

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
 Data File : BG061285.D
 Acq On : 24 May 2024 13:09
 Operator : MA/JU
 Sample : P2548-10
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BH5Y7

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: BG061285.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.078	10	15	39	rVB	1271503	1950059	100.00%	13.836%
2	3.748	124	129	138	rBV2	30496	51860	2.66%	0.368%
3	3.866	143	149	154	rVV2	7146	11663	0.60%	0.083%
4	4.306	220	224	231	rBV	16296	22882	1.17%	0.162%
5	4.718	288	294	299	rBV2	11010	15258	0.78%	0.108%
6	7.062	683	693	700	rBV	94808	160025	8.21%	1.135%
7	7.215	713	719	727	rVB	58822	96799	4.96%	0.687%
8	7.420	748	754	762	rBV	85426	141723	7.27%	1.006%
9	7.884	825	833	840	rBV	80191	130293	6.68%	0.924%
10	8.595	948	954	966	rVB	104334	179014	9.18%	1.270%
11	9.042	1022	1030	1039	rBV	92981	162971	8.36%	1.156%
12	9.759	1146	1152	1159	rBV	84228	148972	7.64%	1.057%
13	10.311	1239	1246	1255	rBV	114390	210280	10.78%	1.492%
14	10.681	1301	1309	1316	rBV	110642	197912	10.15%	1.404%
15	10.816	1325	1332	1340	rBV	88368	159365	8.17%	1.131%
16	13.924	1855	1861	1869	rVV	265280	365562	18.75%	2.594%
17	14.212	1903	1910	1919	rBV2	263500	402505	20.64%	2.856%
18	14.518	1956	1962	1968	rVB	213506	325569	16.70%	2.310%
19	14.741	1992	2000	2017	rBV2	77550	158470	8.13%	1.124%
20	14.923	2025	2031	2036	rVB4	4590	8972	0.46%	0.064%
21	15.511	2125	2131	2137	rBV	369582	540532	27.72%	3.835%
22	15.570	2137	2141	2145	rVB2	6796	9858	0.51%	0.070%
23	15.640	2148	2153	2165	rBV	121724	186223	9.55%	1.321%
24	16.921	2366	2371	2377	rBV2	8392	12996	0.67%	0.092%
25	17.085	2394	2399	2409	rVB2	5429	8924	0.46%	0.063%
26	17.268	2423	2430	2434	rBV	313269	463992	23.79%	3.292%
27	17.309	2434	2437	2441	rVV	147581	199913	10.25%	1.418%
28	17.367	2441	2447	2457	rVB	412852	645523	33.10%	4.580%
29	17.667	2488	2498	2503	rBV4	13498	24789	1.27%	0.176%
30	17.849	2526	2529	2539	rVB5	5448	9157	0.47%	0.065%
31	18.143	2570	2579	2582	rVV	23813	37658	1.93%	0.267%
32	18.190	2582	2587	2592	rVV	35550	60370	3.10%	0.428%
33	18.272	2597	2601	2606	rVV3	7866	13276	0.68%	0.094%
34	18.337	2606	2612	2616	rVV2	31128	60642	3.11%	0.430%
35	18.372	2616	2618	2623	rVB2	17633	23150	1.19%	0.164%
36	18.642	2659	2664	2668	rBV2	19417	26274	1.35%	0.186%

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

37	18.684	2668	2671	2677	rVB2	21471	29801	1.53%	0.211%
38	18.889	2702	2706	2711	rVB2	7568	8046	0.41%	0.057%
39	18.942	2711	2715	2718	rBV2	7671	8701	0.45%	0.062%
40	19.060	2729	2735	2739	rBV3	18231	33720	1.73%	0.239%
41	19.201	2755	2759	2767	rVB5	17897	38268	1.96%	0.272%
42	19.324	2774	2780	2787	rVB	313430	428889	21.99%	3.043%
43	19.418	2793	2796	2800	rVB3	7443	8740	0.45%	0.062%
44	19.524	2809	2814	2816	rBV4	7881	12285	0.63%	0.087%
45	19.565	2818	2821	2823	rBV2	8187	8925	0.46%	0.063%
46	19.659	2831	2837	2840	rBV	555864	868870	44.56%	6.165%
47	19.682	2840	2841	2849	rVB	267834	290111	14.88%	2.058%
48	19.753	2849	2853	2857	rBV4	15432	25506	1.31%	0.181%
49	19.794	2857	2860	2862	rVB2	9836	8881	0.46%	0.063%
50	19.865	2865	2872	2877	rBV2	16234	30805	1.58%	0.219%
51	19.923	2877	2882	2888	rVB2	14481	26841	1.38%	0.190%
52	20.017	2888	2898	2903	rBV3	23690	66028	3.39%	0.468%
53	20.147	2914	2920	2922	rVV7	15902	31206	1.60%	0.221%
54	20.176	2922	2925	2932	rVB2	42391	61020	3.13%	0.433%
55	20.282	2938	2943	2946	rBV	18752	28104	1.44%	0.199%
56	20.335	2946	2952	2956	rVV2	30564	54594	2.80%	0.387%
57	20.376	2956	2959	2962	rVB5	9089	11356	0.58%	0.081%
58	20.411	2963	2965	2970	rVV6	6607	7917	0.41%	0.056%
59	20.476	2972	2976	2980	rBV	21881	31028	1.59%	0.220%
60	20.523	2980	2984	2987	rVV3	19875	29297	1.50%	0.208%
61	20.575	2990	2993	2996	rVV2	21858	27761	1.42%	0.197%
62	20.628	2999	3002	3007	rVB7	12222	17858	0.92%	0.127%
63	20.693	3009	3013	3015	rVB4	9510	12195	0.63%	0.087%
64	20.734	3015	3020	3023	rBV6	17746	26807	1.37%	0.190%
65	20.775	3023	3027	3029	rVV5	10583	15135	0.78%	0.107%
66	20.799	3029	3031	3036	rVB5	6596	9006	0.46%	0.064%
67	20.840	3036	3038	3043	rBV6	6141	9336	0.48%	0.066%
68	20.934	3049	3054	3060	rVB	34637	58653	3.01%	0.416%
69	21.110	3076	3084	3087	rBV3	33181	76688	3.93%	0.544%
70	21.151	3088	3091	3094	rVV	33302	52687	2.70%	0.374%
71	21.192	3094	3098	3104	rVV4	52493	146244	7.50%	1.038%
72	21.257	3106	3109	3115	rVB	41881	70107	3.60%	0.497%
73	21.416	3132	3136	3140	rVB	55588	72570	3.72%	0.515%
74	21.516	3145	3153	3158	rVV3	373012	800022	41.03%	5.676%
75	21.563	3158	3161	3168	rVB	132866	202169	10.37%	1.434%
76	21.709	3182	3186	3192	rBV5	22383	34550	1.77%	0.245%

