

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
 Data File : BP002131.D
 Acq On : 09 Mar 2020 21:12
 Operator : CG/JU
 Sample : L1834-14
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0DK7

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.917	3	5	17	rVB	861265	1085373	12.38%	1.255%
2	3.164	43	47	55	rVB	60700	85925	0.98%	0.099%
3	3.799	152	155	159	rVB3	6976	10107	0.12%	0.012%
4	3.893	168	171	177	rVB3	10997	15190	0.17%	0.018%
5	4.793	319	324	330	rBV6	8855	15616	0.18%	0.018%
6	4.993	352	358	365	rVB	119285	181519	2.07%	0.210%
7	5.340	412	417	420	rBV4	6139	11568	0.13%	0.013%
8	6.058	535	539	542	rBV2	10638	18476	0.21%	0.021%
9	6.816	661	668	671	rBV	212552	395392	4.51%	0.457%
10	6.975	688	695	707	rBV	928172	1422921	16.23%	1.646%
11	7.175	721	729	741	rBV	917390	1542857	17.60%	1.784%
12	7.528	783	789	796	rBV	104047	167832	1.91%	0.194%
13	7.634	801	807	820	rBV	947319	1645206	18.77%	1.903%
14	7.987	858	867	874	rBV	820536	1315573	15.01%	1.522%
15	8.340	917	927	938	rBV	487791	809792	9.24%	0.937%
16	8.775	994	1001	1008	rBV	1027220	1717756	19.59%	1.987%
17	9.263	1078	1084	1091	rVB	61936	95259	1.09%	0.110%
18	9.499	1114	1124	1139	rBV	850513	1422495	16.23%	1.645%
19	10.034	1206	1215	1229	rBV	1316701	2245296	25.61%	2.597%
20	10.193	1235	1242	1244	rBV4	17871	32276	0.37%	0.037%
21	10.275	1250	1256	1259	rBV	175231	313397	3.57%	0.362%
22	10.310	1259	1262	1271	rVB	191364	317996	3.63%	0.368%
23	10.404	1271	1278	1292	rVB	1314721	2237833	25.53%	2.588%
24	10.546	1296	1302	1311	rVB	47727	83498	0.95%	0.097%
25	10.793	1338	1344	1352	rBV3	25766	52527	0.60%	0.061%
26	11.004	1376	1380	1387	rVB3	16440	30093	0.34%	0.035%
27	11.222	1411	1417	1424	rBV2	29437	56640	0.65%	0.066%
