

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
 Data File : BN015054.D
 Acq On : 24 Jun 2021 23:51
 Operator : CG/JU
 Sample : M2682-14
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGD33

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_N\METHODS\SFAM-EPA-BN061821MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN015054.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.287	30	34	38	rVB	17938	21335	0.71%	0.060%
2	3.681	98	101	113	rBV	151500	255581	8.52%	0.718%
3	3.781	114	118	126	rVV	57315	82436	2.75%	0.231%
4	4.005	150	156	164	rVB	30201	46298	1.54%	0.130%
5	4.534	242	246	256	rVB	27958	40006	1.33%	0.112%
6	5.411	390	395	407	rBV	40393	77882	2.60%	0.219%
7	5.775	451	457	463	rBV2	19895	31066	1.04%	0.087%
8	6.893	640	647	655	rVV	302248	469295	15.64%	1.318%
9	7.064	670	676	684	rVB	229042	355696	11.86%	0.999%
10	7.264	703	710	718	rBV	300360	481855	16.06%	1.353%
11	7.728	782	789	796	rBV	550854	865221	28.84%	2.429%
12	8.416	900	906	915	rBV	297299	469714	15.66%	1.319%
13	8.869	977	983	994	rVB	263340	418544	13.95%	1.175%
14	9.587	1097	1105	1111	rBV	179805	288072	9.60%	0.809%
15	10.116	1189	1195	1204	rBV	337177	542968	18.10%	1.524%
16	10.275	1216	1222	1230	rBV	135752	216637	7.22%	0.608%
17	10.499	1253	1260	1266	rBV	748977	1210898	40.37%	3.400%
18	10.546	1266	1268	1274	rVV	19082	25920	0.86%	0.073%
19	10.628	1276	1282	1294	rVV	267627	457892	15.26%	1.286%
20	13.763	1809	1815	1821	rVB	579888	791538	26.39%	2.222%
21	14.040	1856	1862	1874	rVV	677268	1033323	34.45%	2.901%
22	14.351	1908	1915	1921	rBV	1068928	1540362	51.35%	4.325%
23	14.416	1921	1926	1933	rVB	58686	85600	2.85%	0.240%
24	14.540	1939	1947	1956	rBV	193965	279566	9.32%	0.785%
25	14.751	1977	1983	1991	rVB	41010	69848	2.33%	0.196%
26	15.345	2078	2084	2090	rBV	901305	1262838	42.10%	3.545%
27	15.404	2090	2094	2098	rVV2	64571	93345	3.11%	0.262%
28	15.457	2098	2103	2108	rVV	110177	155221	5.17%	0.436%
29	15.698	2140	2144	2150	rBV	18406	25092	0.84%	0.070%
30	15.775	2152	2157	2163	rBV	55175	74822	2.49%	0.210%
31	15.845	2164	2169	2177	rVB	19862	39208	1.31%	0.110%
32	16.081	2204	2209	2214	rVV3	23005	33343	1.11%	0.094%
33	16.751	2318	2323	2331	rVB3	28616	49060	1.64%	0.138%
34	16.839	2331	2338	2339	rBV4	15132	25366	0.85%	0.071%
35	16.916	2347	2351	2357	rVB	45090	60715	2.02%	0.170%
36	17.098	2375	2382	2386	rVV	1285229	1842033	61.41%	5.172%

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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_N\METHODS\SFAM-EPA-BN061821MA.M
 Title : SVOA CALIBRATION

37	17.139	2386	2389	2394	rVV	847995	1158469	38.62%	3.252%
38	17.198	2394	2399	2403	rVV	951377	1412403	47.08%	3.965%
39	17.228	2403	2404	2411	rVV	205531	213263	7.11%	0.599%
40	17.287	2411	2414	2421	rVV4	18559	32218	1.07%	0.090%
41	17.498	2441	2450	2454	rBV2	60447	110350	3.68%	0.310%
42	17.622	2467	2471	2476	rBV3	25266	43190	1.44%	0.121%
43	17.681	2477	2481	2488	rVB2	26183	41147	1.37%	0.116%
44	17.892	2514	2517	2522	rVV5	14767	23884	0.80%	0.067%
45	17.975	2526	2531	2534	rVV	105252	155200	5.17%	0.436%
46	18.022	2534	2539	2545	rVV	161589	274287	9.14%	0.770%
47	18.104	2549	2553	2556	rVV	53931	74128	2.47%	0.208%
48	18.169	2558	2564	2568	rVV3	140212	255319	8.51%	0.717%
49	18.210	2568	2571	2575	rVB	67403	89263	2.98%	0.251%
50	18.316	2585	2589	2595	rVB7	19809	34416	1.15%	0.097%
51	18.481	2612	2617	2620	rBV	80161	107596	3.59%	0.302%
52	18.516	2620	2623	2628	rVB2	42789	58604	1.95%	0.165%
53	18.728	2655	2659	2663	rVV2	29586	37056	1.24%	0.104%
54	18.775	2663	2667	2671	rVV2	30994	41695	1.39%	0.117%
55	18.816	2671	2674	2680	rVB	27231	36152	1.21%	0.101%
56	18.898	2683	2688	2693	rBV2	63252	105721	3.52%	0.297%
57	18.951	2693	2697	2702	rVV4	48192	112730	3.76%	0.316%
58	18.992	2702	2704	2707	rVV2	25892	27414	0.91%	0.077%
59	19.033	2707	2711	2719	rVB2	53667	93091	3.10%	0.261%
60	19.163	2727	2733	2739	rVB	1224769	1633100	54.44%	4.585%
61	19.228	2741	2744	2748	rVV	18930	26059	0.87%	0.073%
62	19.310	2753	2758	2762	rVB	30256	51646	1.72%	0.145%
63	19.357	2762	2766	2769	rBV5	12680	23463	0.78%	0.066%
64	19.404	2772	2774	2778	rVB	22307	23376	0.78%	0.066%
65	19.498	2784	2790	2792	rBV	1098613	1547623	51.59%	4.345%
66	19.528	2792	2795	2803	rVV	741108	1028391	34.28%	2.887%
67	19.598	2803	2807	2814	rVB3	51868	88083	2.94%	0.247%
68	19.710	2821	2826	2830	rBV	59509	87574	2.92%	0.246%
69	19.763	2832	2835	2842	rVB3	31673	50574	1.69%	0.142%
70	19.851	2844	2850	2851	rBV2	75050	105896	3.53%	0.297%
71	19.986	2869	2873	2876	rBV3	55810	99925	3.33%	0.281%
72	20.022	2876	2879	2883	rVB2	110993	155856	5.20%	0.438%
73	20.063	2883	2886	2889	rBV2	28705	31909	1.06%	0.090%
74	20.128	2892	2897	2900	rBV2	55635	77690	2.59%	0.218%
75	20.180	2900	2906	2911	rVV2	99488	162535	5.42%	0.456%
76	20.322	2926	2930	2934	rBV	43541	51384	1.71%	0.144%