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77	21.909	3215	3220	3227	rBV3	22096	44631	2.29%	0.317%
78	21.968	3227	3230	3237	rVV9	9031	15340	0.79%	0.109%
79	22.221	3268	3273	3278	rVB	25655	51962	2.66%	0.369%
80	22.291	3281	3285	3291	rVV5	35037	66208	3.40%	0.470%
81	22.485	3313	3318	3323	rBV7	15002	28262	1.45%	0.201%
82	22.937	3388	3395	3406	rVB6	49012	129807	6.66%	0.921%
83	23.137	3423	3429	3434	rVB7	15657	29983	1.54%	0.213%
84	23.249	3444	3448	3449	rBV4	7596	9452	0.48%	0.067%
85	23.625	3505	3512	3520	rBV3	99816	318910	16.35%	2.263%
86	23.872	3548	3554	3561	rBV3	14478	34849	1.79%	0.247%
87	24.066	3583	3587	3588	rBV4	8532	11467	0.59%	0.081%
88	24.318	3624	3630	3635	rBV2	48910	113994	5.85%	0.809%
89	24.395	3635	3643	3650	rVV2	268402	720764	36.96%	5.114%
90	24.453	3650	3653	3666	rVB	89900	225470	11.56%	1.600%
91	24.606	3669	3679	3688	rBV	202441	581572	29.82%	4.126%
92	25.699	3860	3865	3874	rVB6	24458	66959	3.43%	0.475%
93	25.846	3886	3890	3898	rVB10	14921	35627	1.83%	0.253%
94	26.063	3919	3927	3937	rVB3	52828	167717	8.60%	1.190%
95	26.991	4080	4085	4091	rVB8	11081	18967	0.97%	0.135%
96	27.144	4104	4111	4124	rVB8	27550	90515	4.64%	0.642%
97	27.661	4188	4199	4200	rBV8	7379	20049	1.03%	0.142%
98	28.096	4263	4273	4284	rVV4	36558	153622	7.88%	1.090%
99	28.531	4339	4347	4348	rBV6	17059	31647	1.62%	0.225%
100	29.189	4448	4459	4472	rVB3	28128	122441	6.28%	0.869%

