

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
 Data File : BN014872.D
 Acq On : 10 Jun 2021 21:25
 Operator : CG/JU
 Sample : M2618-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BG5R1

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-BN052521.M
 Title : SVOA CALIBRATION

Signal : TIC: BN014872.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.846	123	129	132	rBB2	36126	56114	1.71%	0.122%
2	4.034	156	161	166	rBB6	17475	34937	1.07%	0.076%
3	4.828	289	296	303	rBB	59142	96044	2.93%	0.208%
4	6.869	635	643	653	rBB	478416	778292	23.73%	1.685%
5	7.034	664	671	682	rBV	323139	528315	16.11%	1.144%
6	7.228	698	704	715	rBB	515291	835742	25.48%	1.810%
7	7.693	778	783	790	rVB	684376	1093722	33.35%	2.368%
8	8.399	894	903	914	rBV	645127	1056102	32.20%	2.287%
9	8.510	918	922	928	rBB5	23888	40249	1.23%	0.087%
10	8.846	972	979	987	rBB	465340	759242	23.15%	1.644%
11	9.563	1093	1101	1112	rBB2	424561	716298	21.84%	1.551%
12	10.104	1185	1193	1206	rBB	656972	1128722	34.41%	2.444%
13	10.481	1250	1257	1270	rBV	886019	1490839	45.45%	3.228%
14	10.616	1272	1280	1290	rBB	546626	955891	29.14%	2.070%
15	13.757	1808	1814	1825	rBV	1154065	1656233	50.50%	3.586%
16	14.034	1852	1861	1869	rBB	1246900	1887054	57.53%	4.086%
17	14.181	1881	1886	1893	rBB10	20001	35238	1.07%	0.076%
18	14.345	1906	1914	1922	rBV	1416547	2152515	65.63%	4.661%
19	14.557	1944	1950	1961	rBB	314980	565501	17.24%	1.225%
20	15.345	2073	2084	2092	rBV	1621845	2400222	73.18%	5.197%
21	15.463	2098	2104	2112	rBB2	429845	596505	18.19%	1.292%
22	15.628	2128	2132	2138	rBB9	35901	41677	1.27%	0.090%
23	16.104	2211	2213	2216	rBB3	47082	40000	1.22%	0.087%
24	16.457	2269	2273	2278	rBB8	19353	38528	1.17%	0.083%
25	16.922	2347	2352	2356	rVB6	83207	122641	3.74%	0.266%
26	17.098	2375	2382	2388	rBB	1786247	2517458	76.75%	5.451%
27	17.198	2393	2399	2408	rBV2	1807356	2566362	78.24%	5.557%
28	17.528	2451	2455	2461	rBB9	63856	111547	3.40%	0.242%
29	18.010	2533	2537	2544	rBB10	86529	156989	4.79%	0.340%
30	18.163	2561	2563	2567	rBB5	36890	40986	1.25%	0.089%
31	18.269	2577	2581	2584	rBB6	45315	44370	1.35%	0.096%
32	18.339	2589	2593	2596	rBB6	31610	53079	1.62%	0.115%
33	18.416	2602	2606	2611	rBB8	64588	97007	2.96%	0.210%
34	18.539	2621	2627	2631	rBB9	65082	75570	2.30%	0.164%
35	18.586	2631	2635	2637	rBB5	29555	46726	1.42%	0.101%
36	18.639	2641	2644	2650	rBB8	23771	35092	1.07%	0.076%

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

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Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
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37	18.716	2652	2657	2659	rBB6	27120	45104	1.38%	0.098%
38	18.775	2663	2667	2670	rBB5	90220	143987	4.39%	0.312%
39	19.022	2705	2709	2711	rBB5	45463	58020	1.77%	0.126%
40	19.057	2713	2715	2717	rBB3	42602	42725	1.30%	0.093%
41	19.116	2719	2725	2728	rBB8	95719	139983	4.27%	0.303%
42	19.163	2728	2733	2739	rBB	184677	257256	7.84%	0.557%
43	19.292	2753	2755	2756	rBB2	41908	37917	1.16%	0.082%
44	19.333	2760	2762	2765	rBB4	63710	38827	1.18%	0.084%
45	19.398	2768	2773	2774	rBB5	46861	69755	2.13%	0.151%
46	19.457	2779	2783	2784	rBB4	53263	46655	1.42%	0.101%
47	19.504	2784	2791	2798	rBB	2049413	3056418	93.18%	6.618%
48	19.704	2821	2825	2829	rBB7	236107	307085	9.36%	0.665%
49	19.769	2831	2836	2839	rBB7	40153	52721	1.61%	0.114%
50	19.839	2844	2848	2853	rBB8	104148	163016	4.97%	0.353%
51	19.916	2860	2861	2867	rBB6	56897	62893	1.92%	0.136%
52	19.969	2867	2870	2872	rBB4	44115	52661	1.61%	0.114%
53	20.033	2878	2881	2883	rBB4	79277	84054	2.56%	0.182%
54	20.069	2883	2887	2889	rBB5	41309	56880	1.73%	0.123%
55	20.127	2889	2897	2900	rBB10	126790	190989	5.82%	0.414%
56	20.257	2913	2919	2920	rBB6	65331	83409	2.54%	0.181%
57	20.292	2923	2925	2927	rBB3	49003	39622	1.21%	0.086%
58	20.369	2934	2938	2939	rBB4	61436	57636	1.76%	0.125%
59	20.410	2943	2945	2949	rBB5	58901	73358	2.24%	0.159%
60	20.480	2955	2957	2959	rBB3	44506	39045	1.19%	0.085%
61	20.645	2980	2985	2990	rBB8	120263	209069	6.37%	0.453%
62	20.733	2998	3000	3002	rBB3	51483	36508	1.11%	0.079%
63	20.816	3012	3014	3015	rBB2	45364	34039	1.04%	0.074%
64	20.851	3017	3020	3024	rBB6	80331	80795	2.46%	0.175%
65	20.892	3024	3027	3031	rBB6	71149	88565	2.70%	0.192%
66	20.998	3040	3045	3047	rBB5	123616	194274	5.92%	0.421%
67	21.021	3047	3049	3051	rBB3	46698	33648	1.03%	0.073%
68	21.039	3051	3052	3056	rBB4	47065	56274	1.72%	0.122%
69	21.110	3061	3064	3066	rBB4	128898	109505	3.34%	0.237%
70	21.157	3070	3072	3074	rBB3	73150	36301	1.11%	0.079%
71	21.227	3078	3084	3089	rBB8	141467	254080	7.75%	0.550%
72	21.298	3091	3096	3100	rBB	2258778	2811013	85.70%	6.087%
73	21.333	3100	3102	3106	rBB5	72039	83654	2.55%	0.181%
74	21.363	3106	3107	3111	rBB4	57558	44901	1.37%	0.097%
75	21.451	3118	3122	3124	rBB5	91087	93818	2.86%	0.203%
76	21.563	3139	3141	3144	rBB4	52033	44592	1.36%	0.097%