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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_N\METHODS\SFAM-EPA-BN061821MA.M
 Title : SVOA CALIBRATION

77	20.369	2934	2938	2942	rBV	59190	86163	2.87%	0.242%
78	20.516	2959	2963	2968	rVB	237673	252390	8.41%	0.709%
79	20.563	2969	2971	2976	rBV4	14413	23021	0.77%	0.065%
80	20.769	3003	3006	3014	rVB	96873	133029	4.43%	0.373%
81	20.939	3029	3035	3039	rBV2	83788	177958	5.93%	0.500%
82	20.980	3039	3042	3045	rVV	95549	119803	3.99%	0.336%
83	21.027	3045	3050	3053	rVV3	228302	343606	11.45%	0.965%
84	21.069	3053	3057	3064	rVB2	109349	189138	6.31%	0.531%
85	21.175	3073	3075	3079	rVB2	31369	33510	1.12%	0.094%
86	21.222	3080	3083	3088	rBV2	78170	92060	3.07%	0.258%
87	21.292	3088	3095	3099	rVV2	1765553	2999804	100.00%	8.422%
88	21.327	3099	3101	3108	rVB	520938	686552	22.89%	1.927%
89	21.463	3118	3124	3127	rBV4	62741	126667	4.22%	0.356%
90	21.845	3187	3189	3193	rVB	98198	106242	3.54%	0.298%
91	21.904	3194	3199	3207	rBV3	212511	397085	13.24%	1.115%
92	22.069	3220	3227	3235	rVB3	129974	283563	9.45%	0.796%
93	22.498	3297	3300	3306	rVB3	52829	74371	2.48%	0.209%
94	22.880	3357	3365	3380	rBV2	610767	1855653	61.86%	5.210%
95	23.039	3389	3392	3398	rVB	79292	129374	4.31%	0.363%
96	23.398	3447	3453	3457	rBV2	396351	741797	24.73%	2.083%
97	23.445	3457	3461	3469	rVB2	245913	479922	16.00%	1.347%
98	23.539	3470	3477	3484	rBV2	1096218	2065651	68.86%	5.799%
99	24.104	3568	3573	3581	rVB	206073	403916	13.46%	1.134%
100	24.857	3696	3701	3708	rVB2	103827	216291	7.21%	0.607%

