

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
 Data File : BG061286.D
 Acq On : 24 May 2024 13:49
 Operator : MA/JU
 Sample : P2548-17
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BH612

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_G\Methods\SFAM-EPA-BG052124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG061286.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.082	11	16	22	rBV	153876	218499	23.45%	2.281%
2	3.341	56	60	64	rVV5	4613	6435	0.69%	0.067%
3	3.611	102	106	112	rBV2	9737	16517	1.77%	0.172%
4	3.746	125	129	139	rBV	54860	88671	9.52%	0.926%
5	7.066	686	694	702	rBV	99605	171532	18.41%	1.790%
6	7.213	712	719	727	rVB	64335	105081	11.28%	1.097%
7	7.424	749	755	762	rBV	90936	151649	16.28%	1.583%
8	7.888	827	834	841	rVB	68071	114302	12.27%	1.193%
9	8.599	949	955	965	rBV	123087	211069	22.66%	2.203%
10	9.040	1021	1030	1037	rBV	102807	178965	19.21%	1.868%
11	9.598	1122	1125	1129	rBB3	3602	4819	0.52%	0.050%
12	9.762	1146	1153	1161	rBV2	96146	169398	18.18%	1.768%
13	10.309	1239	1246	1259	rVB	131993	236299	25.36%	2.466%
14	10.679	1302	1309	1318	rBV	101968	176687	18.97%	1.844%
15	10.814	1324	1332	1343	rBV	131385	232600	24.97%	2.428%
16	11.419	1429	1435	1442	rBV3	7877	13479	1.45%	0.141%
17	13.417	1769	1775	1779	rVB2	2733	4391	0.47%	0.046%
18	13.928	1855	1862	1868	rBV	293735	422717	45.37%	4.412%
19	14.216	1904	1911	1920	rVB2	295560	470373	50.49%	4.910%
20	14.522	1953	1963	1969	rBV	182460	284910	30.58%	2.974%
21	14.745	1993	2001	2020	rBV2	106399	197449	21.19%	2.061%
22	15.515	2126	2132	2138	rBV	424894	614912	66.00%	6.418%
23	15.644	2148	2154	2166	rBV	139958	213069	22.87%	2.224%
24	17.265	2424	2430	2435	rVV	260271	387673	41.61%	4.047%
25	17.307	2435	2437	2441	rVV	19973	25802	2.77%	0.269%
26	17.365	2441	2447	2457	rVB2	484927	693565	74.45%	7.239%
27	18.147	2575	2580	2582	rBV	5239	7524	0.81%	0.079%
28	18.194	2582	2588	2593	rVV4	9517	21095	2.26%	0.220%
29	18.335	2607	2612	2616	rVV5	7345	12909	1.39%	0.135%
30	18.376	2616	2619	2625	rVB4	4876	7735	0.83%	0.081%
31	18.464	2630	2634	2636	rBV2	2131	2890	0.31%	0.030%
32	18.517	2636	2643	2645	rBV3	1291	2814	0.30%	0.029%
33	18.646	2661	2665	2669	rVB4	3988	5779	0.62%	0.060%
34	18.687	2669	2672	2677	rBV5	4775	6446	0.69%	0.067%
35	18.940	2710	2715	2719	rVB3	2260	3608	0.39%	0.038%
36	19.057	2731	2735	2741	rBV4	5353	11962	1.28%	0.125%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
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37	19.122	2743	2746	2749	rVB3	4600	5525	0.59%	0.058%
38	19.222	2759	2763	2768	rVB5	9714	16061	1.72%	0.168%
39	19.292	2768	2775	2776	rBV3	6118	9016	0.97%	0.094%
40	19.322	2776	2780	2787	rVB	59007	80040	8.59%	0.835%

41	19.422	2793	2797	2799	rBV4	2362	2846	0.31%	0.030%
42	19.522	2811	2814	2818	rBV4	3550	5258	0.56%	0.055%
43	19.580	2818	2824	2829	rBV8	5692	12547	1.35%	0.131%
44	19.657	2831	2837	2850	rVB	585109	931646	100.00%	9.724%
45	19.763	2850	2855	2859	rBV6	6264	10270	1.10%	0.107%

46	19.839	2862	2868	2870	rBV6	4553	6322	0.68%	0.066%
47	19.915	2878	2881	2885	rBV6	4367	6314	0.68%	0.066%
48	20.015	2892	2898	2899	rBV5	5696	9554	1.03%	0.100%
49	20.103	2910	2913	2915	rBV4	3092	3227	0.35%	0.034%
50	20.139	2916	2919	2922	rBV5	4933	6653	0.71%	0.069%

51	20.186	2923	2927	2933	rVB4	9417	17494	1.88%	0.183%
52	20.280	2940	2943	2947	rBV3	5644	6588	0.71%	0.069%
53	20.338	2949	2953	2954	rBV4	5902	8005	0.86%	0.084%
54	20.362	2954	2957	2963	rVB4	17984	24790	2.66%	0.259%
55	20.432	2963	2969	2971	rVB5	2757	4645	0.50%	0.048%

56	20.473	2973	2976	2980	rBV3	5619	8573	0.92%	0.089%
57	20.532	2980	2986	2990	rBV	156164	202387	21.72%	2.113%
58	20.573	2990	2993	3000	rVV7	19563	43243	4.64%	0.451%
59	20.626	3000	3002	3007	rVB6	9959	12526	1.34%	0.131%
60	20.697	3011	3014	3017	rVB3	7653	8664	0.93%	0.090%

61	20.932	3050	3054	3059	rBV	7342	12520	1.34%	0.131%
62	21.020	3067	3069	3073	rVB5	4068	5100	0.55%	0.053%
63	21.108	3080	3084	3085	rBV4	4697	6269	0.67%	0.065%
64	21.184	3092	3097	3105	rBV6	40417	98302	10.55%	1.026%
65	21.419	3135	3137	3141	rVB4	5944	5574	0.60%	0.058%