28	11.710	1496	1500	1503	rBV	10763	16492	0.19%	0.019%
29	11.757	1503	1508	1514	rVB	36912	54930	0.63%	0.064%
30	12.004	1546	1550	1557	rVB3	25374	41994	0.48%	0.049%
31	12.098	1561	1566	1573	rBV4	12173	28216	0.32%	0.033%
32	12.216	1583	1586	1596	rVB4	6163	12175	0.14%	0.014%
33	12.298	1596	1600	1607	rVV2	6545	12741	0.15%	0.015%
34	12.381	1610	1614	1620	rVV3	4344	9033	0.10%	0.010%

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 Peak separation: 5

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35	12.457	1622	1627	1632	rVB3	15428	26575	0.30%	0.031%
36	12.840	1688	1692	1696	rBV3	7806	13135	0.15%	0.015%
37	13.069	1727	1731	1737	rBV2	65486	104211	1.19%	0.121%
38	13.293	1760	1769	1781	rVB2	109870	224596	2.56%	0.260%
39	13.463	1793	1798	1799	rBV	7103	10344	0.12%	0.012%
40	13.687	1828	1836	1840	rBV	2760858	3771058	43.02%	4.361%
41	13.734	1840	1844	1859	rVB	250866	388456	4.43%	0.449%
42	13.957	1875	1882	1891	rBV	3017949	4687477	53.47%	5.421%
43	14.028	1891	1894	1900	rVB3	14653	24015	0.27%	0.028%
44	14.269	1928	1935	1947	rVB2	2079229	3143429	35.86%	3.636%
45	14.481	1964	1971	1983	rBV2	92123	228481	2.61%	0.264%
46	14.598	1987	1991	1996	rVB4	8802	15277	0.17%	0.018%
47	15.269	2098	2105	2120	rBV	4170887	6142089	70.06%	7.104%
48	15.387	2120	2125	2139	rVV	868108	1330318	15.18%	1.539%
49	15.510	2142	2146	2158	rVB	51377	89499	1.02%	0.104%
50	15.839	2199	2202	2204	rBV2	7315	10673	0.12%	0.012%
51	16.057	2235	2239	2244	rVB3	17619	24159	0.28%	0.028%
52	16.810	2363	2367	2375	rVB	15740	26727	0.30%	0.031%
53	17.022	2395	2403	2413	rBV	2643675	3806262	43.42%	4.402%
54	17.122	2413	2420	2433	rBV	4346928	6155913	70.22%	7.120%