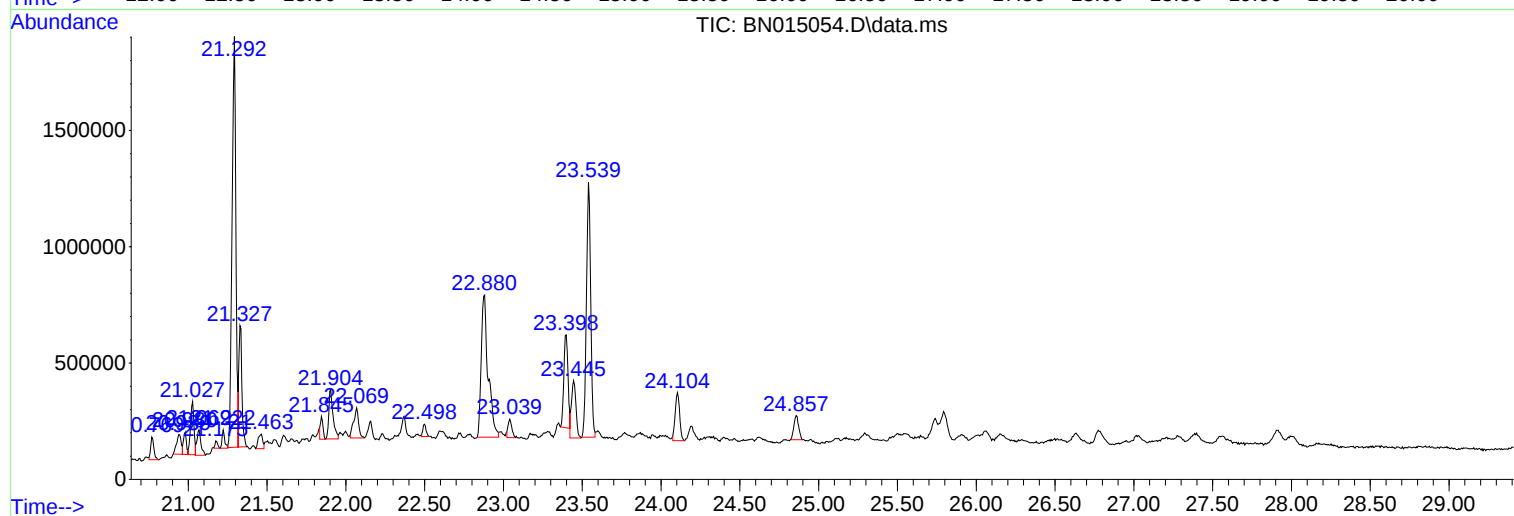
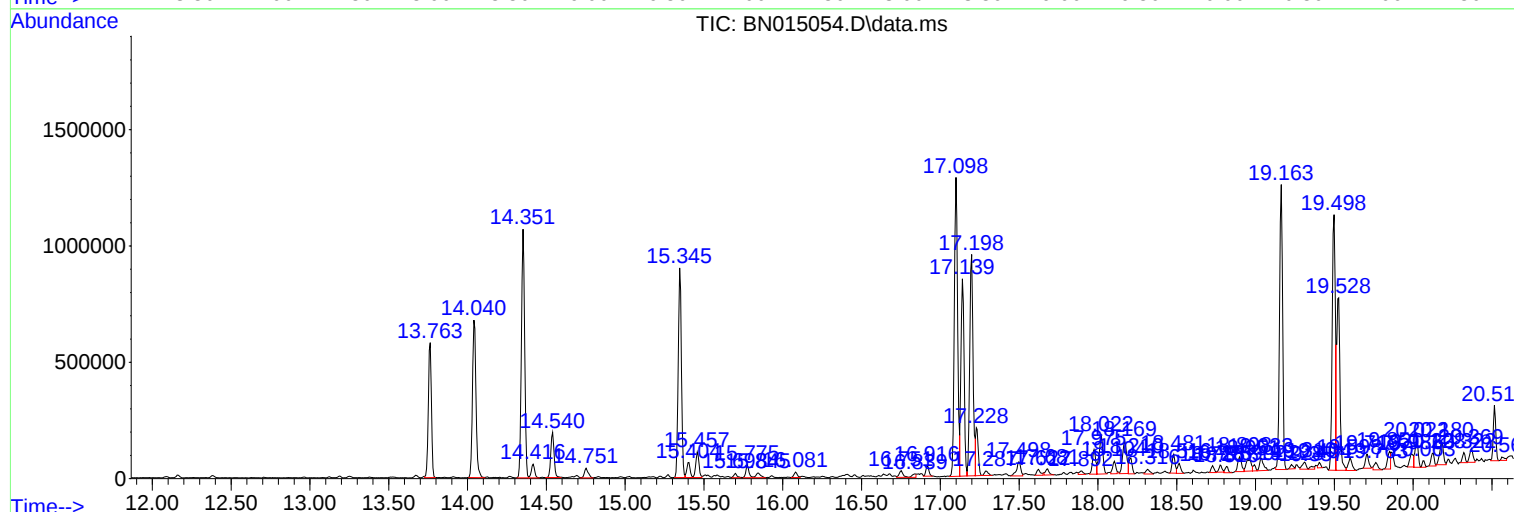
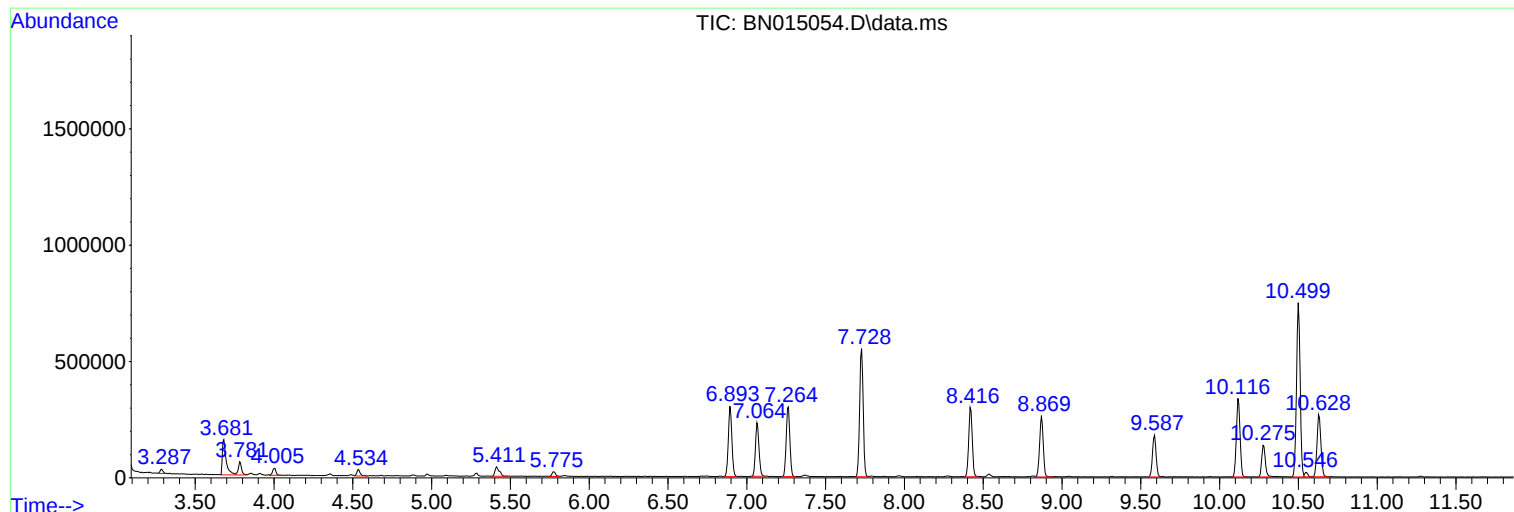
Sum of corrected areas: 35618842

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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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