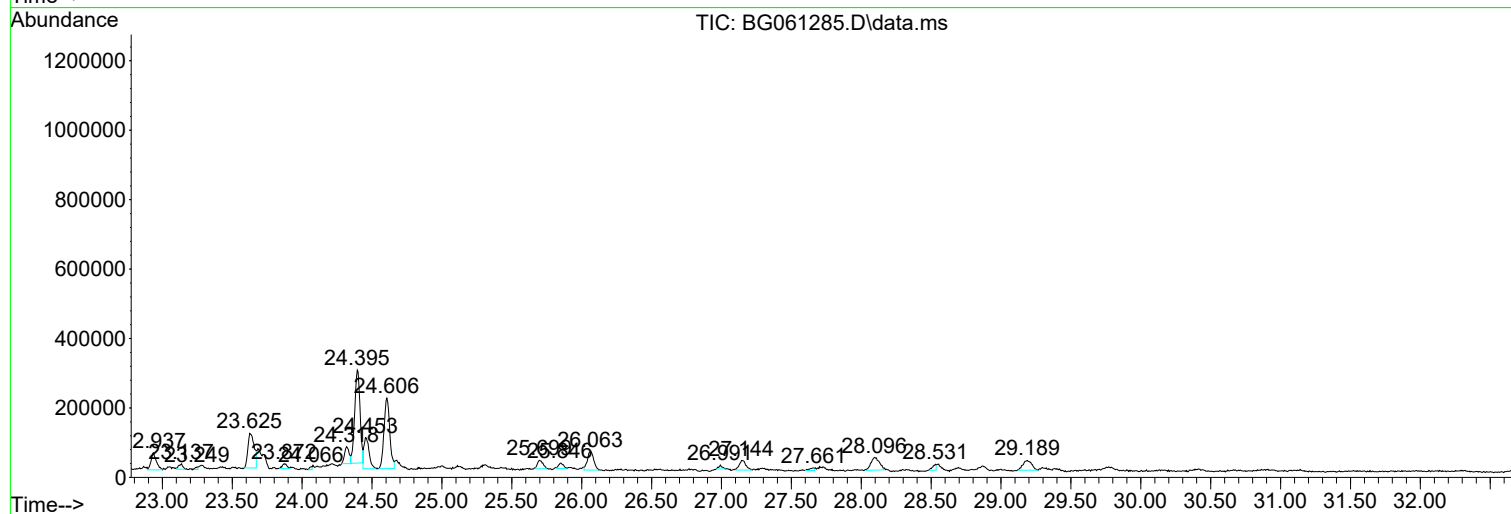
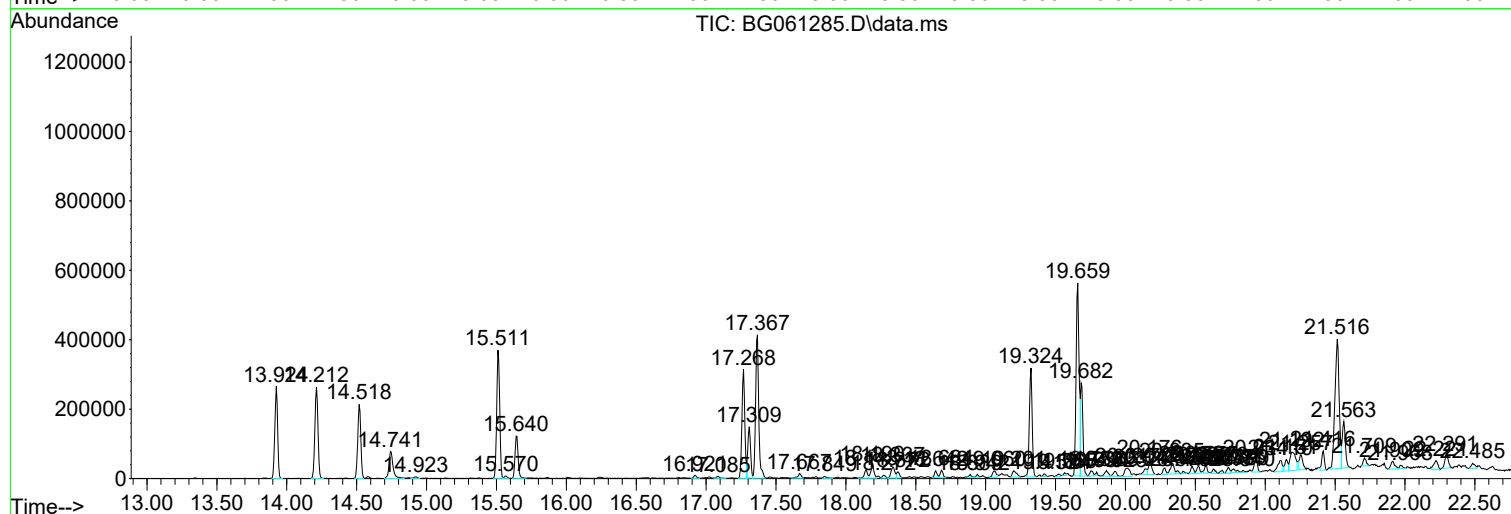
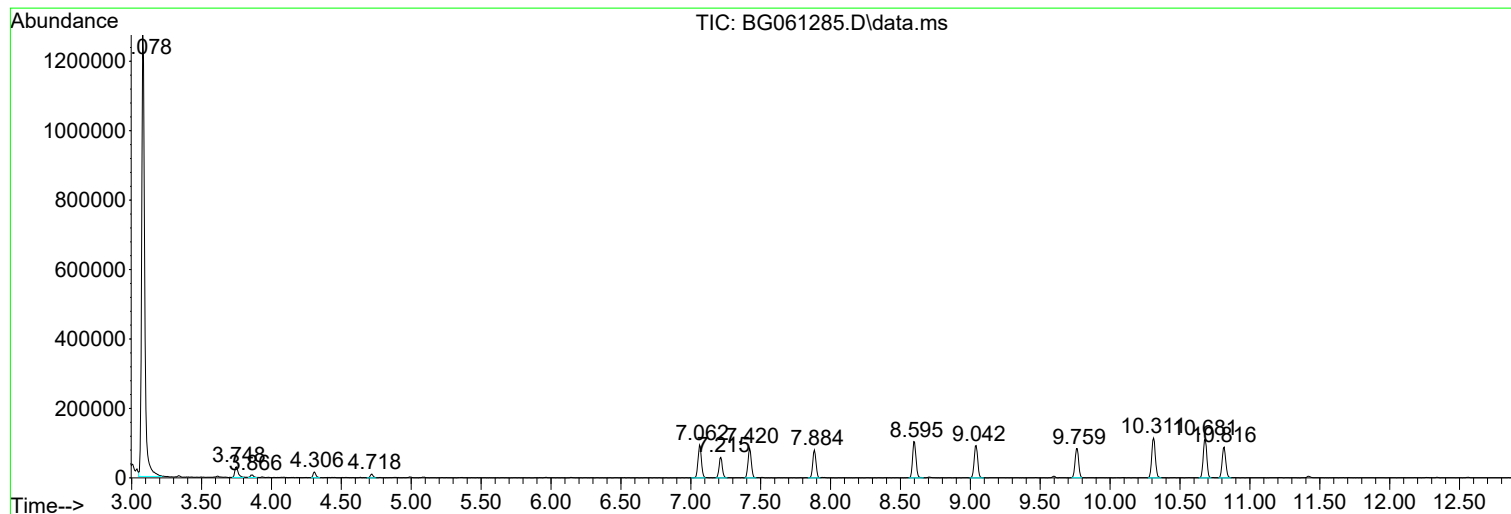
Sum of corrected areas: 14094373