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
 Title : SVOA CALIBRATION

77	21.616	3148	3150	3154	rBB5	47653	35650	1.09%	0.077%
78	21.686	3158	3162	3164	rBB5	90934	83934	2.56%	0.182%
79	21.786	3178	3179	3181	rBB2	65456	44930	1.37%	0.097%
80	21.851	3187	3190	3191	rBB3	46846	52386	1.60%	0.113%
81	21.945	3204	3206	3210	rBB5	58250	39517	1.20%	0.086%
82	22.027	3210	3220	3222	rBB5	59479	116427	3.55%	0.252%
83	22.157	3239	3242	3244	rBB4	64195	76160	2.32%	0.165%
84	22.221	3251	3253	3257	rBB5	53269	47310	1.44%	0.102%
85	22.280	3261	3263	3266	rBB4	66774	45325	1.38%	0.098%
86	22.368	3274	3278	3283	rBB7	138549	165095	5.03%	0.357%
87	22.463	3290	3294	3296	rBB5	64306	72311	2.20%	0.157%
88	22.486	3296	3298	3300	rBB3	52337	46140	1.41%	0.100%
89	22.563	3309	3311	3315	rBB5	41695	42706	1.30%	0.092%
90	22.645	3323	3325	3326	rBB2	40167	37269	1.14%	0.081%
91	22.692	3330	3333	3339	rBB8	48627	74461	2.27%	0.161%
92	22.880	3355	3365	3370	rBB8	234773	571319	17.42%	1.237%
93	22.939	3370	3375	3379	rBB8	91356	111004	3.38%	0.240%
94	23.068	3392	3397	3398	rBB5	46545	77092	2.35%	0.167%
95	23.080	3398	3399	3403	rBB4	46237	52712	1.61%	0.114%
96	23.121	3405	3406	3411	rBB5	37947	54164	1.65%	0.117%
97	23.286	3430	3434	3440	rBB7	117531	193396	5.90%	0.419%
98	23.333	3440	3442	3445	rBB3	36371	32023	0.98%	0.069%
99	23.410	3447	3455	3461	rBB	1564868	2771199	84.49%	6.001%
100	23.457	3461	3463	3466	rBB4	125522	84580	2.58%	0.183%
101	23.557	3473	3480	3488	rBB2	1640970	3279968	100.00%	7.102%
102	23.833	3524	3527	3529	rBB4	38796	35674	1.09%	0.077%
103	23.863	3529	3532	3534	rBB4	52017	44206	1.35%	0.096%
104	24.015	3554	3558	3560	rBB5	37952	49792	1.52%	0.108%
105	24.051	3562	3564	3566	rBB3	39578	35560	1.08%	0.077%
106	24.110	3568	3574	3582	rBB2	303138	619463	18.89%	1.341%
107	24.162	3582	3583	3587	rBB4	42532	34278	1.05%	0.074%
108	24.286	3601	3604	3606	rBB4	24955	40282	1.23%	0.087%
109	24.345	3611	3614	3616	rBB4	40505	31278	0.95%	0.068%
110	24.651	3664	3666	3673	rBB8	33675	34773	1.06%	0.075%
111	24.868	3696	3703	3710	rBB5	318779	725593	22.12%	1.571%
112	25.139	3748	3749	3754	rBB5	29105	36146	1.10%	0.078%
113	25.215	3760	3762	3767	rBB6	37628	45769	1.40%	0.099%
114	25.274	3767	3772	3773	rBB5	24166	35124	1.07%	0.076%
115	25.309	3773	3778	3782	rBB8	76551	137512	4.19%	0.298%
116	25.415	3795	3796	3801	rBB5	40259	49836	1.52%	0.108%

