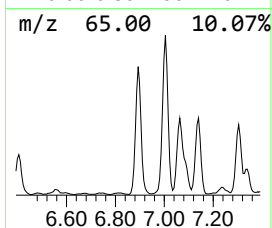
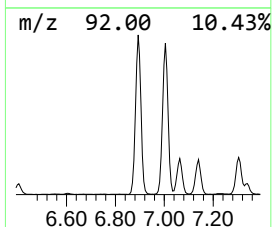
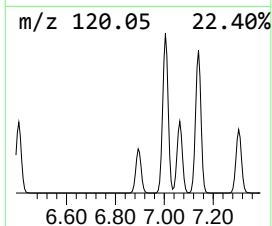
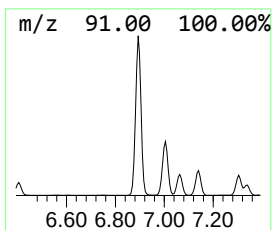
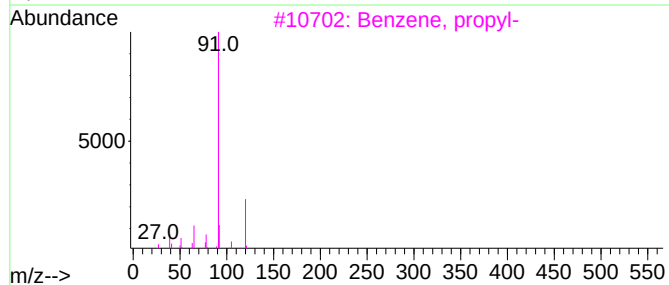
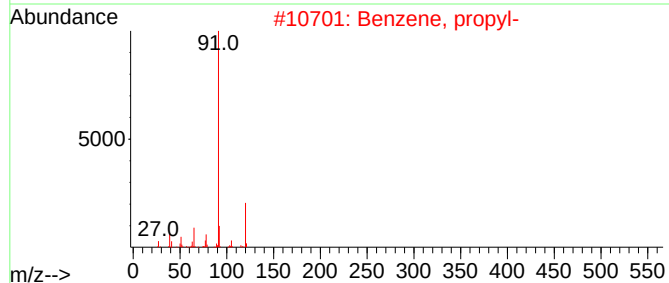
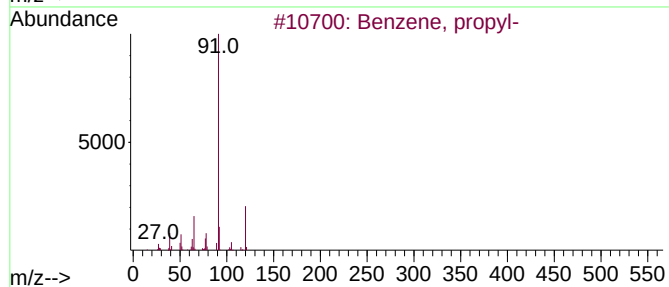
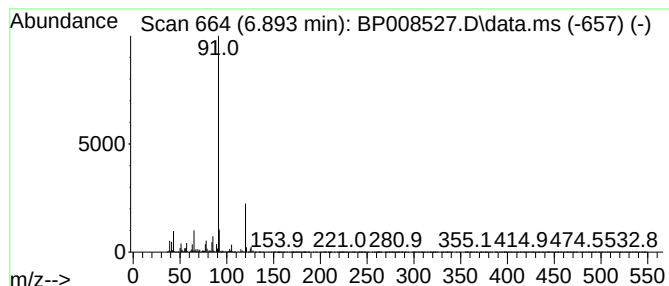
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, propyl- Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.893	15.94 ng/ul	8176940	1,4-Dichlorobenzene-d4	7.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	81
2			Benzene, propyl-	120	C9H12	000103-65-1	76
3			Benzene, propyl-	120	C9H12	000103-65-1	76
4			Benzene, propyl-	120	C9H12	000103-65-1	68
5			1-(Benzylamino)propan-2-ol	165	C10H15NO	1000486-84-7	64



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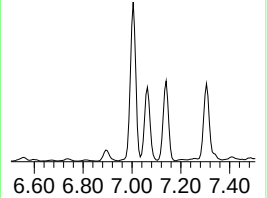
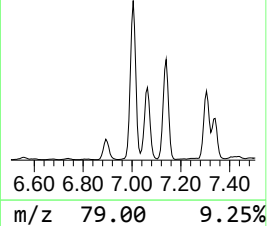
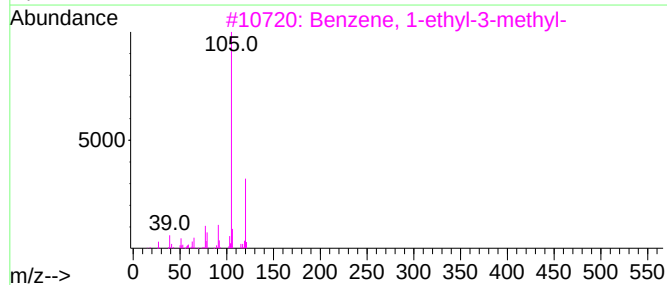
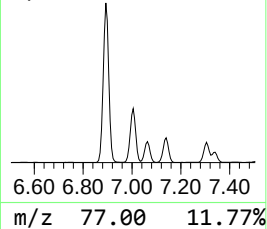
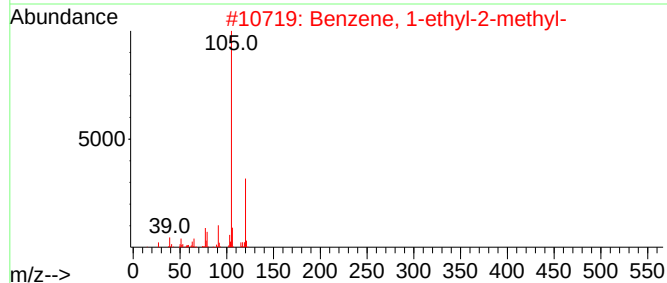
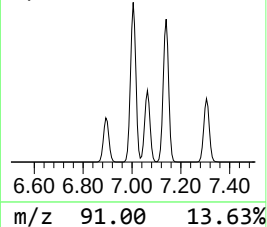
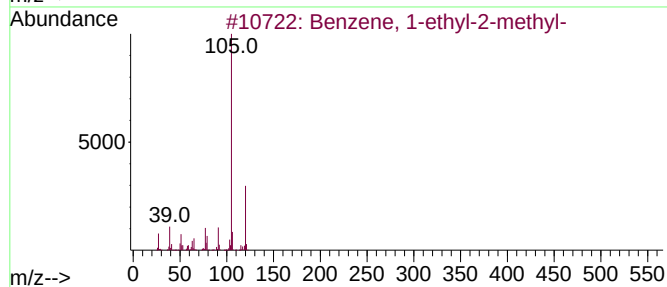
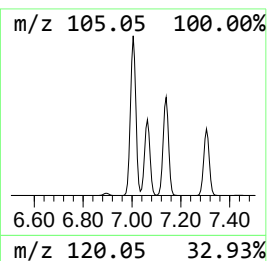
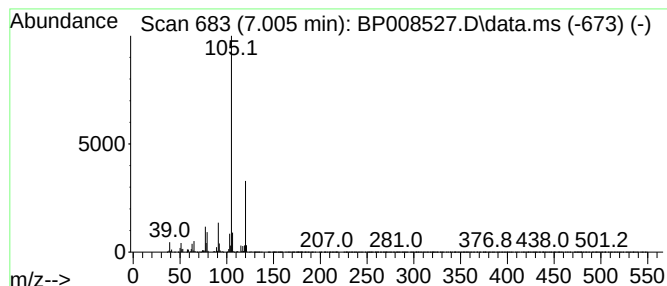
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.005	40.73 ng/ul	20887100	1,4-Dichlorobenzene-d4	7.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGLA5ME

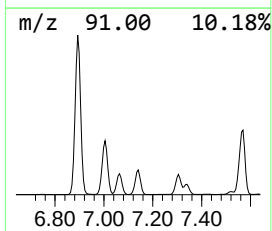
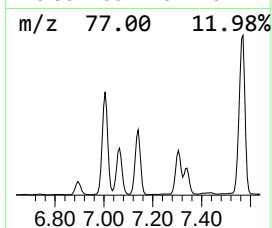
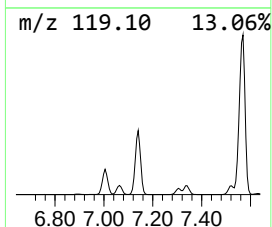
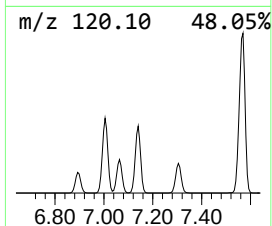
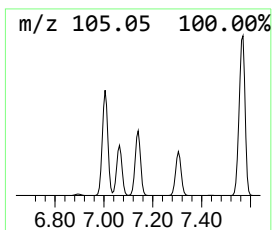
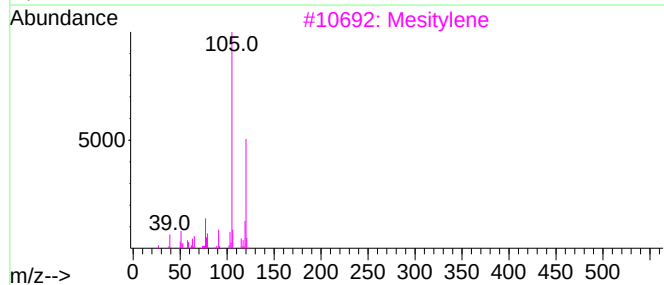
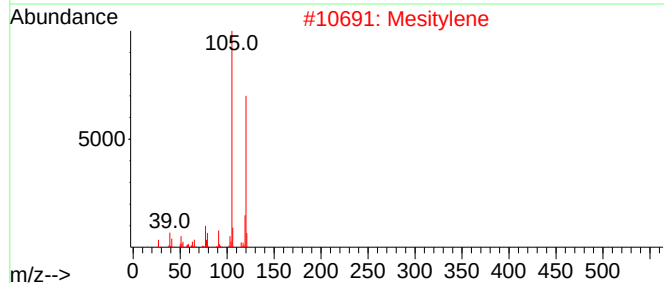
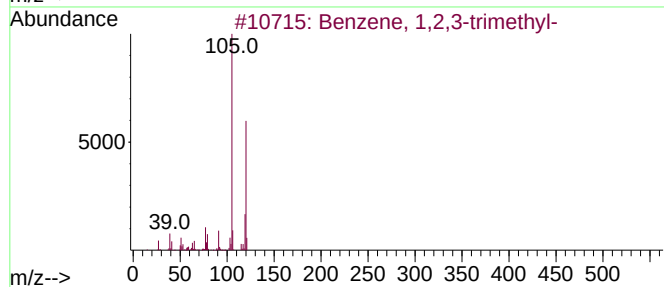
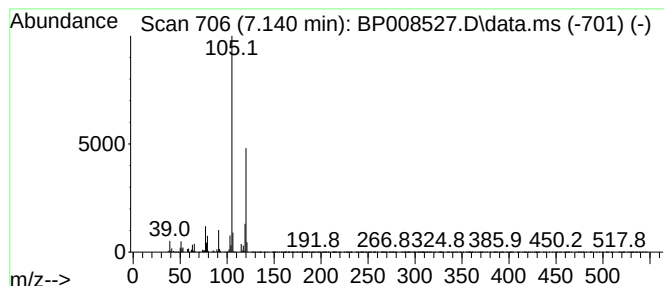
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.140	21.45 ng/ul	10999500	1,4-Dichlorobenzene-d4	7.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2			Mesitylene	120	C9H12	000108-67-8	97
3			Mesitylene	120	C9H12	000108-67-8	97
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5			Mesitylene	120	C9H12	000108-67-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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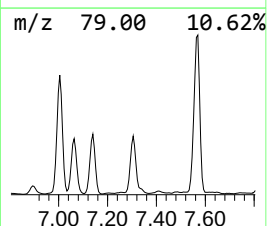
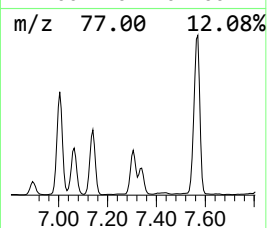
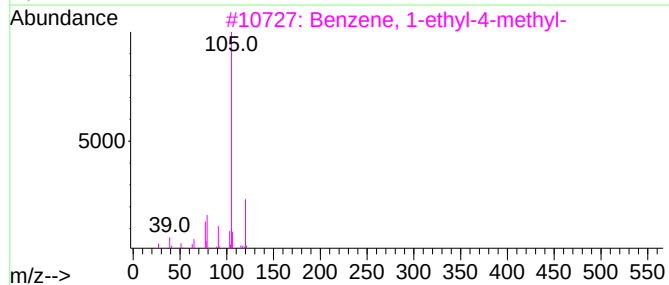
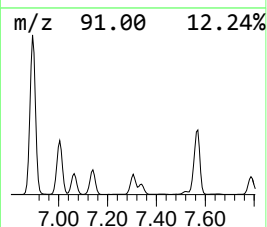
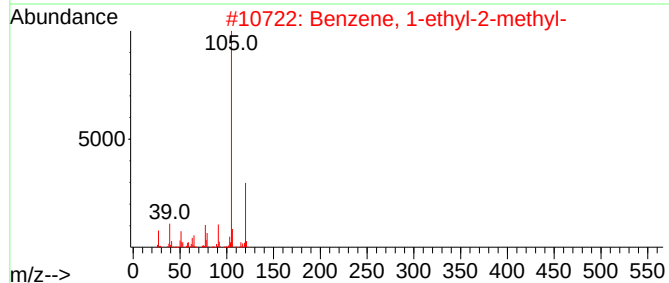
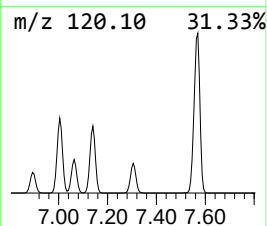
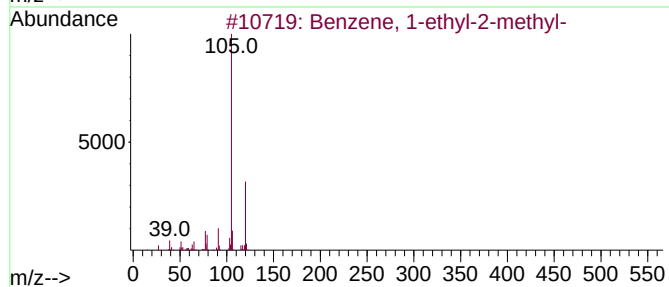
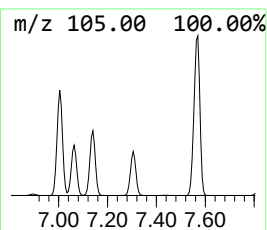
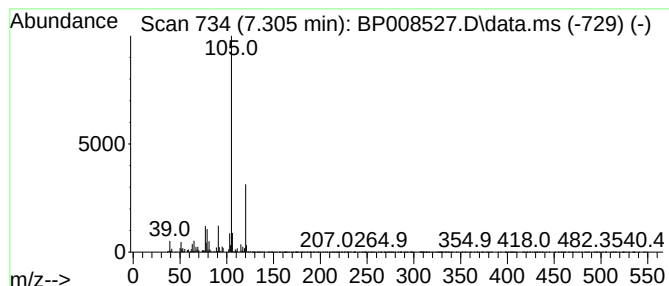
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-ethyl-4-methyl- Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.305	13.28 ng/ul	6808870	1,4-Dichlorobenzene-d4	7.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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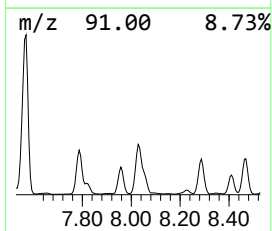
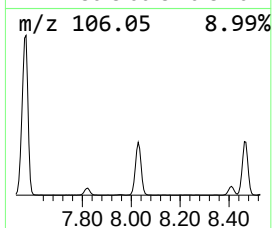
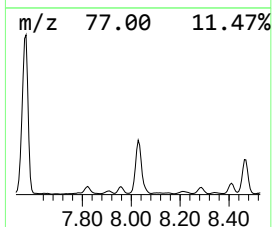
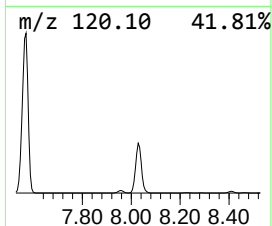
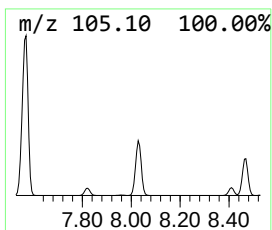
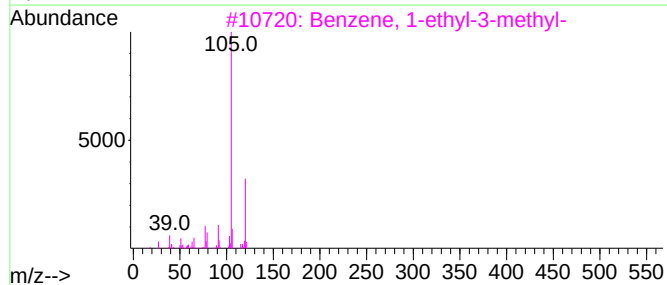
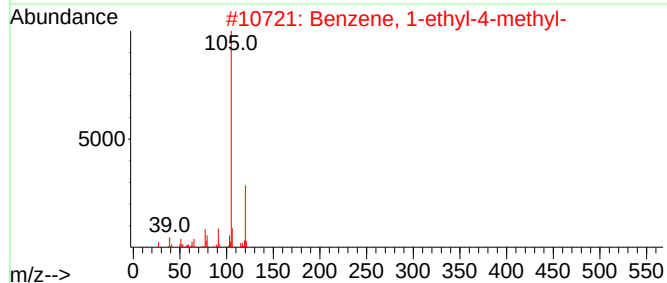
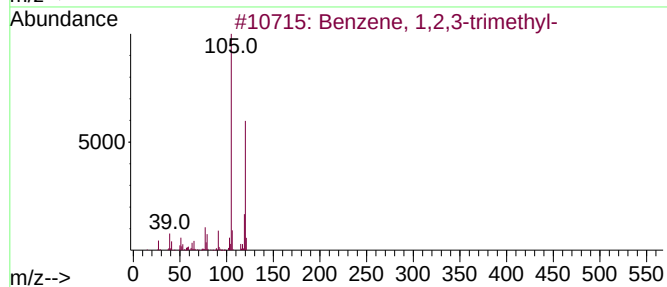
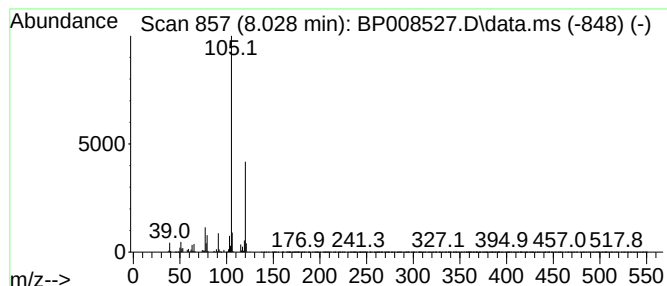
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzene, 1-ethyl-3-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.028	23.64 ng/ul	12125500	1,4-Dichlorobenzene-d4	7.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
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Sample : M5171-09ME
Misc :
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Instrument :
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ClientSampleId :
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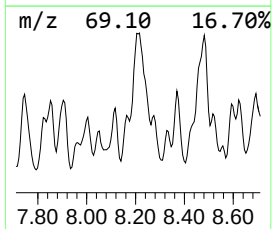
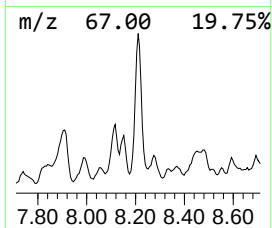
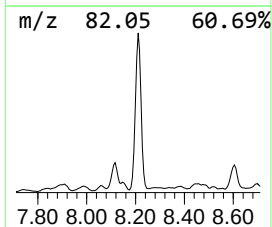
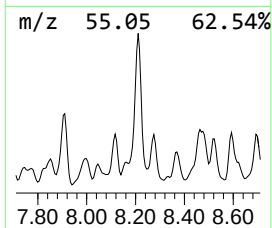
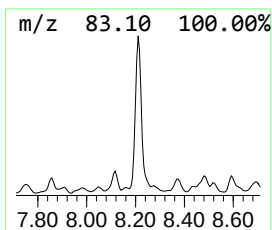
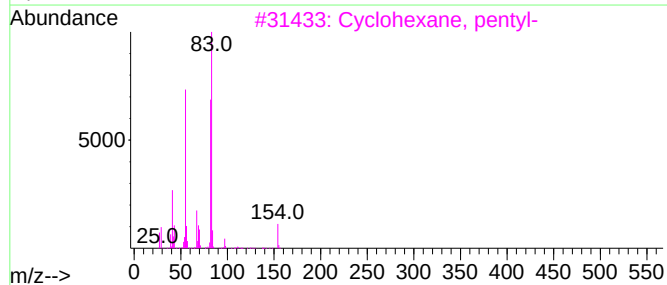
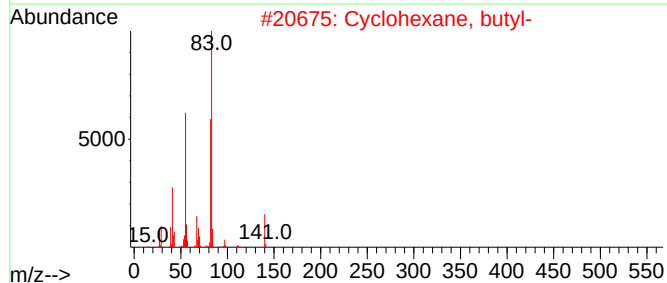
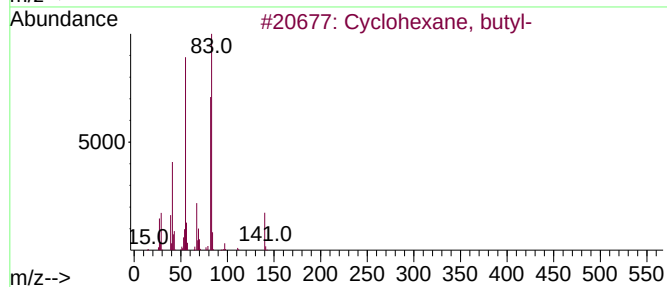
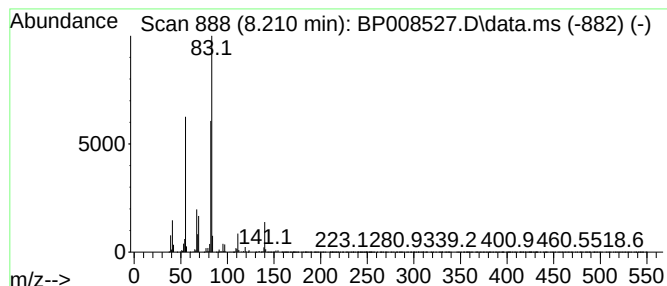
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Cyclic8.210 Concentration Rank 62