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ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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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



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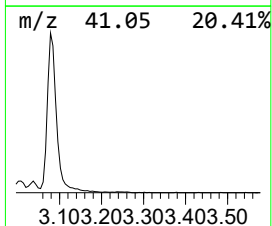
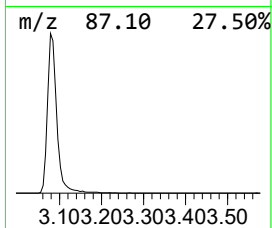
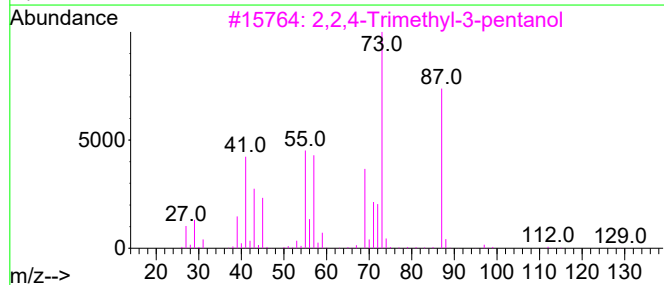
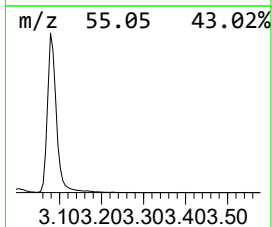
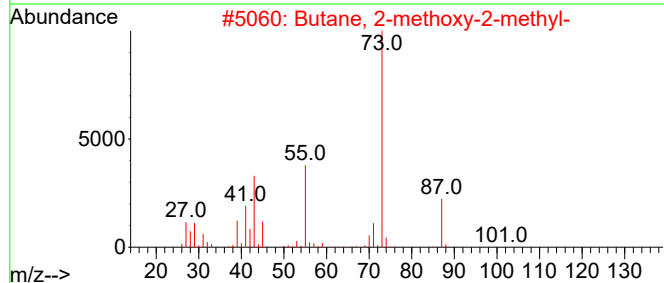
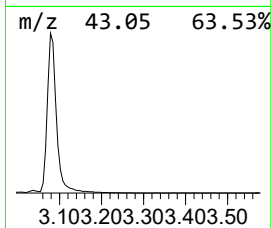
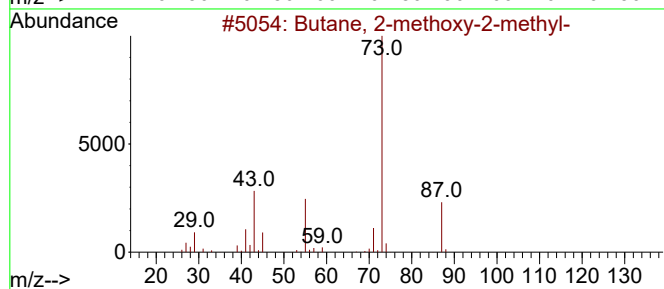
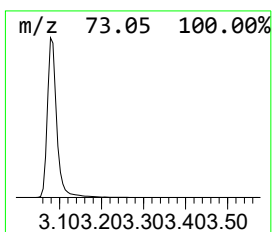
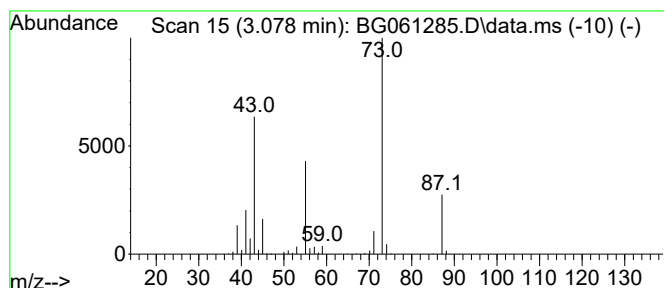
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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.
3.078	299.33 ng/ul	1950060	1,4-Dichlorobenzene-d4	7.884

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	32
5			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	25