55	17.945	2555	2560	2565	rBV2	119154	178669	2.04%	0.207%
56	17.998	2565	2569	2575	rVB2	57795	88109	1.01%	0.102%
57	18.239	2606	2610	2616	rBV4	19252	29752	0.34%	0.034%
58	18.386	2633	2635	2640	rVB	13742	17082	0.19%	0.020%
59	19.016	2737	2742	2747	rBV	61928	91646	1.05%	0.106%
60	19.180	2766	2770	2777	rBV2	65320	105418	1.20%	0.122%
61	19.257	2779	2783	2789	rVB	53861	80411	0.92%	0.093%
62	19.369	2796	2802	2811	rBV	5700562	7843260	89.47%	9.071%
63	19.439	2811	2814	2824	rVB4	42802	80430	0.92%	0.093%
64	19.780	2869	2872	2876	rBV3	18712	28043	0.32%	0.032%
65	19.933	2894	2898	2900	rBV3	20336	26114	0.30%	0.030%
66	20.410	2976	2979	2983	rVB	52939	58439	0.67%	0.068%
67	20.863	3052	3056	3061	rBV	205038	243522	2.78%	0.282%
68	21.116	3094	3099	3109	rBV	3726258	4572598	52.16%	5.288%
69	21.292	3125	3129	3135	rBV	523510	560215	6.39%	0.648%
70	21.721	3197	3202	3209	rBV	975928	1144257	13.05%	1.323%
71	22.180	3275	3280	3293	rVB	1465663	1868032	21.31%	2.160%

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Smoothing : OFF Filtering: 5
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72	22.686	3360	3366	3379	rBV	1512027	2232385	25.47%	2.582%
73	23.180	3442	3450	3457	rBV	4831491	8766380	100.00%	10.139%
74	23.251	3457	3462	3467	rVV	1384970	2373767	27.08%	2.745%
75	23.321	3467	3474	3489	rVB	2554967	4907899	55.99%	5.676%
76	23.898	3565	3572	3582	rBV	941568	1875505	21.39%	2.169%
77	24.645	3692	3699	3713	rVB	498747	1135400	12.95%	1.313%
78	25.510	3840	3846	3858	rVB2	172972	429120	4.90%	0.496%

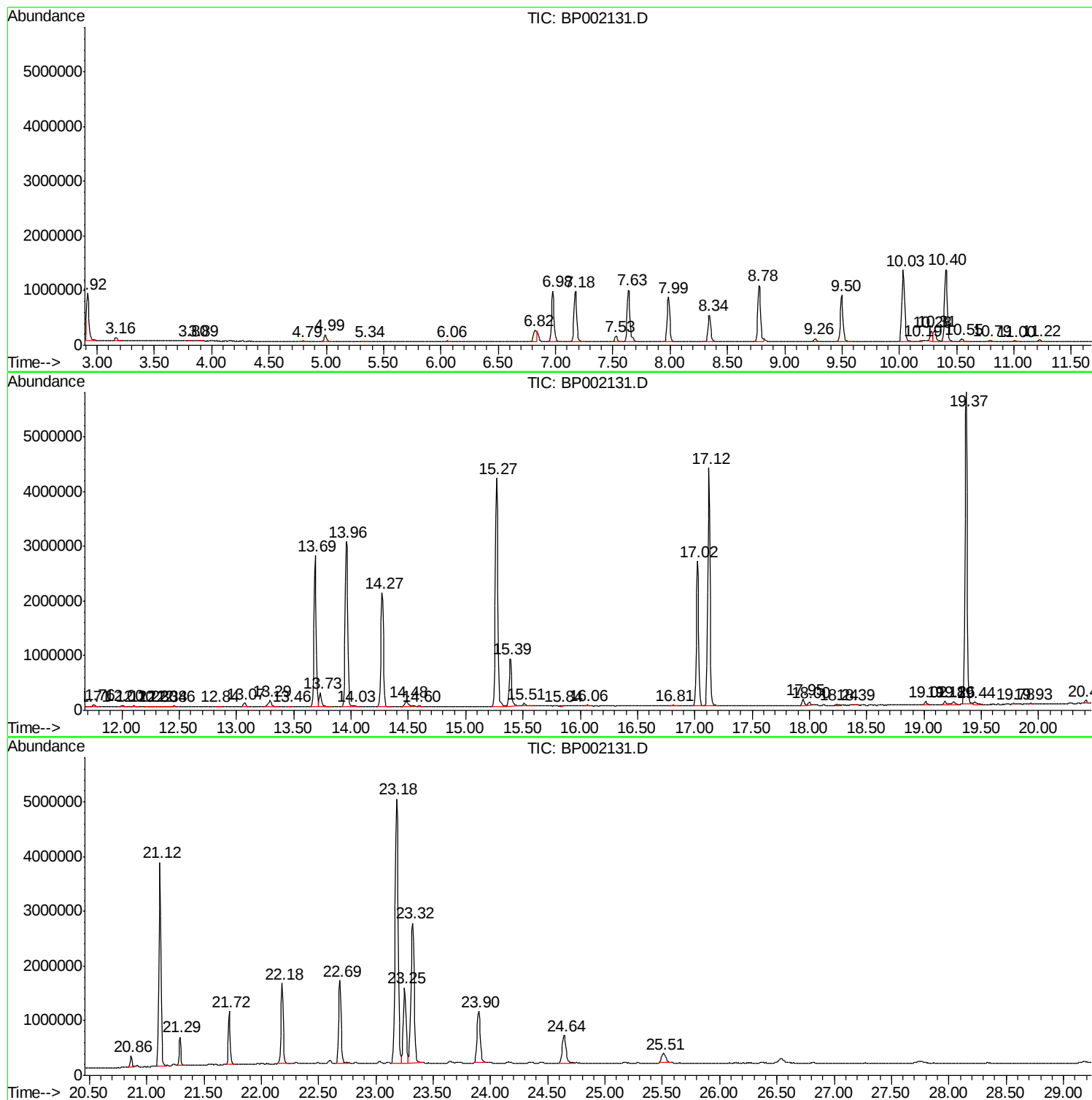
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Instrument :
BNA_P
ClientSampled :