66	21.519	3146	3154	3159	rVV2	294050	517007	55.49%	5.397%
67	21.560	3159	3161	3166	rVB2	33945	46405	4.98%	0.484%
68	21.719	3183	3188	3193	rBV8	14496	29445	3.16%	0.307%
69	21.925	3221	3223	3228	rVB6	4416	5587	0.60%	0.058%
70	22.224	3270	3274	3280	rVB6	18485	32357	3.47%	0.338%

71	22.301	3282	3287	3291	rBV5	18781	31226	3.35%	0.326%
72	22.483	3315	3318	3321	rBV5	5509	8844	0.95%	0.092%
73	22.959	3391	3399	3409	rVB5	21310	50203	5.39%	0.524%
74	23.629	3507	3513	3520	rBV2	24536	67459	7.24%	0.704%
75	23.728	3526	3530	3537	rVB2	27523	55277	5.93%	0.577%

76	24.063	3586	3587	3590	rBV3	4390	4732	0.51%	0.049%
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 Title : SVOA CALIBRATION

77	24.316	3625	3630	3631	rBV3	11658	17430	1.87%	0.182%
78	24.398	3635	3644	3652	rVV2	319698	826831	88.75%	8.630%
79	24.610	3670	3680	3692	rBV	181317	510274	54.77%	5.326%
80	24.986	3742	3744	3749	rBV6	3509	6814	0.73%	0.071%
81	25.708	3857	3867	3874	rBV5	26675	77927	8.36%	0.813%
82	25.867	3891	3894	3895	rBV2	4001	3888	0.42%	0.041%
83	26.249	3956	3959	3961	rBV3	5035	5564	0.60%	0.058%
84	26.931	4073	4075	4076	rBV	3413	2929	0.31%	0.031%
85	26.983	4079	4084	4094	rVV5	23138	68629	7.37%	0.716%
86	27.277	4132	4134	4135	rBV2	3453	3034	0.33%	0.032%
87	27.436	4158	4161	4163	rBV4	2869	3295	0.35%	0.034%
88	28.000	4255	4257	4260	rBV4	2879	3543	0.38%	0.037%
89	28.100	4267	4274	4275	rBV5	8368	15261	1.64%	0.159%
90	28.247	4297	4299	4302	rVB3	3350	3434	0.37%	0.036%
91	28.394	4322	4324	4326	rBV3	3375	2822	0.30%	0.029%
92	28.493	4338	4341	4342	rBV3	5859	5779	0.62%	0.060%
93	28.535	4342	4348	4358	rVB8	17834	56711	6.09%	0.592%
94	28.658	4366	4369	4370	rBV3	3334	3148	0.34%	0.033%
95	29.622	4531	4533	4535	rBV3	3091	3055	0.33%	0.032%
96	29.798	4561	4563	4568	rVB5	3905	3630	0.39%	0.038%
97	30.109	4614	4616	4618	rBV3	3170	3048	0.33%	0.032%
98	30.215	4632	4634	4637	rBV3	3068	3472	0.37%	0.036%
99	30.409	4657	4667	4670	rBV10	11239	33088	3.55%	0.345%
100	31.425	4837	4840	4842	rBV3	3492	4669	0.50%	0.049%