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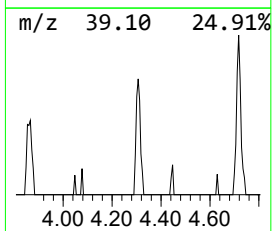
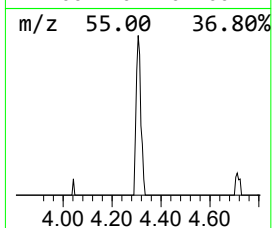
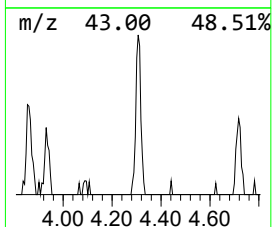
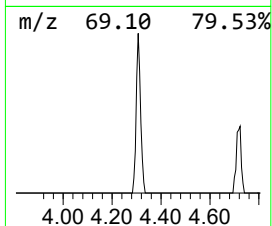
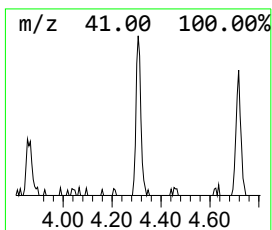
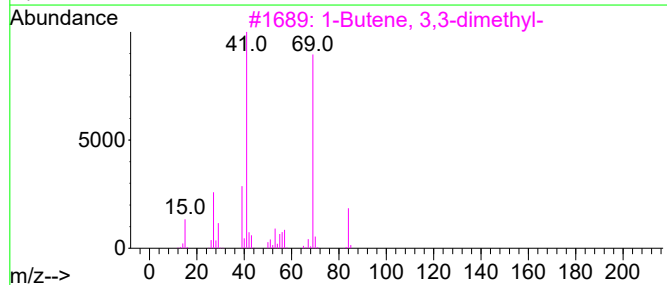
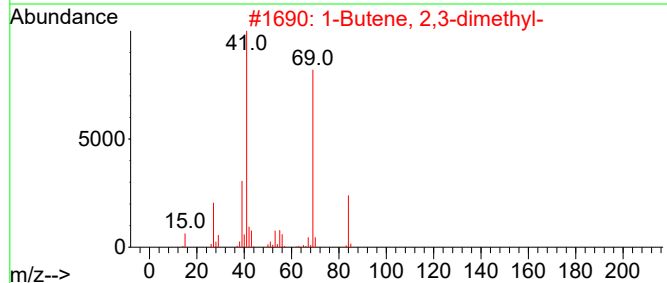
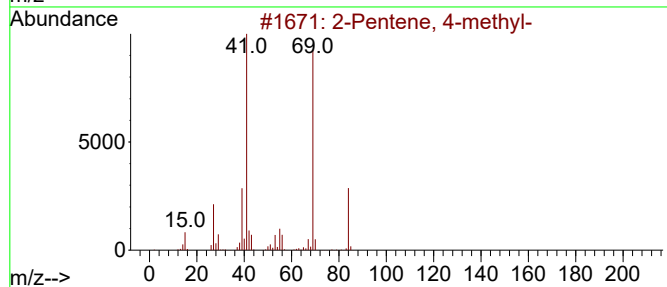
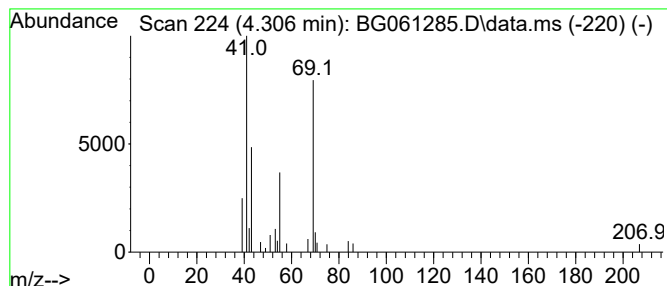
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	3.51 ng/ul	22882	1,4-Dichlorobenzene-d4	7.884

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentene, 4-methyl-	84	C6H12	004461-48-7	72
2		1-Butene, 2,3-dimethyl-	84	C6H12	000563-78-0	72
3		1-Butene, 3,3-dimethyl-	84	C6H12	000558-37-2	72
4		Butane, 2,3-dimethyl-2-nitro-	131	C6H13NO2	034075-28-0	56
5		2-Pentene, 3-methyl-, (E)-	84	C6H12	000616-12-6	50



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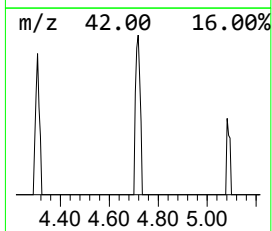
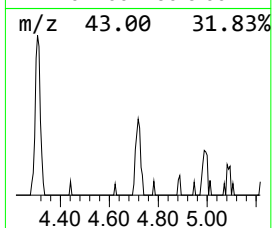
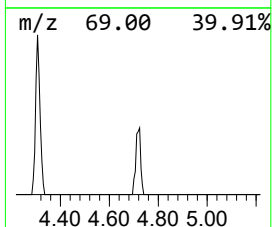
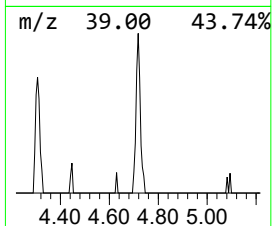
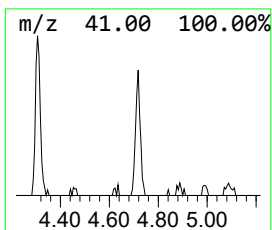
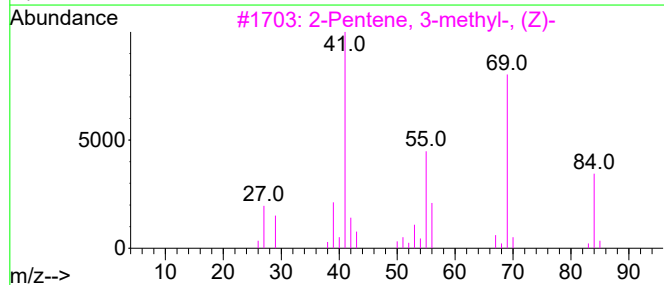
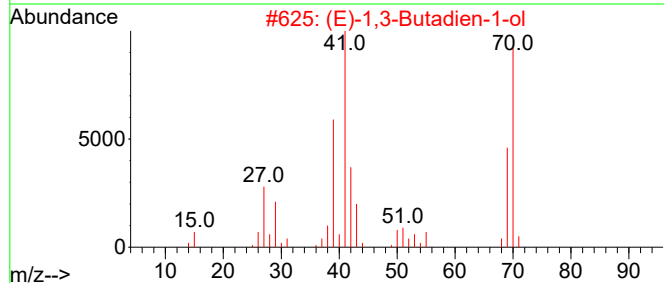
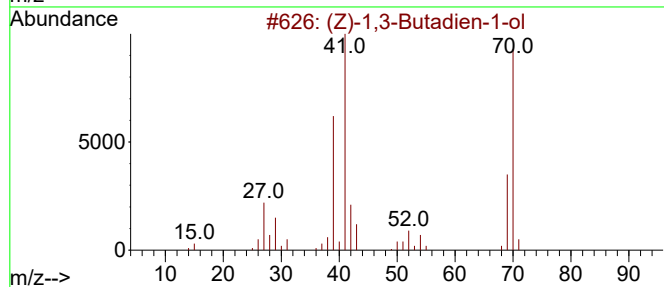
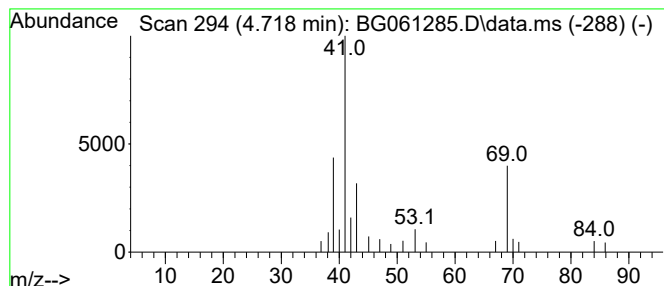
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.718	2.34 ng/ul	15258	1,4-Dichlorobenzene-d4	7.884