C0DK7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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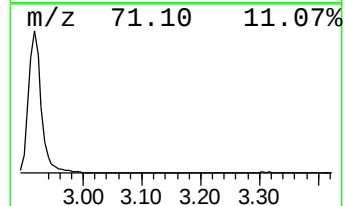
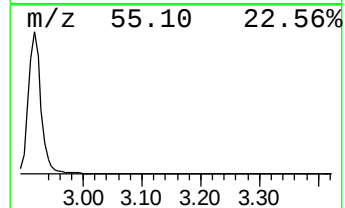
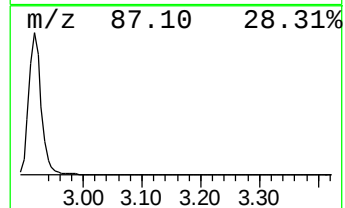
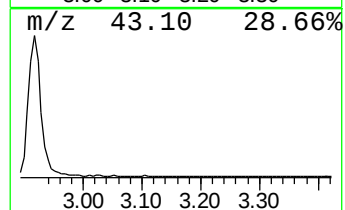
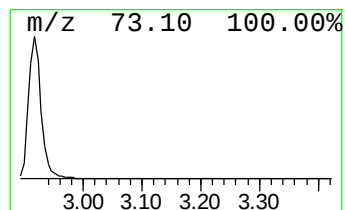
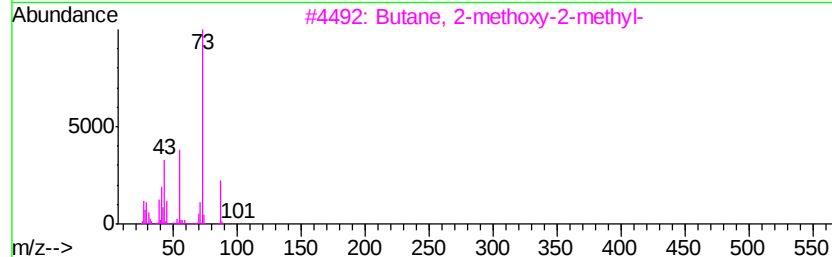
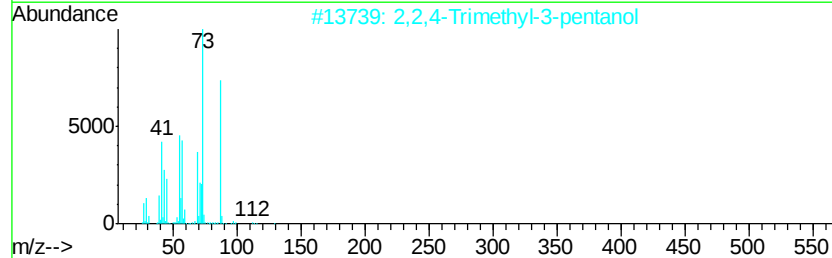
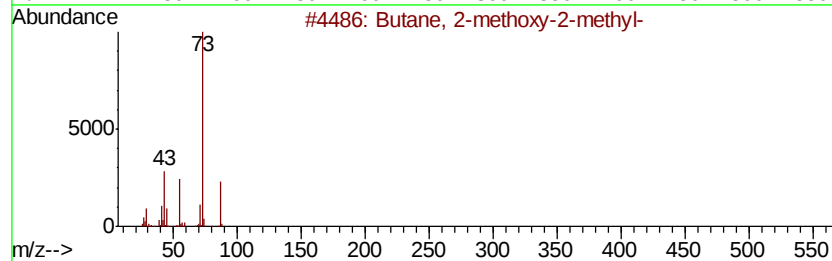
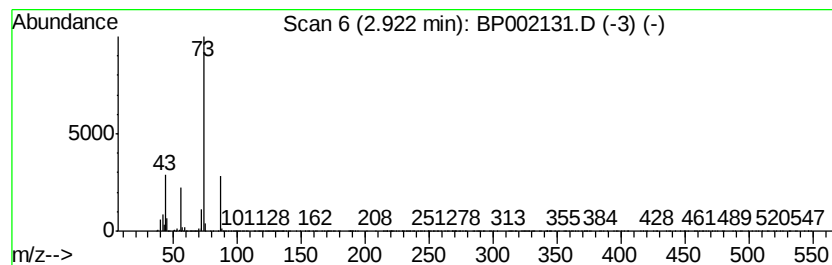
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	13.19 ng/ul	1085370	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
3		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	39
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	12



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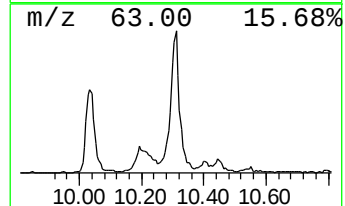
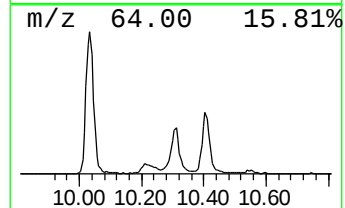
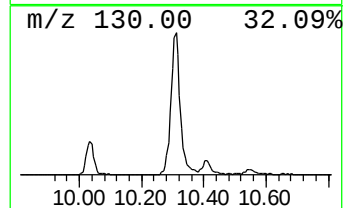
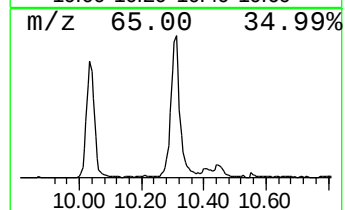
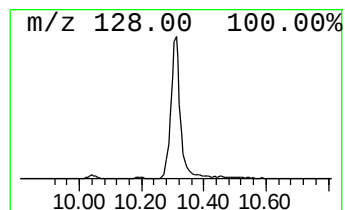
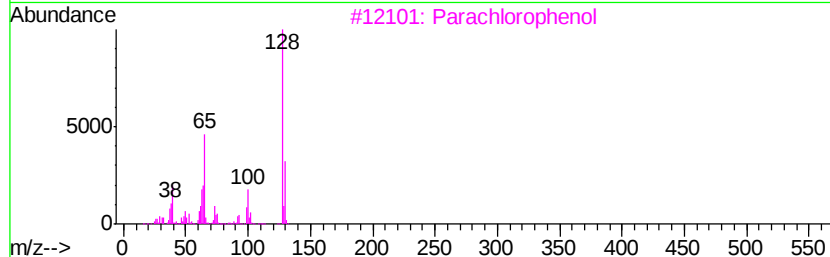
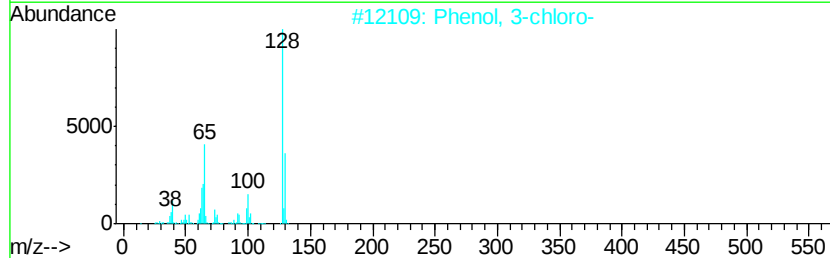
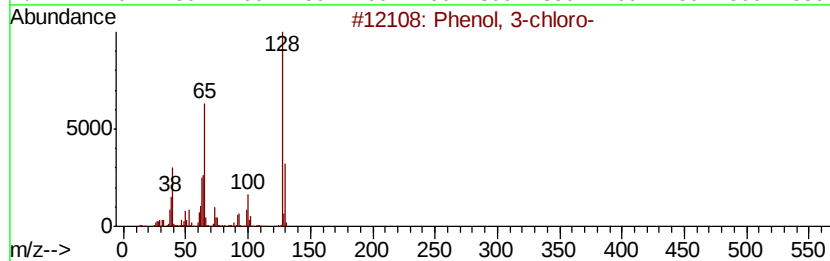
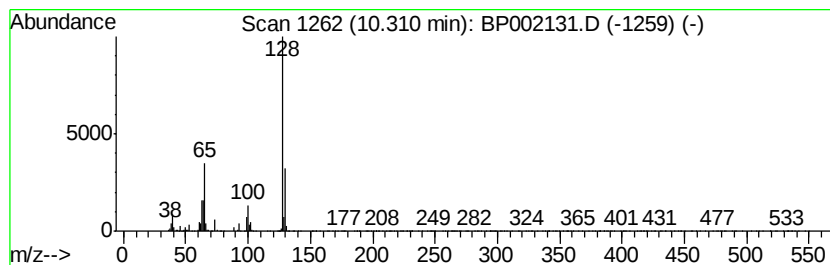
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Phenol, 3-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.31	2.84 ng/ul	317996	Naphthalene-d8	10.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3-chloro-		128	C6H5ClO	000108-43-0	94
2	Phenol, 3-chloro-		128	C6H5ClO	000108-43-0	94
3	Parachlorophenol		128	C6H5ClO	000106-48-9	94
4	Parachlorophenol		128	C6H5ClO	000106-48-9	93
5	Phenol, 3-chloro-		128	C6H5ClO	000108-43-0	91



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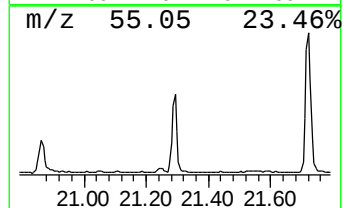
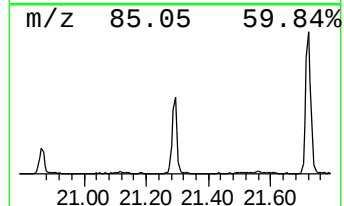