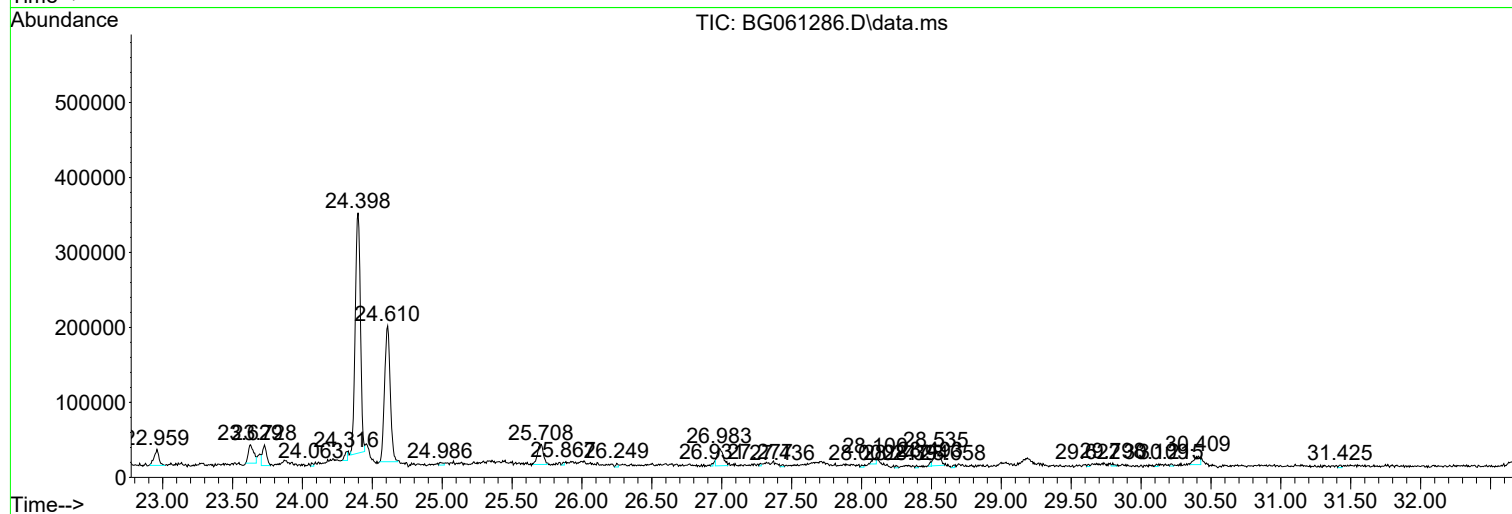
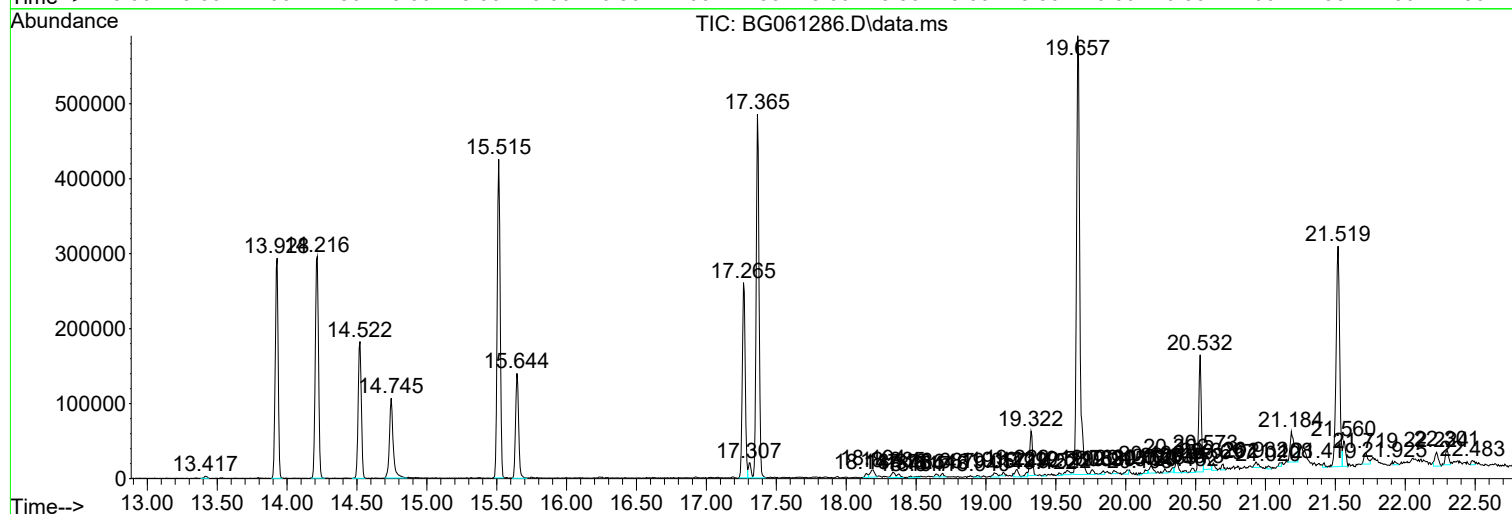
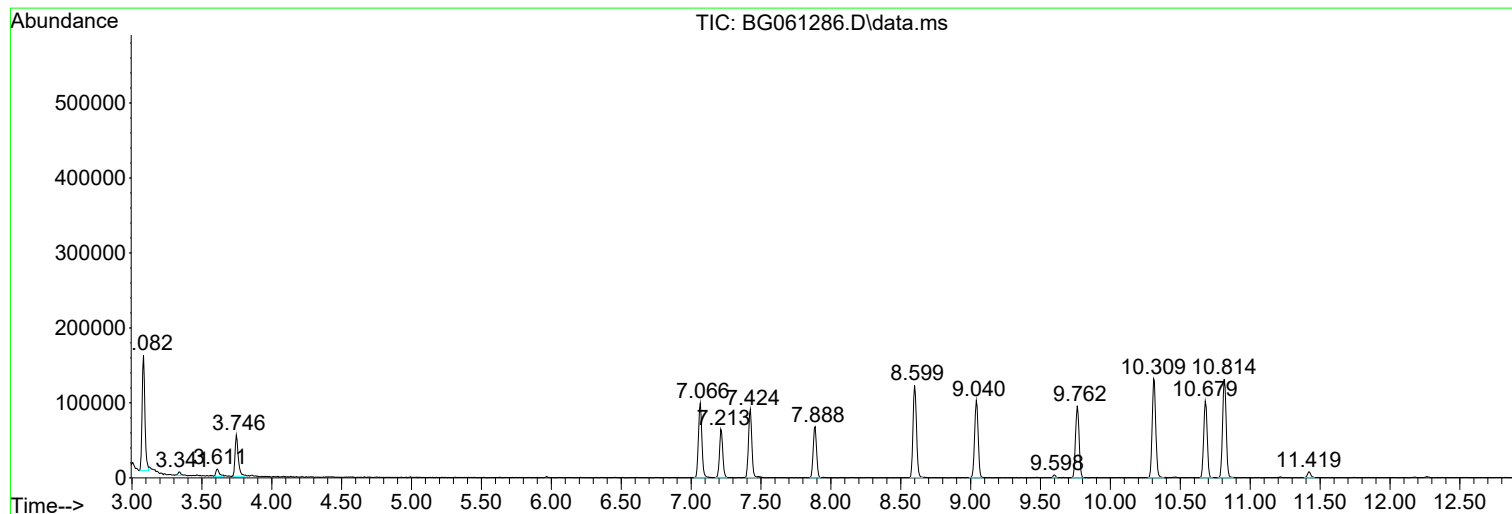
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
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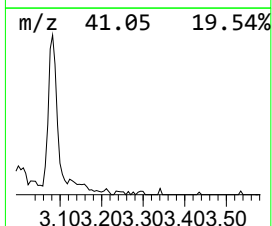
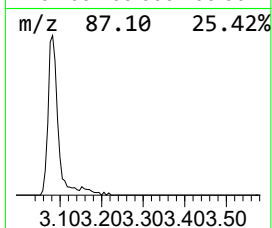
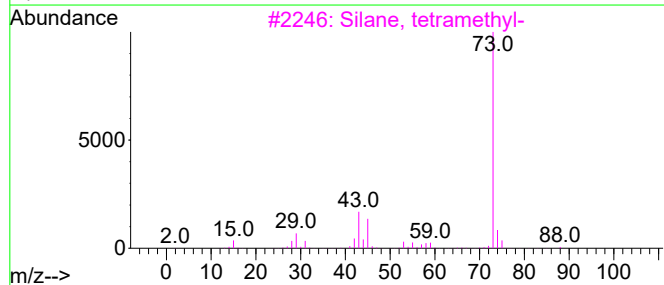
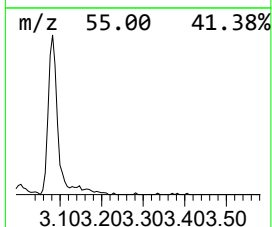
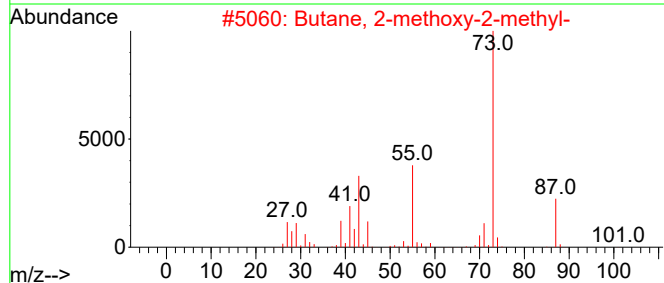
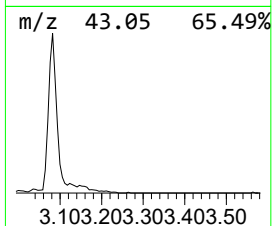
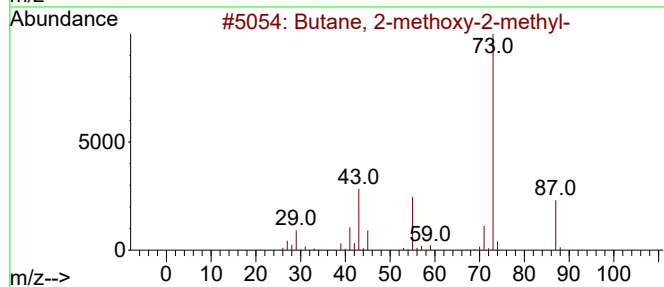
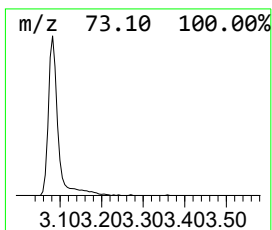
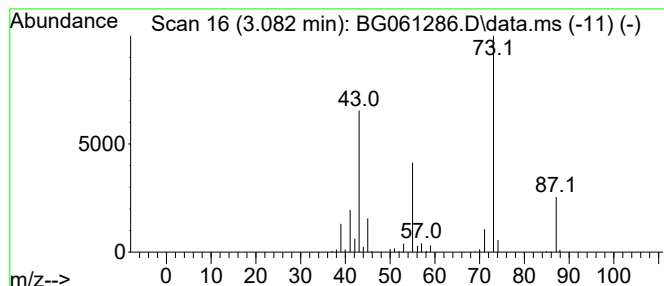
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.082	38.23 ng/ul	218499	1,4-Dichlorobenzene-d4	7.888