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(Z)-1,3-Butadien-1-ol			70	C4H6O	070415-58-6	36
2	(E)-1,3-Butadien-1-ol			70	C4H6O	070411-98-2	36
3	2-Pentene, 3-methyl-, (Z)-			84	C6H12	000922-62-3	35
4	3-Penten-2-one			84	C5H8O	000625-33-2	33
5	2-Pentene, 3-methyl-, (Z)-			84	C6H12	000922-62-3	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061285.D
Acq On : 24 May 2024 13:09
Operator : MA/JU
Sample : P2548-10
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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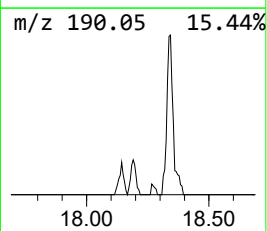
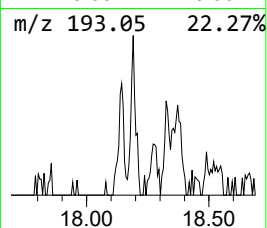
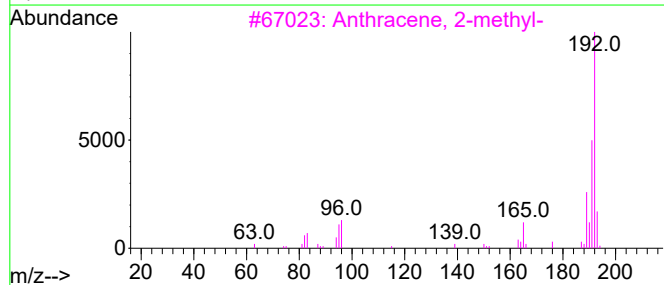
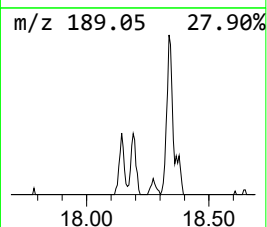
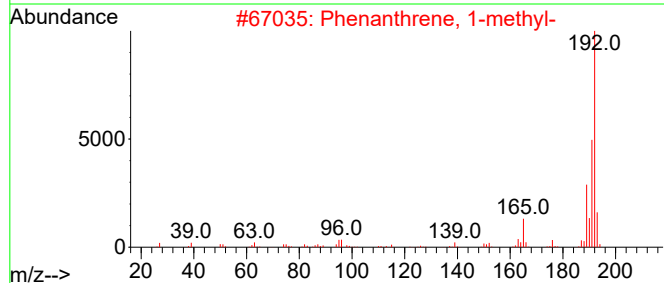
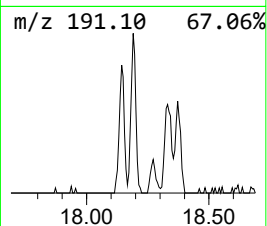
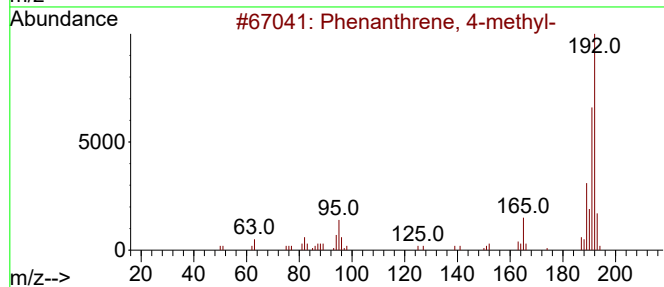
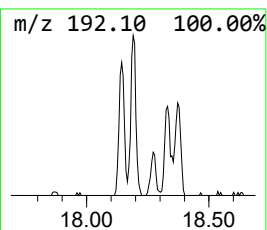
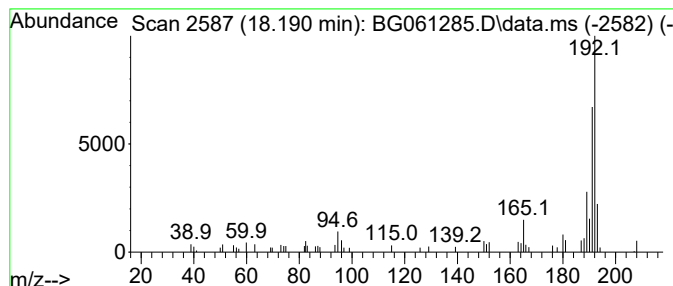
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenanthrene, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.190	2.60 ng/ul	60370	Phenanthrene-d10	17.268