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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-BN052521.M
 Title : SVOA CALIBRATION

117	25.504	3808	3811	3815	rBB6	35895	44448	1.36%	0.096%
118	25.598	3823	3827	3829	rBB5	32988	38756	1.18%	0.084%
119	25.680	3838	3841	3844	rBB5	34537	36623	1.12%	0.079%
120	25.751	3844	3853	3862	rBB7	207283	532331	16.23%	1.153%
121	26.086	3907	3910	3913	rBB5	42045	44135	1.35%	0.096%
122	26.121	3913	3916	3918	rBV4	45326	51427	1.57%	0.111%
123	26.380	3956	3960	3965	rBV8	40810	54859	1.67%	0.119%
124	26.433	3967	3969	3972	rBB4	36896	40865	1.25%	0.088%
125	26.504	3978	3981	3982	rBB3	49994	36018	1.10%	0.078%
126	26.798	4021	4031	4039	rBB3	167985	491475	14.98%	1.064%
127	26.898	4046	4048	4052	rBB5	27213	35471	1.08%	0.077%
128	27.151	4087	4091	4094	rBV6	35851	35899	1.09%	0.078%
129	27.662	4176	4178	4181	rBB4	40758	40024	1.22%	0.087%
130	27.692	4181	4183	4186	rBB4	31051	37103	1.13%	0.080%
131	28.021	4237	4239	4248	rBB10	30371	51396	1.57%	0.111%
132	28.080	4248	4249	4257	rBB8	25429	43278	1.32%	0.094%
133	28.356	4293	4296	4298	rBB4	42022	37494	1.14%	0.081%
134	28.974	4397	4401	4404	rBB6	32757	38901	1.19%	0.084%
135	29.092	4416	4421	4423	rBB6	24978	32145	0.98%	0.070%
136	29.233	4441	4445	4448	rBB6	20854	33535	1.02%	0.073%