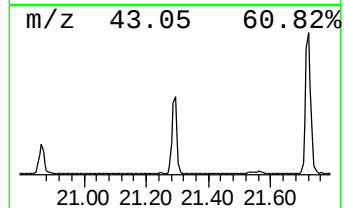
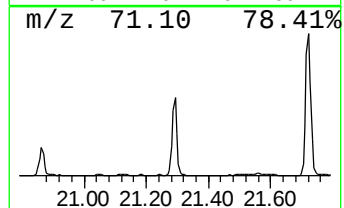
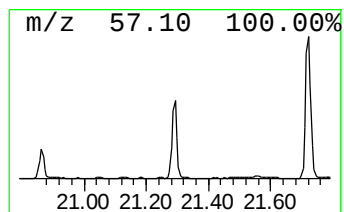
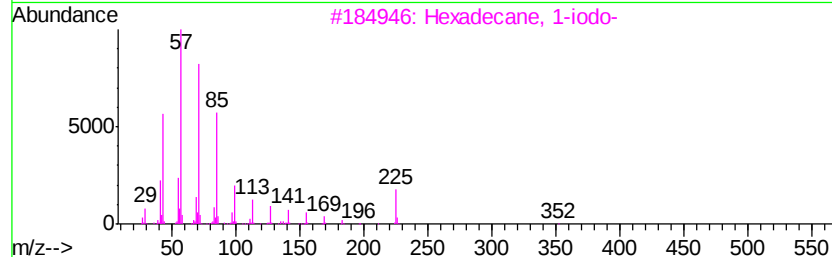
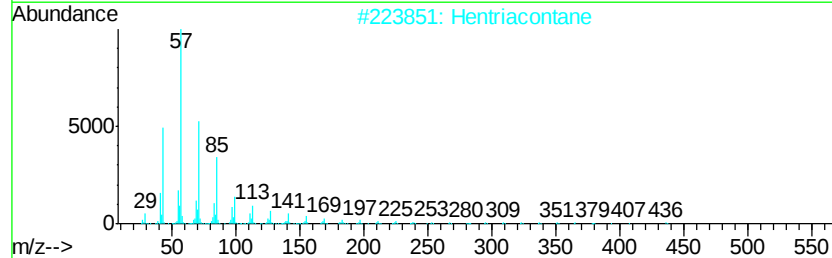
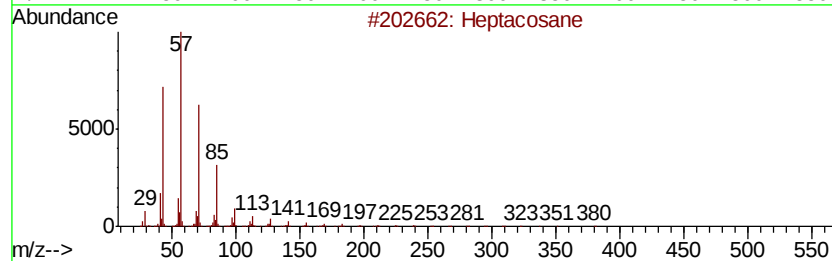
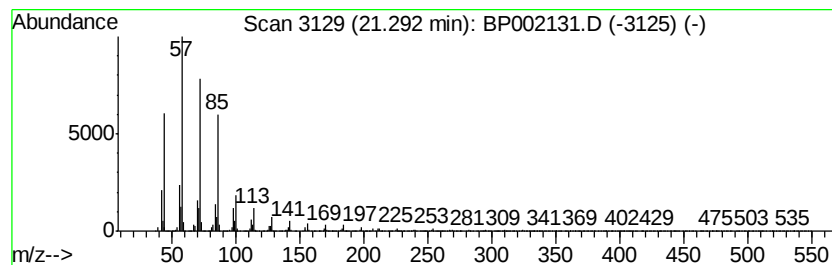
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.29	2.45 ng/ul	560215	Chrysene-d12	21.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	98
2		Hentriacontane	437	C31H64	000630-04-6	97
3		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	95
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
5		Triacontane	422	C30H62	000638-68-6	94



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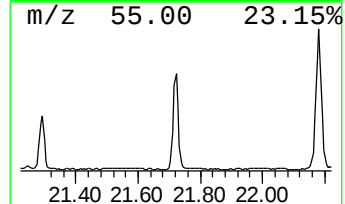
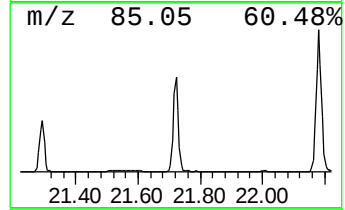
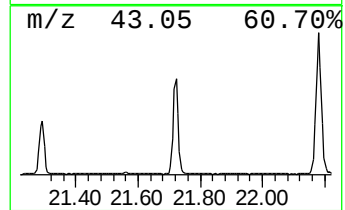
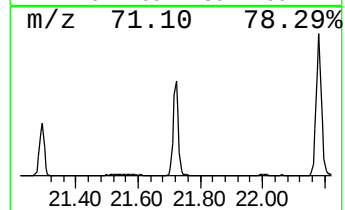
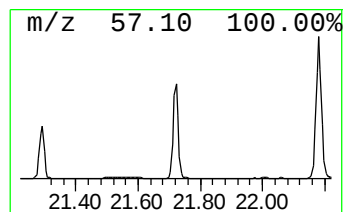
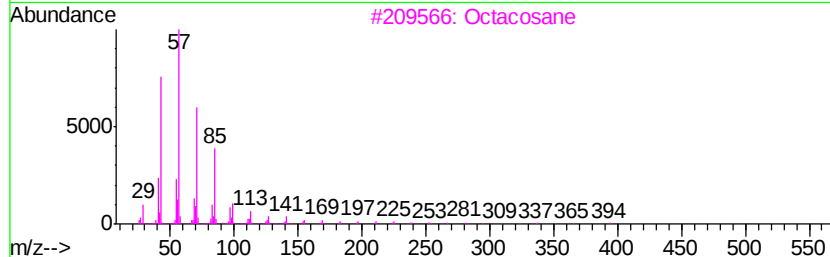
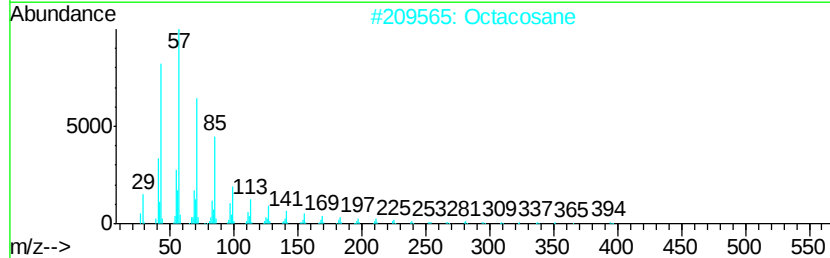
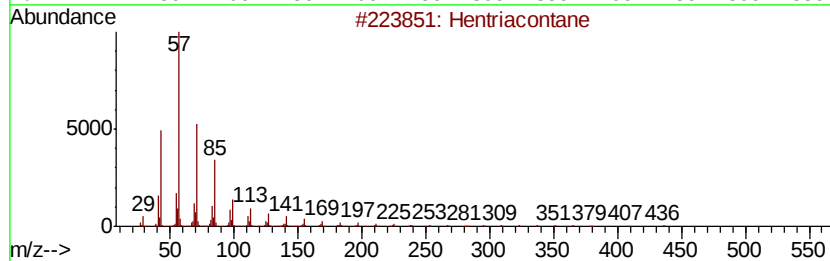
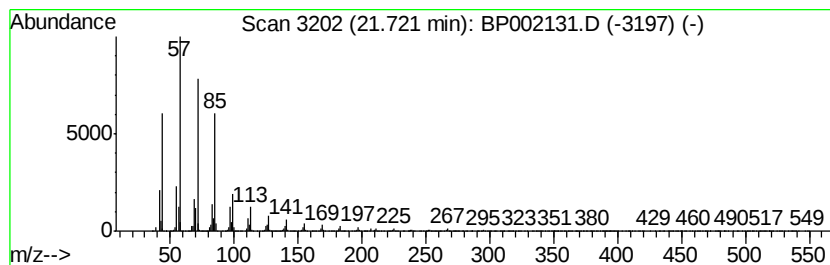
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.72	5.00 ng/ul	1144260	Chrysene-d12	21.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hentriacontane	437	C31H64	000630-04-6	97
2			Octacosane	394	C28H58	000630-02-4	97
3			Octacosane	394	C28H58	000630-02-4	95
4			2-Bromo dodecane	248	C12H25Br	013187-99-0	95
5			Tetratriacontane	479	C34H70	014167-59-0	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002131.D
Acq On : 09 Mar 2020 21:12
Operator : CG/JU
Sample : L1834-14
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0DK7

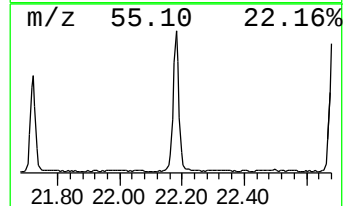
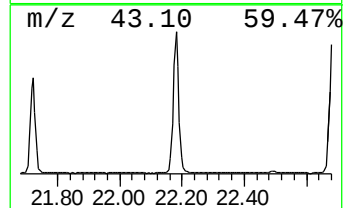