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.210	9.37 ng/ul	4804420	1,4-Dichlorobenzene-d4	7.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, butyl-	140	C10H20	001678-93-9	90
2			Cyclohexane, butyl-	140	C10H20	001678-93-9	86
3			Cyclohexane, pentyl-	154	C11H22	004292-92-6	78
4			Cyclohexane, propyl-	126	C9H18	001678-92-8	78
5			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
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Sample : M5171-09ME
Misc :
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Instrument :
BNA_P
ClientSampleId :
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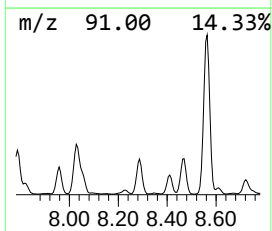
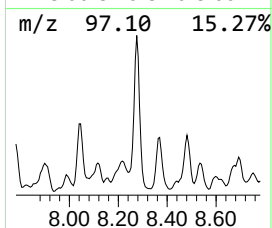
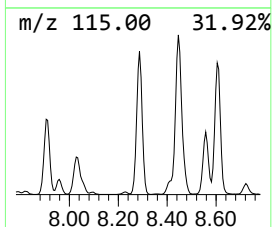
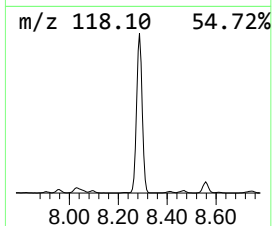
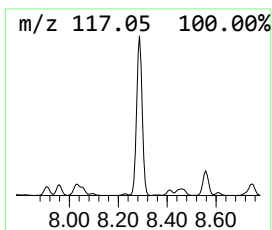
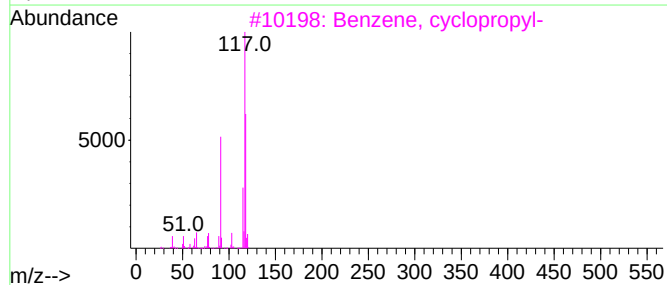
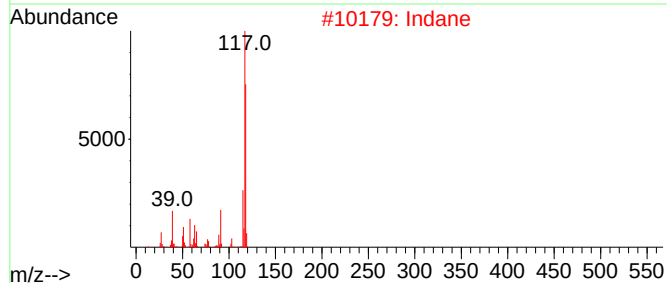
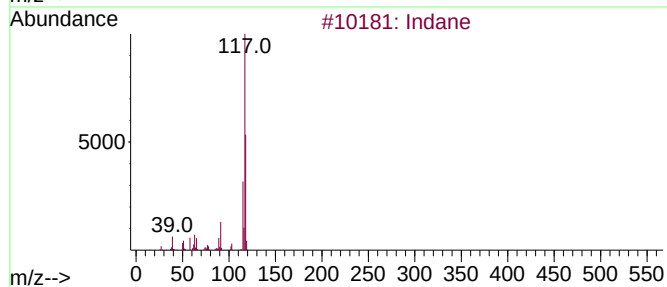
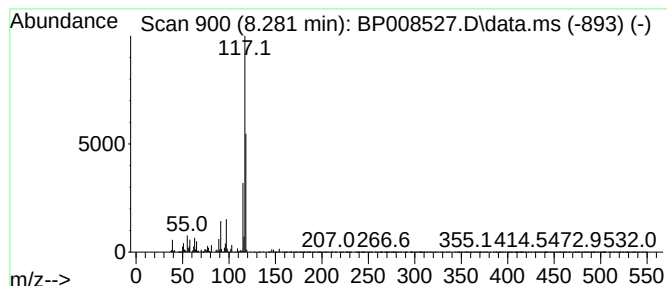
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Indane Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.281	13.94 ng/ul	7148730	1,4-Dichlorobenzene-d4	7.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	76
2	Indane			118	C9H10	000496-11-7	76
3	Benzene, cyclopropyl-			118	C9H10	000873-49-4	64
4	Indane			118	C9H10	000496-11-7	59
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-			118	C9H10	1000191-13-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
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Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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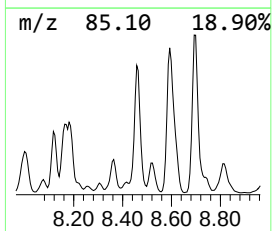
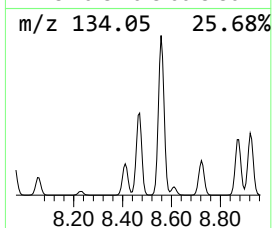
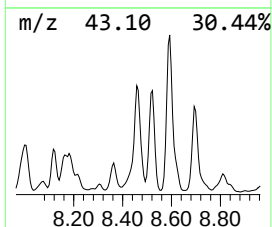
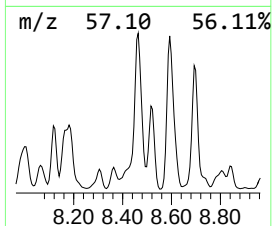
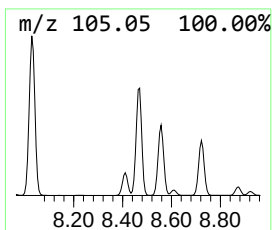
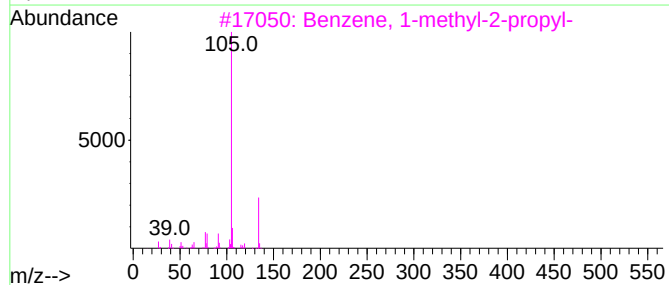
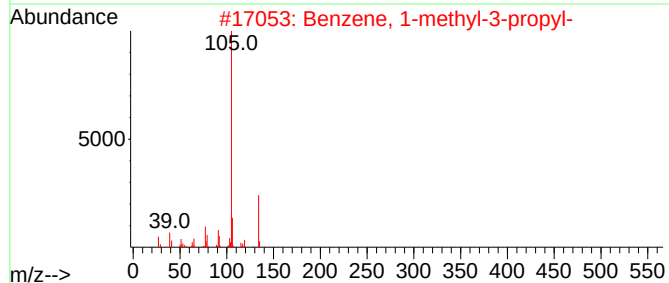
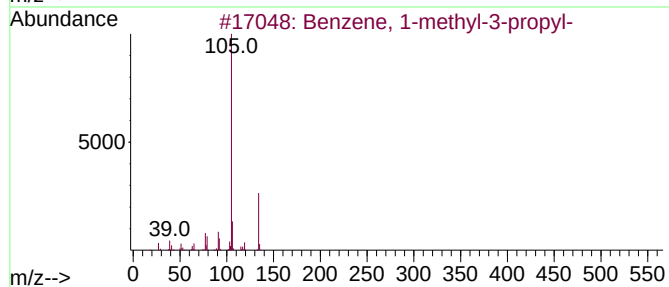
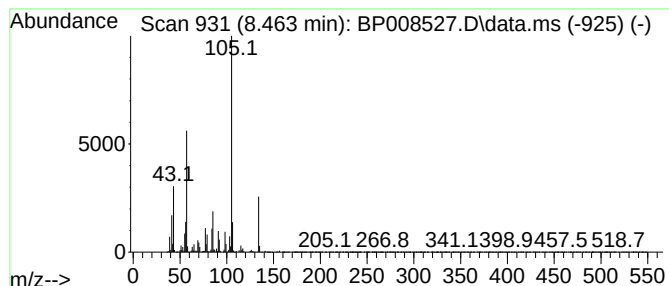
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Benzene, 1-methyl-3-propyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.463	23.24 ng/ul	11916800	1,4-Dichlorobenzene-d4	7.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	64
2			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	64
3			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	64
4			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	64
5			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
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Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
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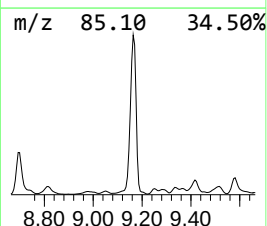
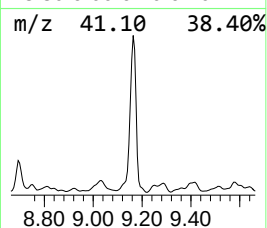
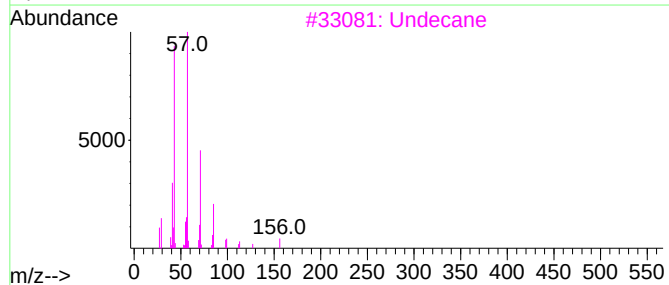
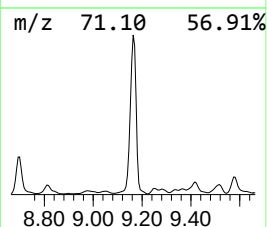
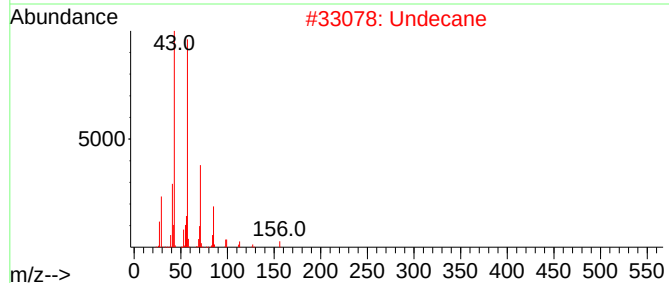
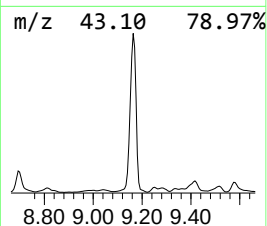
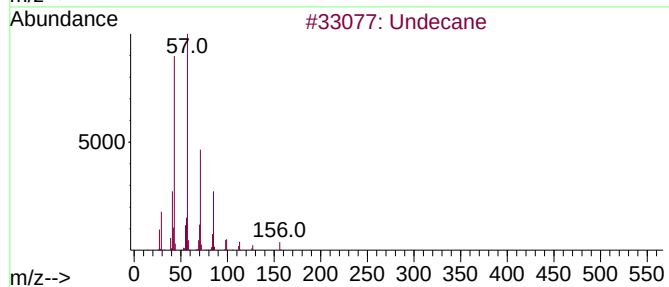
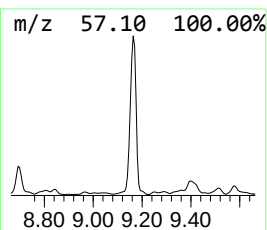
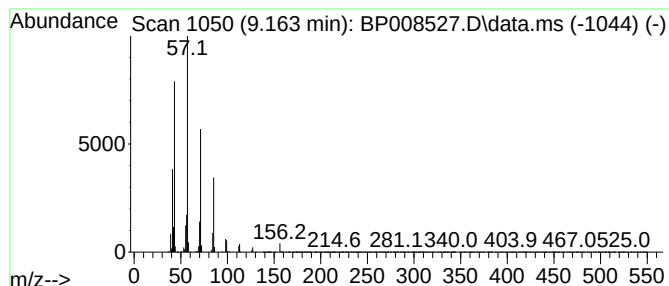
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.163	61.39 ng/ul	31481000	1,4-Dichlorobenzene-d4	7.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	97
2	Undecane			156	C11H24	001120-21-4	91
3	Undecane			156	C11H24	001120-21-4	90
4	Decane			142	C10H22	000124-18-5	90
5	Tetradecane			198	C14H30	000629-59-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

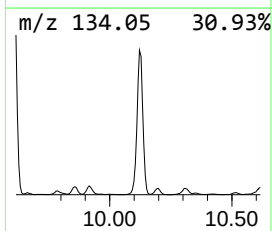
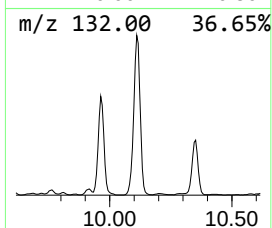
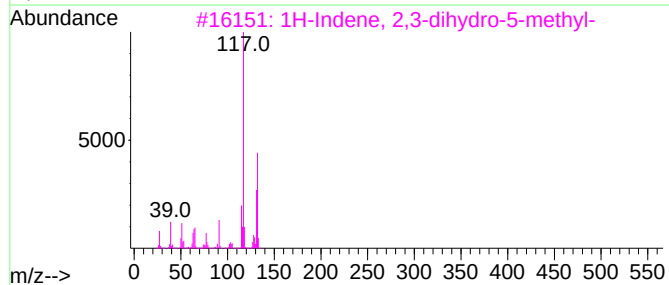
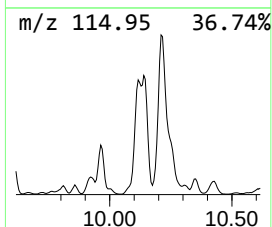
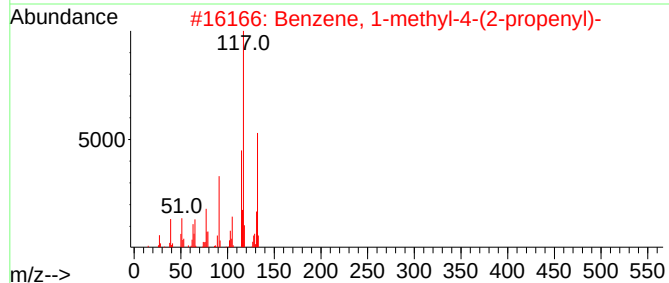
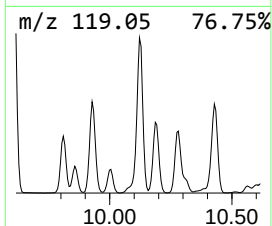
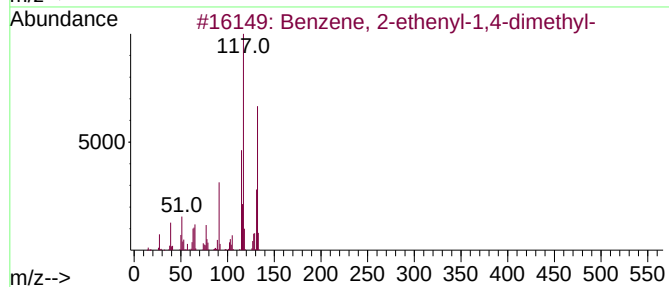
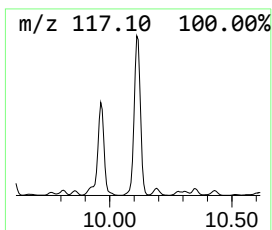
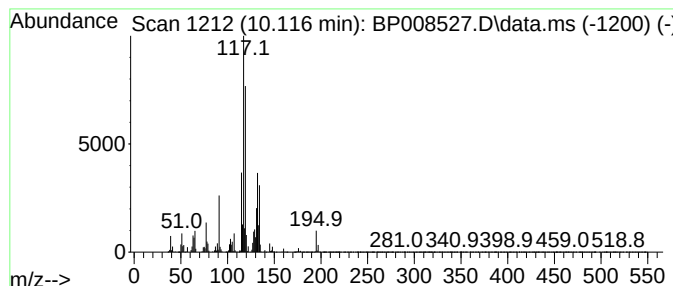
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.116	15.09 ng/ul	17459400	Naphthalene-d8	10.716	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	74
2	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	60
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	50
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	50
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
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Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

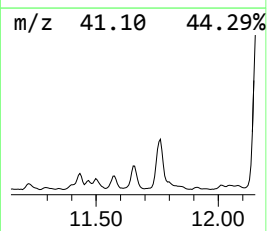
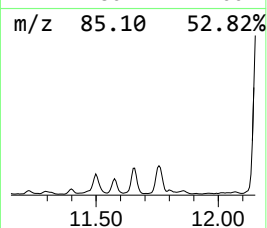
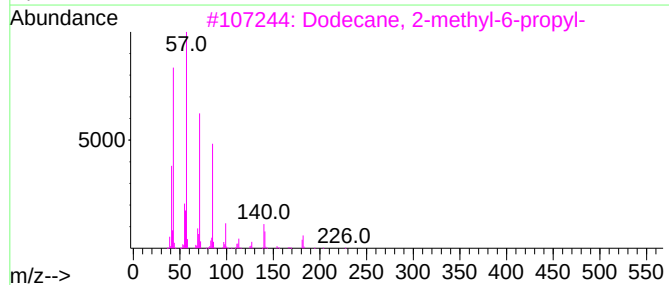
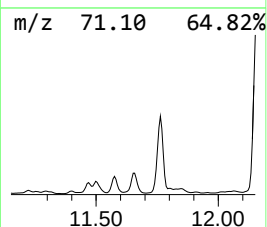
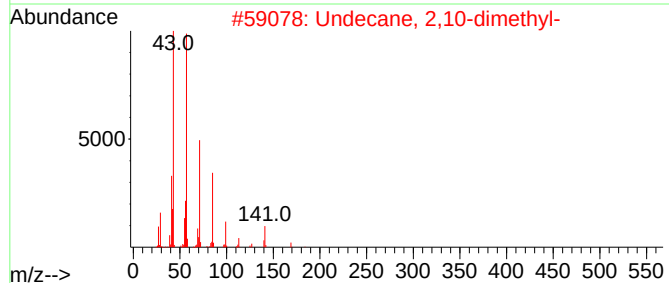
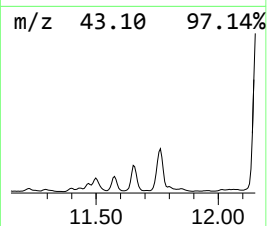
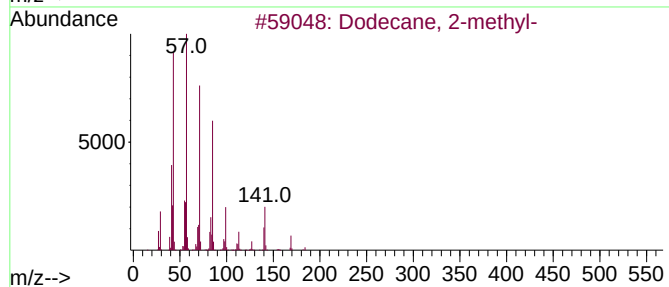
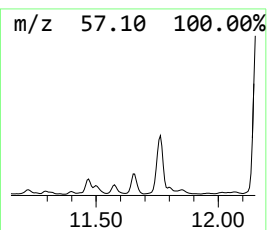
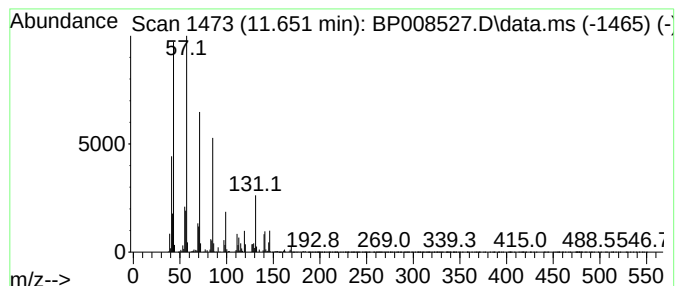
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 66

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.651	3.61 ng/ul	4178510	Naphthalene-d8	10.716	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2-methyl-	184	C13H28	001560-97-0	93
2	Undecane, 2,10-dimethyl-	184	C13H28	017301-27-8	70
3	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	64
4	Tridecane, 7-propyl-	226	C16H34	055045-09-5	64
5	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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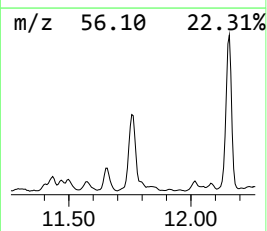
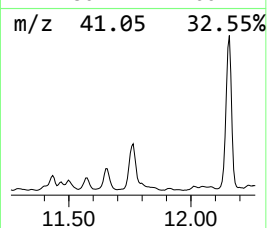
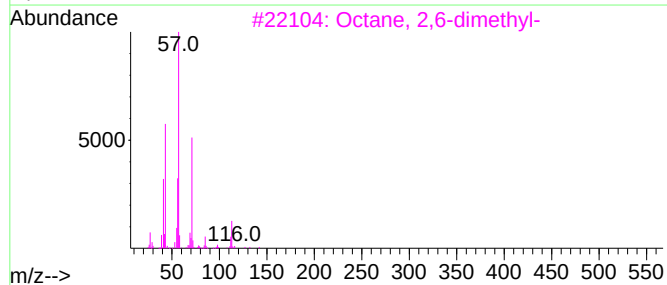
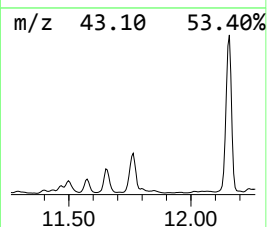
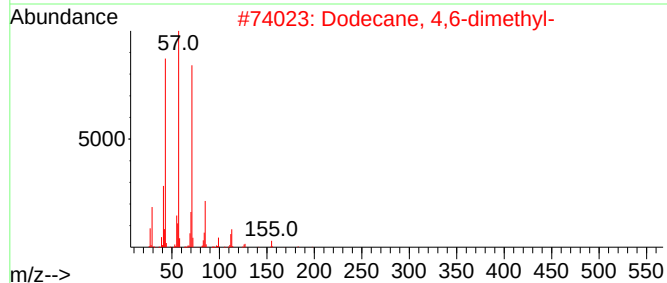
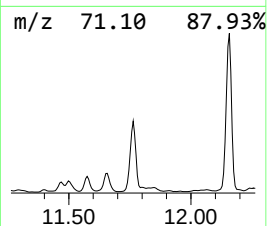
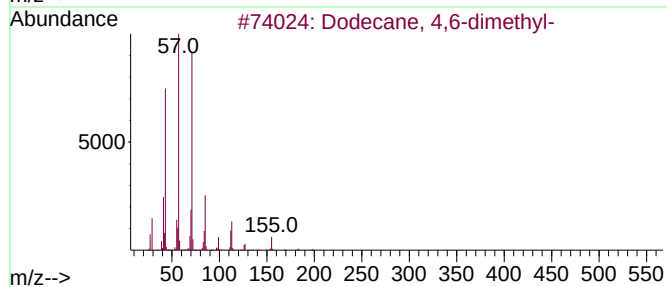
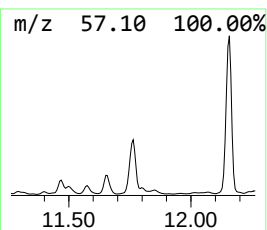
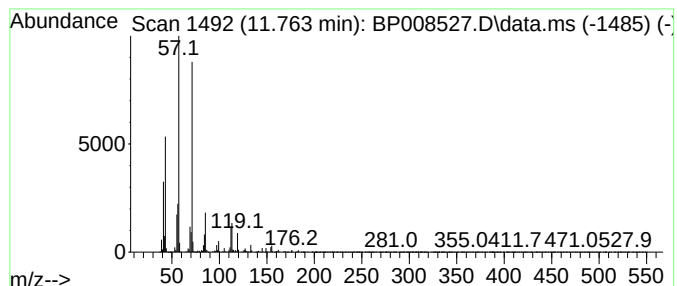
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 63

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.763	7.27 ng/ul	8415130	Naphthalene-d8	10.716

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	80
2			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	80
3			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
4			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	72
5			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
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Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