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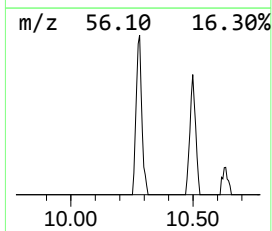
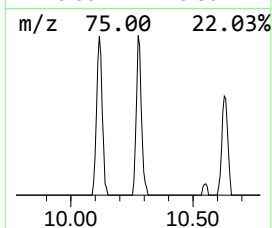
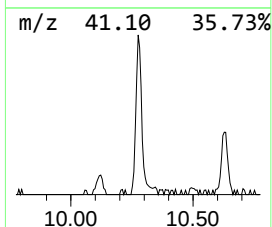
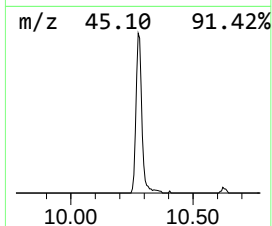
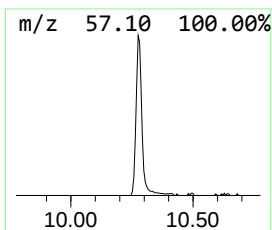
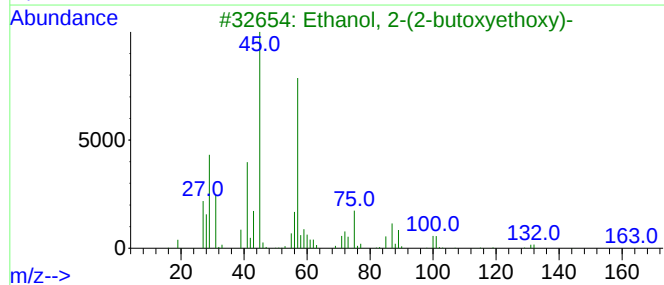
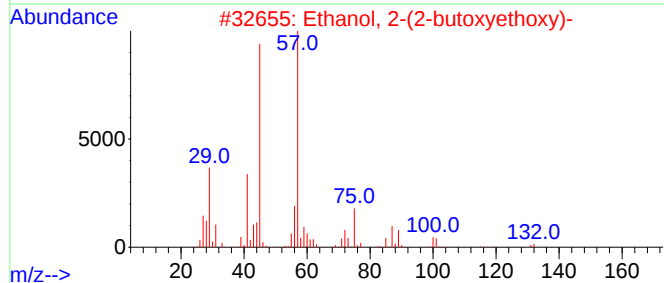
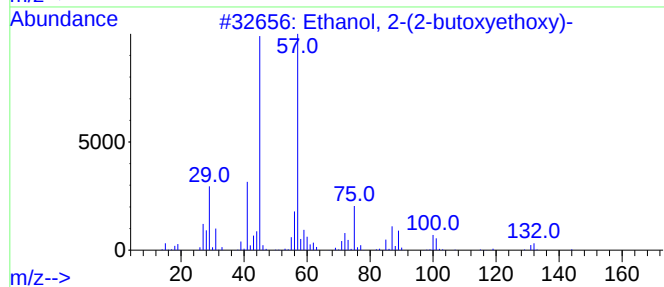
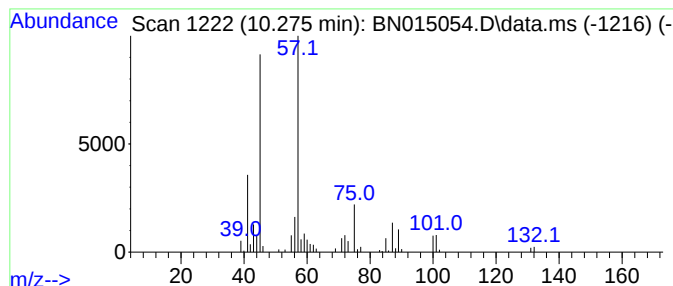
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.275	3.58 ng/ul	216637	Naphthalene-d8	10.499

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	64
5			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64



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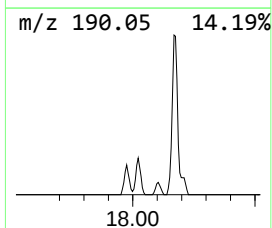
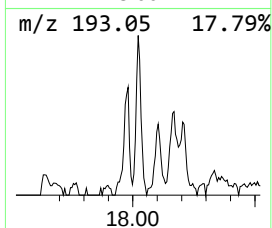
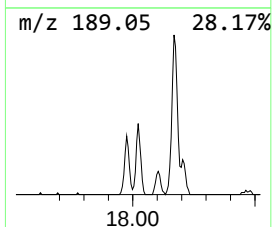
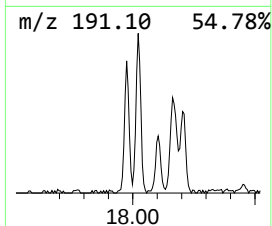
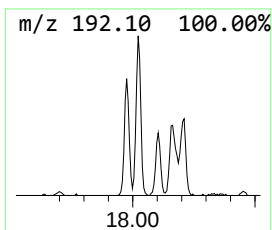
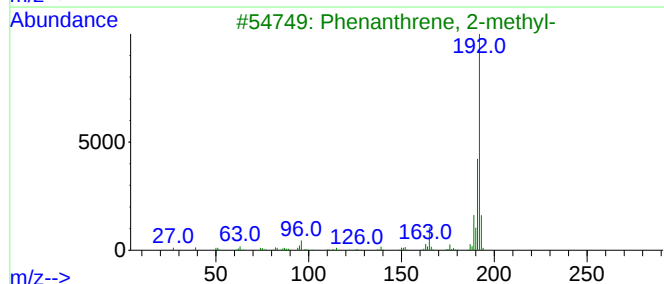
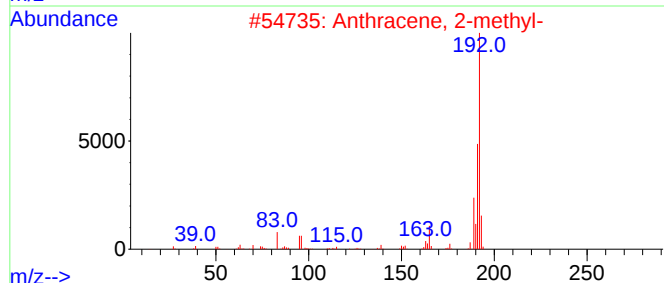
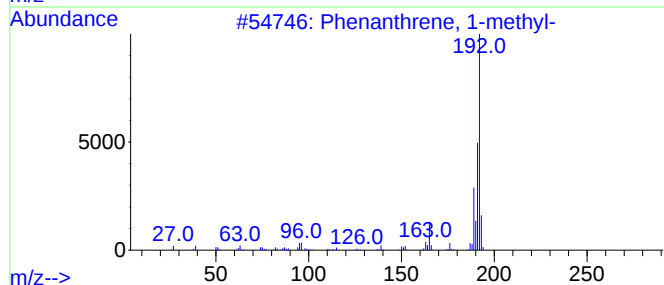
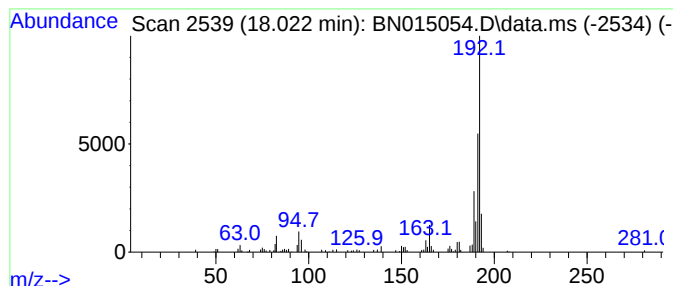
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenanthrene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.022	2.98 ng/ul	274287	Phenanthrene-d10	17.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95