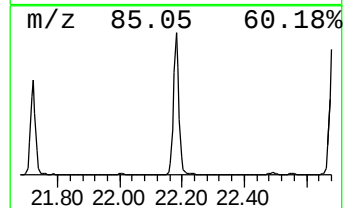
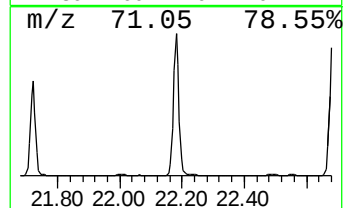
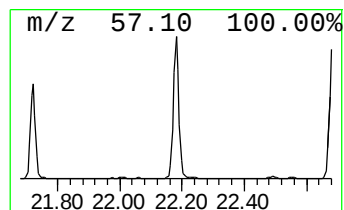
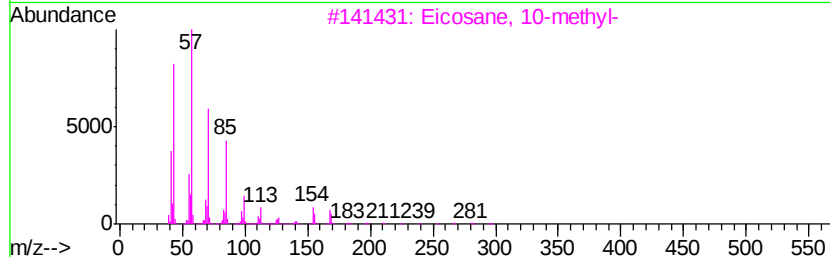
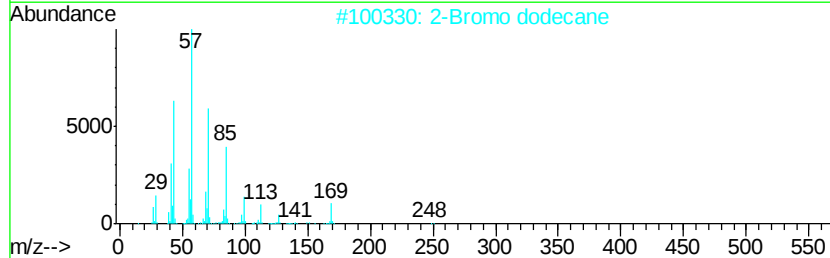
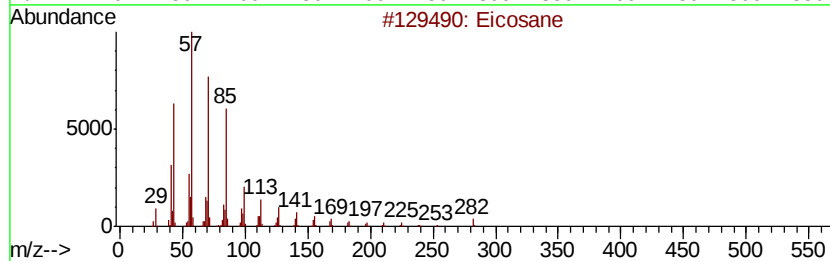
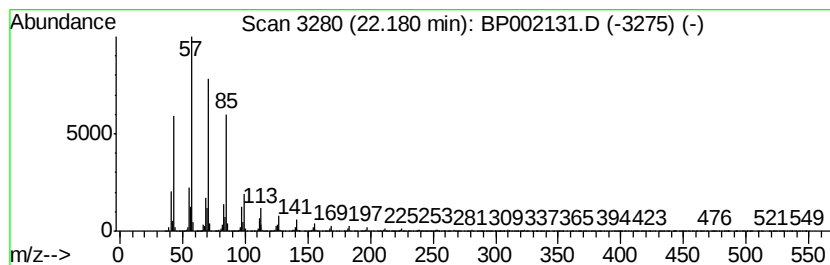
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.18	8.17 ng/ul	1868030	Chrysene-d12	21.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		2-Bromo dodecane	248	C12H25Br	013187-99-0	94
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
5		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002131.D
Acq On : 09 Mar 2020 21:12
Operator : CG/JU
Sample : L1834-14
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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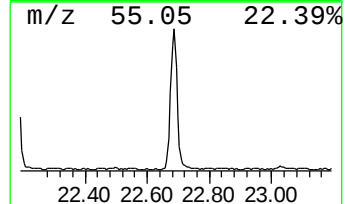
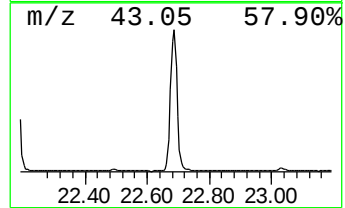
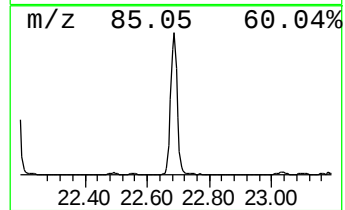
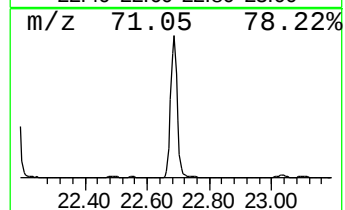
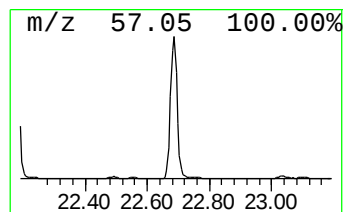
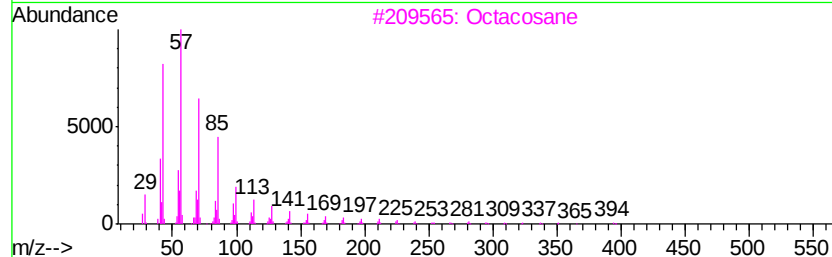
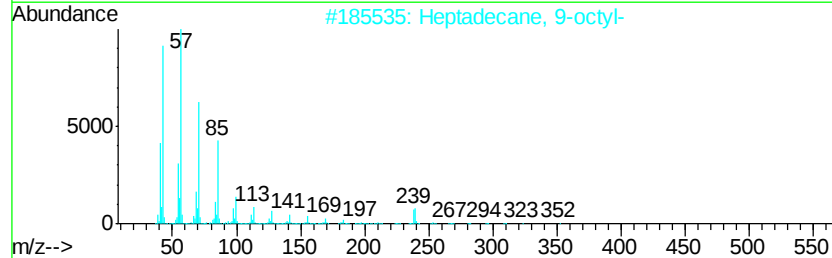
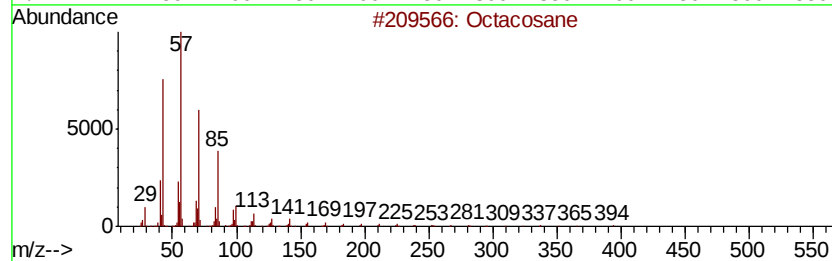
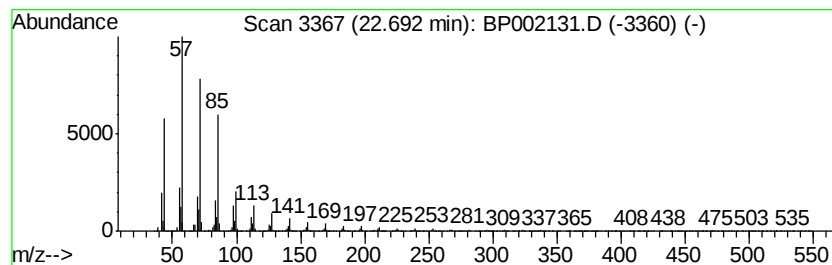
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.69	9.10 ng/ul	2232390	Perylene-d12	23.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	99
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
3		Octacosane	394	C28H58	000630-02-4	94
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Tetratriacontane	479	C34H70	014167-59-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002131.D
Acq On : 09 Mar 2020 21:12
Operator : CG/JU
Sample : L1834-14