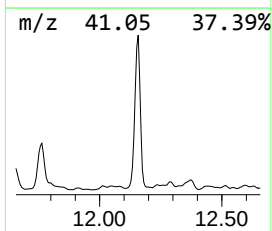
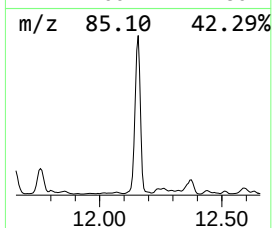
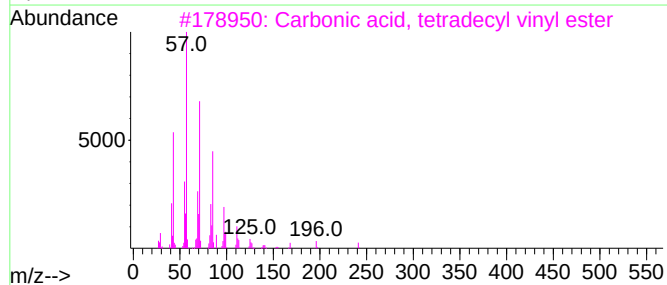
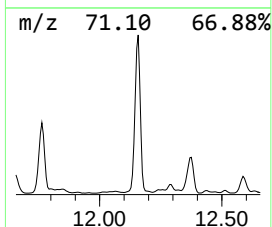
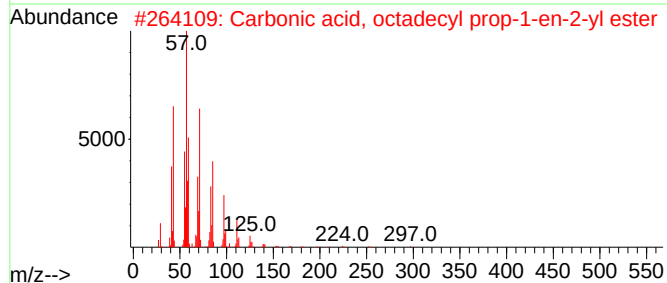
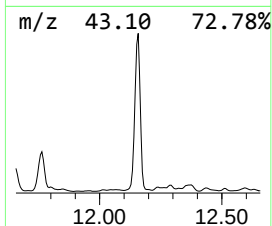
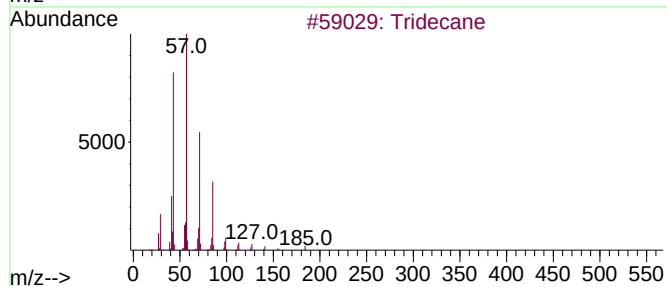
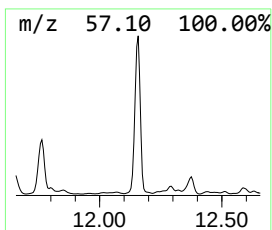
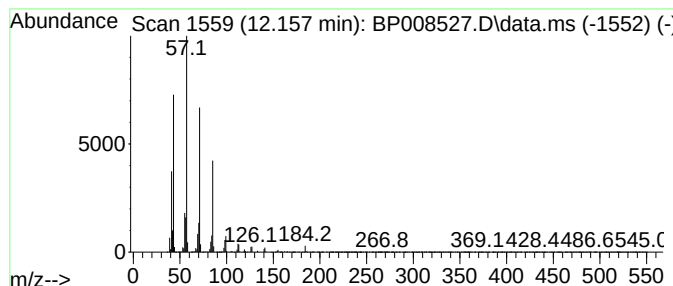
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.157	16.30 ng/ul	18859400	Naphthalene-d8	10.716	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	96
2	Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	90
3	Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	90
4	Tetradecane	198	C14H30	000629-59-4	90
5	Tridecane	184	C13H28	000629-50-5	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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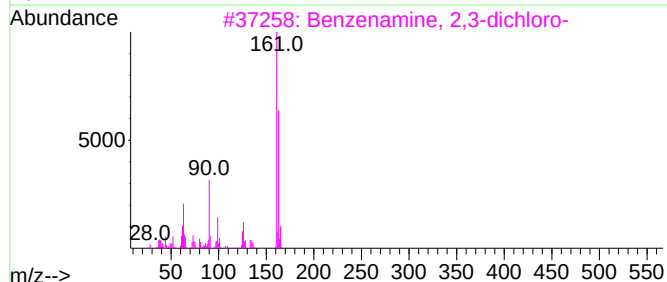
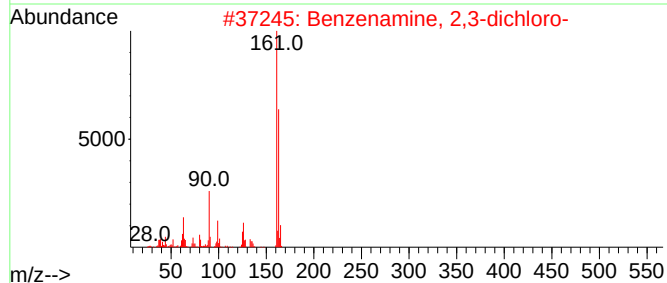
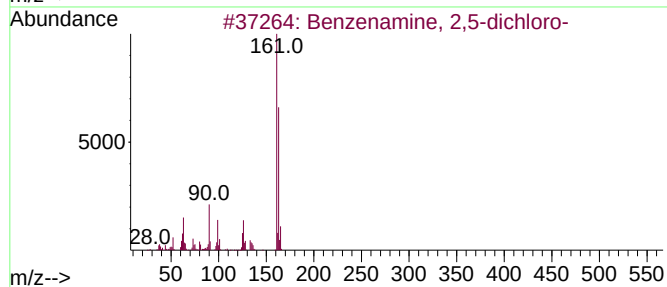
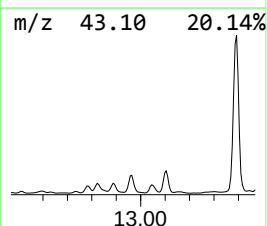
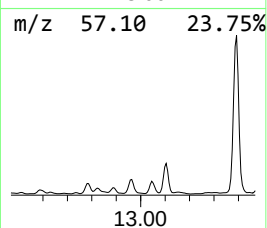
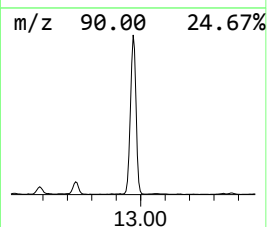
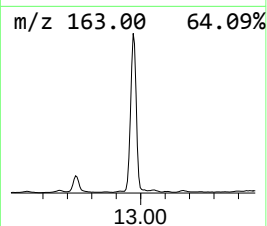
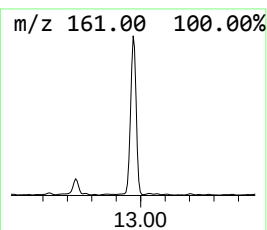
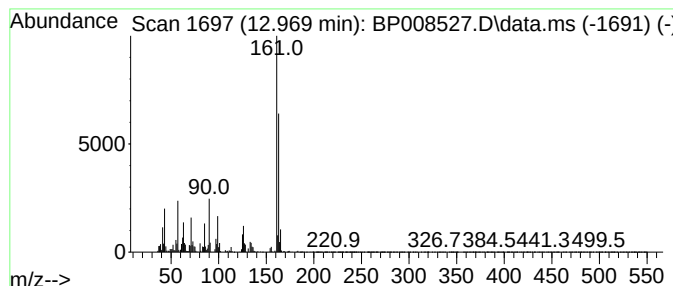
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Benzenamine, 2,5-dichloro- Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.969	17.86 ng/ul	8641670	Acenaphthene-d10	14.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	97
2			Benzenamine, 2,3-dichloro-	161	C6H5Cl2N	000608-27-5	96
3			Benzenamine, 2,3-dichloro-	161	C6H5Cl2N	000608-27-5	96
4			Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	96
5			Benzenamine, 3,4-dichloro-	161	C6H5Cl2N	000095-76-1	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
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Misc :
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Instrument :
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ClientSampleId :
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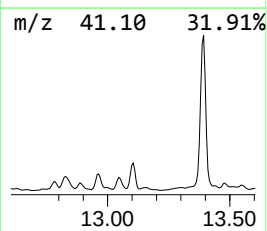
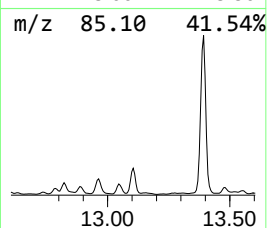
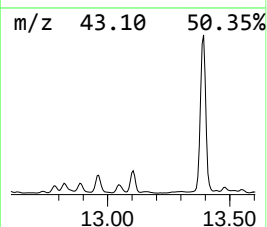
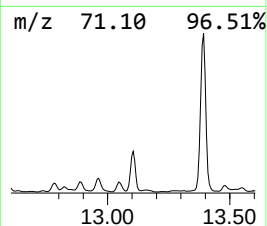
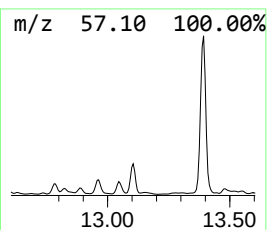
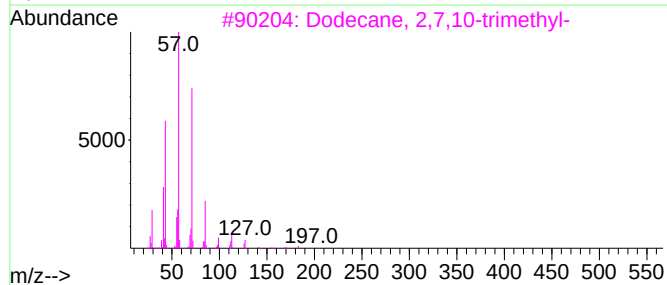
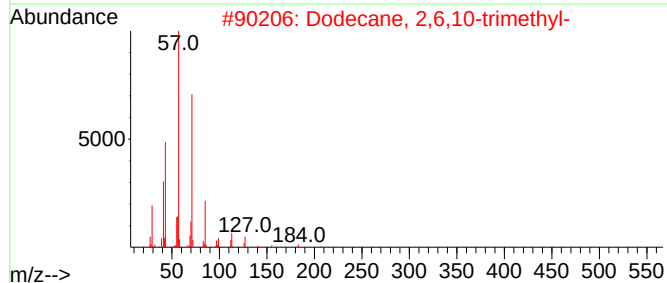
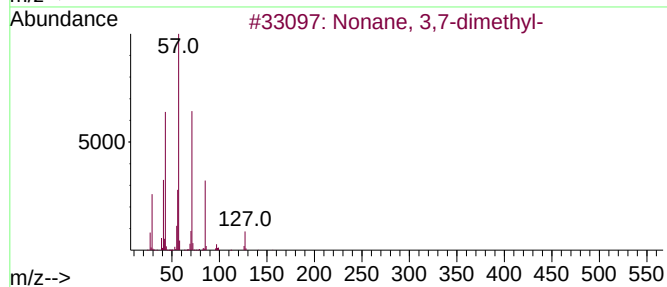
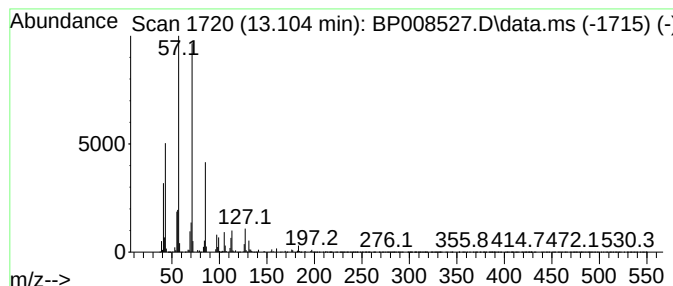
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.104	10.14 ng/ul	4905870	Acenaphthene-d10	14.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	76
2			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	72
3			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72
4			Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	72
5			Decane, 5-propyl-	184	C13H28	017312-62-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
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Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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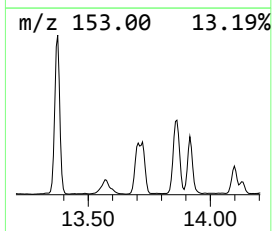
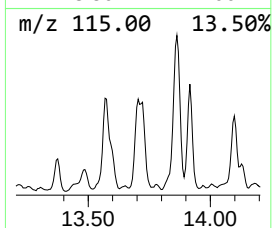
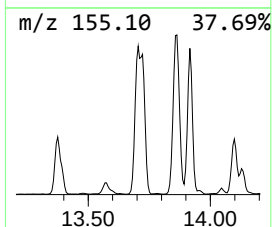
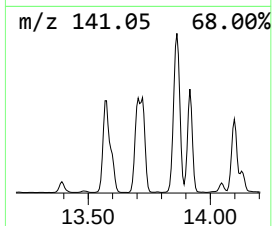
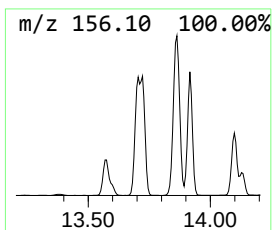
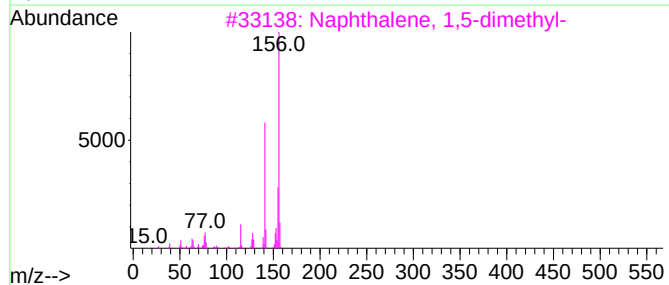
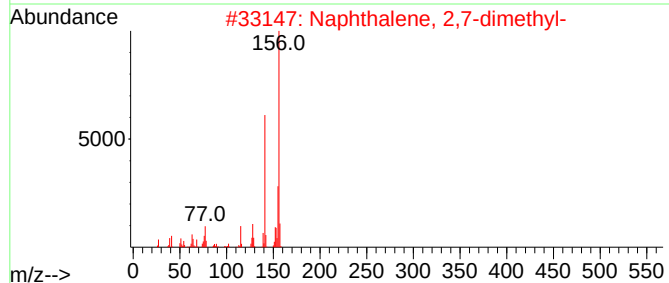
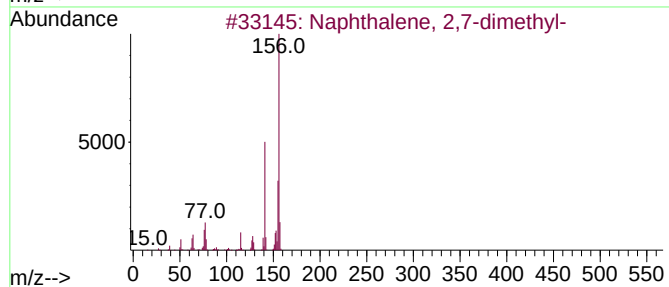
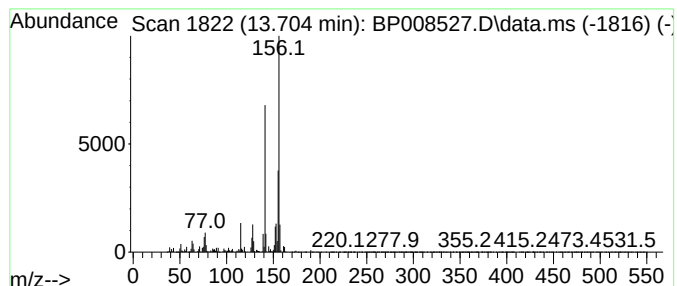
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 Naphthalene, 2,7-dimethyl- Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.704	13.47 ng/ul	6518170	Acenaphthene-d10	14.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
4			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

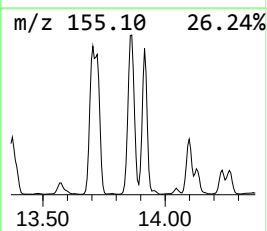
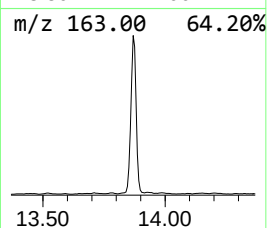
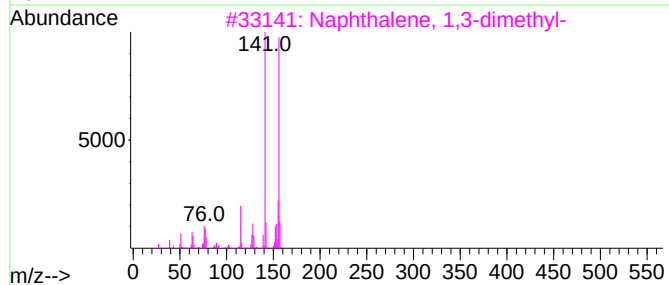
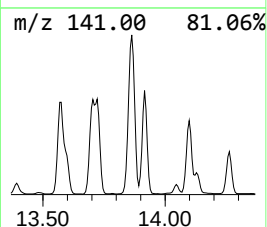
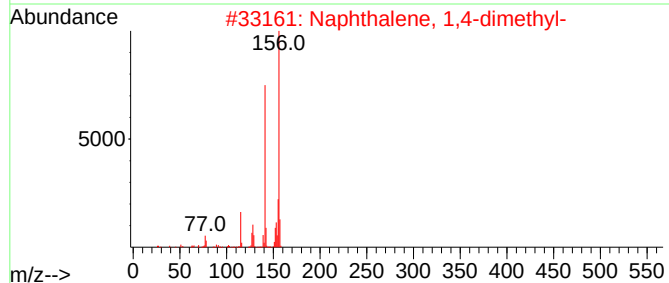
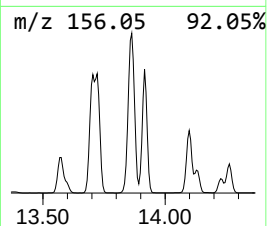
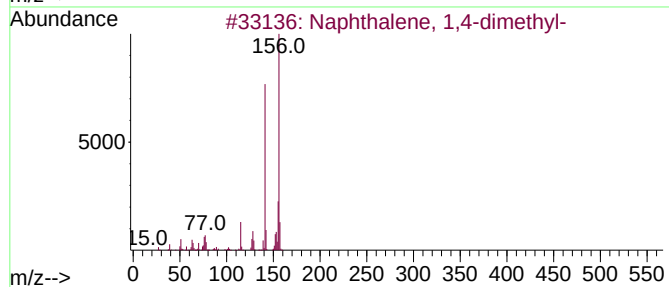
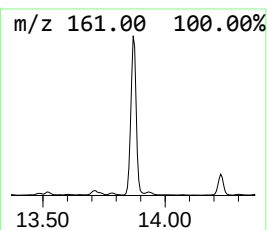
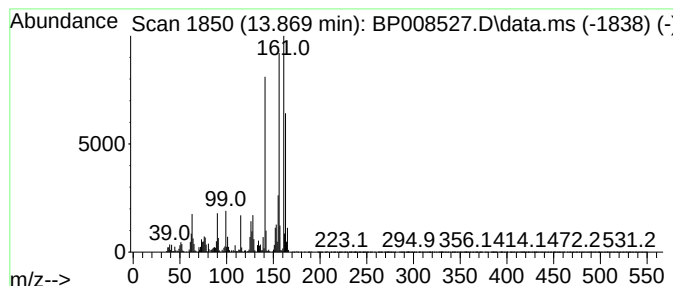
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 26 Naphthalene, 1,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.869	42.82 ng/ul	20722600	Acenaphthene-d10	14.534	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
2	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
3	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

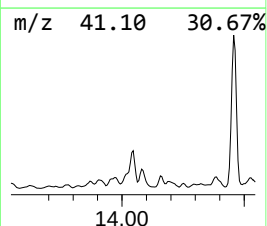
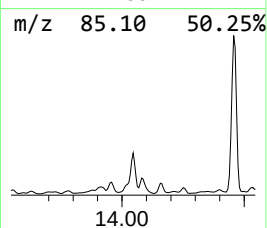
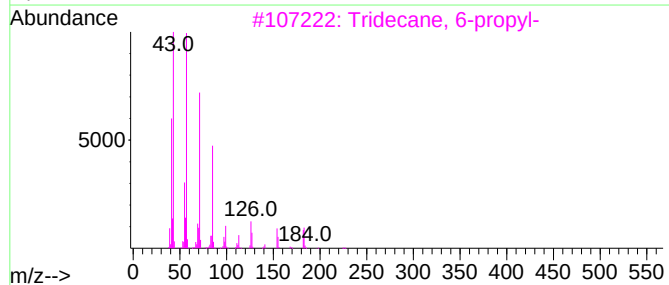
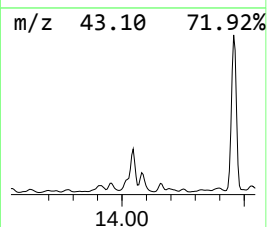
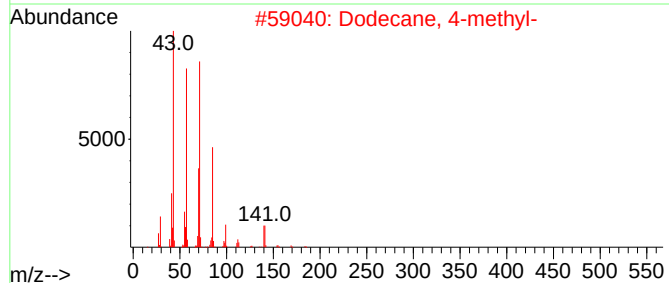
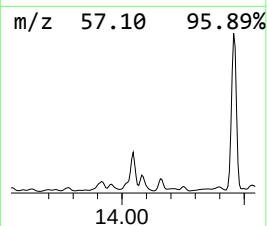
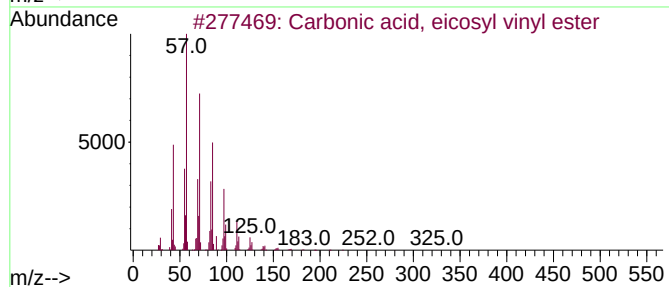
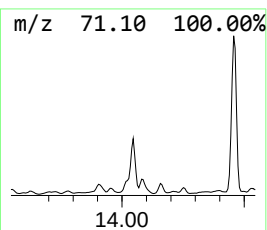
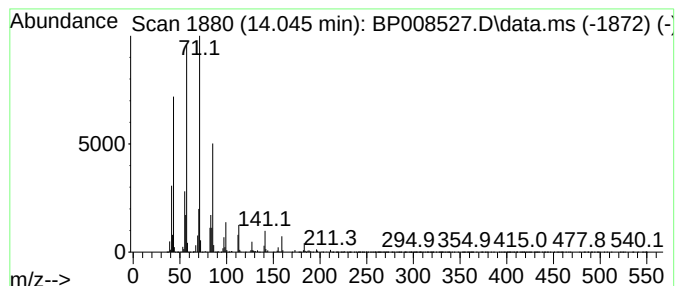
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 27 Carbonic acid, eicosyl vinyl... Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.045	14.13 ng/ul	6837260	Acenaphthene-d10	14.534	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	86
2	Dodecane, 4-methyl-	184	C13H28	006117-97-1	81
3	Tridecane, 6-propyl-	226	C16H34	055045-10-8	80
4	Dodecane, 4-methyl-	184	C13H28	006117-97-1	72
5	Hexadecane	226	C16H34	000544-76-3	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGLA5ME

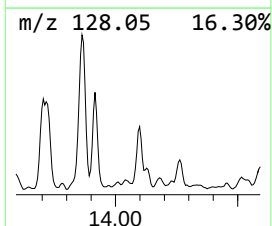
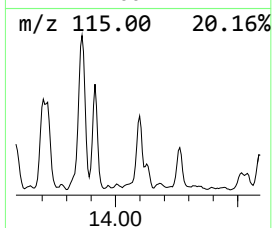
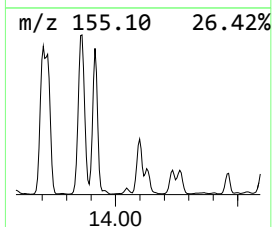
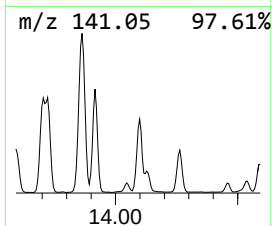
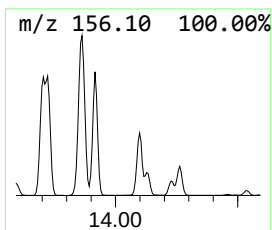
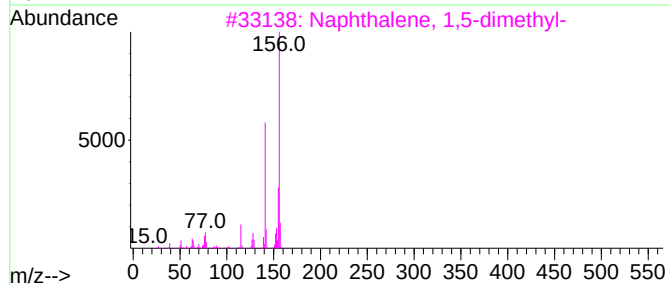
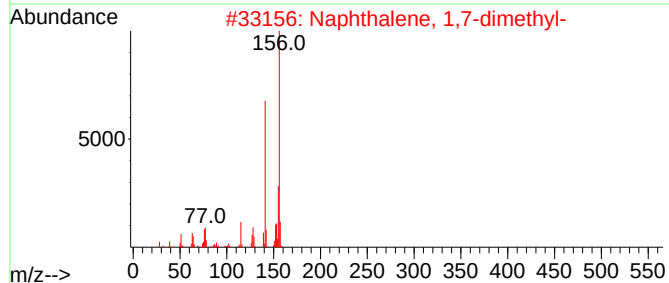
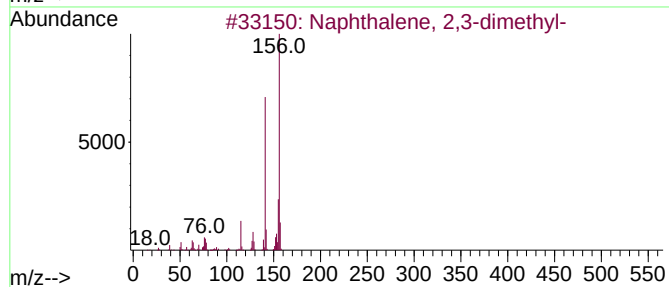
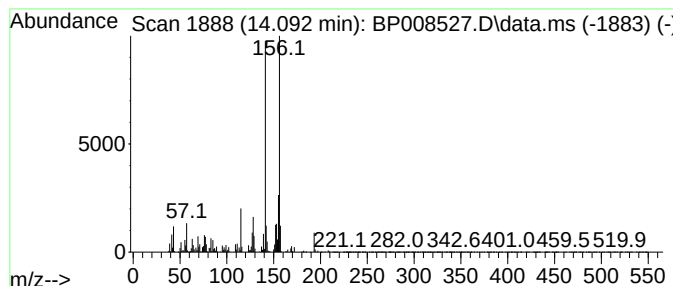
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 Naphthalene, 2,3-dimethyl- Concentration Rank 61

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.092	9.69 ng/ul	4687950	Acenaphthene-d10	14.534

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
4		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