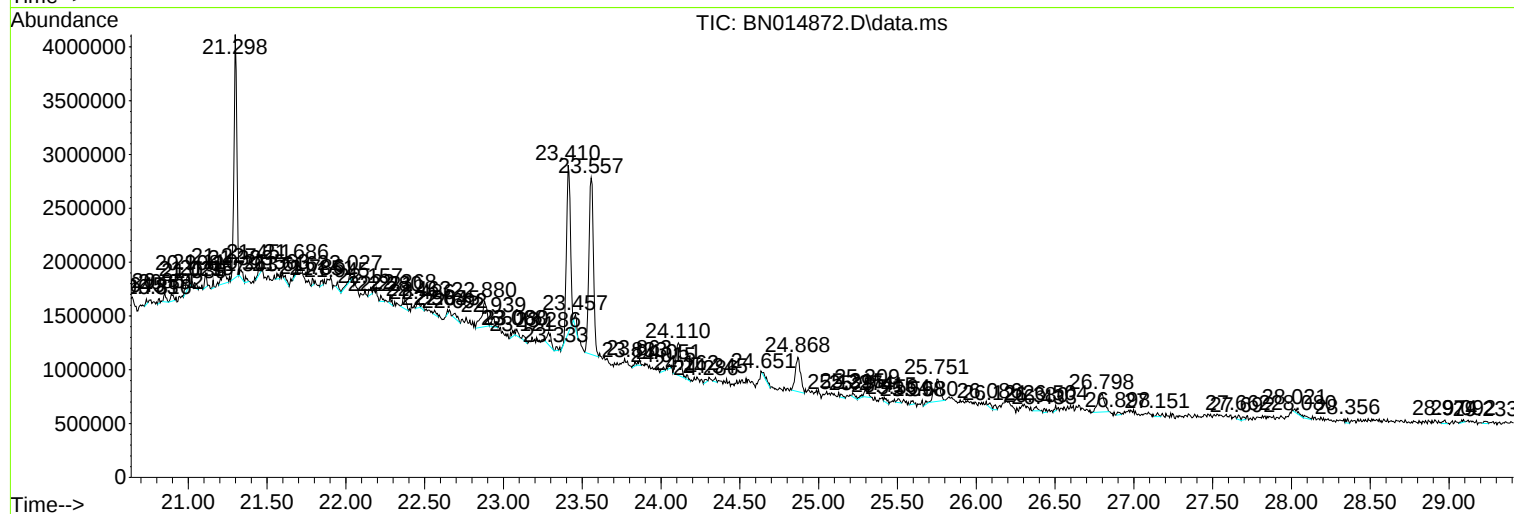
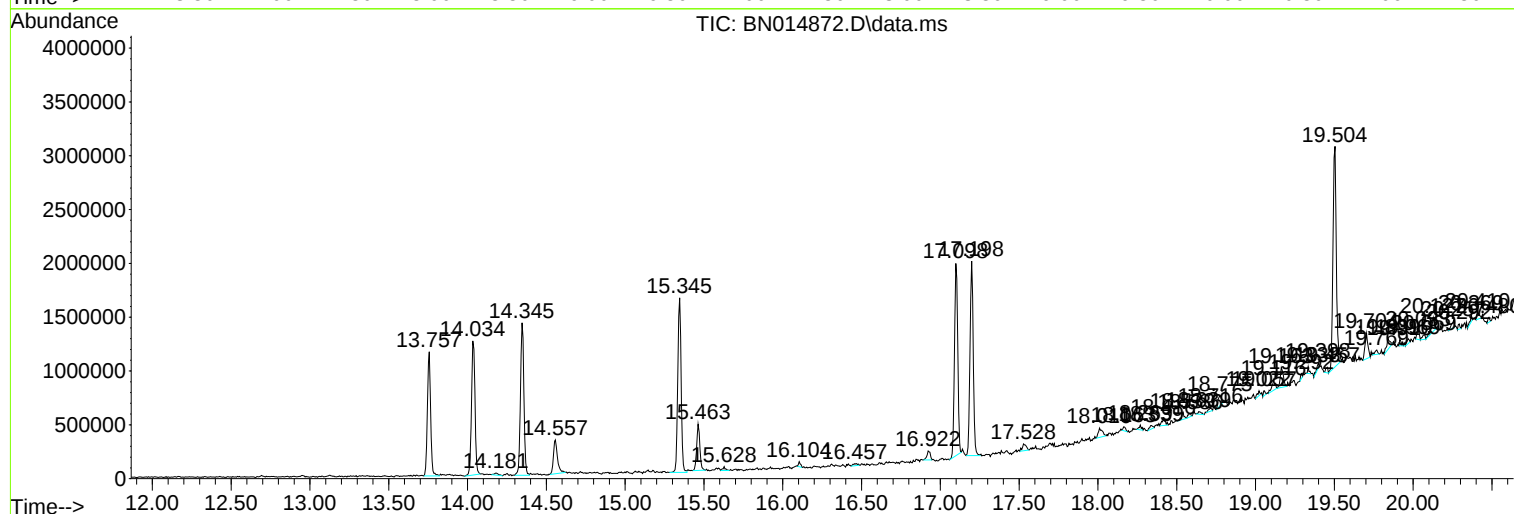
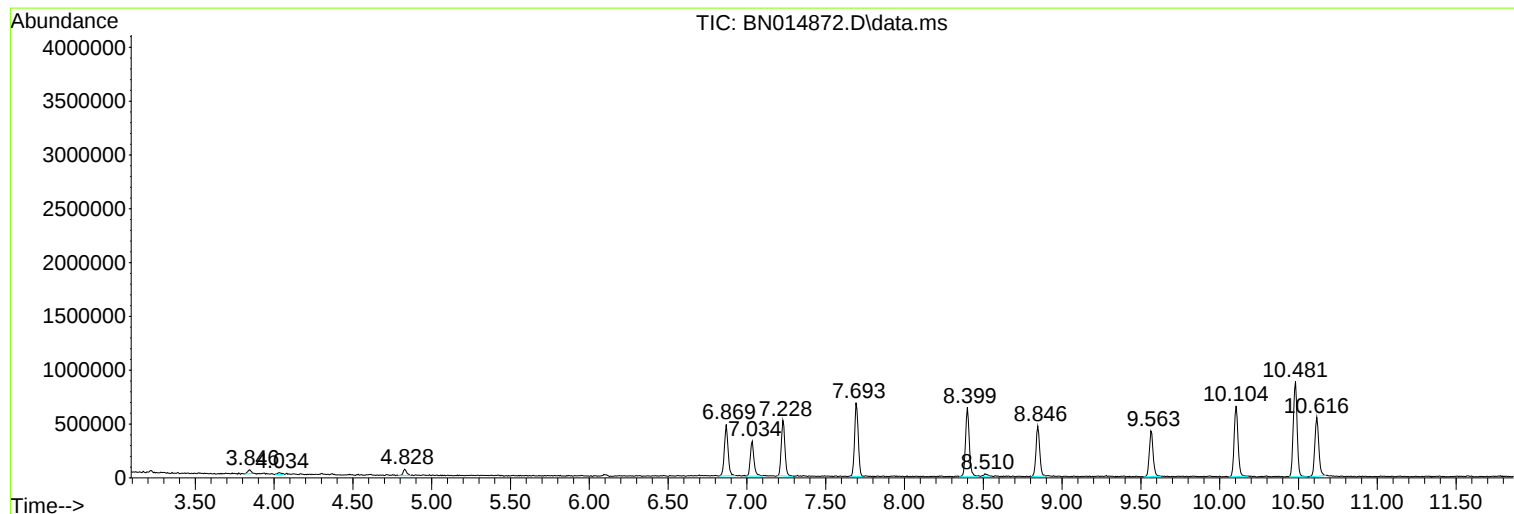
Sum of corrected areas: 46181408