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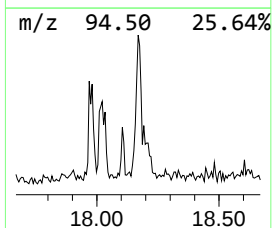
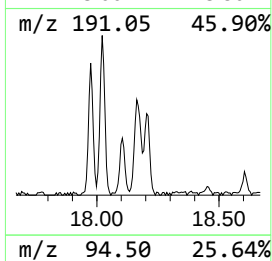
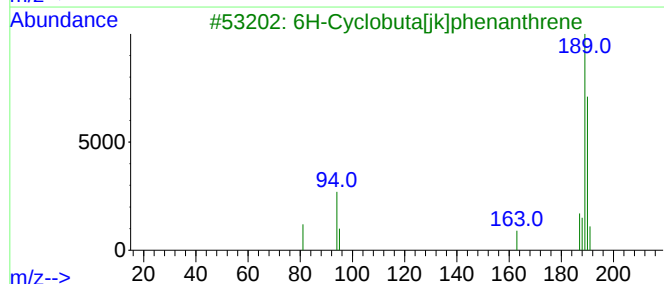
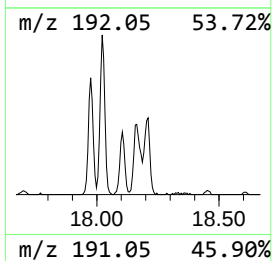
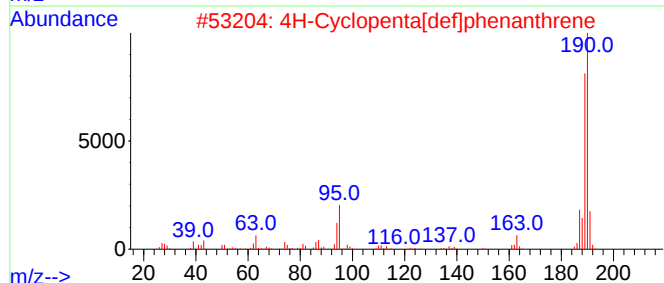
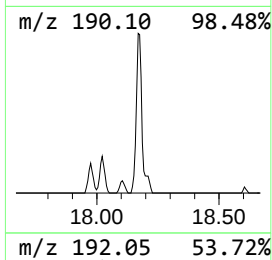
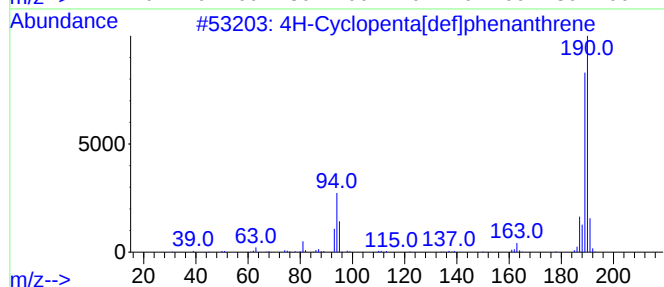
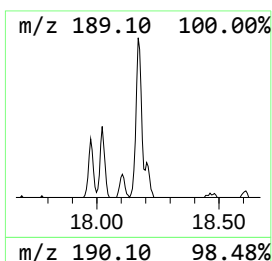
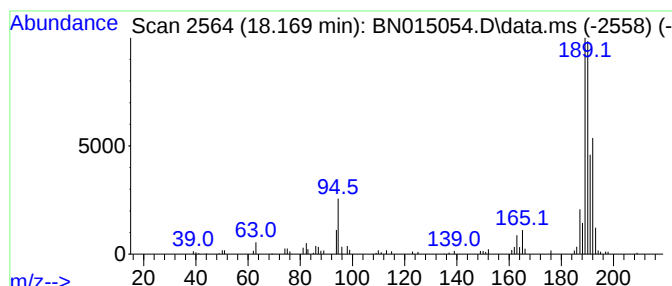
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.169	2.77 ng/ul	255319	Phenanthrene-d10	17.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
3		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	58
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	38
5		Methyl diselenide	190	C2H6Se2	007101-31-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015054.D
Acq On : 24 Jun 2021 23:51
Operator : CG/JU
Sample : M2682-14
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGD33