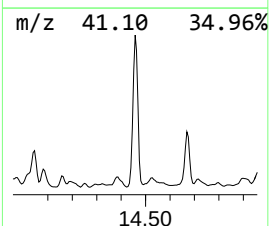
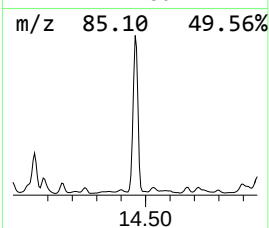
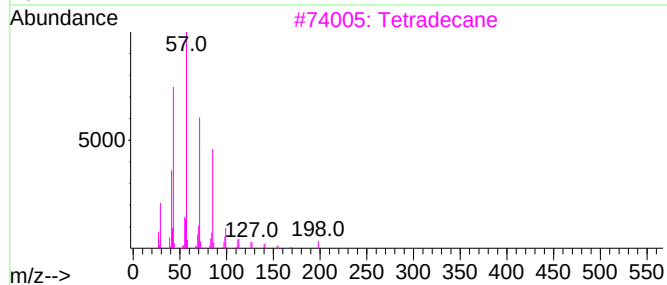
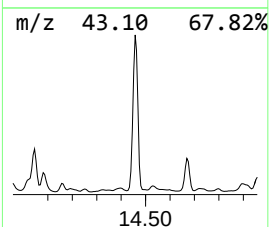
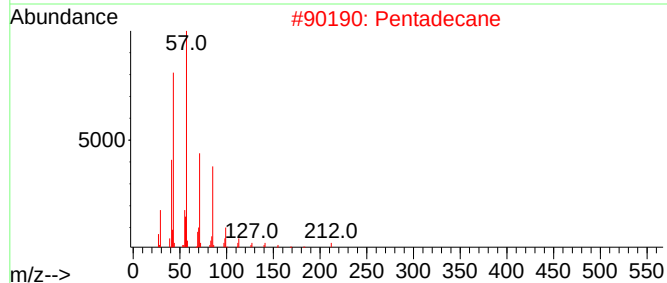
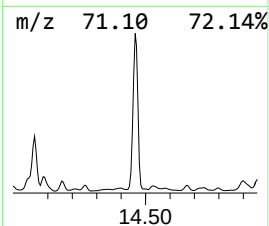
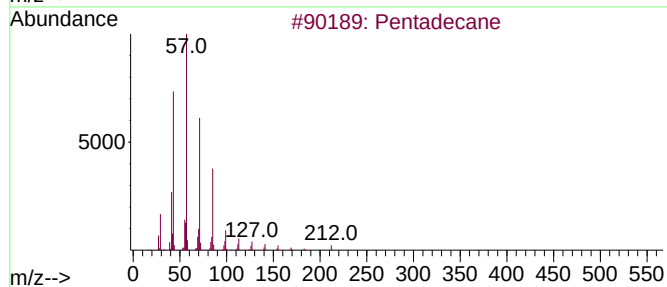
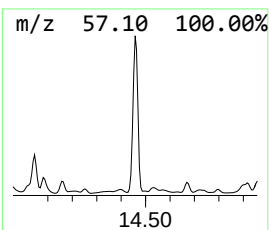
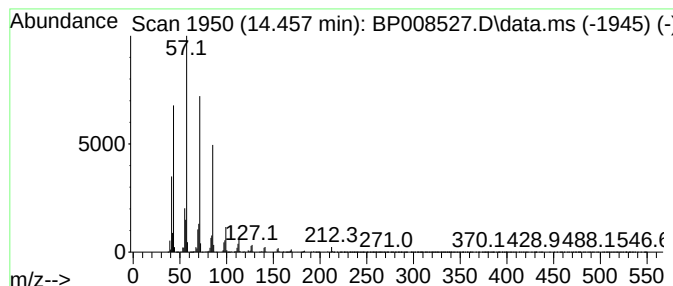
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.457	44.28 ng/ul	21431600	Acenaphthene-d10		14.534	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecane		212	C15H32	000629-62-9	93
2	Pentadecane		212	C15H32	000629-62-9	92
3	Tetradecane		198	C14H30	000629-59-4	91
4	Hexadecane		226	C16H34	000544-76-3	91
5	Carbonic acid, eicosyl vinyl ester		368	C23H44O3	1000382-54-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

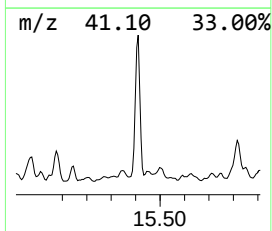
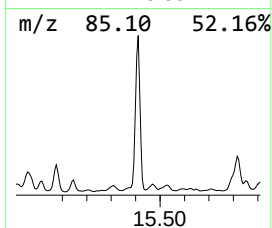
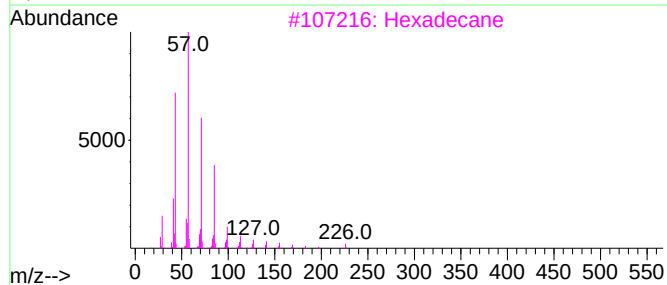
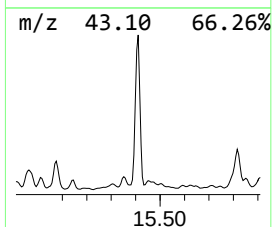
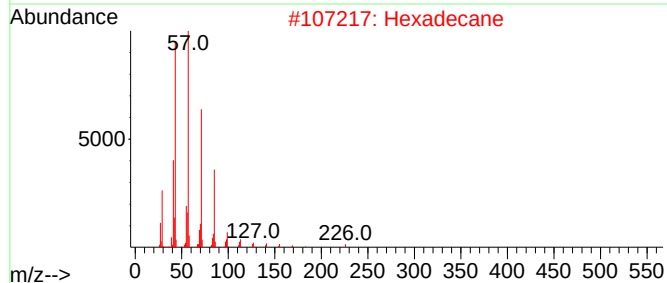
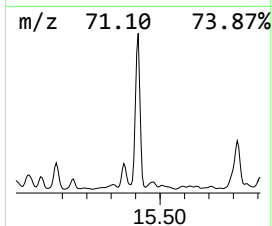
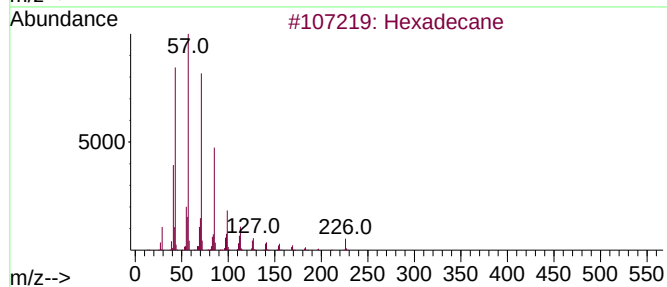
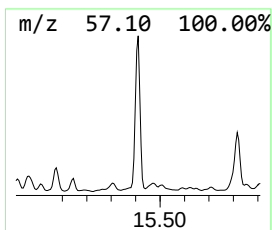
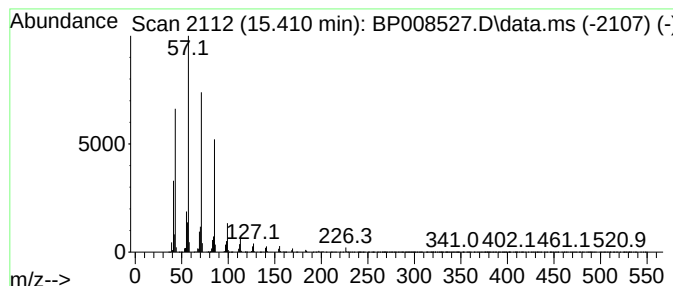
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.410	29.29 ng/ul	14172800	Acenaphthene-d10		14.534	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	96
2	Hexadecane		226	C16H34	000544-76-3	95
3	Hexadecane		226	C16H34	000544-76-3	92
4	Tetradecane		198	C14H30	000629-59-4	91
5	Carbonic acid, eicosyl vinyl ester		368	C23H44O3	1000382-54-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

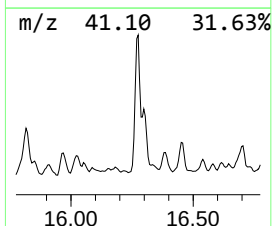
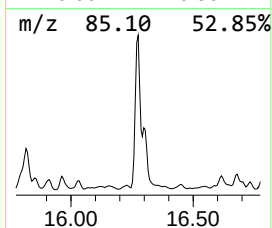
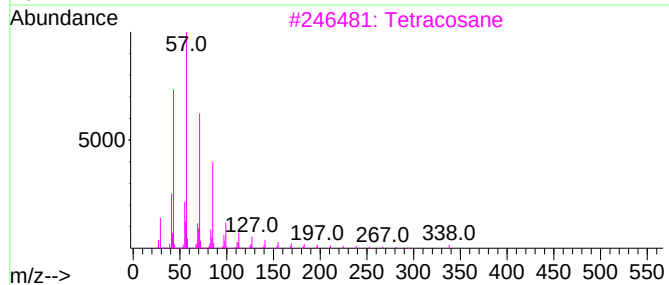
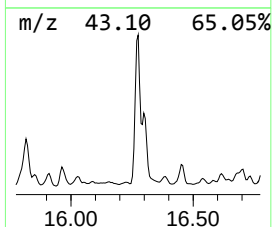
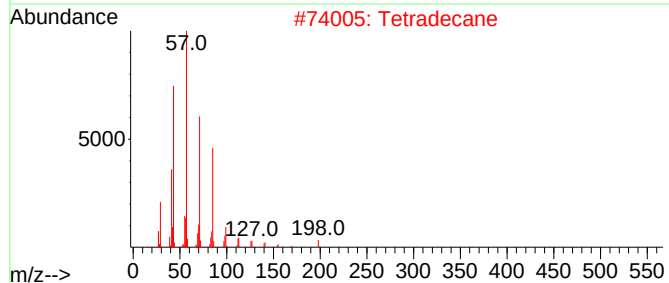
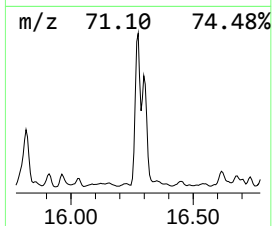
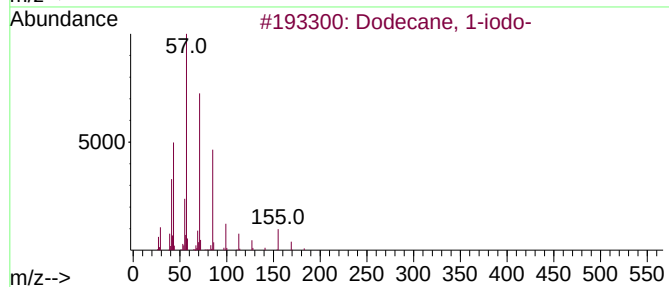
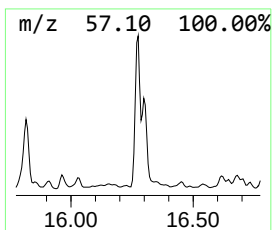
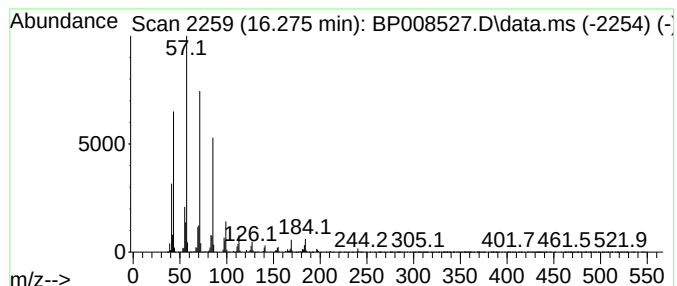
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 34 Dodecane, 1-iodo- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.275	53.25 ng/ul	14271100	Phenanthrene-d10		17.286	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	91
2	Tetradecane		198	C14H30	000629-59-4	87
3	Tetracosane		338	C24H50	000646-31-1	87
4	Heptadecane		240	C17H36	000629-78-7	87
5	Heptadecane		240	C17H36	000629-78-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGLA5ME

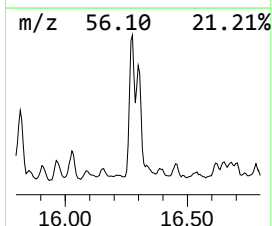
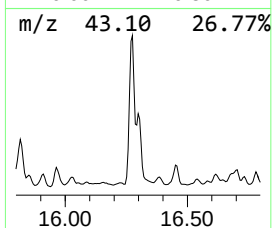
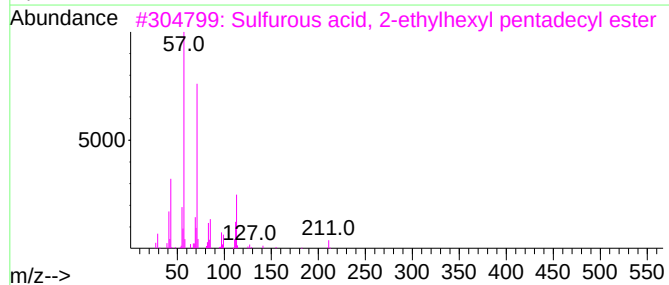
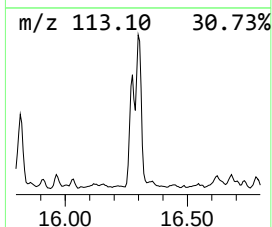
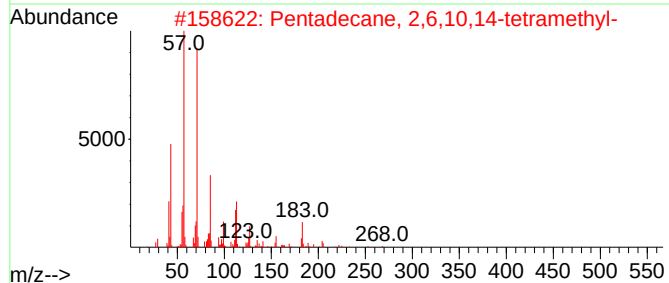
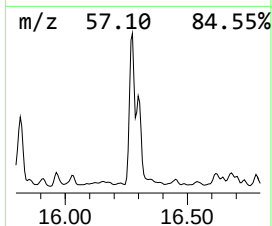
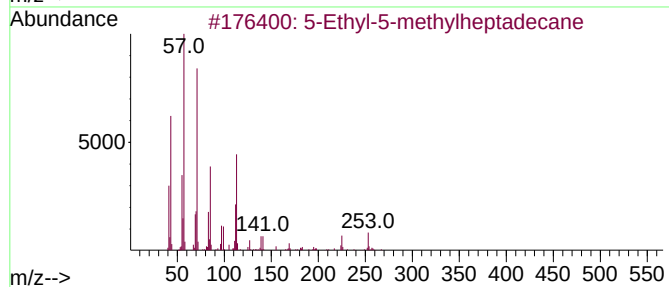
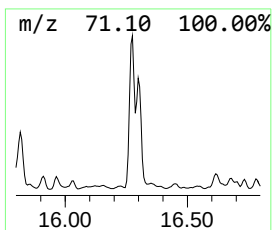
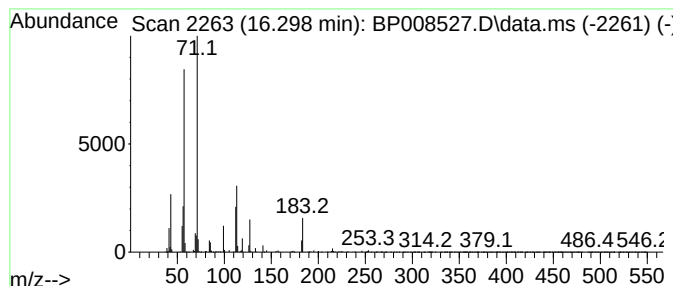
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Quant Title : SVOA CALIBRATION

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

Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.298	23.61 ng/ul	6325910	Phenanthrene-d10	17.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Ethyl-5-methylheptadecane	282	C20H42	1000360-42-5	59
2			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	56
3			Sulfurous acid, 2-ethylhexyl pen...	404	C23H48O3S	1000309-19-8	53
4			Sulfurous acid, 2-ethylhexyl tet...	390	C22H46O3S	1000309-19-7	53
5			Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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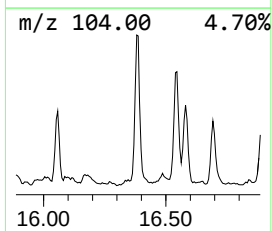
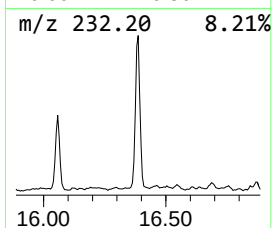
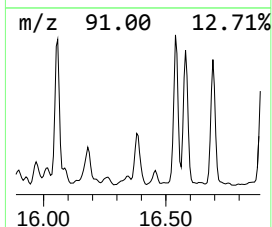
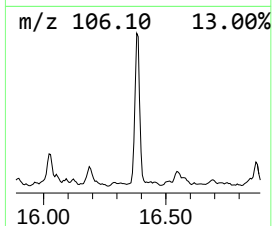
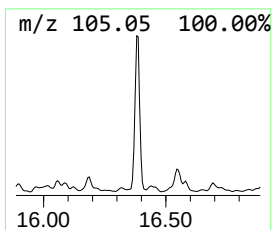
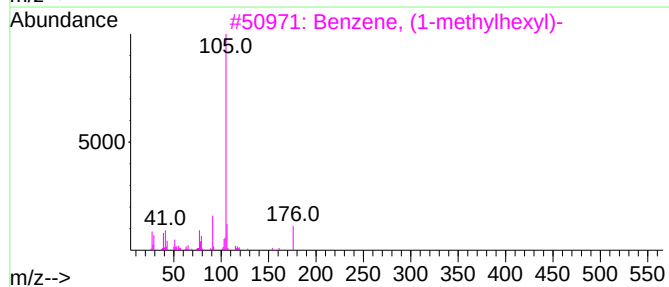
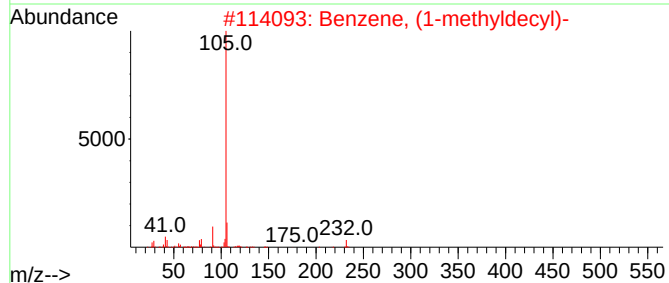
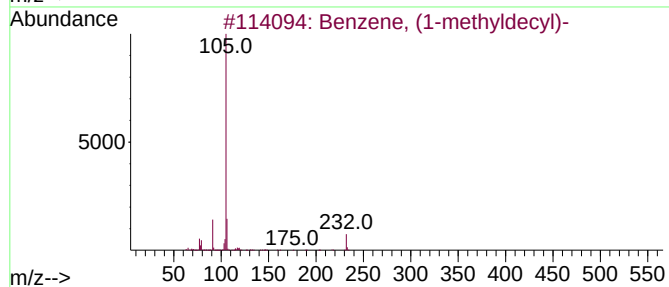
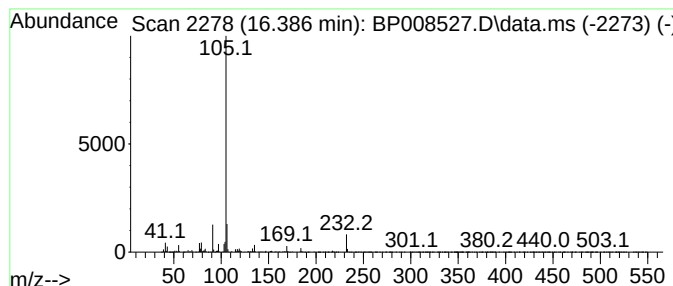
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 Benzene, (1-methyldecyl)- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.387	21.22 ng/ul	5685880	Phenanthrene-d10	17.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyldecyl)-	232	C17H28	004536-88-3	94
2			Benzene, (1-methyldecyl)-	232	C17H28	004536-88-3	83
3			Benzene, (1-methylhexyl)-	176	C13H20	002132-84-5	68
4			Benzene, (1-methylnonyl)-	218	C16H26	004537-13-7	59
5			Benzene, (1-methylnonyl)-	218	C16H26	004537-13-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

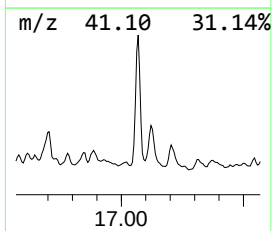
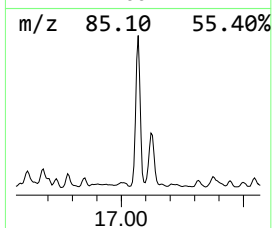
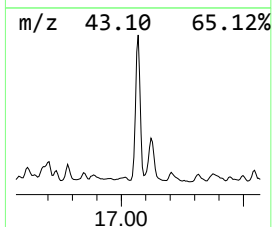
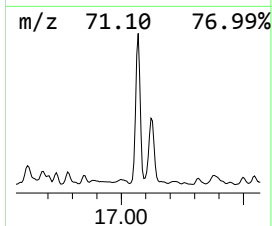
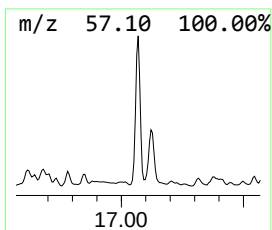
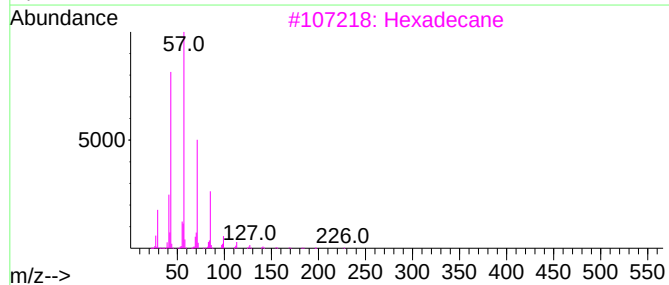
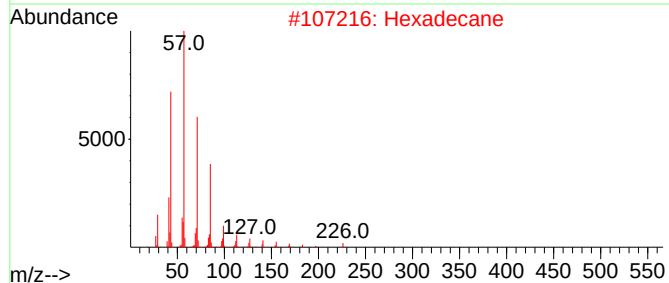
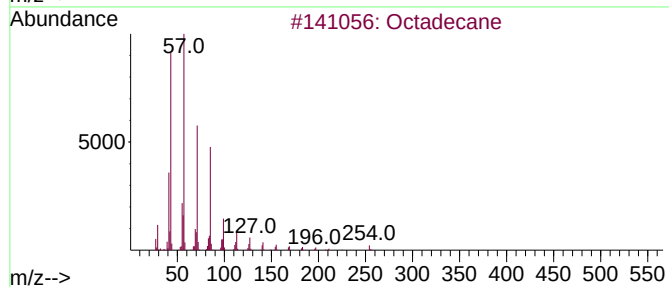
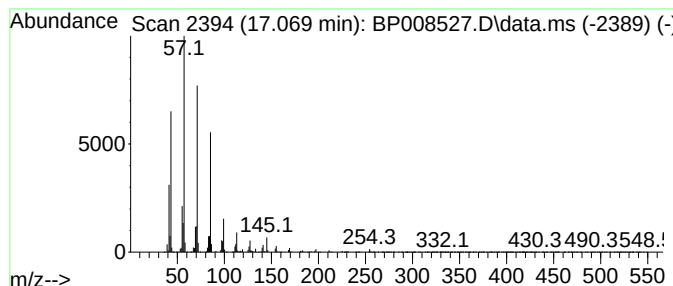
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.069	35.58 ng/ul	9535770	Phenanthrene-d10	17.286	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	94
2	Hexadecane	226	C16H34	000544-76-3	90
3	Hexadecane	226	C16H34	000544-76-3	90
4	Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	89
5	Tetradecane	198	C14H30	000629-59-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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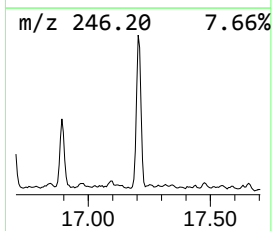
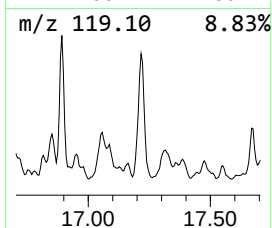
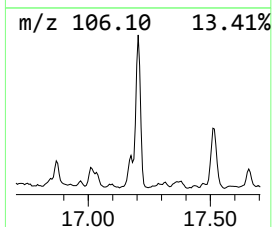
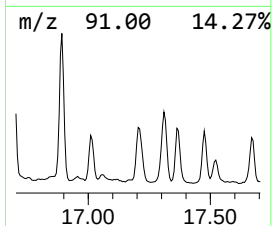
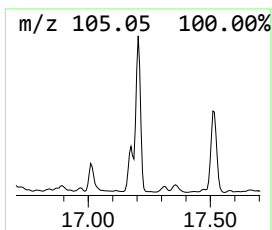
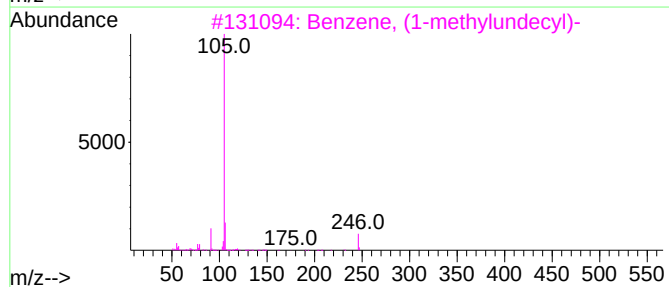
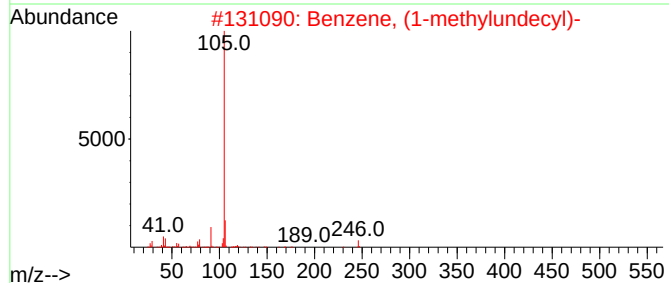
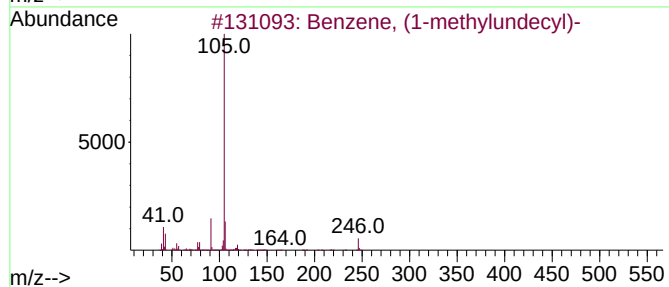
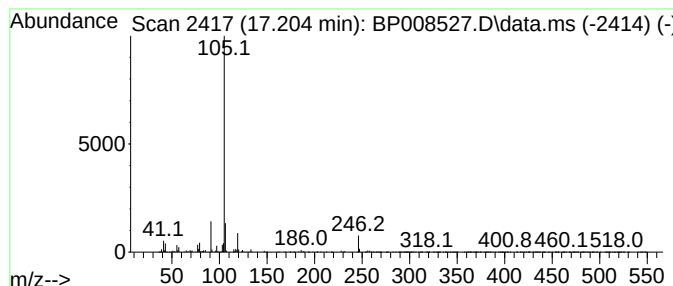
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 Benzene, (1-methylundecyl)- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.204	18.67 ng/ul	5004220	Phenanthrene-d10	17.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylundecyl)-	246	C18H30	002719-61-1	91
2			Benzene, (1-methylundecyl)-	246	C18H30	002719-61-1	87
3			Benzene, (1-methylundecyl)-	246	C18H30	002719-61-1	81
4			3-Phenyl-1-butanol	150	C10H14O	002722-36-3	64
5			Benzene, (1-methylundecyl)-	246	C18H30	002719-61-1	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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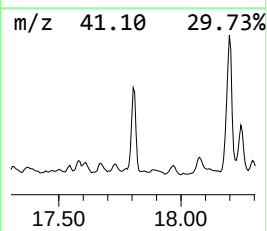
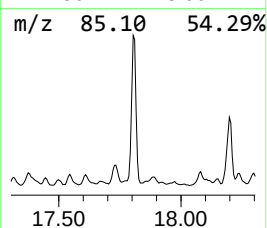
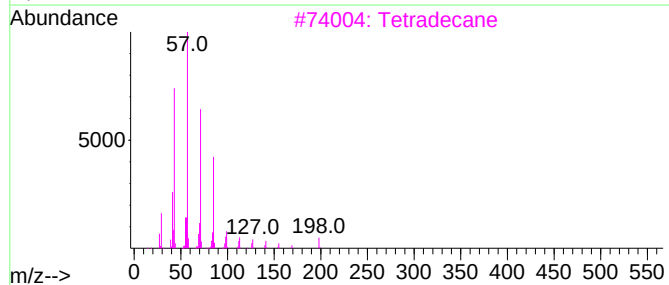
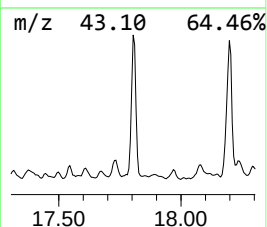
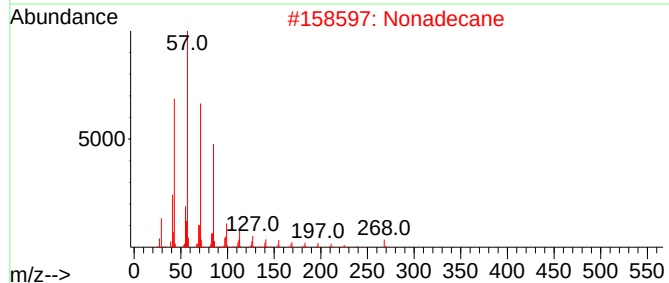
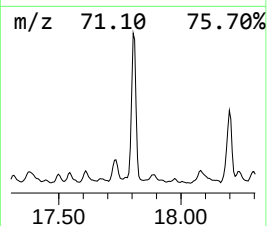
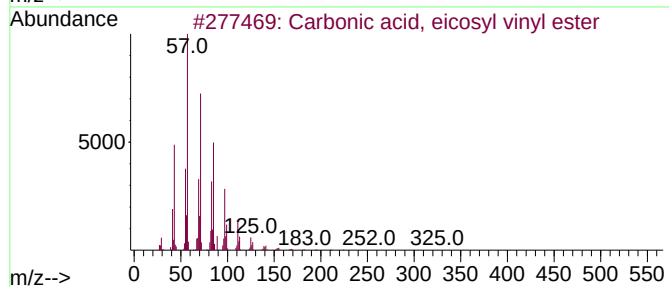
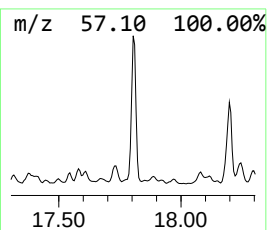
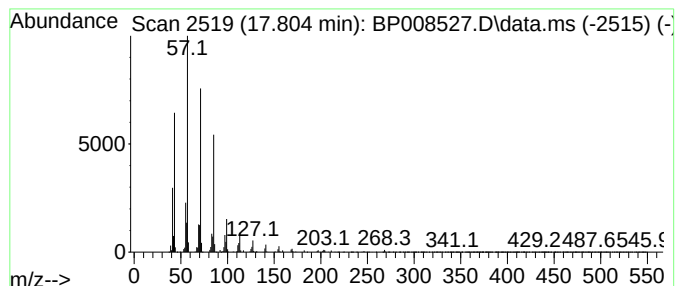
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 Nonadecane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.804	27.88 ng/ul	7470920	Phenanthrene-d10	17.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	91
2			Nonadecane	268	C19H40	000629-92-5	91
3			Tetradecane	198	C14H30	000629-59-4	91
4			Nonadecane	268	C19H40	000629-92-5	90
5			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
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Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