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061285.D
Acq On : 24 May 2024 13:09
Operator : MA/JU
Sample : P2548-10
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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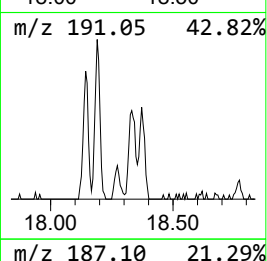
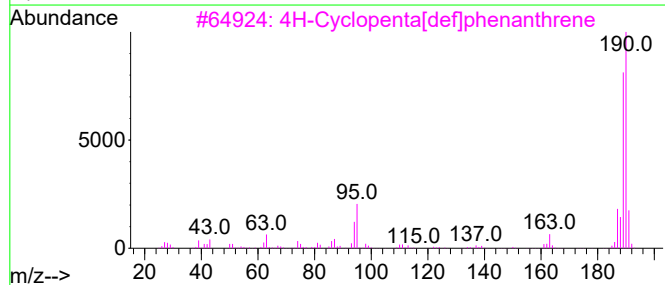
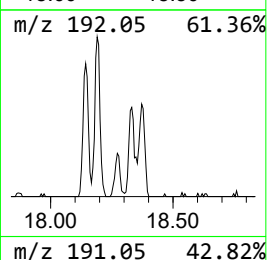
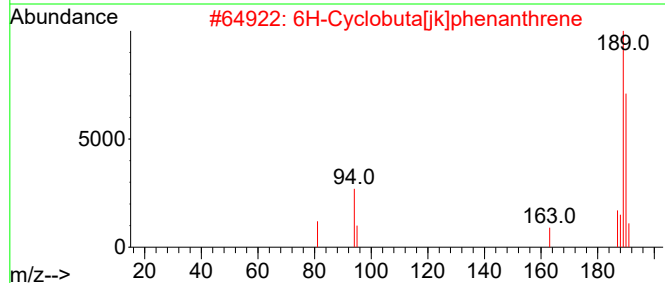
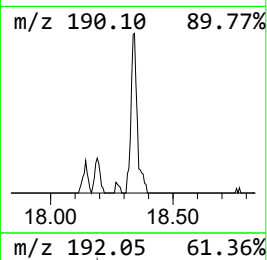
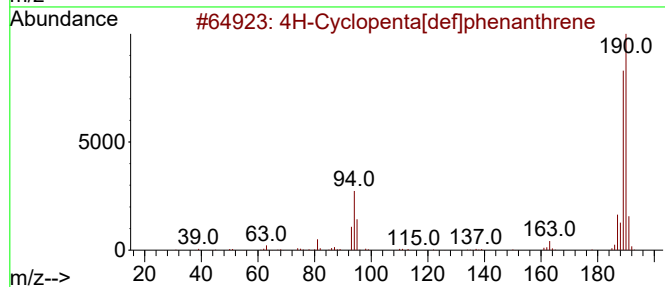
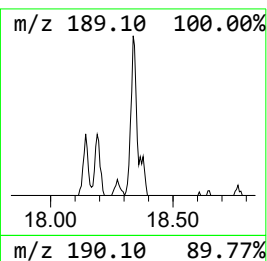
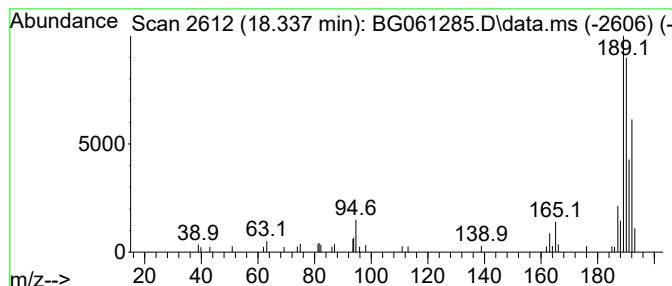
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.337	2.61 ng/ul	60642	Phenanthrene-d10	17.268