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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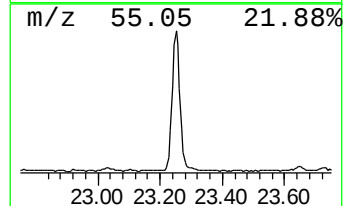
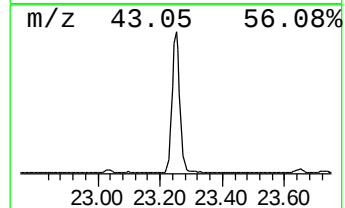
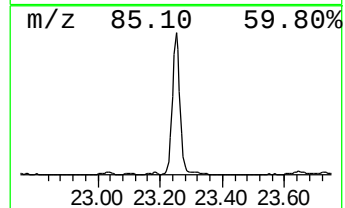
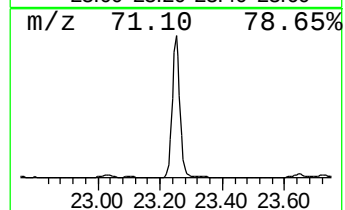
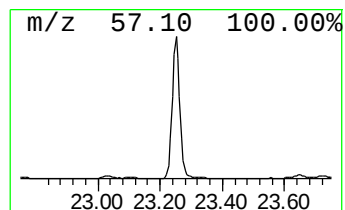
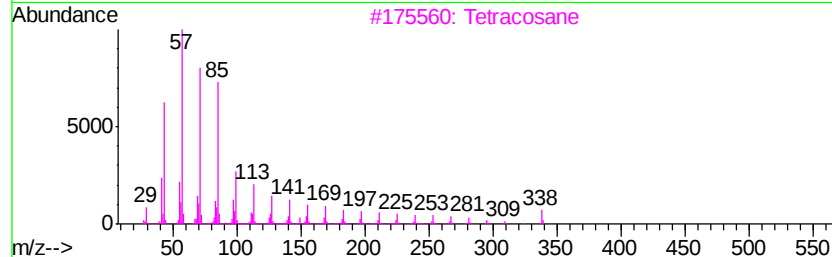
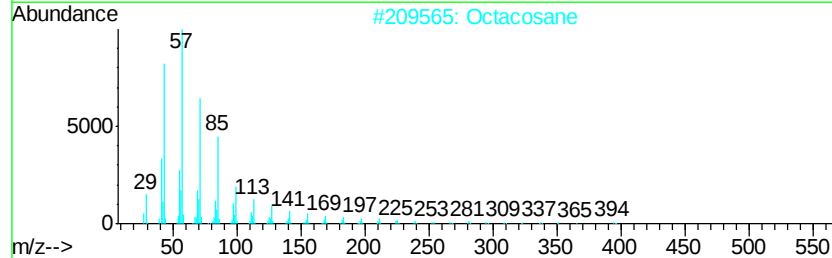
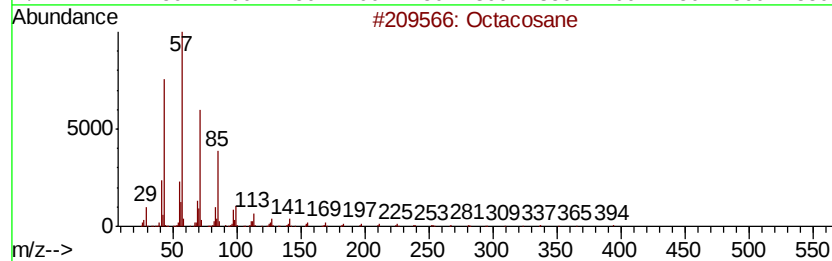
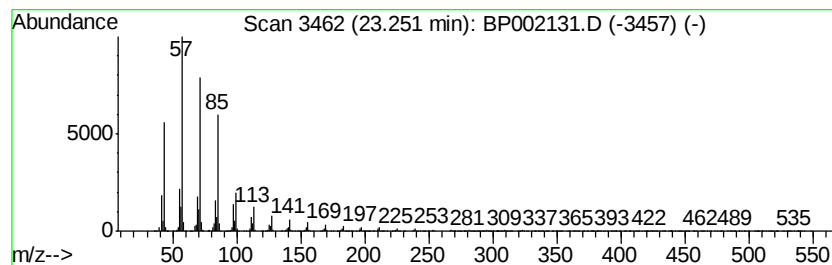
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.25	9.67 ng/ul	2373770	Perylene-d12	23.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	97
2			Octacosane	394	C28H58	000630-02-4	96
3			Tetracosane	338	C24H50	000646-31-1	95
4			Hexatriacontane	507	C36H74	000630-06-8	93
5			Eicosane, 10-methyl-	296	C21H44	054833-23-7	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002131.D
Acq On : 09 Mar 2020 21:12
Operator : CG/JU
Sample : L1834-14
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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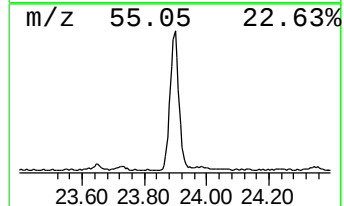
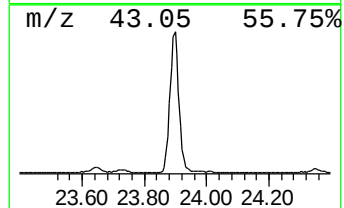
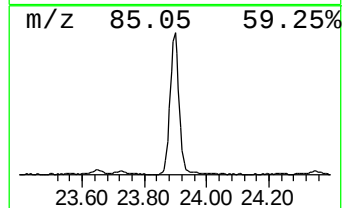
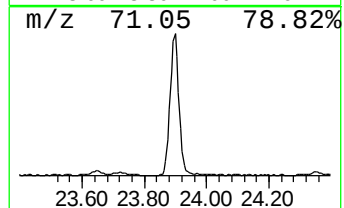
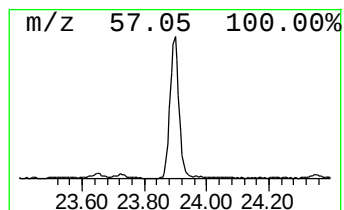
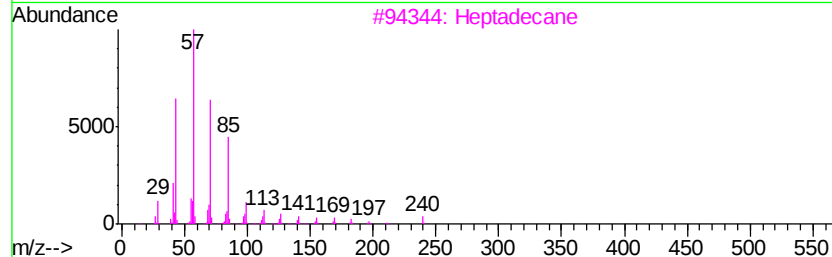
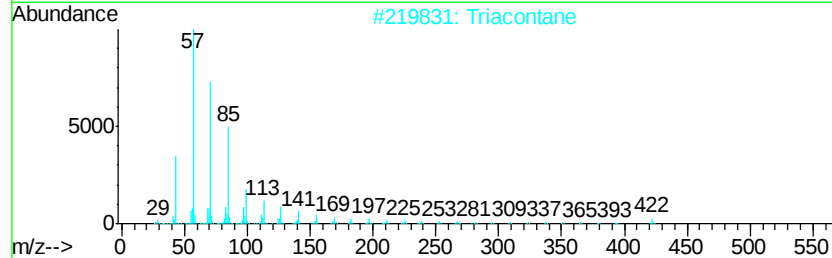
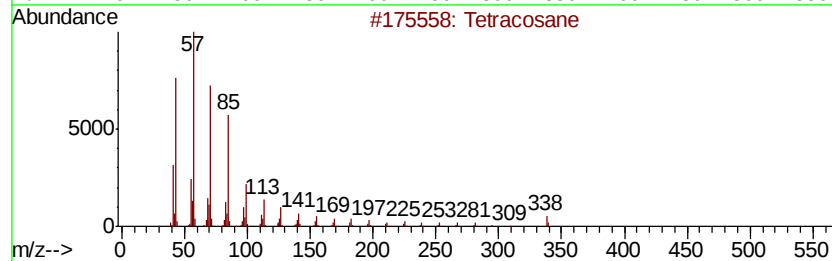
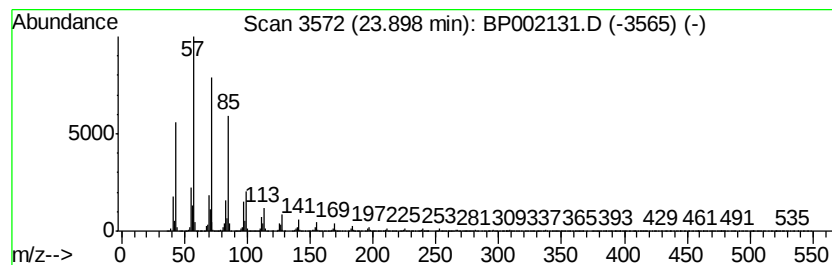
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.90	7.64 ng/ul	1875510	Perylene-d12	23.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	98
2		triacontane	422	C30H62	000638-68-6	94
3		heptadecane	240	C17H36	000629-78-7	94
4		hexatriacontane	507	C36H74	000630-06-8	93
5		heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002131.D