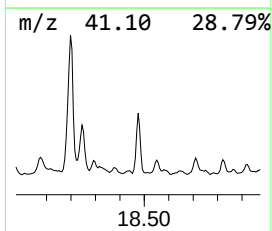
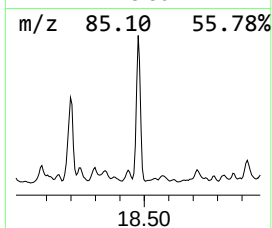
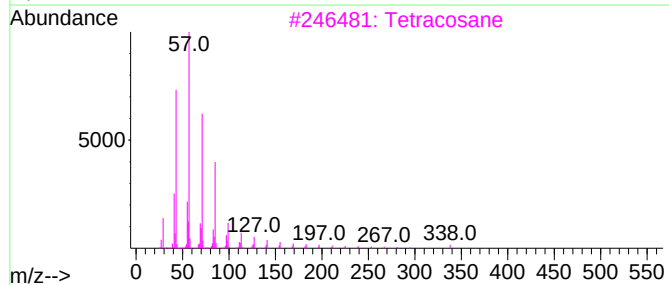
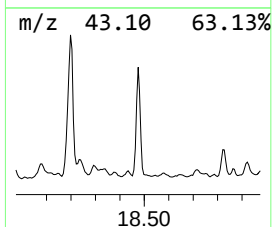
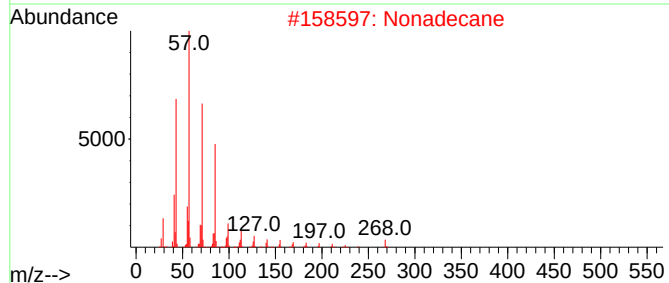
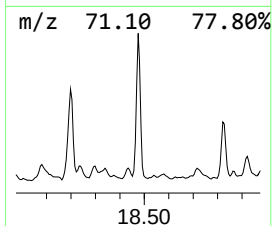
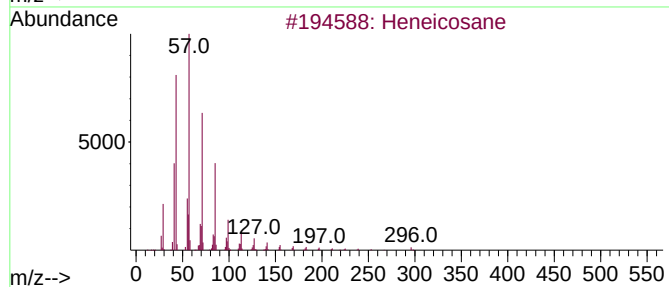
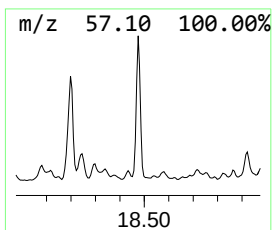
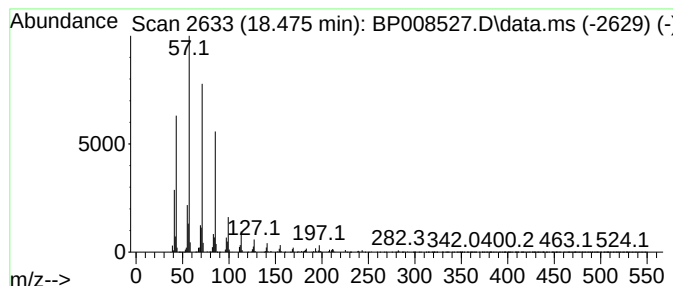
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 (DEL) Alkane: Straight-Chai... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.475	19.03 ng/ul	5101160	Phenanthrene-d10		17.286	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Nonadecane		268	C19H40	000629-92-5	91
3	Tetracosane		338	C24H50	000646-31-1	91
4	Octadecane		254	C18H38	000593-45-3	91
5	Nonadecane		268	C19H40	000629-92-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
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Operator : CG/JU
Sample : M5171-09ME
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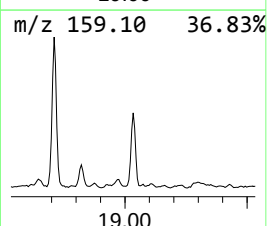
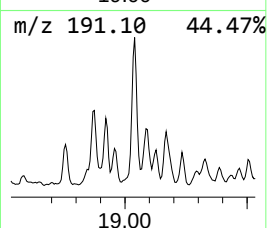
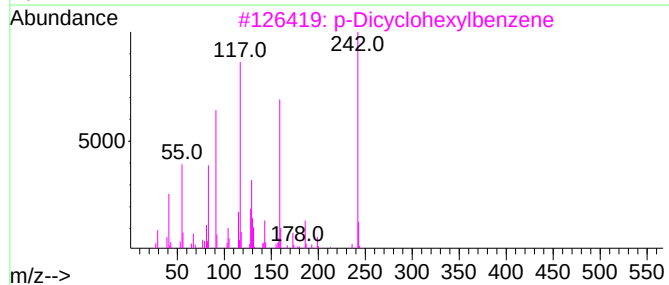
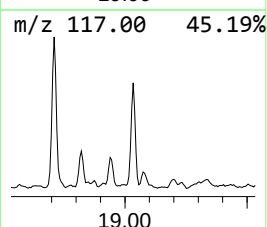
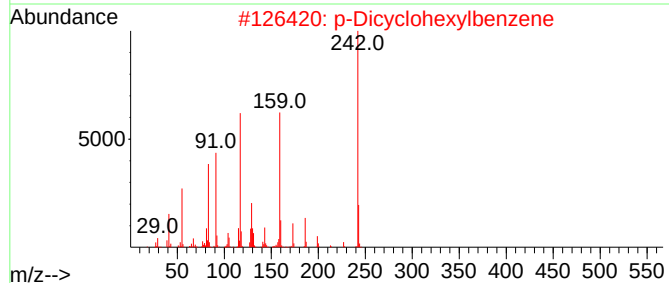
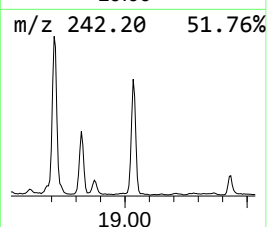
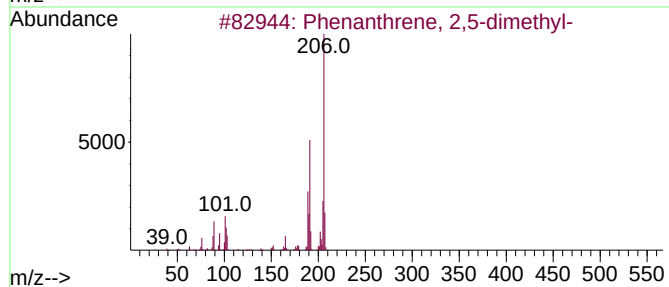
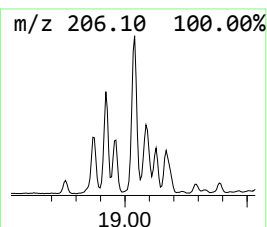
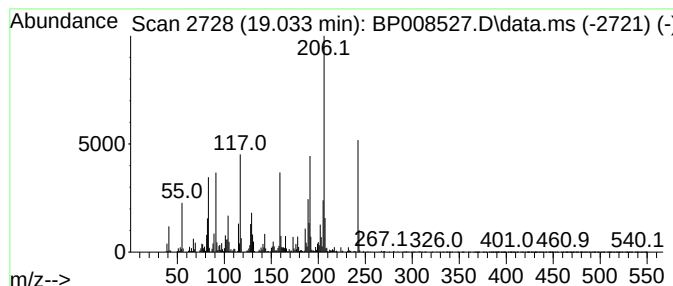
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Quant Title : SVOA CALIBRATION

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

Peak Number 42 Phenanthrene, 2,5-dimethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD			R.T.
19.033	17.84 ng/ul	4781460	Phenanthrene-d10			17.286
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-		206	C16H14	003674-66-6	95
2	p-Dicyclohexylbenzene		242	C18H26	001087-02-1	94
3	p-Dicyclohexylbenzene		242	C18H26	001087-02-1	93
4	p-Dicyclohexylbenzene		242	C18H26	001087-02-1	93
5	Phenanthrene, 1,7-dimethyl-		206	C16H14	000483-87-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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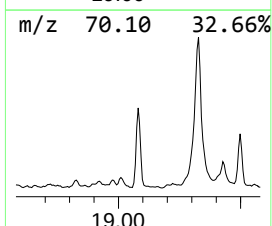
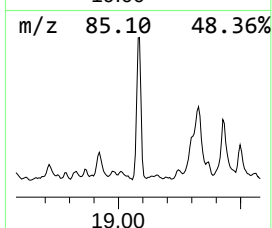
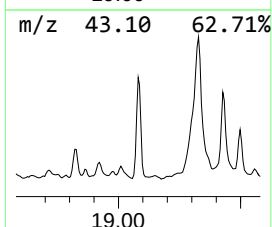
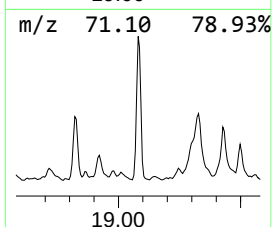
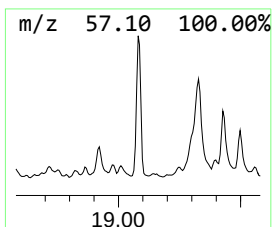
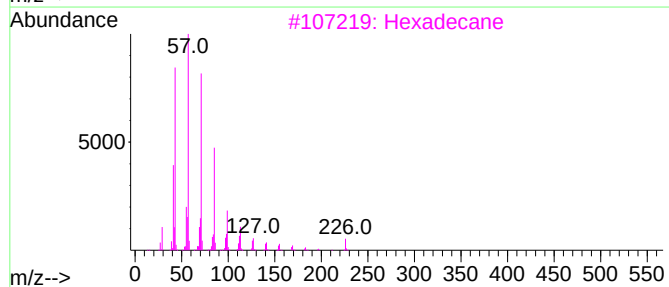
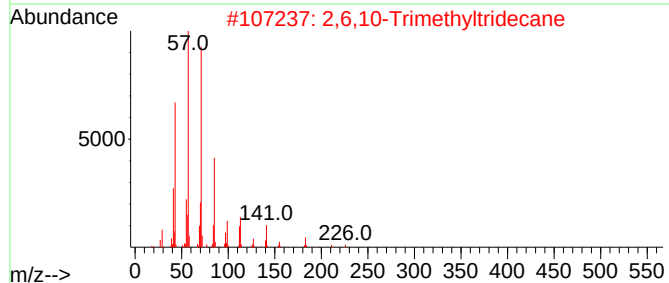
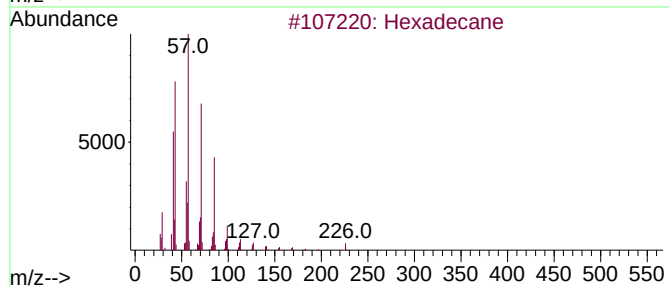
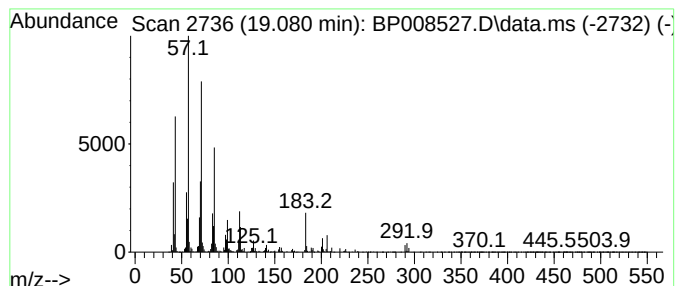
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.080	26.91 ng/ul	7212290	Phenanthrene-d10	17.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	94
2			2,6,10-Trimethyltridecane	226	C16H34	003891-99-4	91
3			Hexadecane	226	C16H34	000544-76-3	90
4			3-Ethyl-2,6,10-trimethylundecane	226	C16H34	1000432-25-9	90
5			Pentadecane, 7-methyl-	226	C16H34	006165-40-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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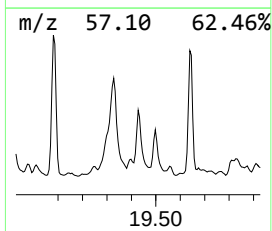
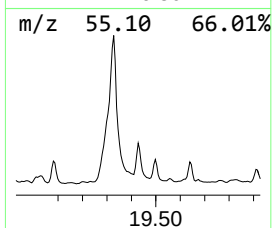
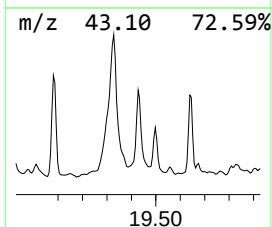
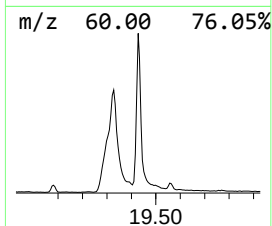
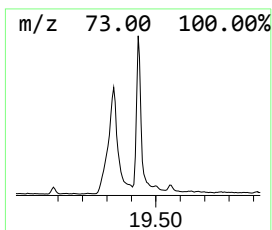
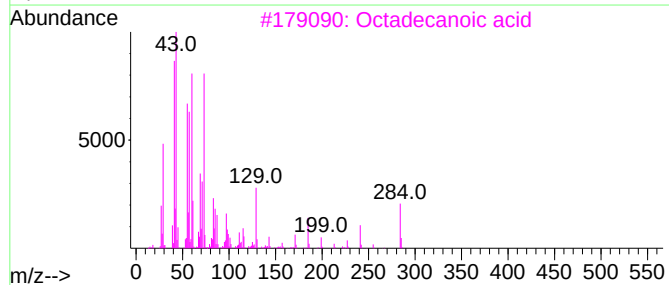
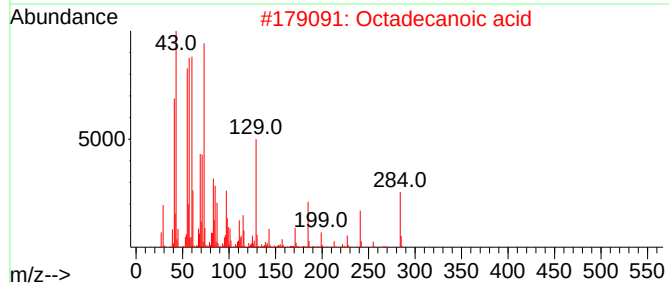
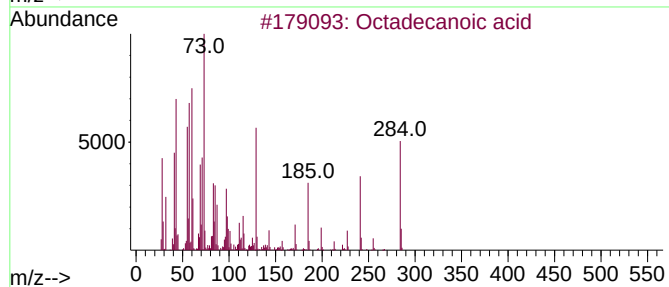
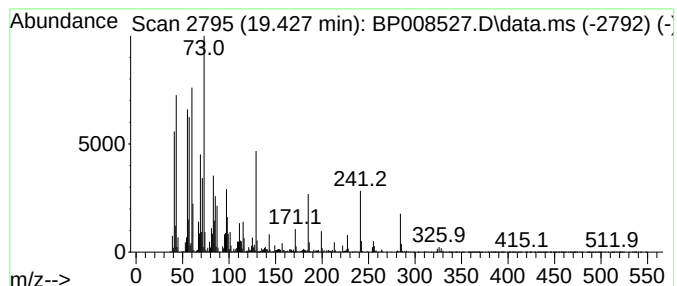
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Quant Title : SVOA CALIBRATION

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

Peak Number 44 Octadecanoic acid Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.427	26.63 ng/ul	9917370	Chrysene-d12	21.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	97
5			Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGLA5ME

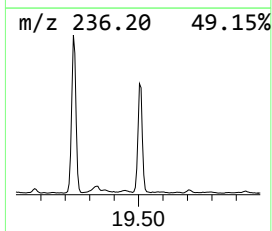
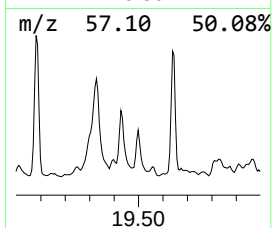
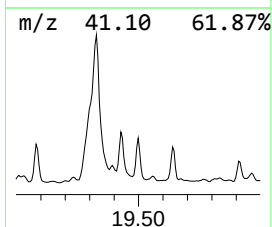
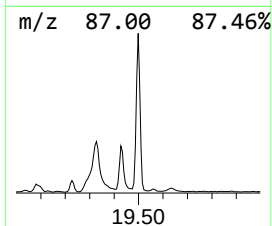
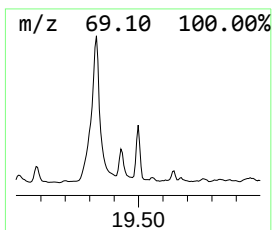
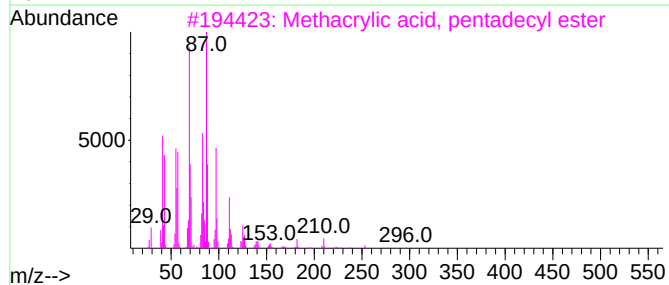
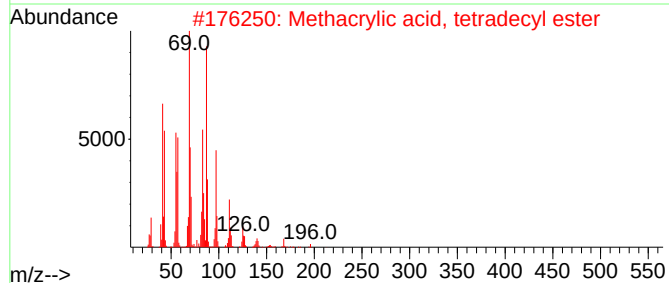
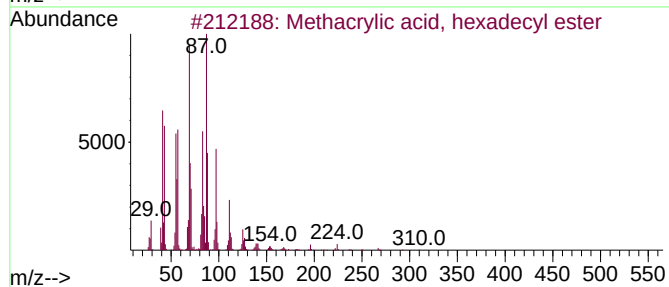
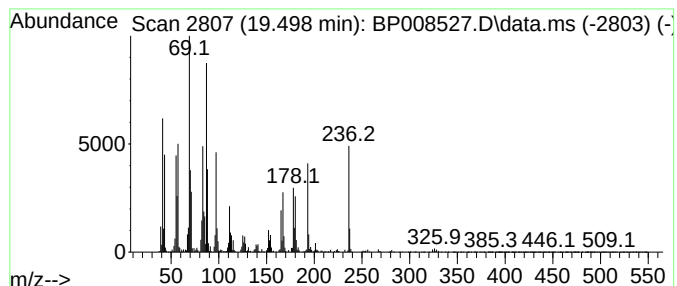
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 Methacrylic acid, hexadecyl... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.498	19.72 ng/ul	7346160	Chrysene-d12	21.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylic acid, hexadecyl ester	310	C20H38O2	1000340-29-2	78
2			Methacrylic acid, tetradecyl ester	282	C18H34O2	1000340-29-0	55
3			Methacrylic acid, pentadecyl ester	296	C19H36O2	1000340-29-1	53
4			Octadecyl methacrylate	338	C22H42O2	000112-08-3	49
5			n-Dodecyl methacrylate	254	C16H30O2	000142-90-5	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