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	43
4			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	40
5			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	28



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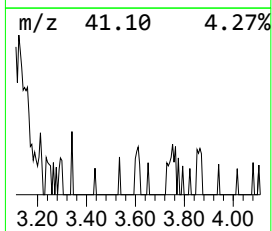
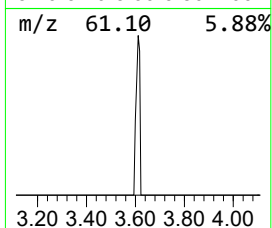
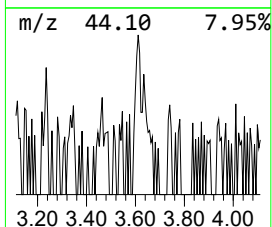
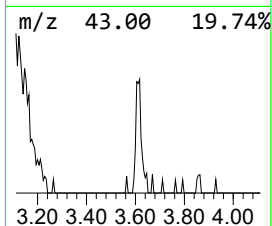
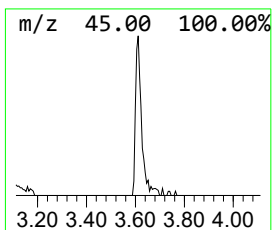
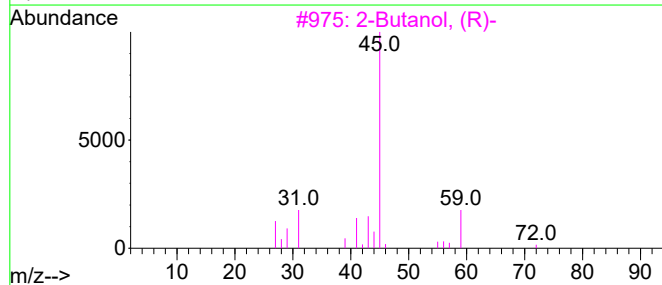
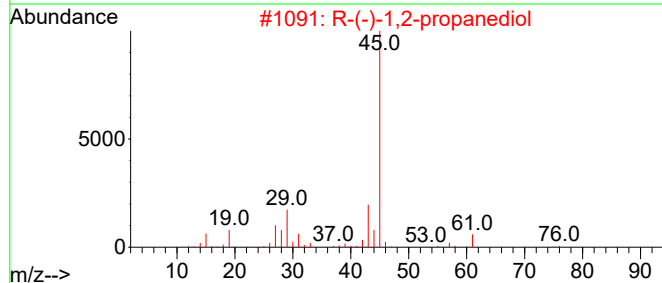
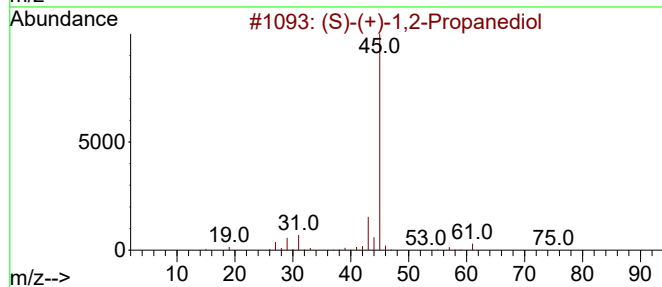
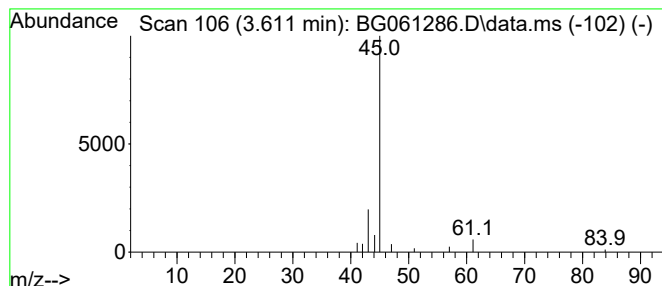
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.611	2.89 ng/ul	16517	1,4-Dichlorobenzene-d4	7.888