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION

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



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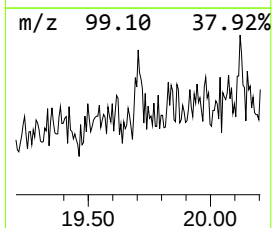
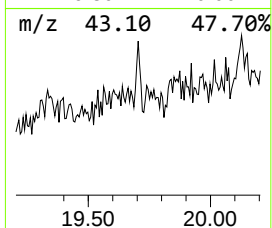
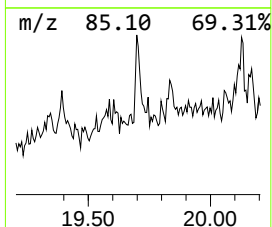
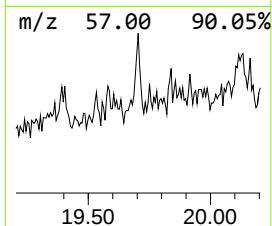
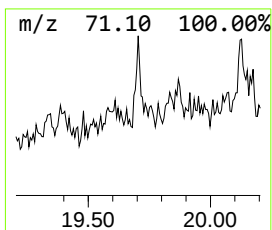
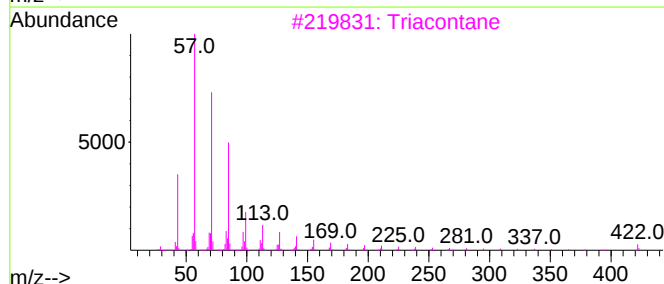
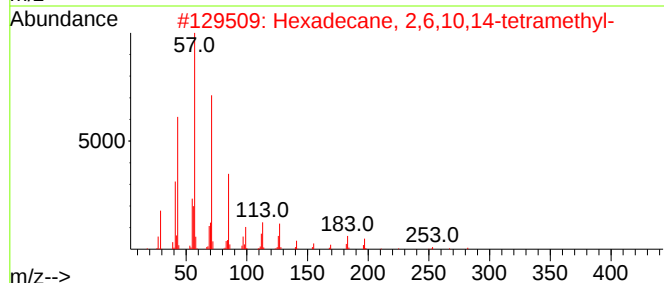
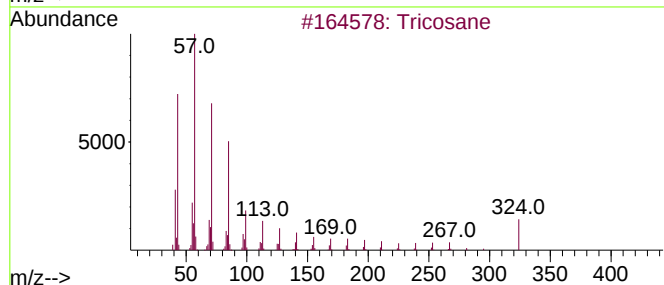
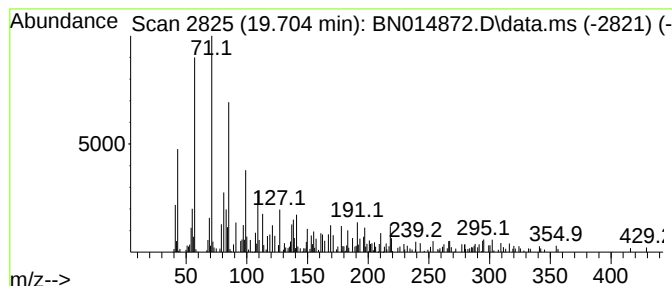
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.704	2.18 ng/ul	307085	Chrysene-d12	21.298	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricosane	324	C23H48	000638-67-5	81
2	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	76
3	Triacontane	422	C30H62	000638-68-6	72
4	Tetracosane	338	C24H50	000646-31-1	72
5	Tetracosane	338	C24H50	000646-31-1	64