Acq On : 09 Mar 2020 21:12
Operator : CG/JU
Sample : L1834-14
Misc :
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ClientSampleId :
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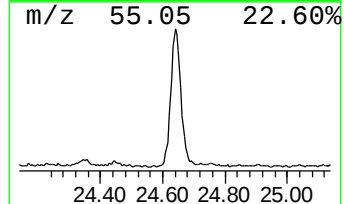
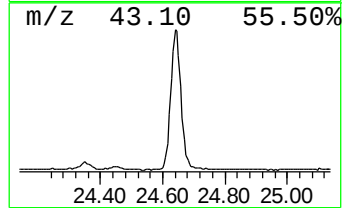
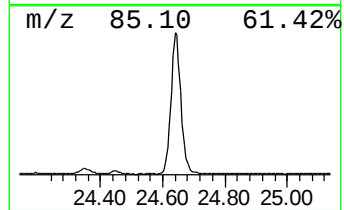
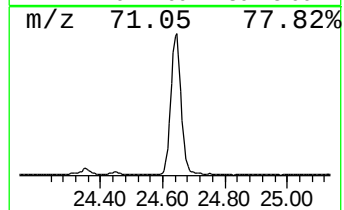
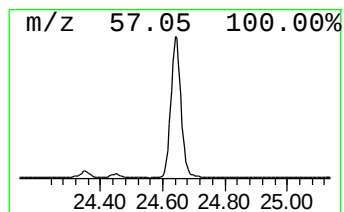
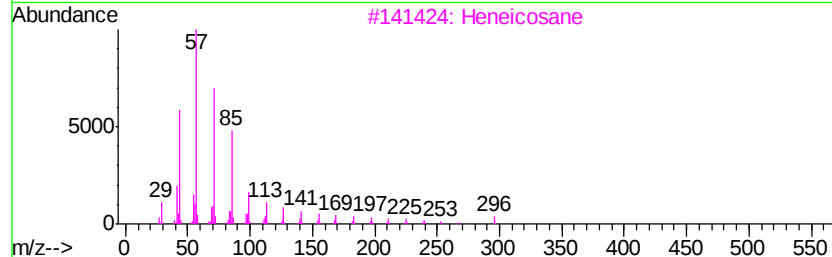
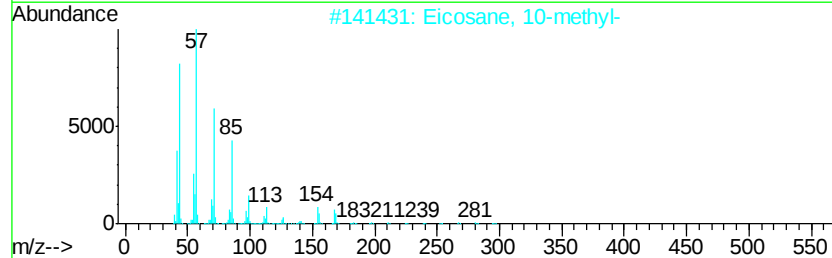
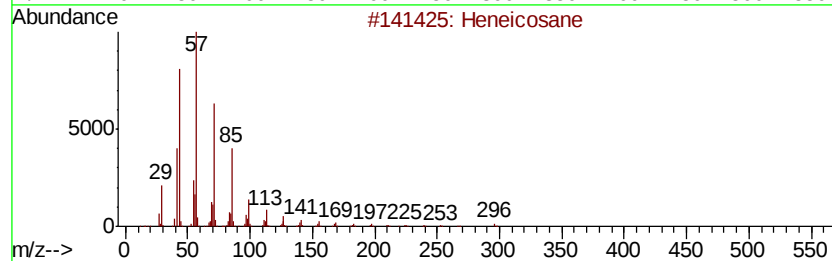
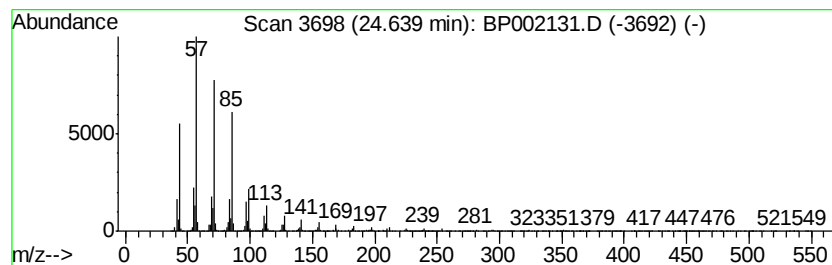
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.64	4.63 ng/ul	1135400	Perylene-d12	23.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	97
2			Eicosane, 10-methyl-	296	C21H44	054833-23-7	97
3			Heneicosane	296	C21H44	000629-94-7	97
4			Heneicosane	296	C21H44	000629-94-7	96
5			Docosane, 7-hexyl-	394	C28H58	055373-86-9	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030920\
Data File : BP002131.D
Acq On : 09 Mar 2020 21:12
Operator : CG/JU
Sample : L1834-14
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0DK7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Phenol, 3-chloro-	10.31	2.8	ng/ul	317996	2	10.40	2237830	20.0
(DEL) Alkane: Str...	21.29	2.5	ng/ul	560215	5	21.12	4572600	20.0
(DEL) Alkane: Str...	21.72	5.0	ng/ul	1144260	5	21.12	4572600	20.0
(DEL) Alkane: Str...	22.18	8.2	ng/ul	1868030	5	21.12	4572600	20.0
(DEL) Alkane: Str...	22.69	9.1	ng/ul	2232390	6	23.32	4907900	20.0
(DEL) Alkane: Str...	23.25	9.7	ng/ul	2373770	6	23.32	4907900	20.0
(DEL) Alkane: Str...	23.90	7.6	ng/ul	1875510	6	23.32	4907900	20.0
(DEL) Alkane: Str...	24.64	4.6	ng/ul	1135400	6	23.32	4907900	20.0