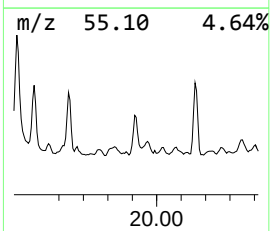
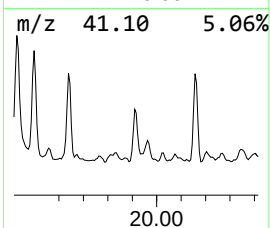
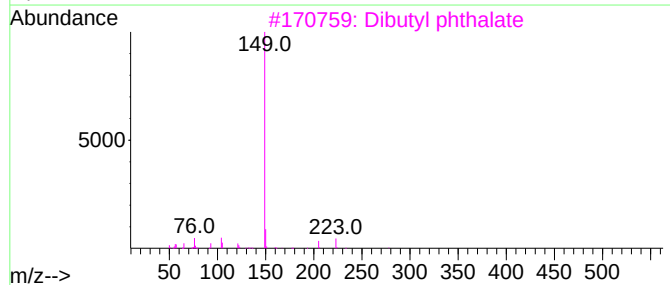
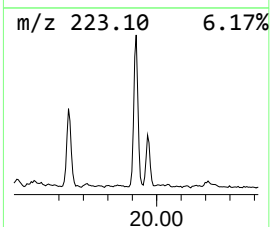
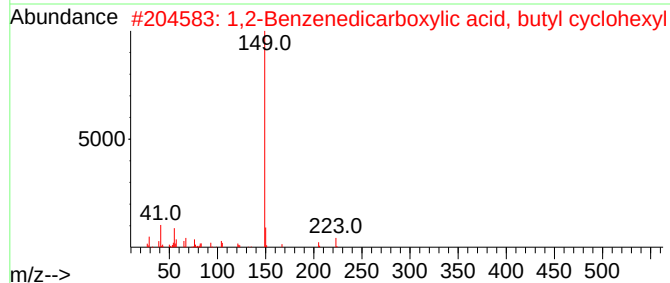
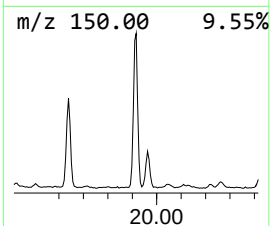
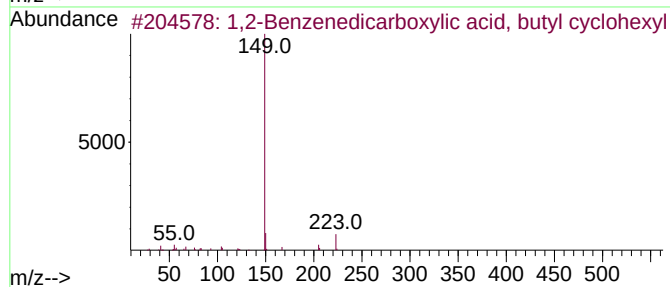
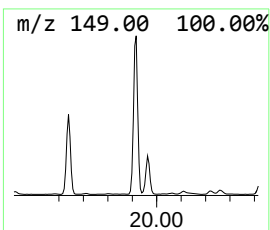
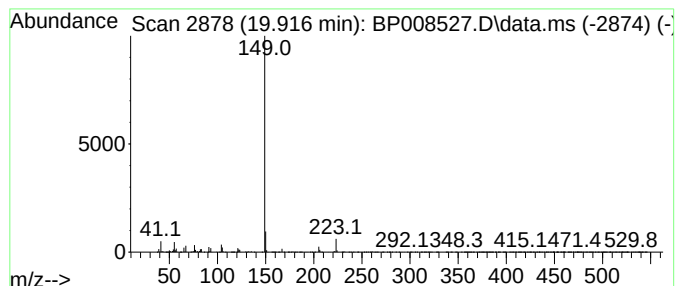
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 46 1,2-Benzenedicarboxylic aci... Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.916	15.18 ng/ul	5652390	Chrysene-d12		21.369	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1,2-Benzenedicarboxylic acid, bu...		304	C18H24O4	000084-64-0	90
2	1,2-Benzenedicarboxylic acid, bu...		304	C18H24O4	000084-64-0	90
3	Dibutyl phthalate		278	C16H22O4	000084-74-2	86
4	1,2-Benzenedicarboxylic acid, bu...		334	C20H30O4	000084-78-6	78
5	Phthalic acid, 8-chlorooctyl iso...		368	C20H29ClO4	1000356-85-8	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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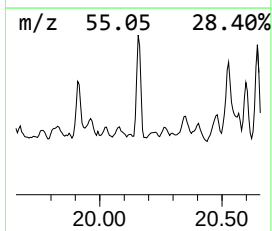
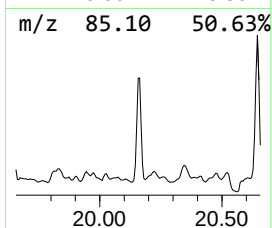
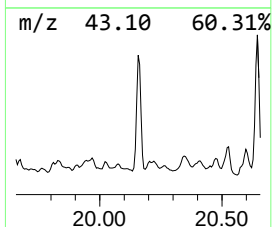
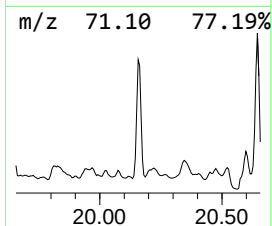
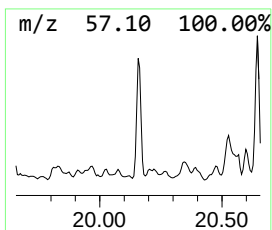
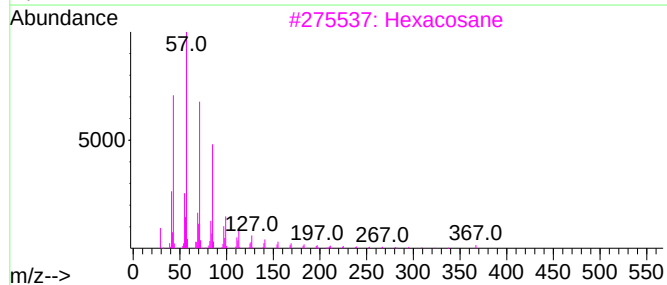
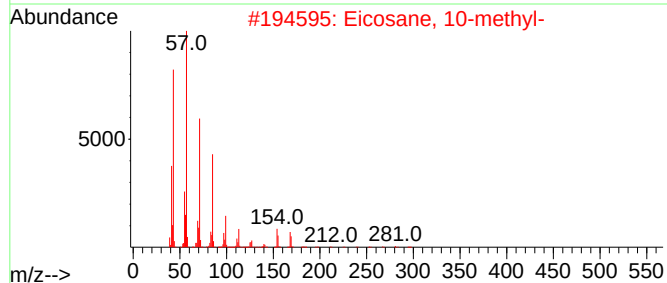
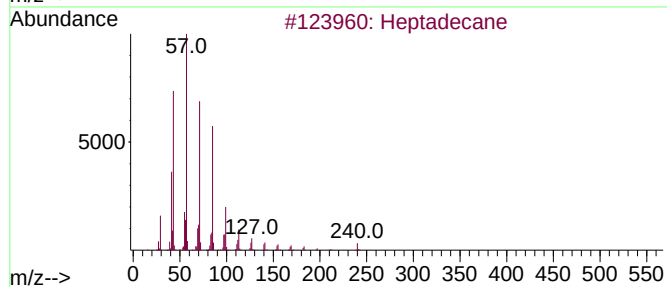
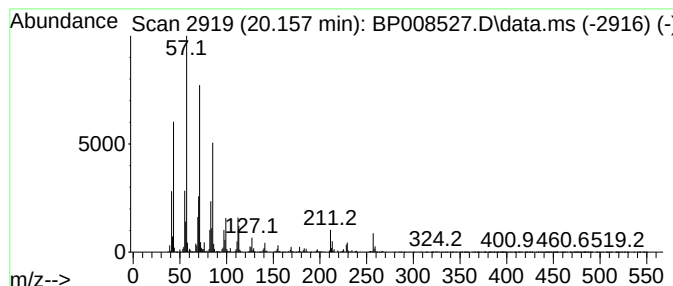
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Quant Title : SVOA CALIBRATION

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

Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.157	18.22 ng/ul	6787000	Chrysene-d12	21.369

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	94
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	89
3		Hexacosane	366	C26H54	000630-01-3	87
4		Heptadecane, 3-methyl-	254	C18H38	006418-44-6	81
5		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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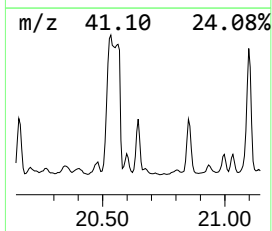
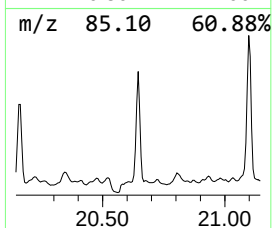
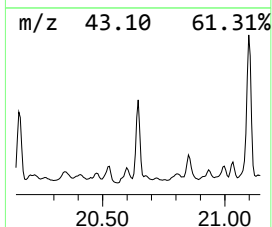
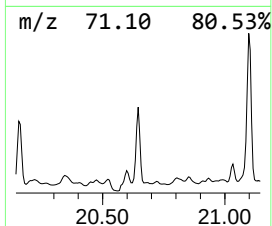
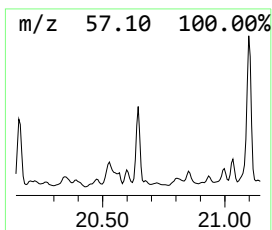
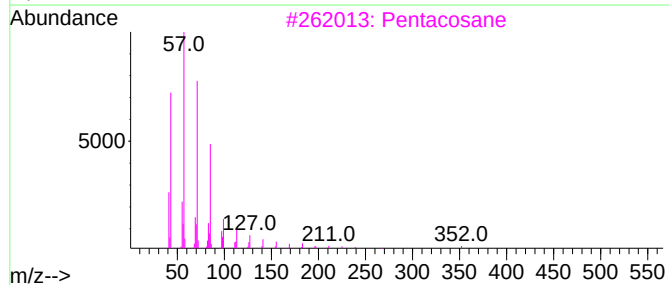
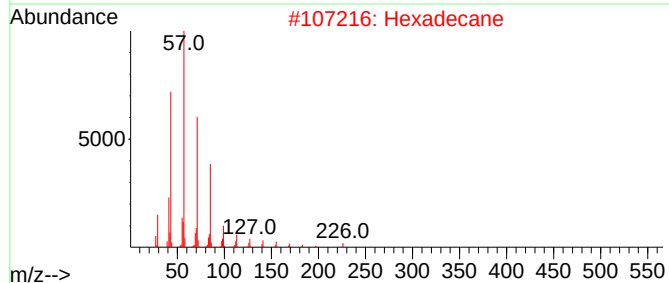
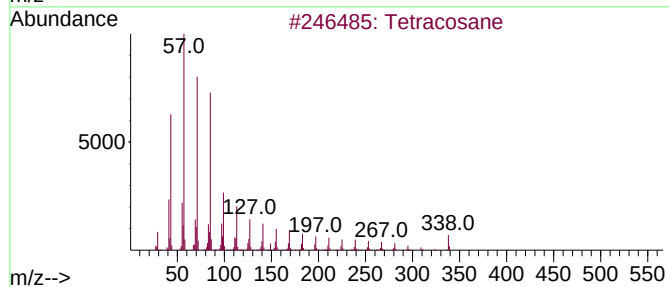
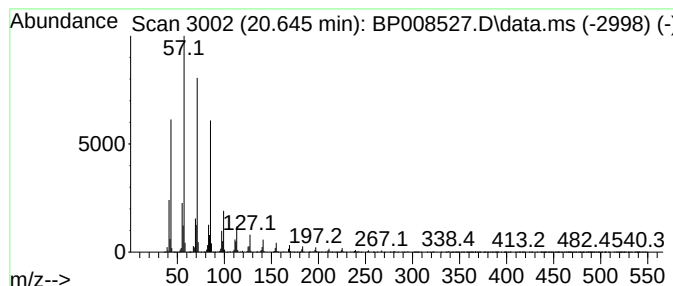
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Quant Title : SVOA CALIBRATION

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

Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.645	15.44 ng/ul	5748980	Chrysene-d12	21.369

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	97
2		Hexadecane	226	C16H34	000544-76-3	94
3		Pentacosane	352	C25H52	000629-99-2	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
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Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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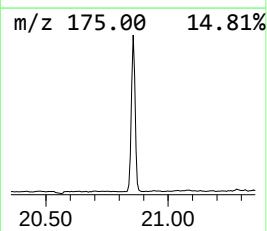
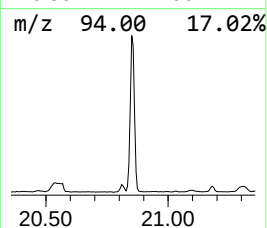
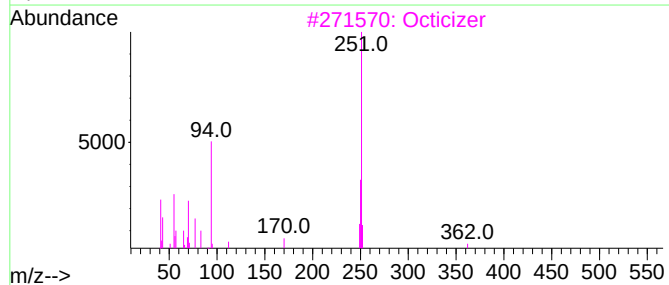
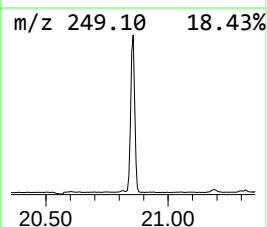
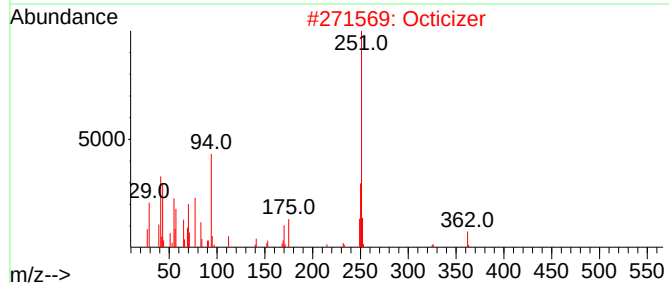
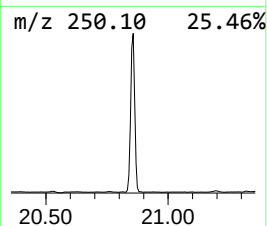
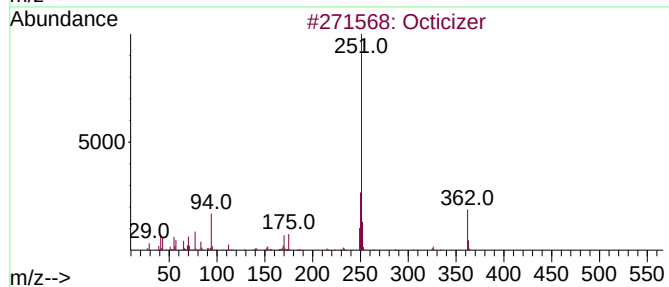
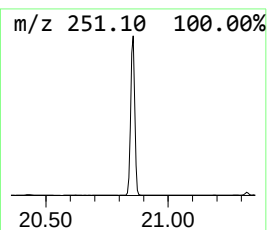
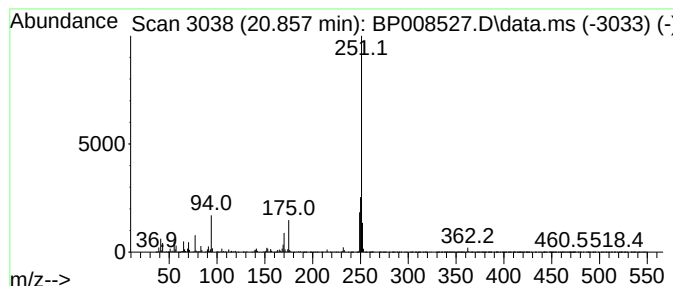
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Quant Title : SVOA CALIBRATION

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

Peak Number 49 Octicizer Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.857	32.48 ng/ul	12096800	Chrysene-d12	21.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octicizer			362	C20H27O4P	001241-94-7	97
2	Octicizer			362	C20H27O4P	001241-94-7	94
3	Octicizer			362	C20H27O4P	001241-94-7	56
4	trans-4-(Dimethylamino)chalcone			251	C17H17NO	022965-98-6	53
5	Phosphoric acid, isodecyl diphen...			390	C22H31O4P	1346599-15-2	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
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Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

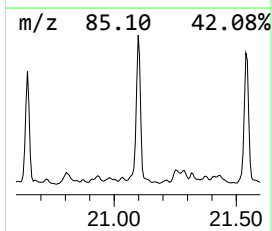
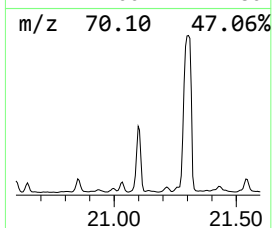
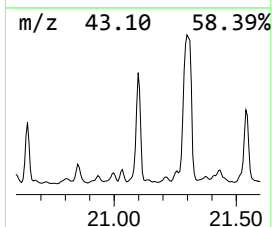
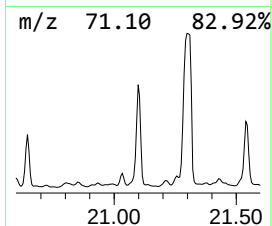
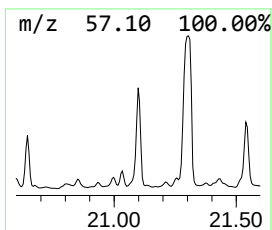
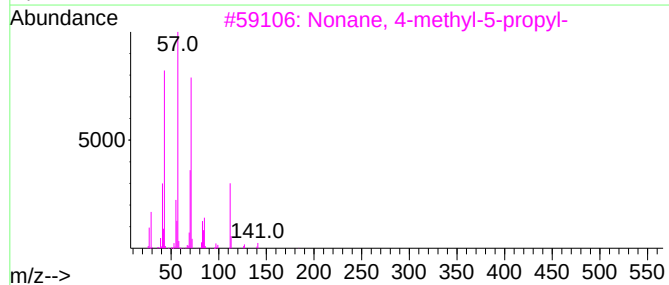
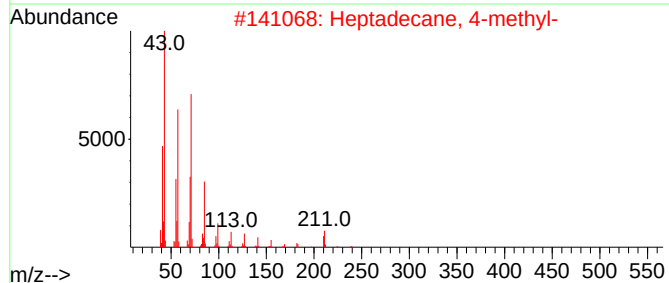
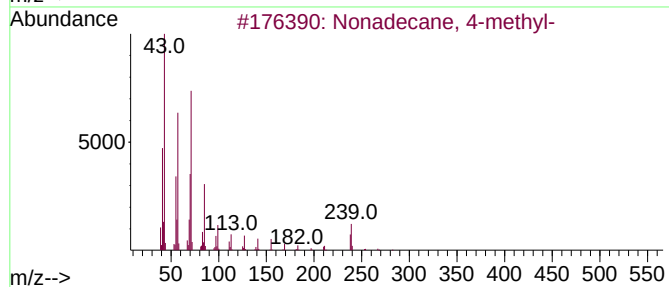
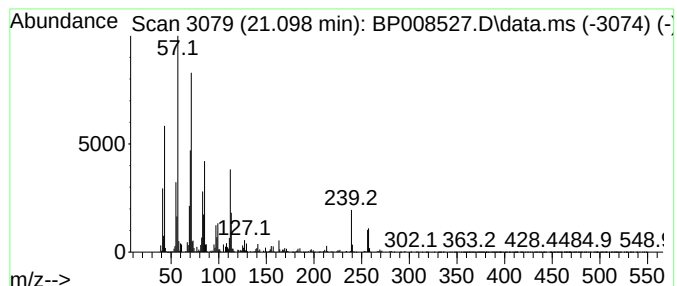
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.098	49.67 ng/ul	18501300	Chrysene-d12	21.369	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane, 4-methyl-	282	C20H42	025117-27-5	86
2	Heptadecane, 4-methyl-	254	C18H38	026429-11-8	64
3	Nonane, 4-methyl-5-propyl-	184	C13H28	062185-55-1	58
4	Heptadecane	240	C17H36	000629-78-7	55
5	Octadecane, 4-methyl-	268	C19H40	010544-95-3	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGLA5ME

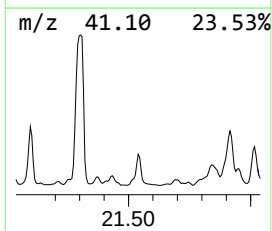
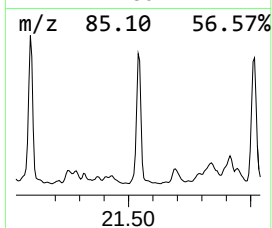
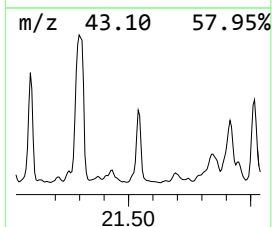
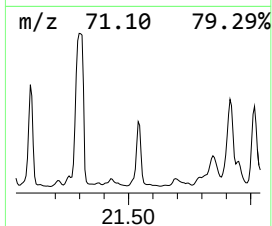
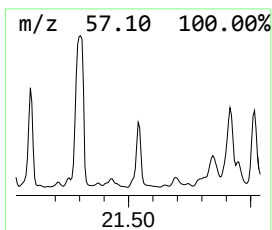
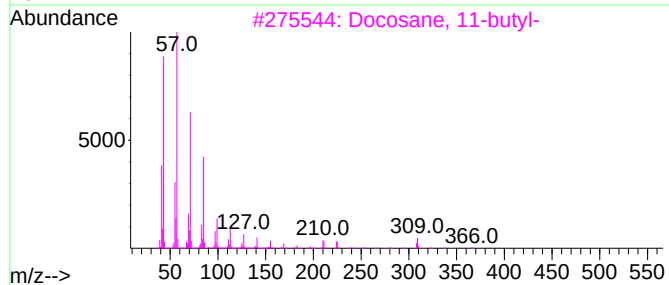
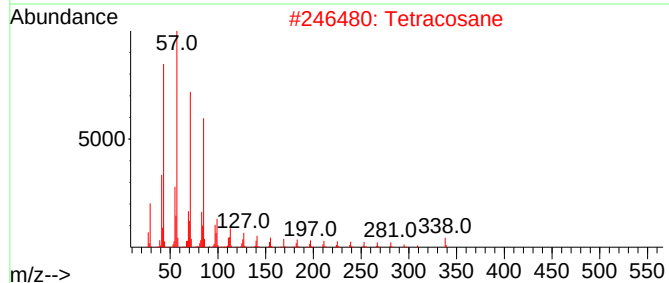
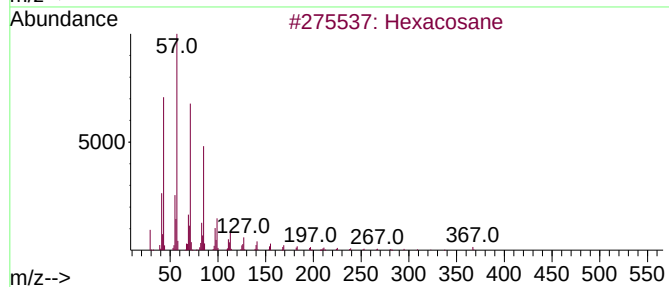
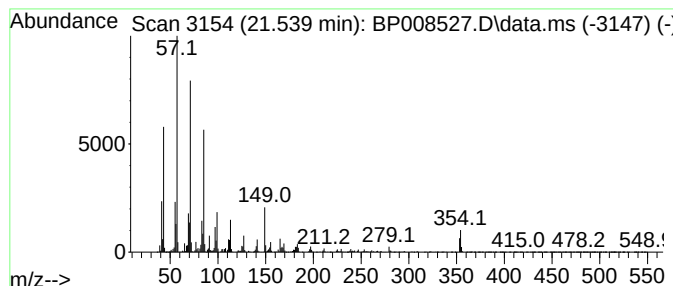
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Quant Title : SVOA CALIBRATION