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(S)-(+)-1,2-Propanediol			76	C3H8O2	004254-15-3	9
2	R-(-)-1,2-propanediol			76	C3H8O2	004254-14-2	9
3	2-Butanol, (R)-			74	C4H10O	014898-79-4	4
4	2-Hexanol, (R)-			102	C6H14O	026549-24-6	4
5	2,3-Butanediol			90	C4H10O2	000513-85-9	4



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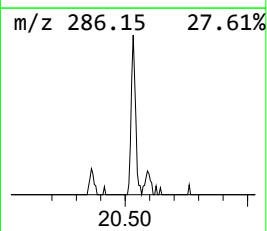
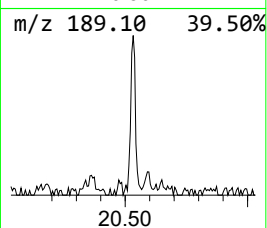
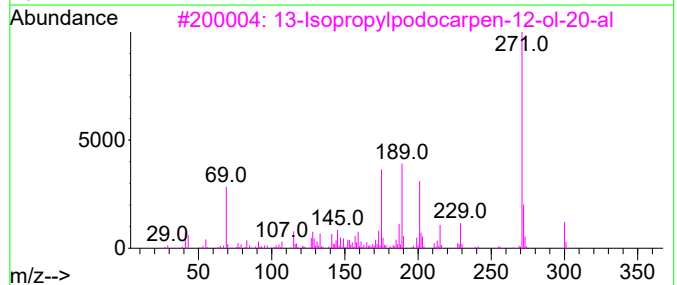
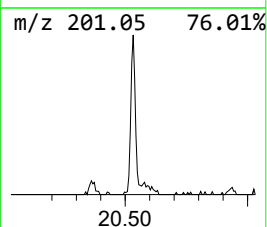
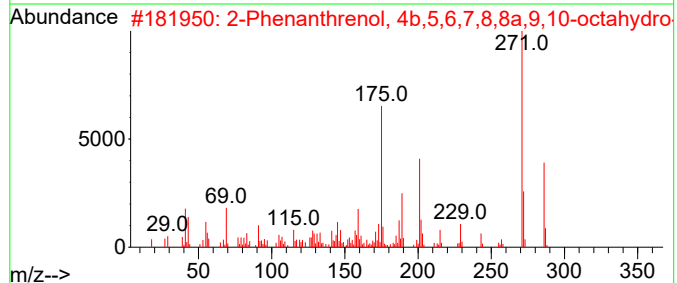
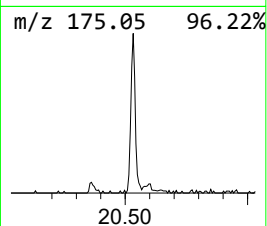
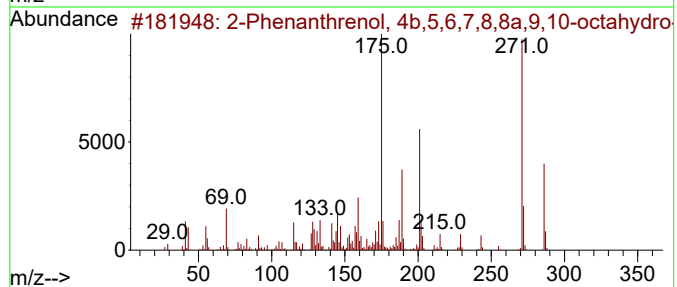
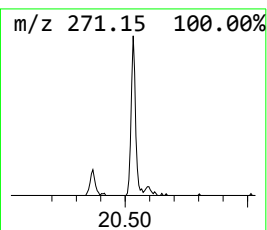
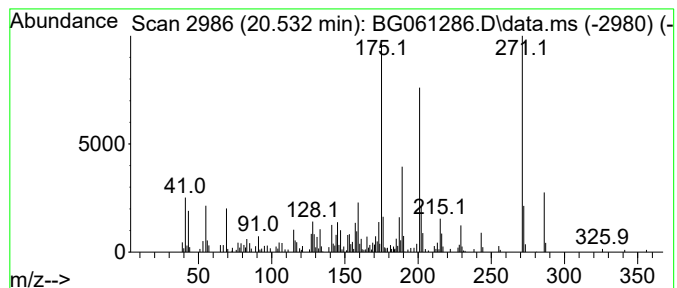
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.532	7.83 ng/ul	202387	Chrysene-d12	21.519