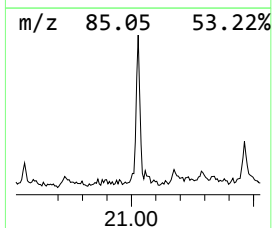
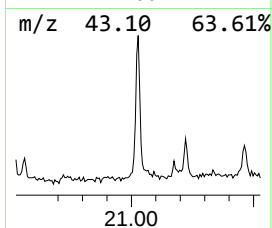
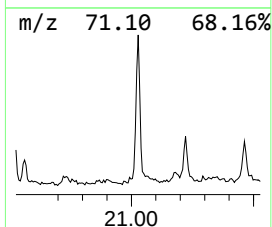
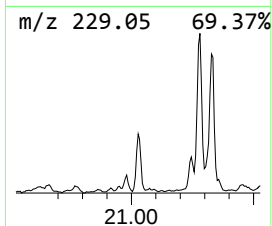
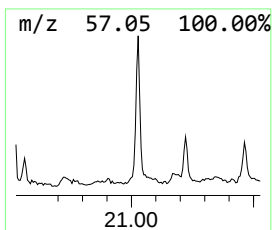
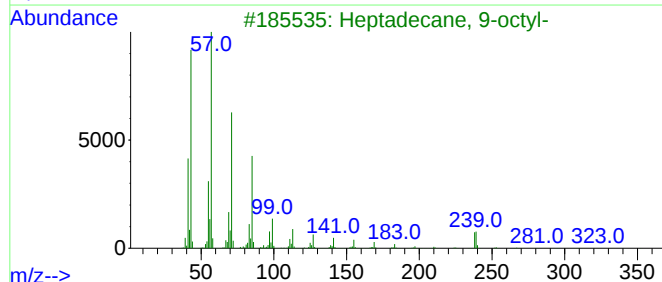
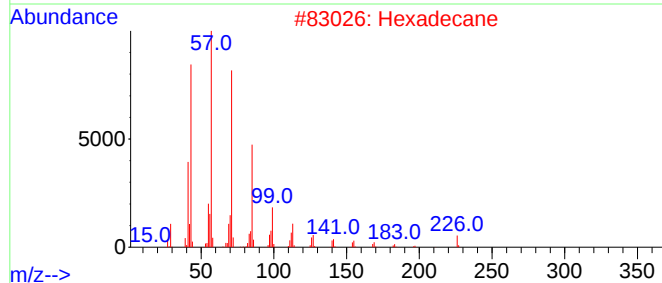
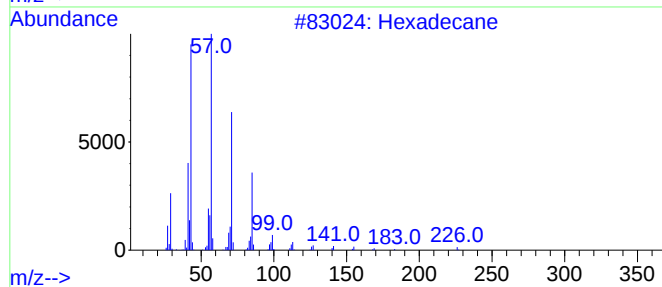
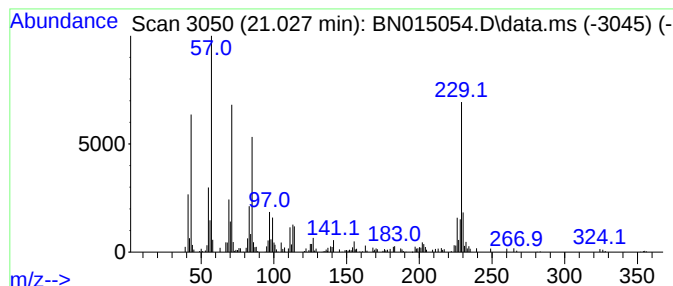
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.028	2.29 ng/ul	343606	Chrysene-d12	21.292

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	90
2		Hexadecane	226	C16H34	000544-76-3	70
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	64
4		Tritetracontane	605	C43H88	007098-21-7	55
5		Hexadecane	226	C16H34	000544-76-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015054.D
Acq On : 24 Jun 2021 23:51
Operator : CG/JU
Sample : M2682-14
Misc :
ALS Vial : 18 Sample Multiplier: 1