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

Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.539	38.18 ng/ul	14220300	Chrysene-d12	21.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	93
2			Tetracosane	338	C24H50	000646-31-1	87
3			Docosane, 11-butyl-	366	C26H54	013475-76-8	86
4			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	83
5			Tridecanol, 2-ethyl-2-methyl-	242	C16H34O	1010115-66-1	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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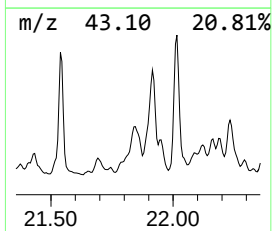
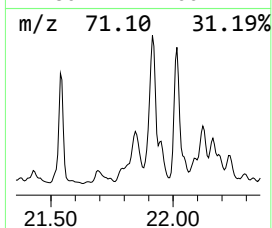
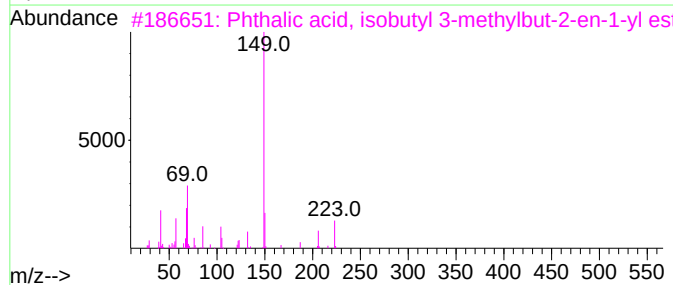
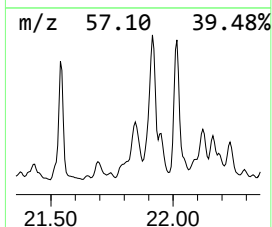
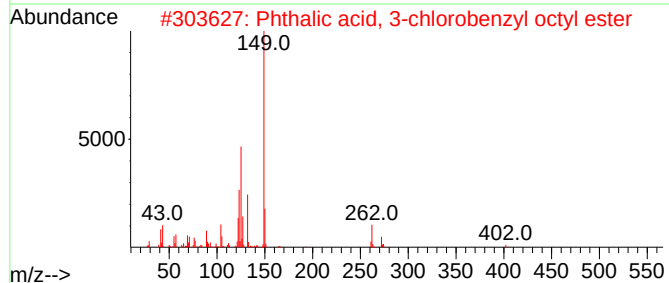
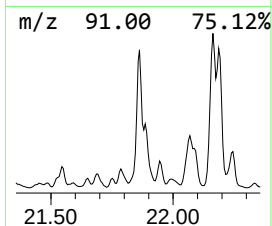
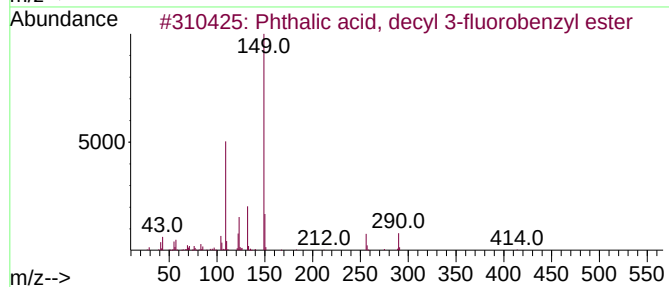
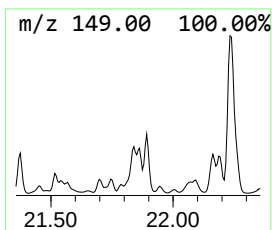
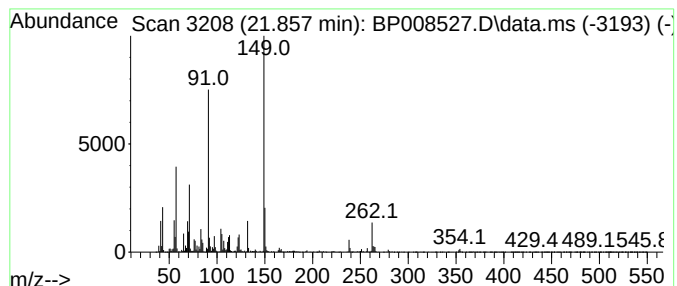
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Quant Title : SVOA CALIBRATION

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

Peak Number 52 Phthalic acid, decyl 3-fluo... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.857	57.24 ng/ul	21319600	Chrysene-d12	21.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, decyl 3-fluoroben...	414	C25H31FO4	1000377-89-4	53
2			Phthalic acid, 3-chlorobenzyl oc...	402	C23H27ClO4	1010371-15-2	50
3			Phthalic acid, isobutyl 3-methyl...	290	C17H22O4	1000315-45-0	50
4			Phthalic acid, hexyl 4-methyl-3-...	399	C22H25NO6	1000382-58-4	50
5			2-(Neopentylloxycarbonyl)benzoic ...	236	C13H16O4	083646-70-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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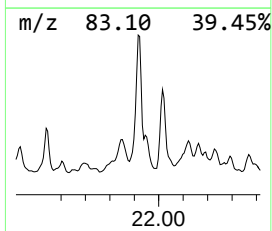
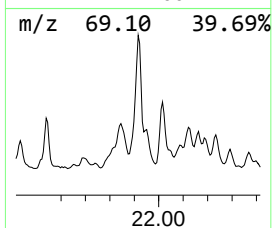
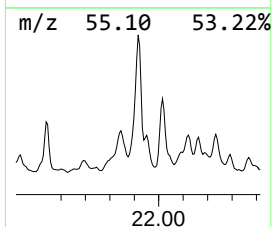
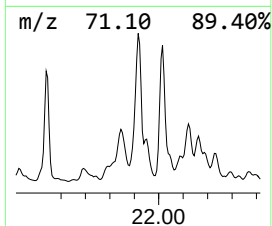
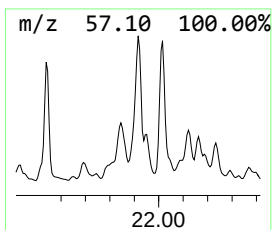
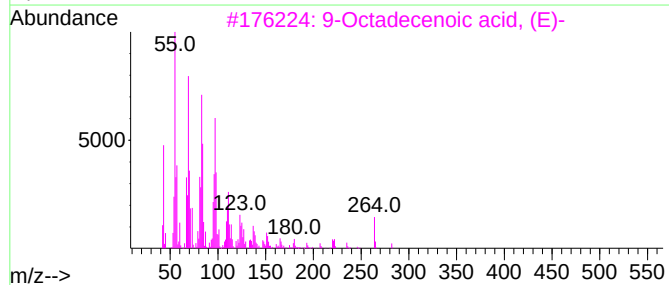
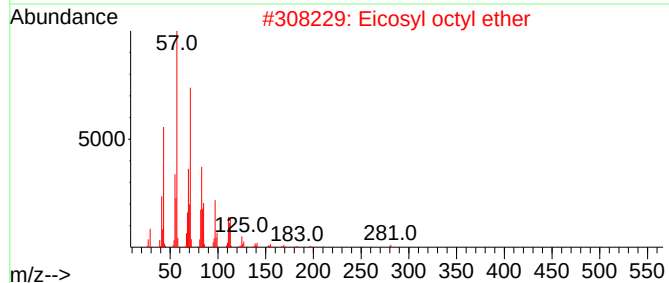
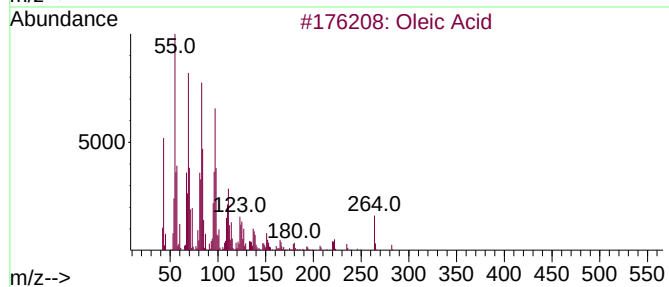
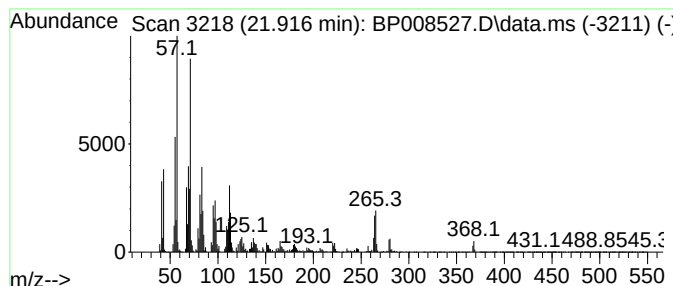
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Quant Title : SVOA CALIBRATION

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

Peak Number 53 Oleic Acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.916	60.55 ng/ul	22551800	Chrysene-d12	21.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oleic Acid	282	C18H34O2	000112-80-1	72
2			Eicosyl octyl ether	410	C28H58O	1000406-38-8	62
3			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	60
4			Octyl tetracosyl ether	467	C32H66O	1000406-39-0	58
5			Oleic Acid	282	C18H34O2	000112-80-1	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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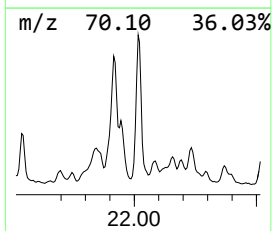
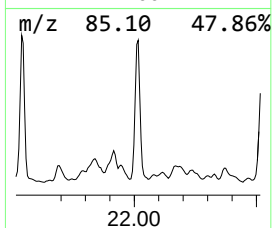
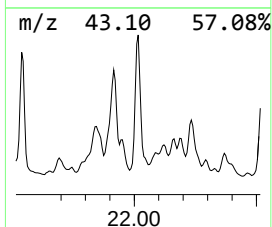
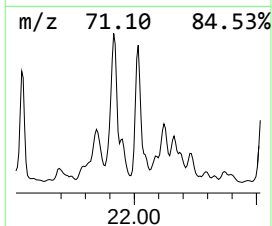
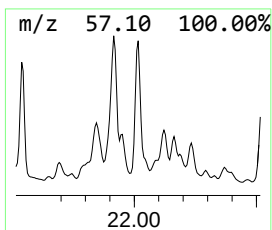
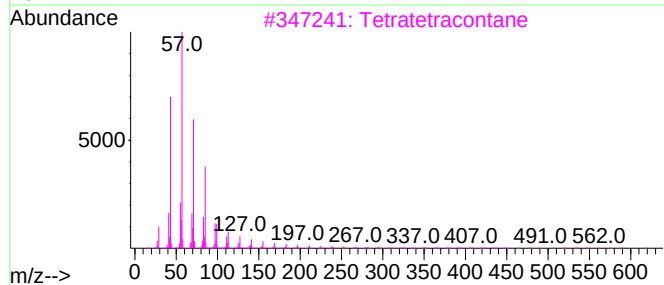
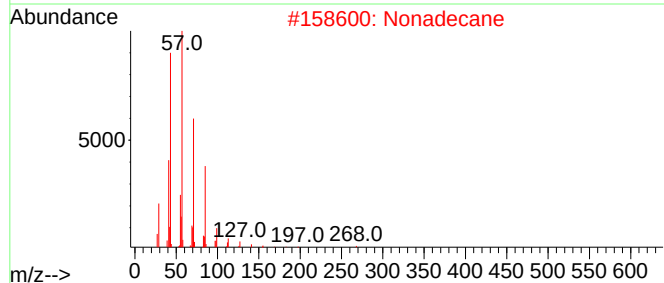
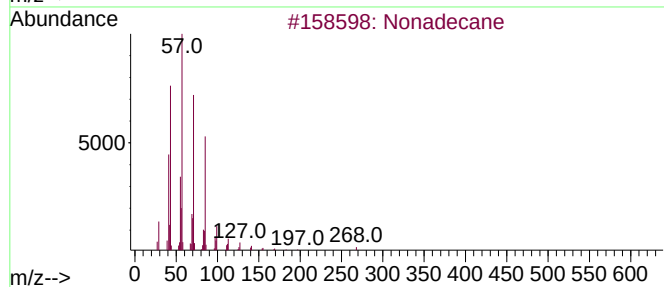
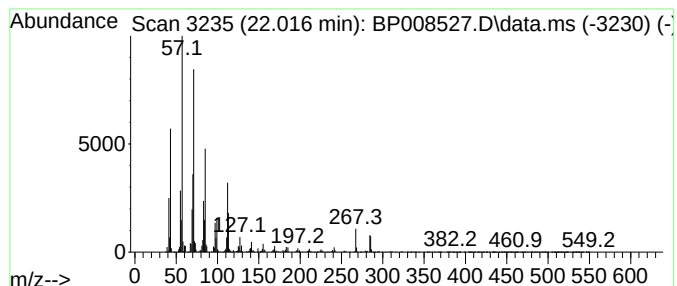
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Quant Title : SVOA CALIBRATION

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

Peak Number 54 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.016	40.13 ng/ul	14945300	Chrysene-d12	21.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	89
2			Nonadecane	268	C19H40	000629-92-5	86
3			Tetratetracontane	619	C44H90	007098-22-8	76
4			Hexacosane	366	C26H54	000630-01-3	76
5			Pentacosane	352	C25H52	000629-99-2	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
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Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

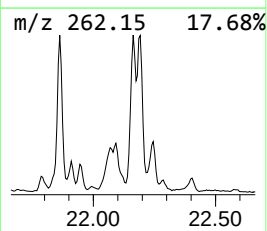
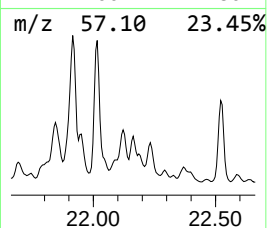
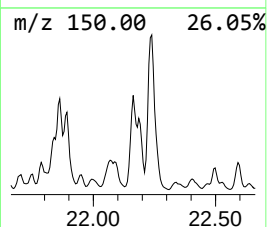
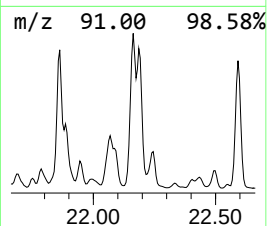
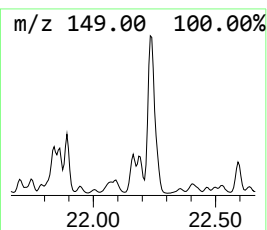
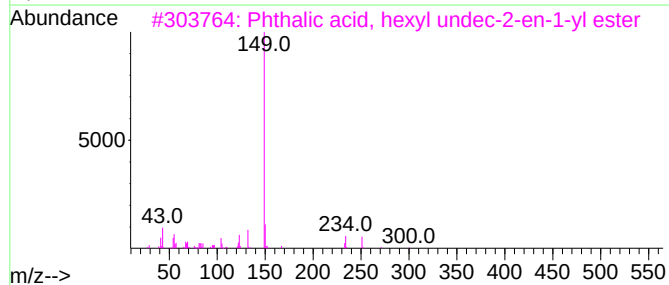
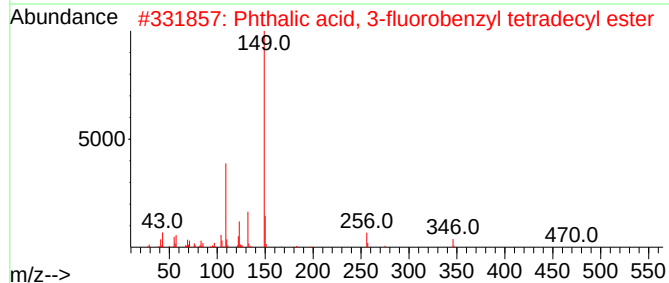
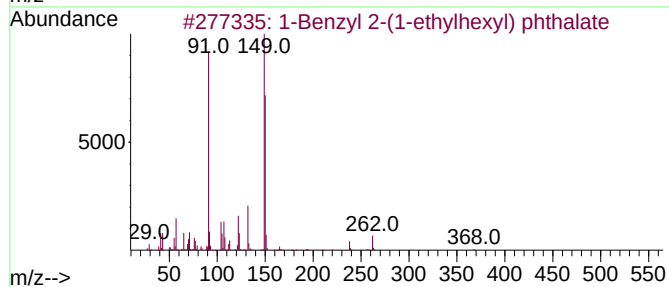
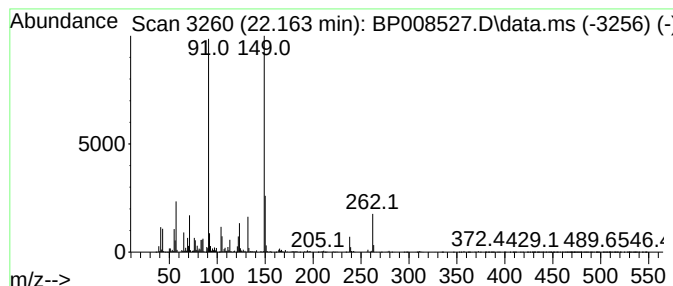
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 55 1-Benzyl 2-(1-ethylhexyl) p... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.163	18.78 ng/ul	6994540	Chrysene-d12	21.369	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Benzyl 2-(1-ethylhexyl) phthalate	368	C23H28O4	1000499-09-5	64
2	Phthalic acid, 3-fluorobenzyl te...	470	C29H39FO4	1010377-89-8	53
3	Phthalic acid, hexyl undec-2-en-...	402	C25H38O4	1000315-41-7	53
4	Phthalic acid, heptyl 4-methyl-3...	413	C23H27NO6	1000382-58-5	53
5	Phthalic acid, 2-methylallyl pen...	290	C17H22O4	1000315-82-4	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
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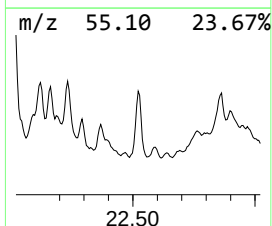
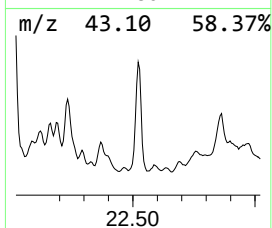
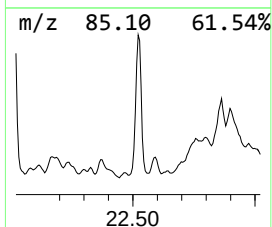
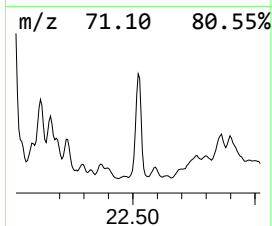
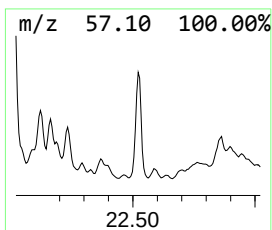
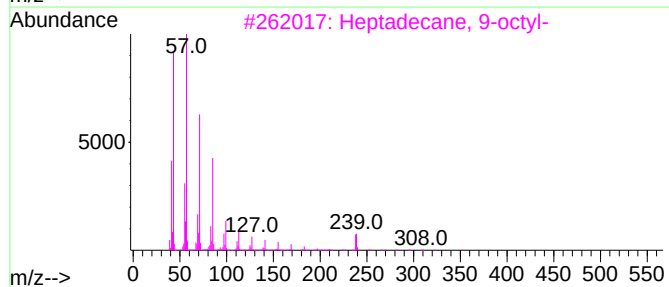
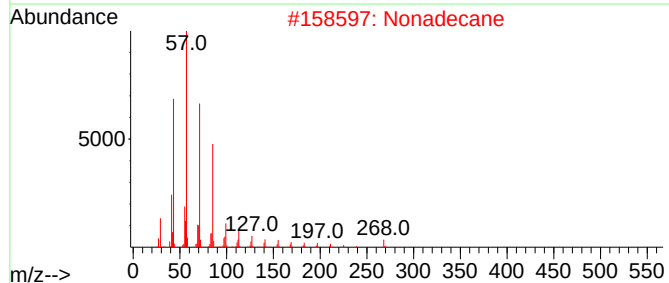
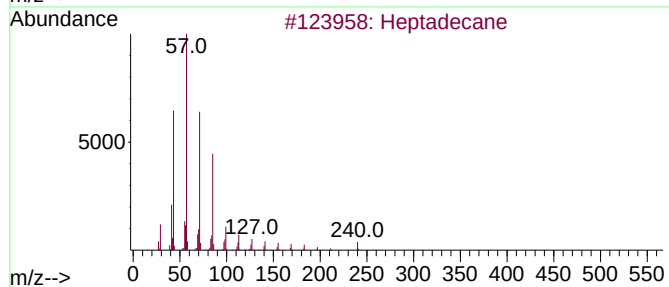
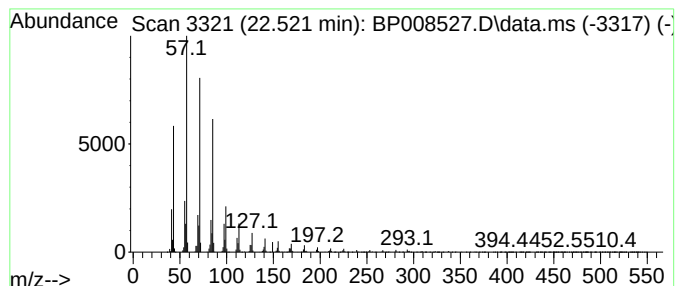
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 56 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.521	21.21 ng/ul	7900110	Chrysene-d12	21.369	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	94
2	Nonadecane	268	C19H40	000629-92-5	94
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
4	Heptacosane	380	C27H56	000593-49-7	91
5	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGLA5ME

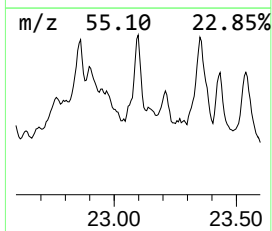
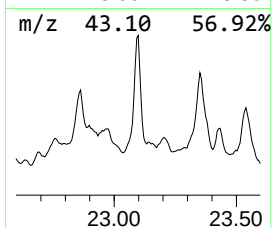
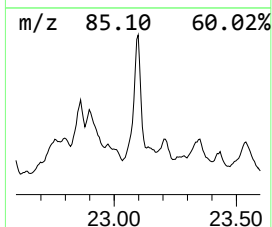
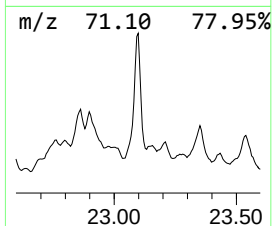
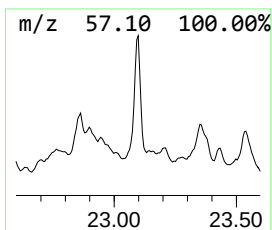
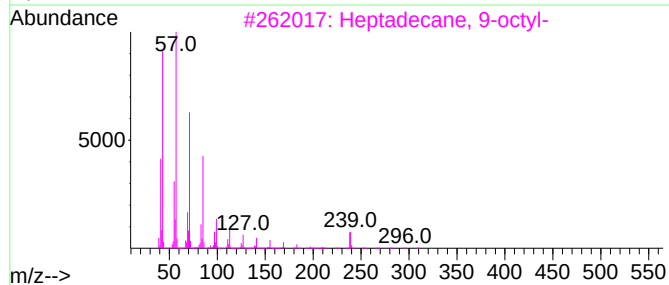
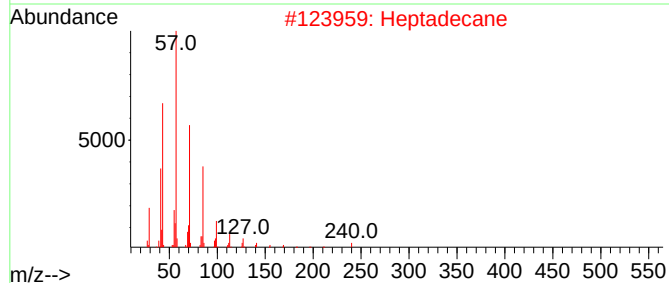
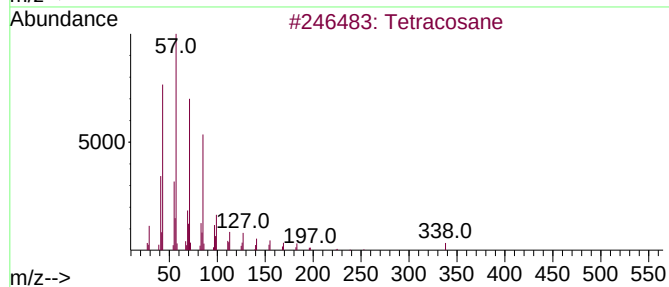
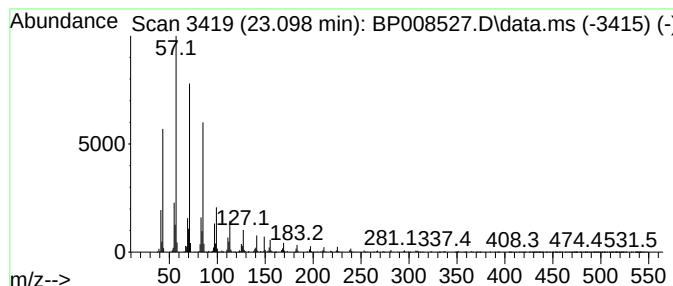
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Quant Title : SVOA CALIBRATION

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

Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.098	14.77 ng/ul	6033740	Perylene-d12	23.804

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	95
2		Heptadecane	240	C17H36	000629-78-7	94
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5		11-Methylpentacosane	366	C26H54	015689-71-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGLA5ME