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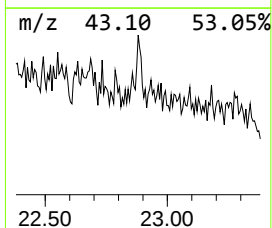
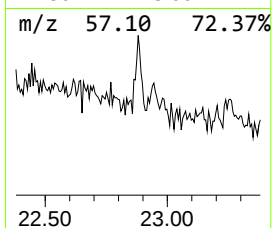
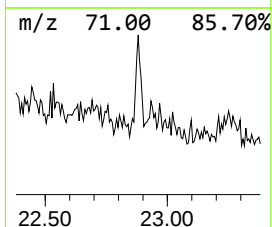
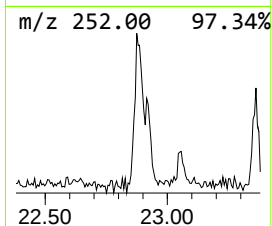
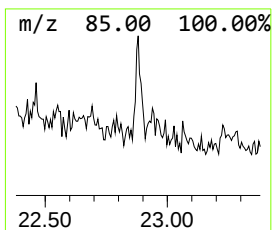
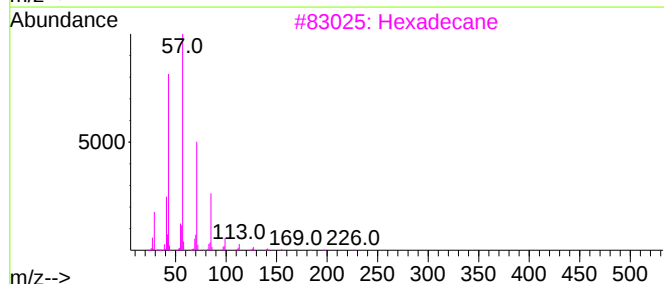
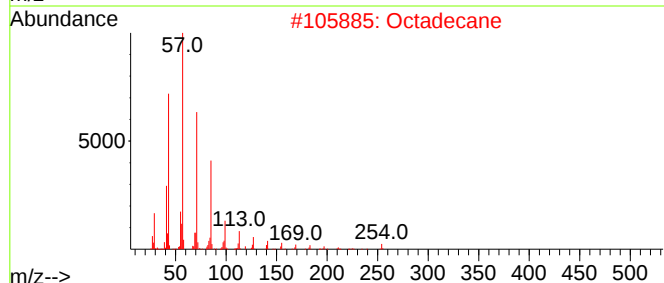
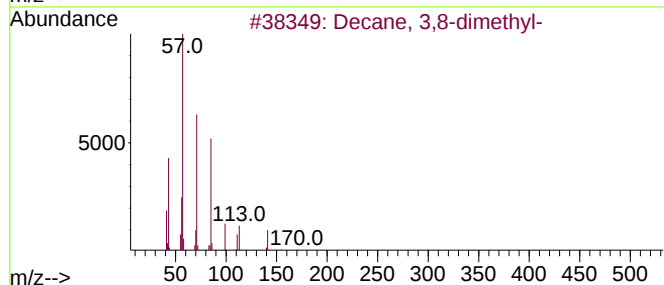
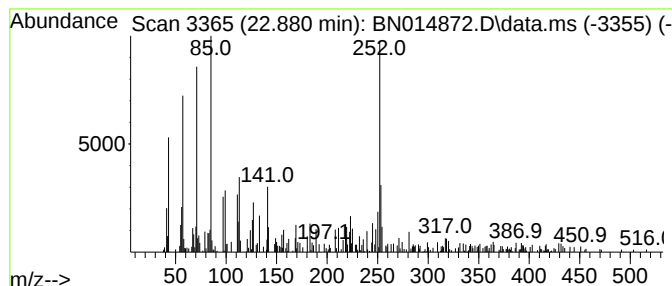
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.880	3.48 ng/ul	571319	Perylene-d12	23.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	60
2			Octadecane	254	C18H38	000593-45-3	49
3			Hexadecane	226	C16H34	000544-76-3	47
4			1-Iodoundecane	282	C11H23I	004282-44-4	47
5			Decane, 1-iodo-	268	C10H21I	002050-77-3	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014872.D
Acq On : 10 Jun 2021 21:25
Operator : CG/JU
Sample : M2618-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BG5R1

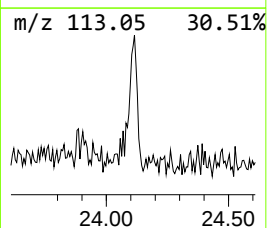
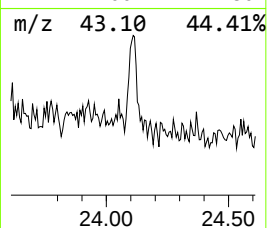
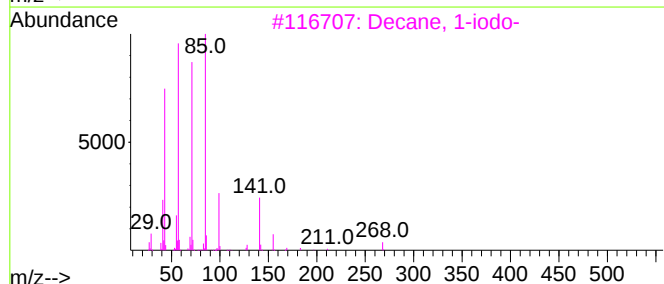
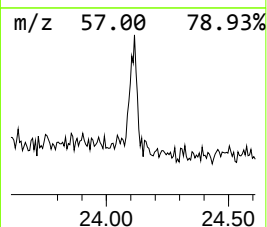
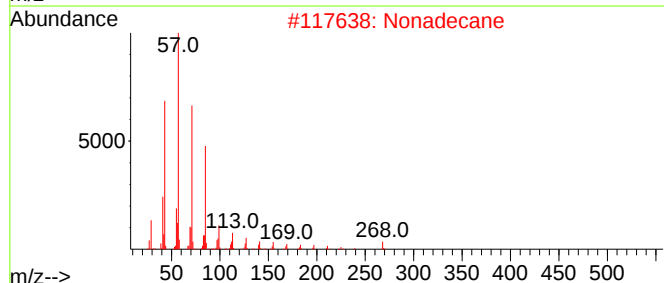
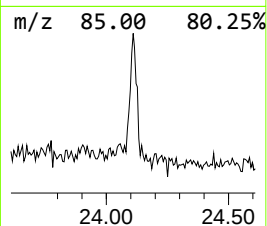
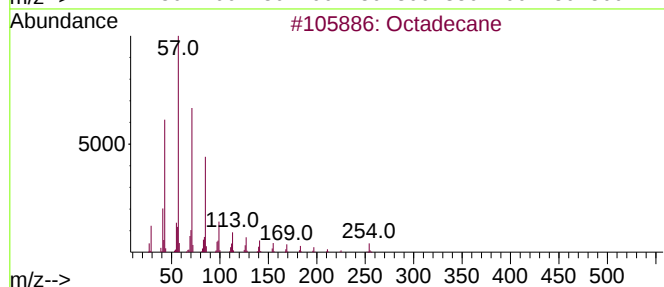
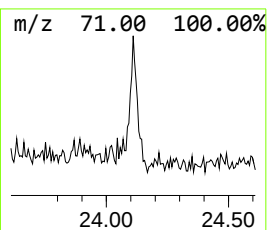
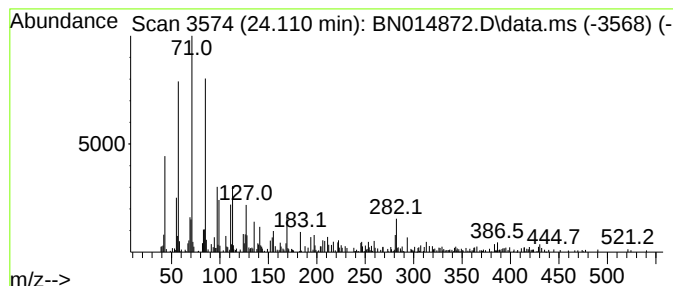
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.110	3.78 ng/ul	619463	Perylene-d12	23.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	93
2			Nonadecane	268	C19H40	000629-92-5	86
3			Decane, 1-iodo-	268	C10H21I	002050-77-3	78
4			Hexacosane, 1,26-dibromo-	522	C26H52Br2	143389-26-8	68
5			Eicosane	282	C20H42	000112-95-8	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014872.D
Acq On : 10 Jun 2021 21:25
Operator : CG/JU
Sample : M2618-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BG5R1