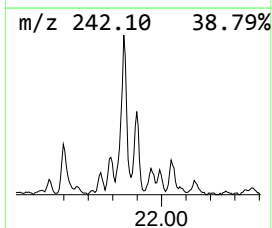
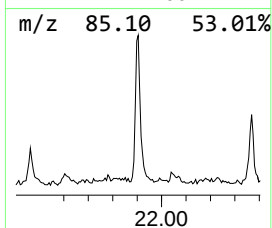
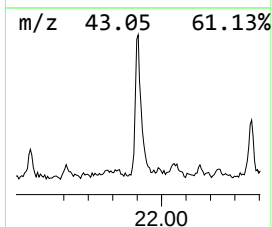
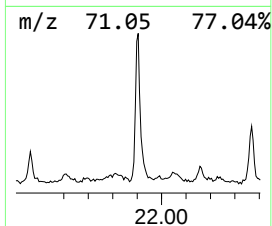
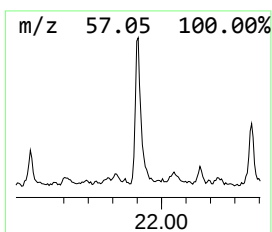
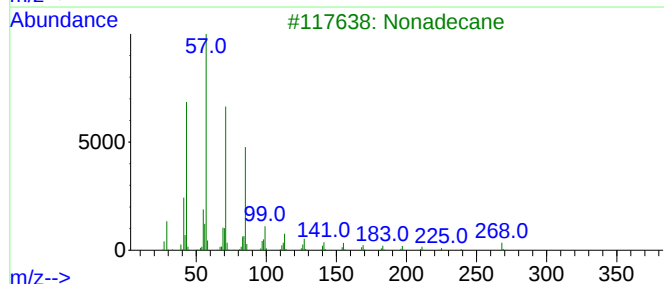
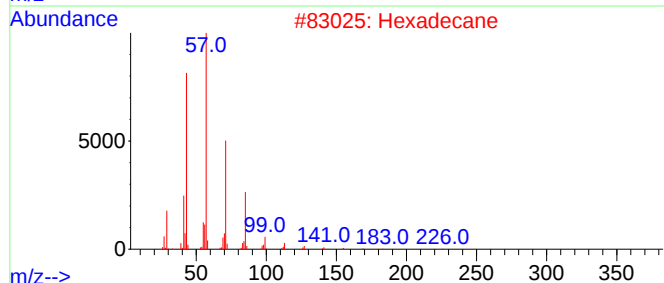
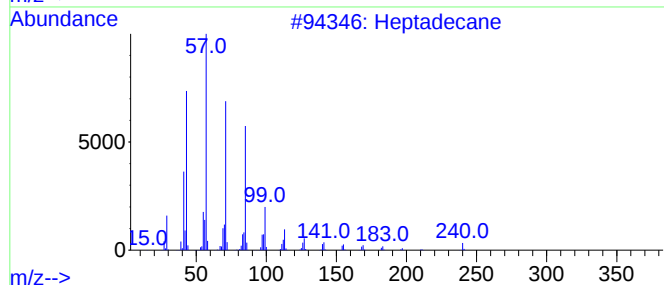
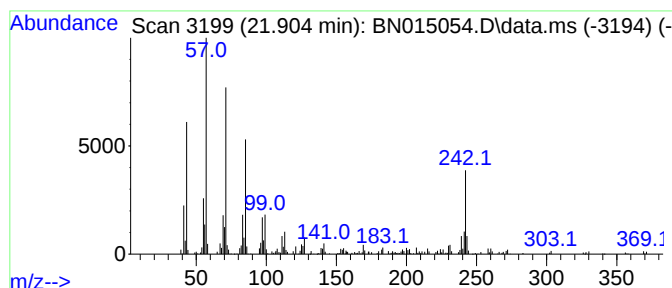
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.904	2.65 ng/ul	397085	Chrysene-d12	21.292	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	95
2	Hexadecane	226	C16H34	000544-76-3	95
3	Nonadecane	268	C19H40	000629-92-5	93
4	Hexadecane	226	C16H34	000544-76-3	93
5	Hexadecane	226	C16H34	000544-76-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015054.D
Acq On : 24 Jun 2021 23:51
Operator : CG/JU
Sample : M2682-14
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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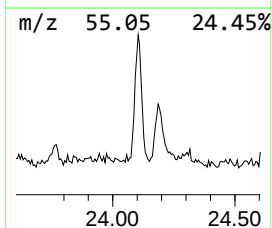
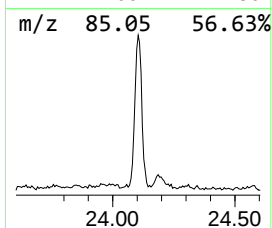
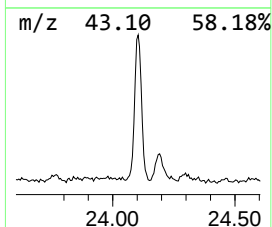
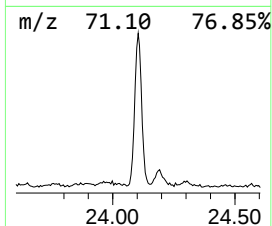
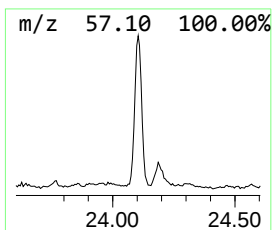
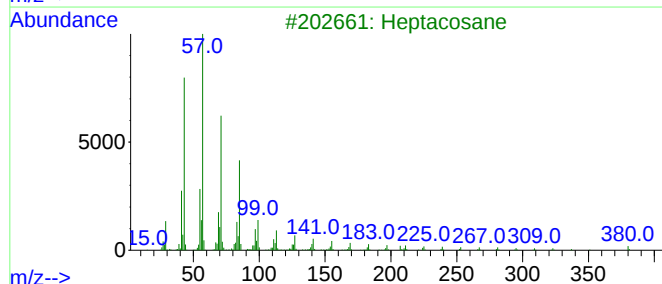
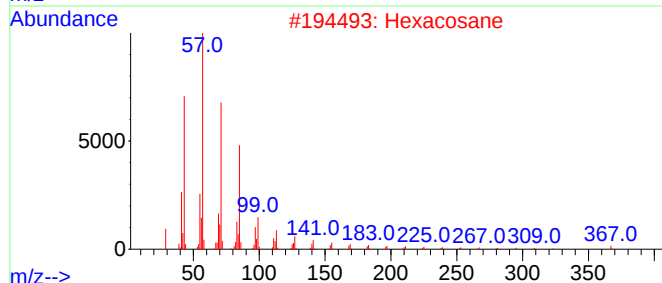
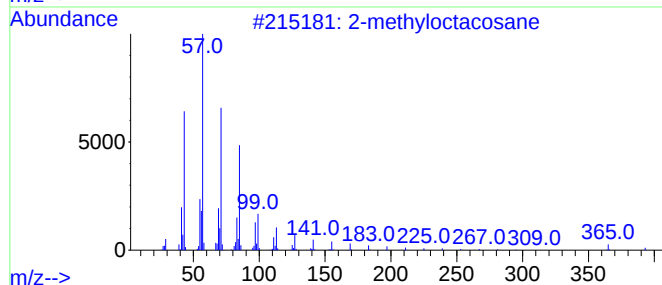
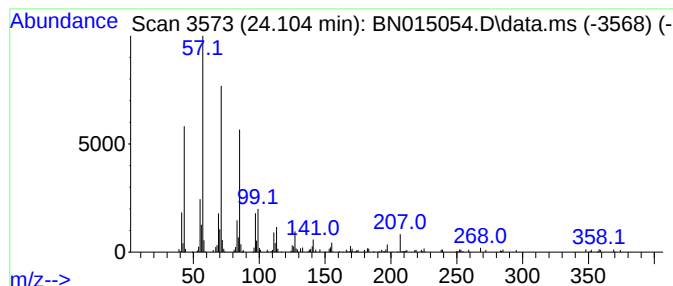
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.104	3.91 ng/ul	403916	Perylene-d12	23.539