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	87		
2	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	86		
3	13-Isopropylpodocarpene-12-ol-20-al	300	C20H28O2	024035-37-8	43		
4	Pisiferol	302	C20H30O2	024035-36-7	43		
5	13-Isopropylpodocarpene-12-ol-20-al	300	C20H28O2	024035-37-8	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061286.D
Acq On : 24 May 2024 13:49
Operator : MA/JU
Sample : P2548-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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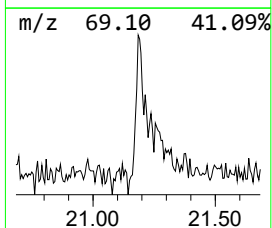
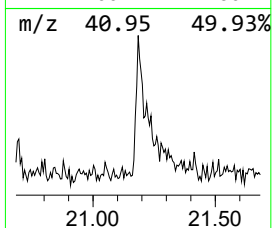
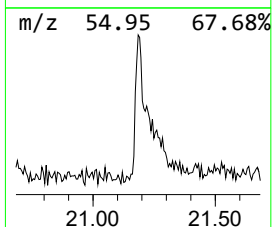
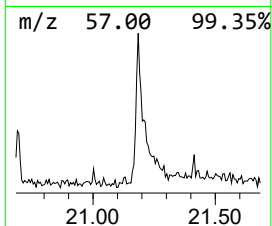
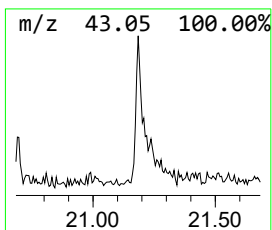
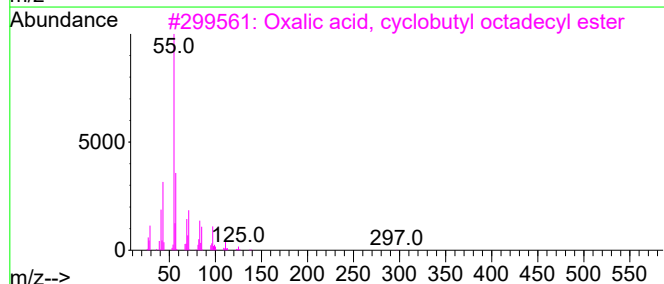
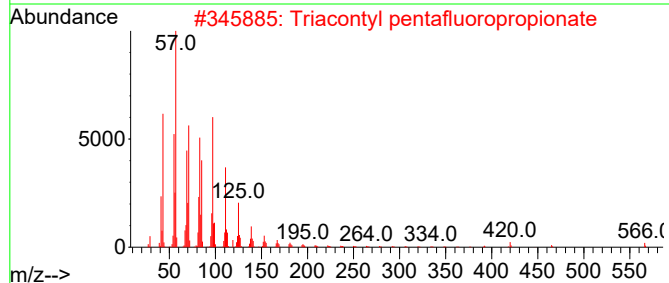
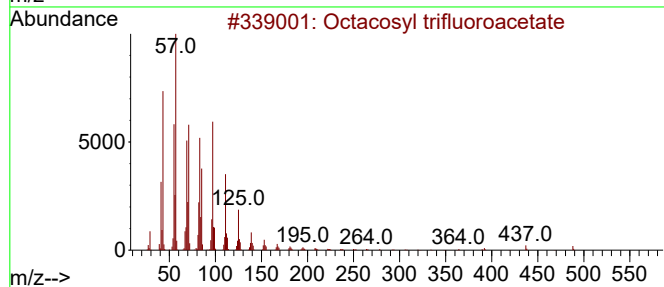
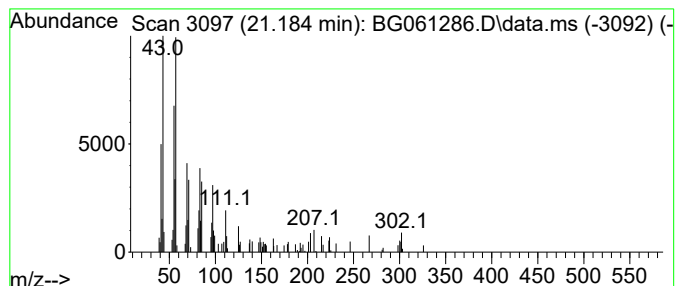
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octacosyl trifluoroacetate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.184	3.80 ng/ul	98302	Chrysene-d12	21.519