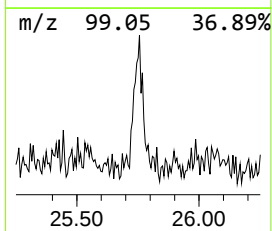
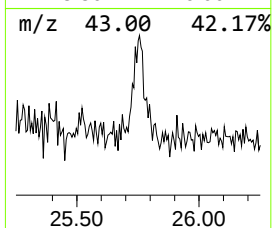
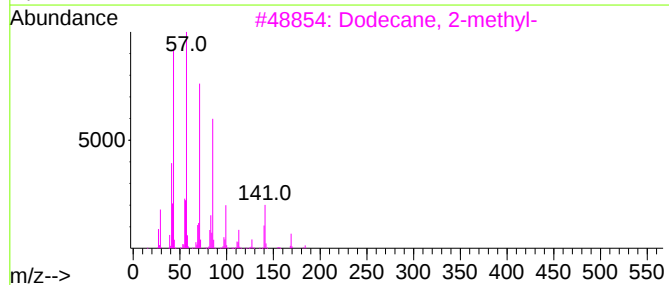
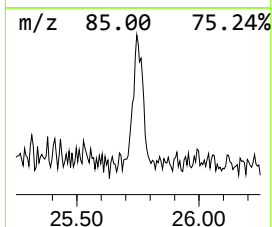
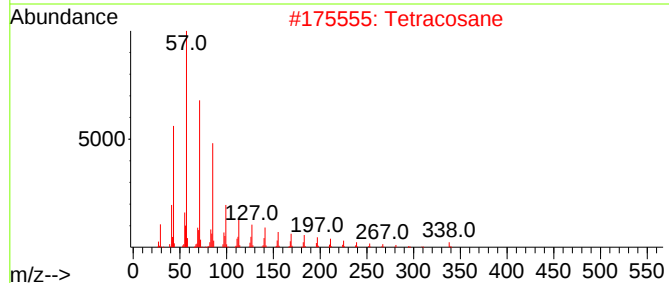
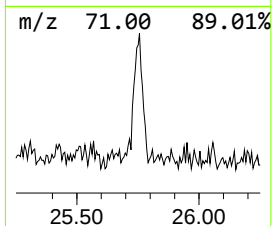
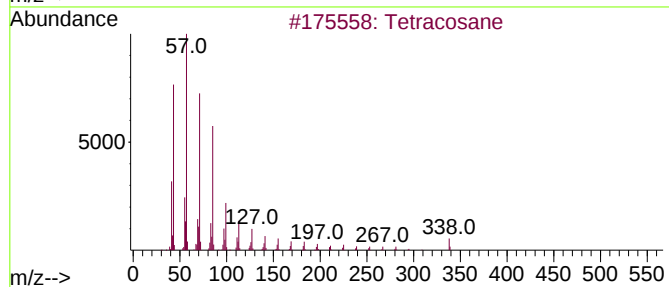
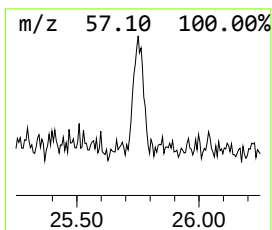
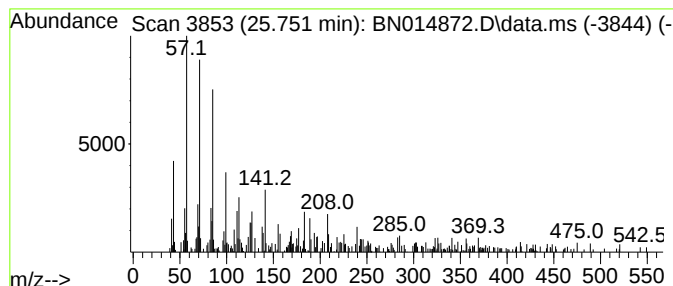
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.751	3.25 ng/ul	532331	Perylene-d12	23.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	91
2			Tetracosane	338	C24H50	000646-31-1	91
3			Dodecane, 2-methyl-	184	C13H28	001560-97-0	87
4			Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	87
5			Behenyl chloride	344	C22H45Cl	042217-03-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014872.D
Acq On : 10 Jun 2021 21:25
Operator : CG/JU
Sample : M2618-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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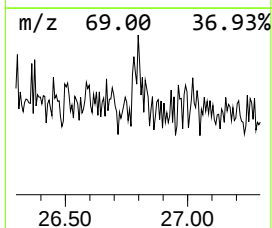
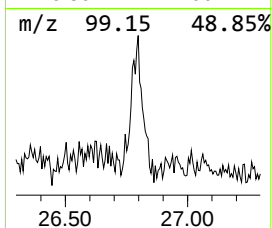
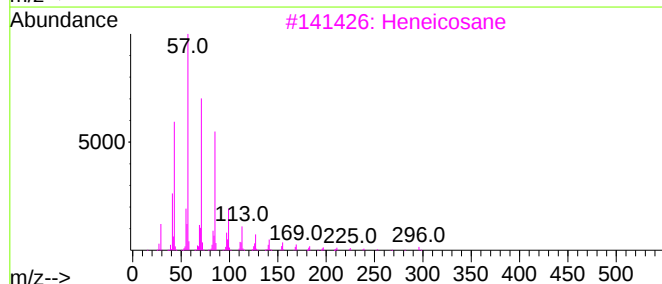
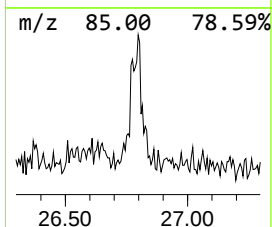
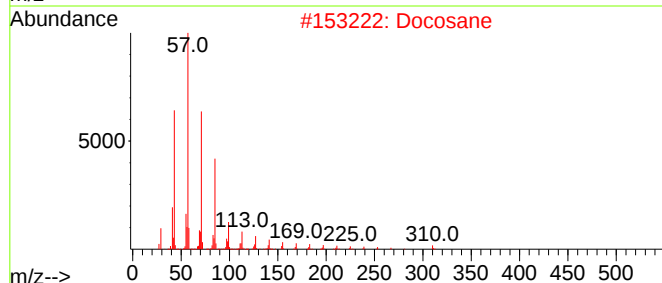