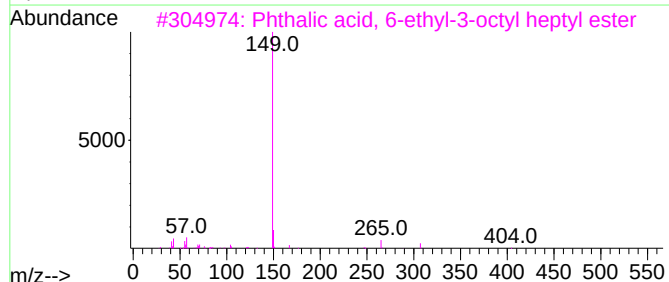
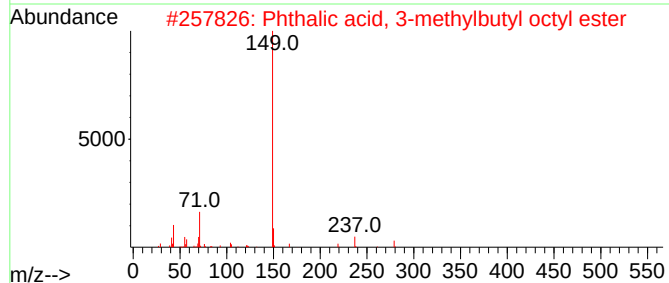
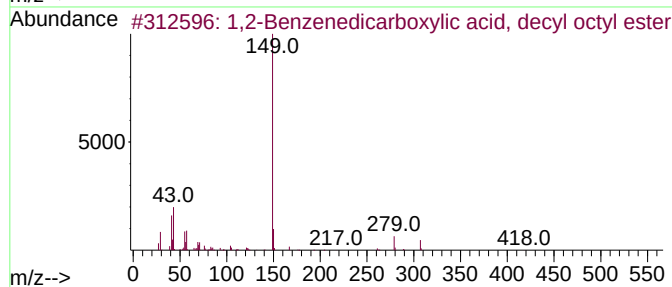
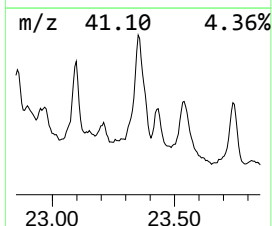
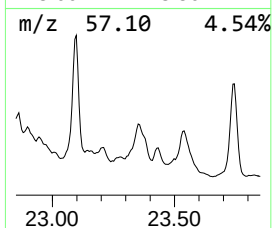
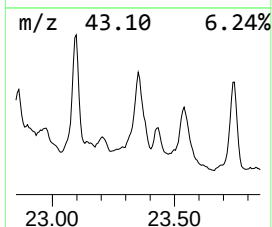
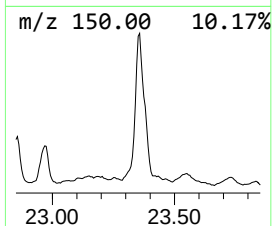
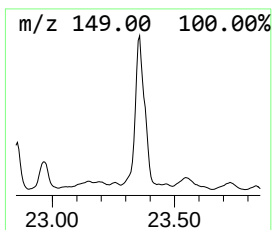
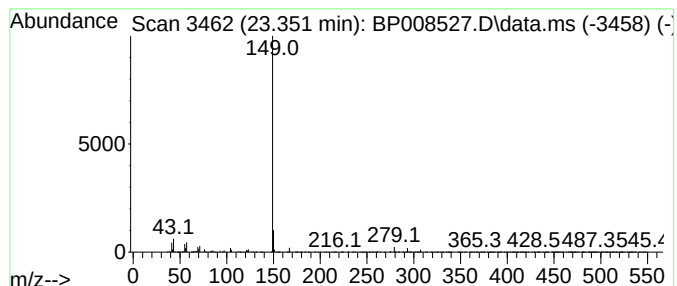
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Quant Title : SVOA CALIBRATION

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

Peak Number 60 1,2-Benzenedicarboxylic aci... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.351	35.61 ng/ul	14546600	Perylene-d12	23.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenedicarboxylic acid, de...	418	C26H42O4	000119-07-3	83
2			Phthalic acid, 3-methylbutyl oct...	348	C21H32O4	1000356-80-9	80
3			Phthalic acid, 6-ethyl-3-octyl h...	404	C25H40O4	1000315-17-7	72
4			Phthalic acid, hept-4-yl pentyl ...	334	C20H30O4	1000356-78-5	72
5			Di-n-octyl phthalate	390	C24H38O4	000117-84-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGLA5ME

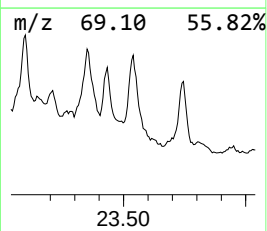
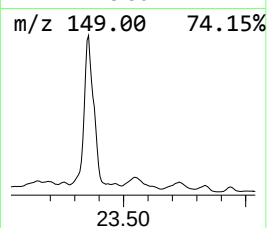
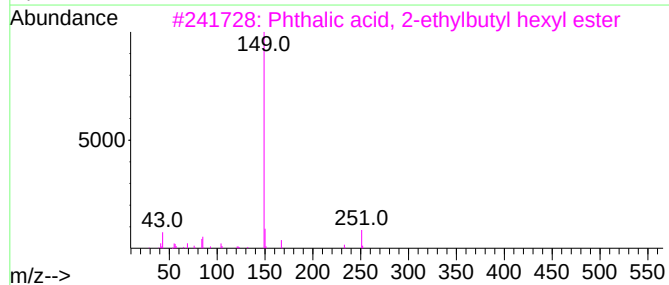
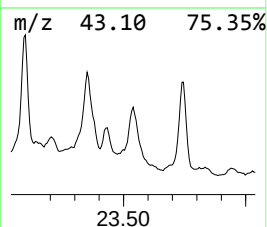
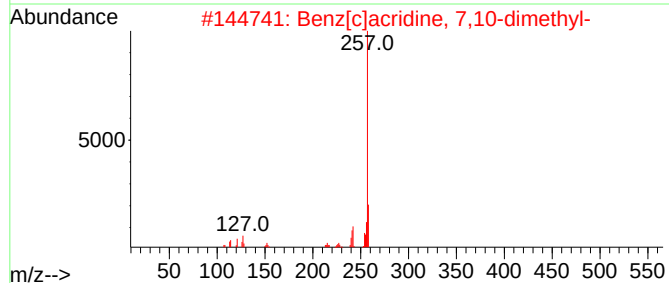
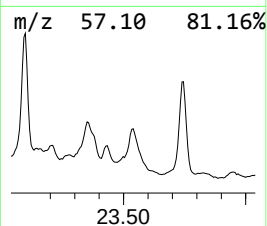
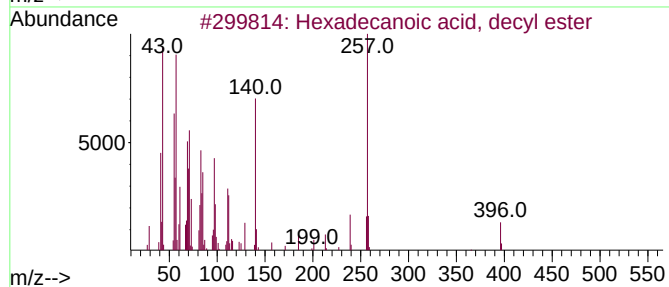
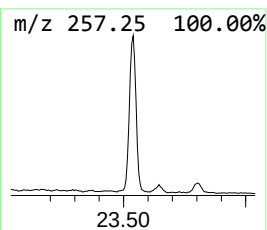
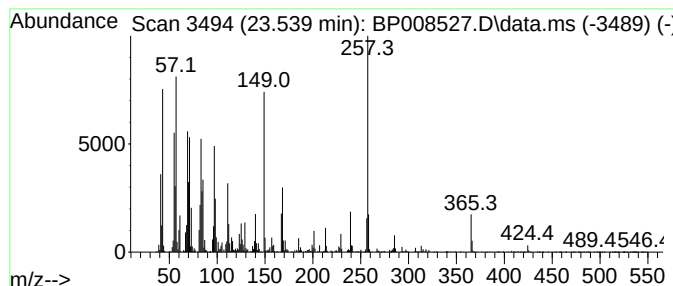
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Quant Title : SVOA CALIBRATION

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

Peak Number 61 Hexadecanoic acid, decyl ester Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.539	22.33 ng/ul	9122500	Perylene-d12	23.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, decyl ester	396	C26H52O2	042232-27-9	64
2			Benz[c]acridine, 7,10-dimethyl-	257	C19H15N	002381-40-0	30
3			Phthalic acid, 2-ethylbutyl hexy...	334	C20H30O4	1000356-89-1	22
4			1-Hexadecanethiol	258	C16H34S	002917-26-2	18
5			Pyrrole, 1-methyl-2-phenyl-4-(ph...	257	C19H15N	066463-26-1	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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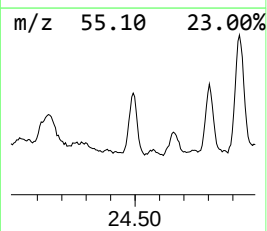
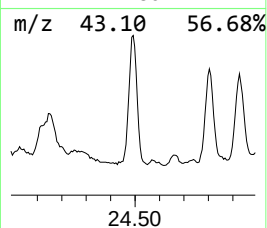
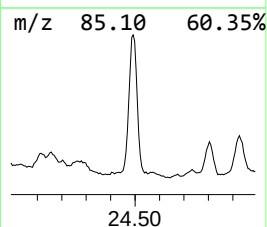
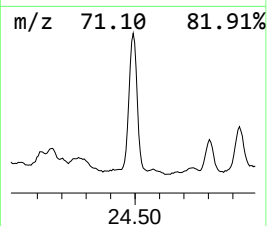
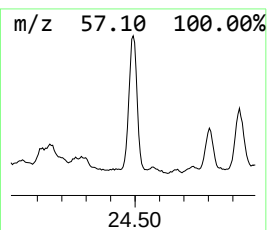
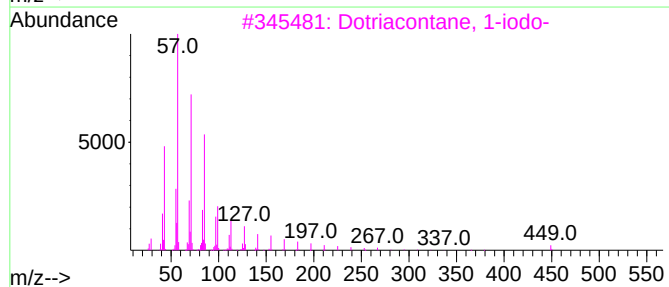
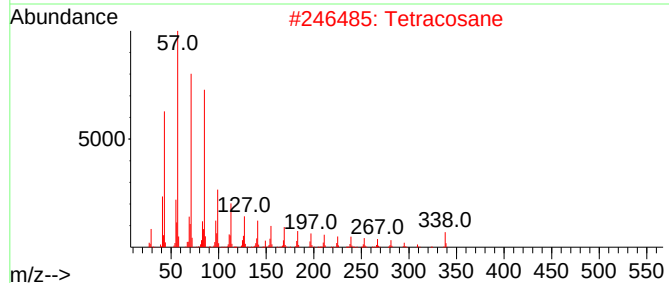
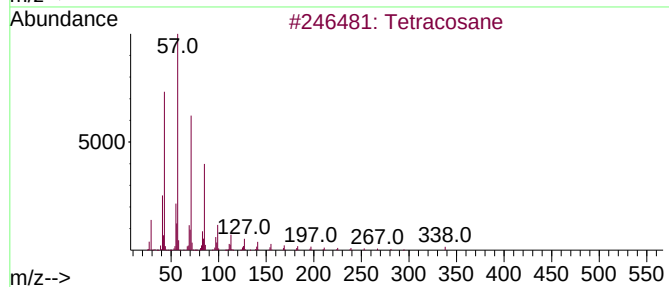
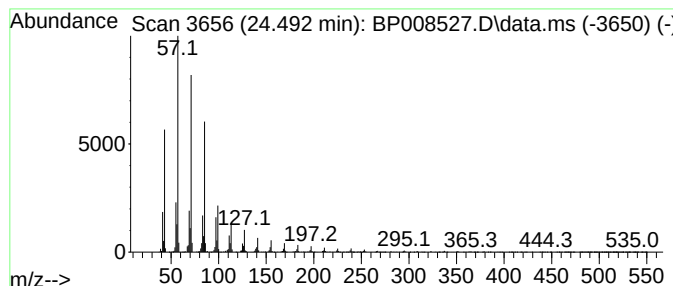
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Quant Title : SVOA CALIBRATION

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

Peak Number 63 (DEL) Alkane: Straight-Chai... Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.492	12.73 ng/ul	5200490	Perylene-d12	23.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	97
2			Tetracosane	338	C24H50	000646-31-1	95
3			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	94
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
5			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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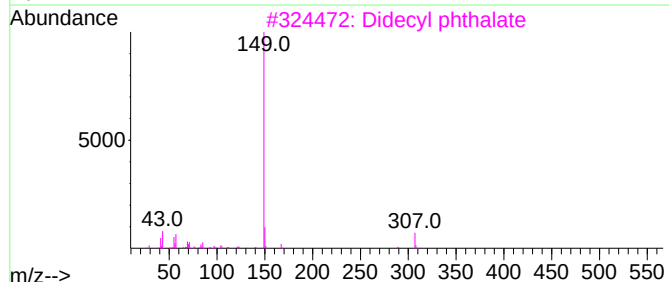
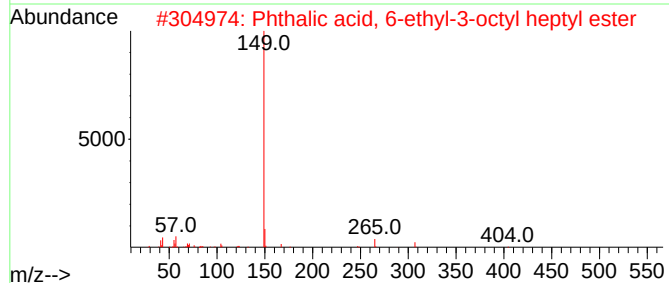
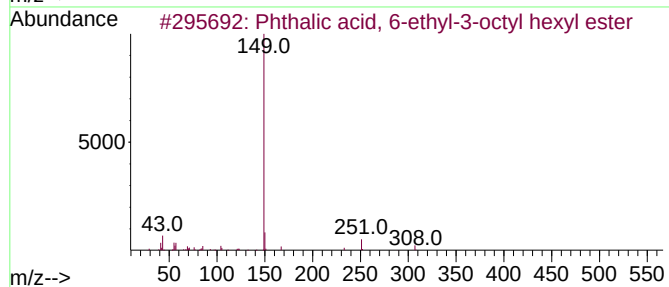
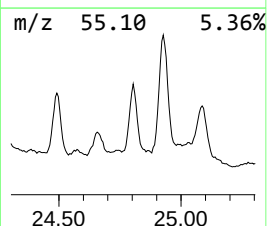
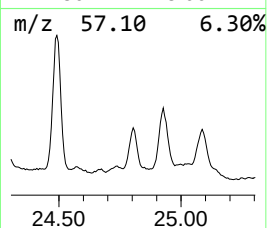
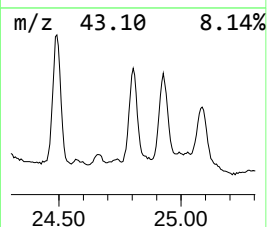
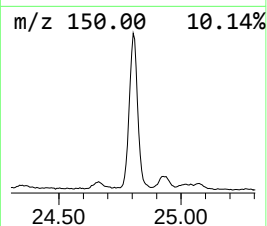
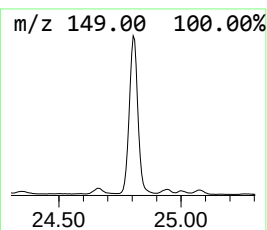
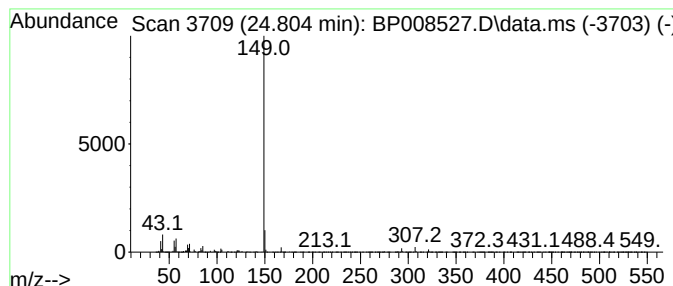
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Quant Title : SVOA CALIBRATION

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

Peak Number 64 Phthalic acid, 6-ethyl-3-oc... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.804	19.21 ng/ul	7847570	Perylene-d12	23.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, 6-ethyl-3-octyl h...	390	C24H38O4	1000315-17-6	86
2			Phthalic acid, 6-ethyl-3-octyl h...	404	C25H40O4	1000315-17-7	86
3			Didecyl phthalate	446	C28H46O4	000084-77-5	80
4			Phthalic acid, decyl pentyl ester	376	C23H36O4	1000309-00-9	78
5			Phthalic acid, 3-methylbutyl oct...	488	C31H52O4	1010356-81-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
Data File : BP008527.D
Acq On : 23 Dec 2021 16:34
Operator : CG/JU
Sample : M5171-09ME
Misc :
ALS Vial : 13 Sample Multiplier: 1

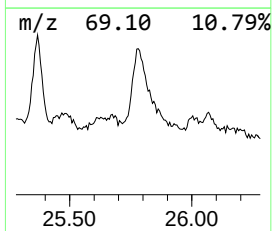
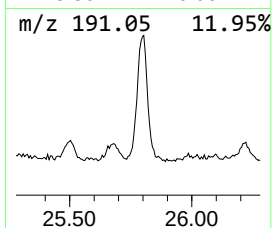
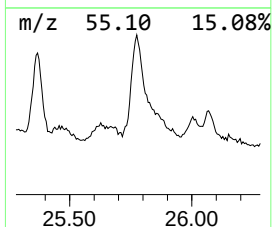
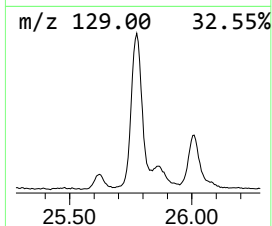
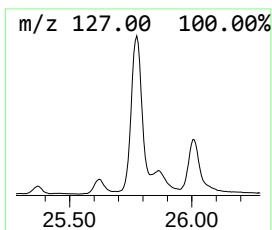
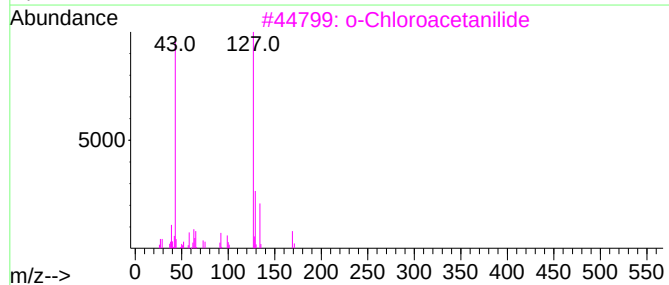
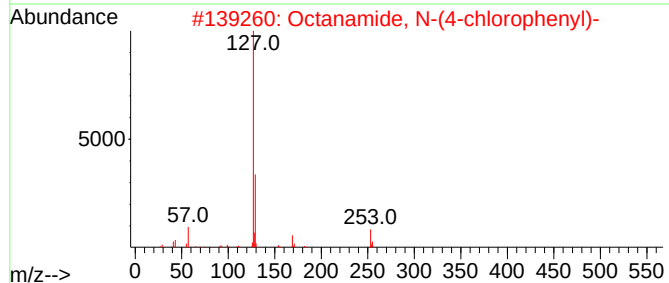
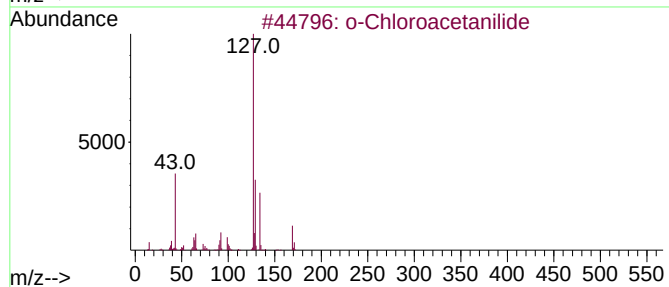
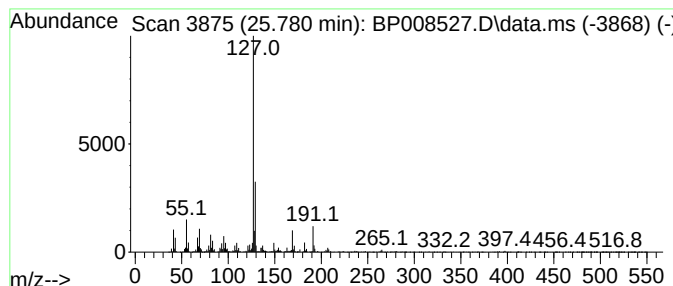
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 66 o-Chloroacetanilide Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.	
25.780	17.09 ng/ul	6981480	Perylene-d12	23.804	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Chloroacetanilide	169	C8H8ClNO	000533-17-5	80
2	Octanamide, N-(4-chlorophenyl)-	253	C14H20ClNO	1000340-16-6	72
3	o-Chloroacetanilide	169	C8H8ClNO	000533-17-5	72
4	o-Chloroacetanilide	169	C8H8ClNO	000533-17-5	72
5	Octanamide, N-(3-chlorophenyl)-	253	C14H20ClNO	1000306-92-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
 Data File : BP008527.D
 Acq On : 23 Dec 2021 16:34
 Operator : CG/JU
 Sample : M5171-09ME
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGLA5ME

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.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
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Benzene, propyl-	6.893	15.9	ng/ul	8176940	1	7.910	10256700	20.0
Benzene, 1-ethy...	7.005	40.7	ng/ul	20887100	1	7.910	10256700	20.0
Benzene, 1,2,3-...	7.140	21.4	ng/ul	10999500	1	7.910	10256700	20.0
Benzene, 1-ethy...	7.305	13.3	ng/ul	6808870	1	7.910	10256700	20.0
Benzene, 1-ethy...	8.028	23.6	ng/ul	12125500	1	7.910	10256700	20.0
(DEL) Alkane: C...	8.210	9.4	ng/ul	4804420	1	7.910	10256700	20.0
Indane	8.281	13.9	ng/ul	7148730	1	7.910	10256700	20.0
Benzene, 1-meth...	8.463	23.2	ng/ul	11916800	1	7.910	10256700	20.0
(DEL) Alkane: S...	9.163	61.4	ng/ul	31481000	1	7.910	10256700	20.0
Benzene, 2-ethe...	10.116	15.1	ng/ul	17459400	2	10.716	23141700	20.0
(DEL) Alkane: S...	11.651	3.6	ng/ul	4178510	2	10.716	23141700	20.0
(DEL) Alkane: S...	11.763	7.3	ng/ul	8415130	2	10.716	23141700	20.0
(DEL) Alkane: S...	12.157	16.3	ng/ul	18859400	2	10.716	23141700	20.0
Benzenamine, 2,...	12.969	17.9	ng/ul	8641670	3	14.534	9678980	20.0
(DEL) Alkane: S...	13.104	10.1	ng/ul	4905870	3	14.534	9678980	20.0
Naphthalene, 2,...	13.704	13.5	ng/ul	6518170	3	14.534	9678980	20.0
Naphthalene, 1,...	13.869	42.8	ng/ul	20722600	3	14.534	9678980	20.0
Carbonic acid, ...	14.045	14.1	ng/ul	6837260	3	14.534	9678980	20.0
Naphthalene, 2,...	14.092	9.7	ng/ul	4687950	3	14.534	9678980	20.0
(DEL) Alkane: S...	14.457	44.3	ng/ul	21431600	3	14.534	9678980	20.0
(DEL) Alkane: S...	15.410	29.3	ng/ul	14172800	3	14.534	9678980	20.0
Dodecane, 1-iodo-	16.275	53.3	ng/ul	14271100	4	17.286	5359780	20.0
(DEL) Alkane: S...	16.298	23.6	ng/ul	6325910	4	17.286	5359780	20.0
Benzene, (1-met...	16.387	21.2	ng/ul	5685880	4	17.286	5359780	20.0
(DEL) Alkane: S...	17.069	35.6	ng/ul	9535770	4	17.286	5359780	20.0
Benzene, (1-met...	17.204	18.7	ng/ul	5004220	4	17.286	5359780	20.0
Nonadecane	17.804	27.9	ng/ul	7470920	4	17.286	5359780	20.0
(DEL) Alkane: S...	18.475	19.0	ng/ul	5101160	4	17.286	5359780	20.0
Phenanthrene, 2...	19.033	17.8	ng/ul	4781460	4	17.286	5359780	20.0
(DEL) Alkane: S...	19.080	26.9	ng/ul	7212290	4	17.286	5359780	20.0
Octadecanoic acid	19.427	26.6	ng/ul	9917370	5	21.369	7448990	20.0
Methacrylic aci...	19.498	19.7	ng/ul	7346160	5	21.369	7448990	20.0
1,2-Benzenedica...	19.916	15.2	ng/ul	5652390	5	21.369	7448990	20.0
(DEL) Alkane: S...	20.157	18.2	ng/ul	6787000	5	21.369	7448990	20.0
(DEL) Alkane: S...	20.645	15.4	ng/ul	5748980	5	21.369	7448990	20.0
Octicizer	20.857	32.5	ng/ul	12096800	5	21.369	7448990	20.0
(DEL) Alkane: S...	21.098	49.7	ng/ul	18501300	5	21.369	7448990	20.0
(DEL) Alkane: S...	21.539	38.2	ng/ul	14220300	5	21.369	7448990	20.0
Phthalic acid, ...	21.857	57.2	ng/ul	21319600	5	21.369	7448990	20.0
Oleic Acid	21.916	60.5	ng/ul	22551800	5	21.369	7448990	20.0
(DEL) Alkane: S...	22.016	40.1	ng/ul	14945300	5	21.369	7448990	20.0
1-Benzyl 2-(1-e...	22.163	18.8	ng/ul	6994540	5	21.369	7448990	20.0
(DEL) Alkane: S...	22.521	21.2	ng/ul	7900110	5	21.369	7448990	20.0
(DEL) Alkane: S...	23.098	14.8	ng/ul	6033740	6	23.804	8170610	20.0
1,2-Benzenedica...	23.351	35.6	ng/ul	14546600	6	23.804	8170610	20.0
Hexadecanoic ac...	23.539	22.3	ng/ul	9122500	6	23.804	8170610	20.0
(DEL) Alkane: S...	24.492	12.7	ng/ul	5200490	6	23.804	8170610	20.0
Phthalic acid, ...	24.804	19.2	ng/ul	7847570	6	23.804	8170610	20.0
o-Chloroacetani...	25.780	17.1	ng/ul	6981480	6	23.804	8170610	20.0