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	62
2			Triacontyl pentafluoropropionate	584	C33H61F5O2	1000351-80-0	62
3			Oxalic acid, cyclobutyl octadecy...	396	C24H44O4	1000309-70-8	62
4			Tetratriacontyl pentafluoropropi...	641	C37H69F5O2	1000351-81-5	62
5			1-Hentetracontanol	593	C41H84O	040710-42-7	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061286.D
Acq On : 24 May 2024 13:49
Operator : MA/JU
Sample : P2548-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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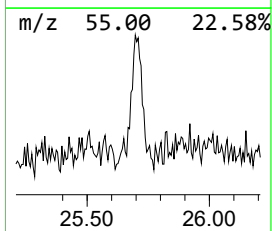
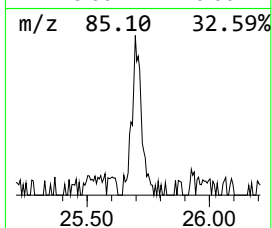
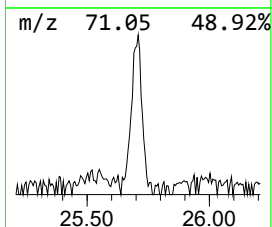
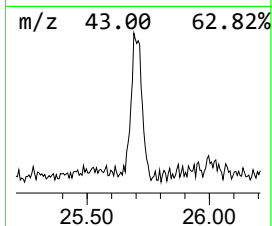
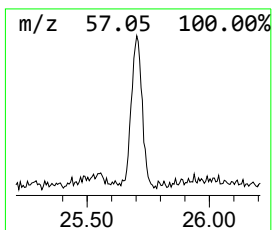
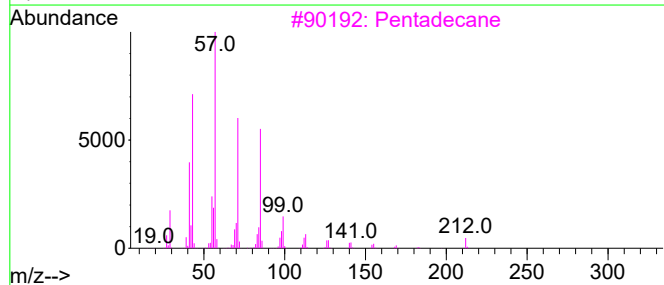
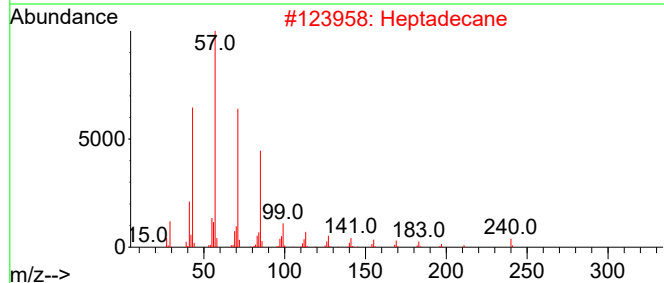
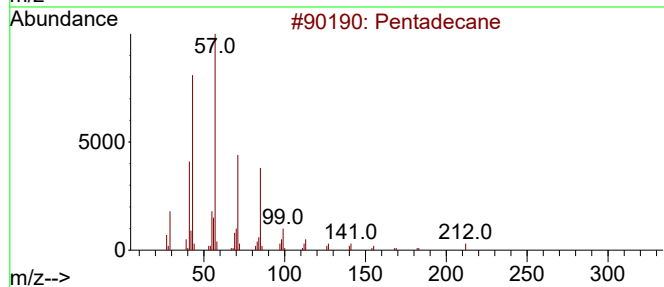
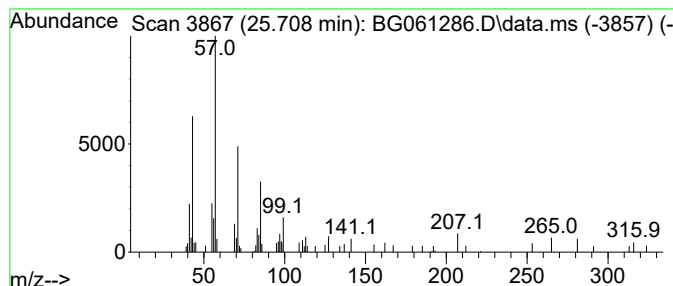
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.709	3.05 ng/ul	77927	Perylene-d12	24.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	74
2			Heptadecane	240	C17H36	000629-78-7	72
3			Pentadecane	212	C15H32	000629-62-9	70
4			Eicosane, 2-methyl-	296	C21H44	001560-84-5	68
5			Hexatriacontane	507	C36H74	000630-06-8	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061286.D
Acq On : 24 May 2024 13:49
Operator : MA/JU
Sample : P2548-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH612