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Hexacosane	366	C26H54	000630-01-3	90
3		Heptacosane	380	C27H56	000593-49-7	90
4		Heneicosane	296	C21H44	000629-94-7	87
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015054.D
Acq On : 24 Jun 2021 23:51
Operator : CG/JU
Sample : M2682-14
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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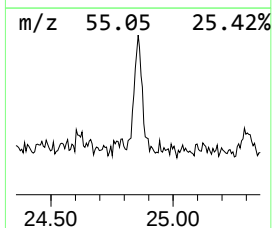
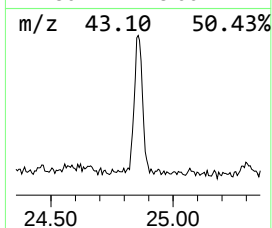
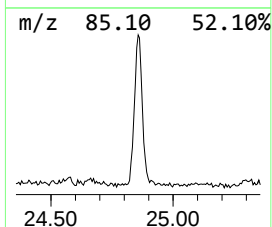
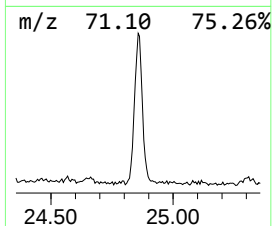
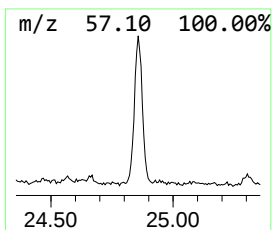
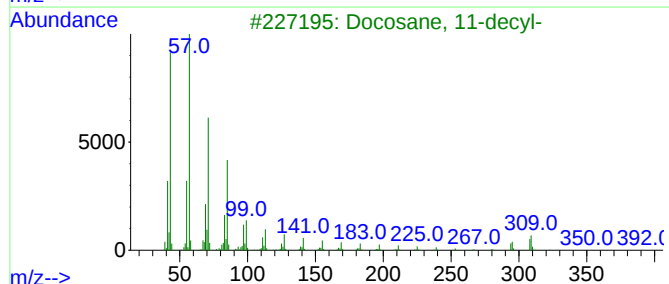
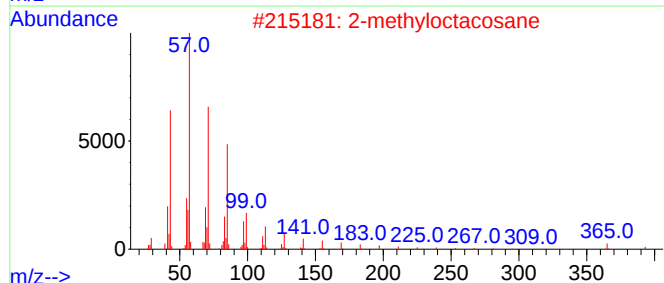
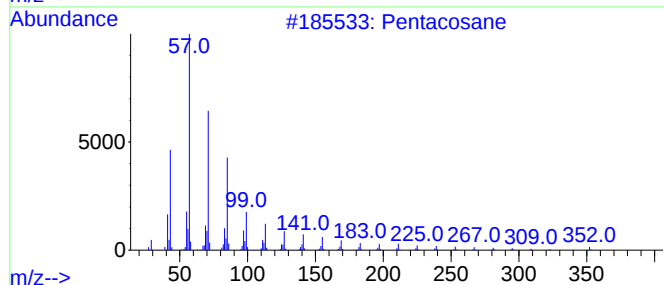
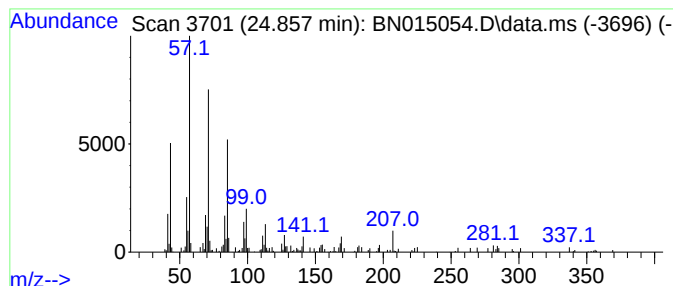
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.857	2.09 ng/ul	216291	Perylene-d12	23.539

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentacosane	352	C25H52	000629-99-2	90
2		2-methyloctacosane	408	C29H60	1000376-72-8	87
3		Docosane, 11-decyl-	451	C32H66	055401-55-3	87
4		Hentriacontane	437	C31H64	000630-04-6	87
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062521\
Data File : BN015054.D
Acq On : 24 Jun 2021 23:51
Operator : CG/JU
Sample : M2682-14
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGD33

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.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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Phenanthrene, 1...	18.022	3.0	ng/u1	274287	4	17.098	1842030	20.0
4H-Cyclopenta[d...	18.169	2.8	ng/u1	255319	4	17.098	1842030	20.0
(DEL) Alkane: S...	21.028	2.3	ng/u1	343606	5	21.292	2999800	20.0
(DEL) Alkane: S...	21.904	2.6	ng/u1	397085	5	21.292	2999800	20.0
(DEL) Alkane: S...	24.104	3.9	ng/u1	403916	6	23.539	2065650	20.0
(DEL) Alkane: S...	24.857	2.1	ng/u1	216291	6	23.539	2065650	20.0