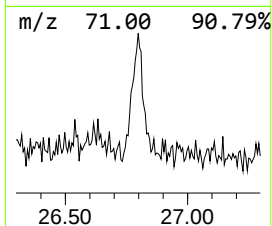
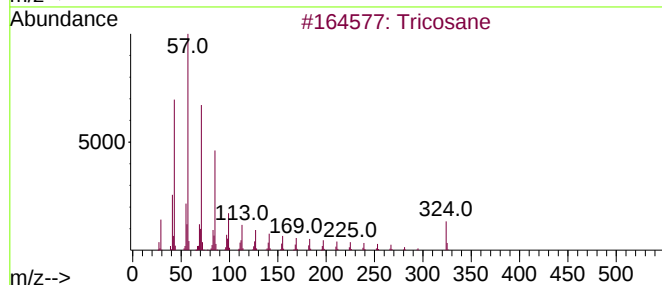
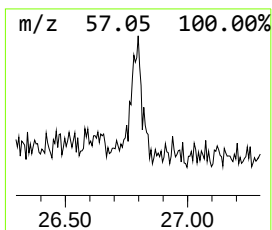
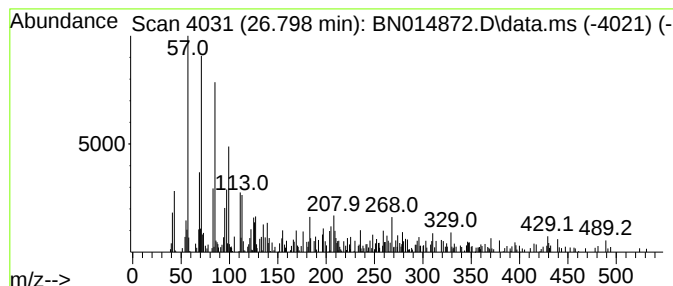
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.798	3.00 ng/ul	491475	Perylene-d12	23.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricosane	324	C23H48	000638-67-5	89
2			Docosane	310	C22H46	000629-97-0	76
3			Heneicosane	296	C21H44	000629-94-7	76
4			Tetratriacontane	479	C34H70	014167-59-0	76
5			Docosane	310	C22H46	000629-97-0	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014872.D
Acq On : 10 Jun 2021 21:25
Operator : CG/JU
Sample : M2618-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BG5R1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.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
(DEL) Alkane: S...	19.704	2.2	ng/ul	307085	5	21.298	2811010	20.0
(DEL) Alkane: S...	22.880	3.5	ng/ul	571319	6	23.557	3279970	20.0
(DEL) Alkane: S...	24.110	3.8	ng/ul	619463	6	23.557	3279970	20.0
(DEL) Alkane: S...	24.868	4.4	ng/ul	725593	6	23.557	3279970	20.0
(DEL) Alkane: S...	25.751	3.3	ng/ul	532331	6	23.557	3279970	20.0
(DEL) Alkane: S...	26.798	3.0	ng/ul	491475	6	23.557	3279970	20.0