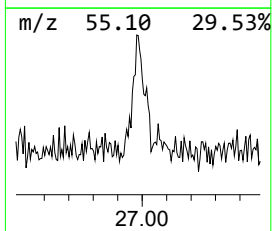
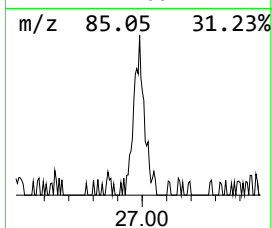
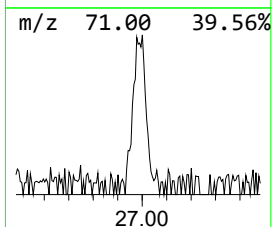
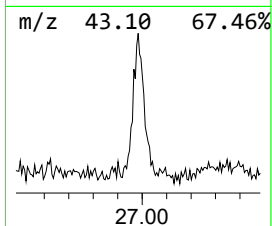
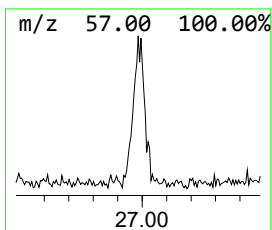
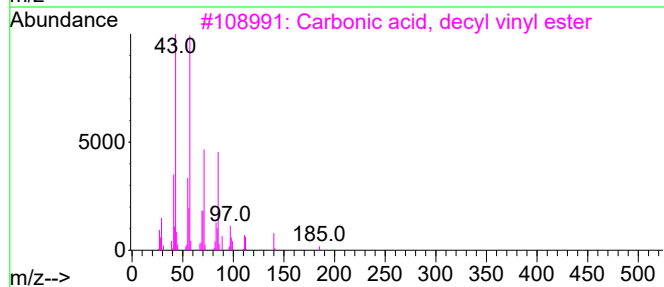
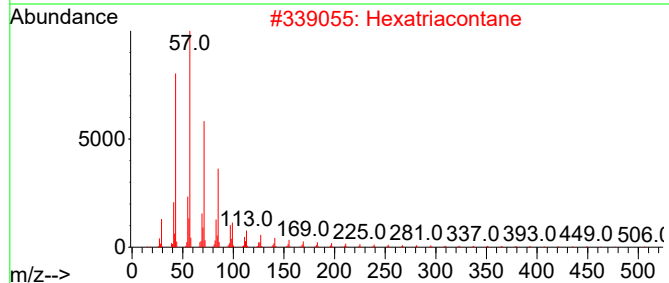
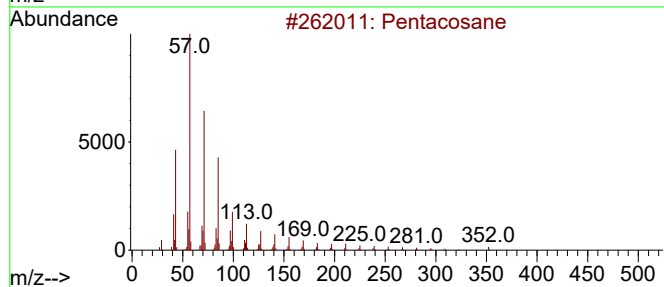
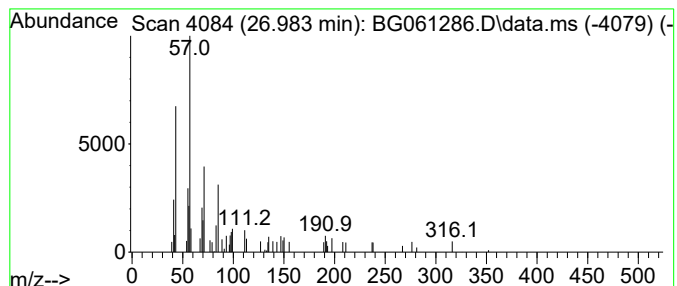
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.983	2.69 ng/ul	68629	Perylene-d12	24.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentacosane	352	C25H52	000629-99-2	52
2			Hexatriacontane	507	C36H74	000630-06-8	50
3			Carbonic acid, decyl vinyl ester	228	C13H24O3	1000383-25-7	50
4			Octacosane	394	C28H58	000630-02-4	50
5			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061286.D
Acq On : 24 May 2024 13:49
Operator : MA/JU
Sample : P2548-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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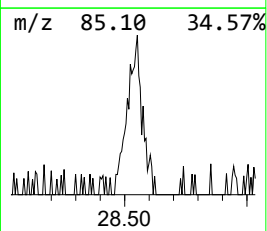
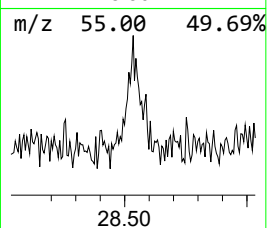
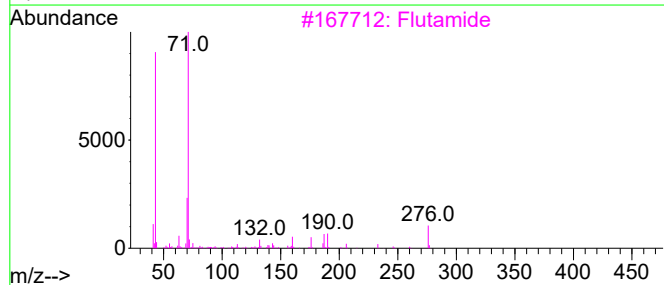
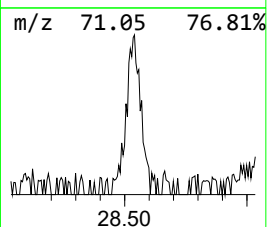
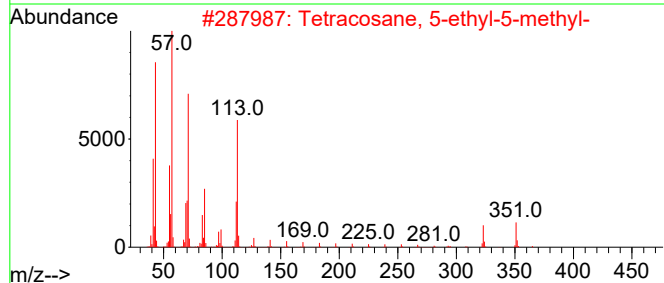
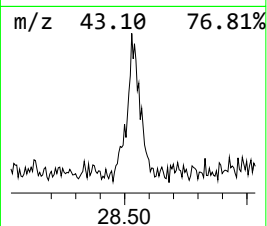
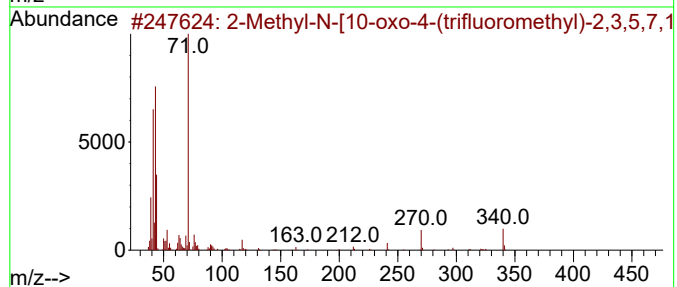
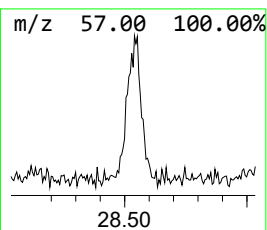
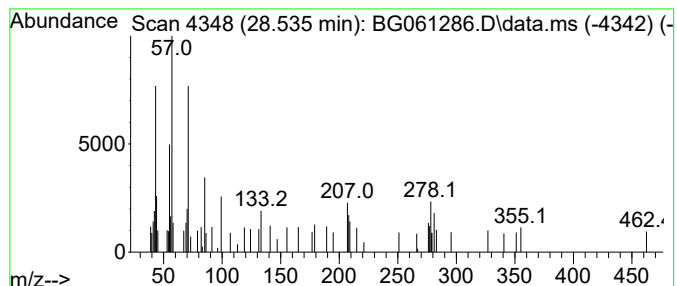
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.535	2.22 ng/ul	56711	Perylene-d12	24.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-N-[10-oxo-4-(trifluoromethyl)-2,3,5,7,1	340	C13H11F3N6O2	1000435-27-1	35		
2	Tetracosane, 5-ethyl-5-methyl-	380	C27H56	055282-30-9	32		
3	Flutamide	276	C11H11F3N2O3	013311-84-7	25		
4	Flutamide	276	C11H11F3N2O3	013311-84-7	25		
5	1,9,12,15-Octadecatetraene, 1-me...	276	C19H32O	056847-00-8	25		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061286.D
Acq On : 24 May 2024 13:49
Operator : MA/JU
Sample : P2548-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH612

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown-01	3.611	2.9	ng/ul	16517	1	7.888	114302	20.0
2-Phenanthrenol...	20.532	7.8	ng/ul	202387	5	21.519	517007	20.0
Octacosyl trifl...	21.184	3.8	ng/ul	98302	5	21.519	517007	20.0
(DEL) Alkane: S...	25.709	3.0	ng/ul	77927	6	24.604	510274	20.0
(DEL) Alkane: S...	26.983	2.7	ng/ul	68629	6	24.604	510274	20.0
unknown-02	28.535	2.2	ng/ul	56711	6	24.604	510274	20.0