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	53
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
4			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	43
5			Methyl diselenide	190	C2H6Se2	007101-31-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061285.D
Acq On : 24 May 2024 13:09
Operator : MA/JU
Sample : P2548-10
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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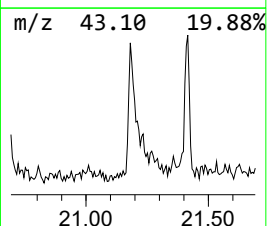
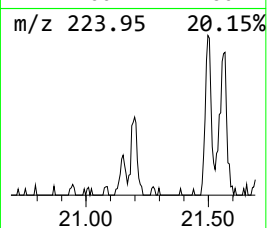
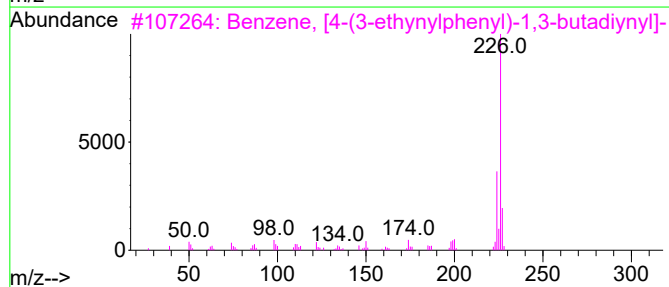
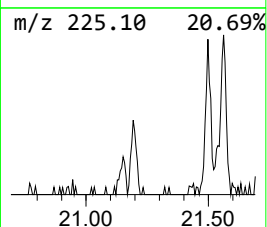
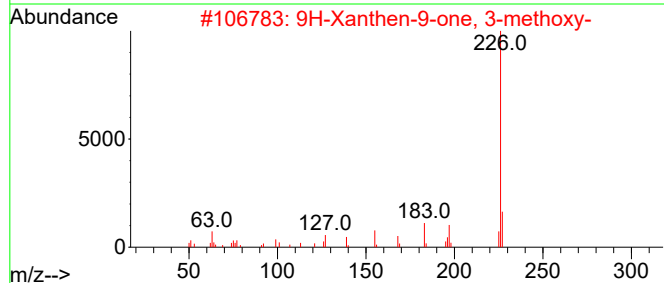
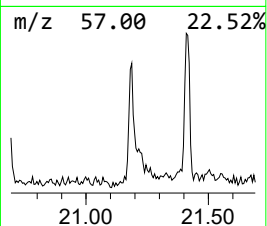
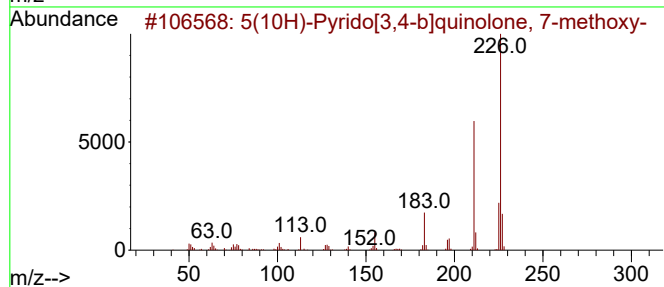
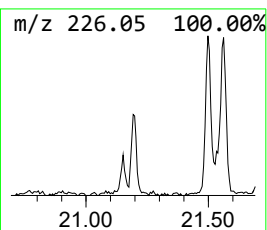
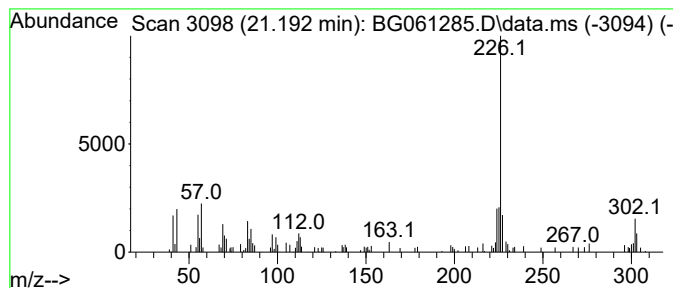
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 5(10H)-Pyrido[3,4-b]quinolo... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.192	3.66 ng/ul	146244	Chrysene-d12	21.515

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5(10H)-Pyrido[3,4-b]quinolone, 7...	226	C13H10N2O2	1000256-99-5	52
2			9H-Xanthen-9-one, 3-methoxy-	226	C14H10O3	003722-52-9	46
3			Benzene, [4-(3-ethynylphenyl)-1,...	226	C18H10	1000115-87-3	43
4			Silane, dimethyl(4-methoxyphenox...	226	C11H18O3Si	1000347-13-9	43
5			9H-Xanthen-9-one, 2-methoxy-	226	C14H10O3	001214-20-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061285.D
Acq On : 24 May 2024 13:09
Operator : MA/JU
Sample : P2548-10
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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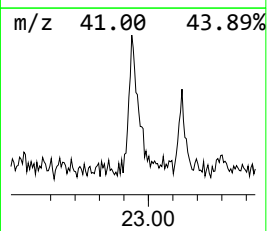
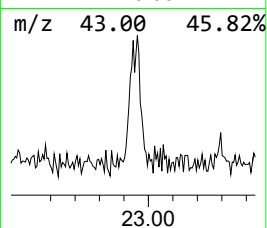
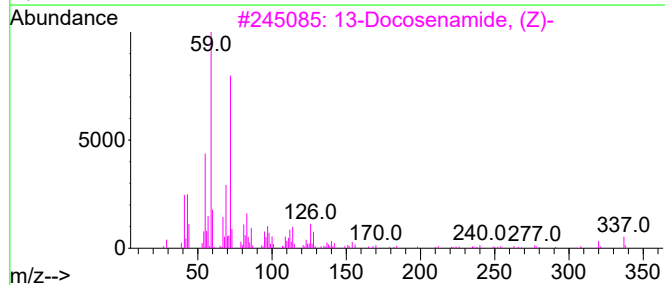
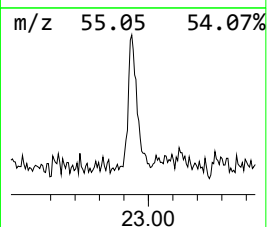
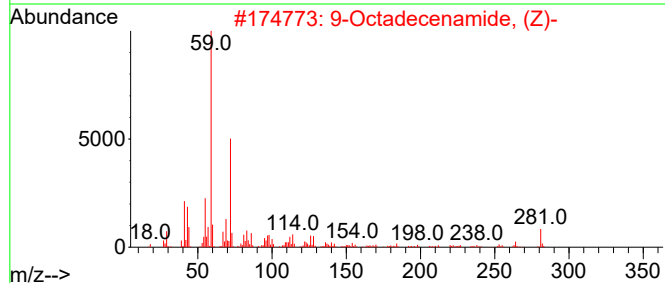
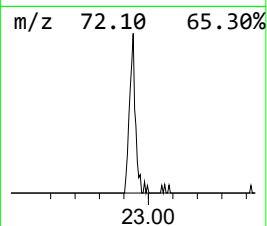
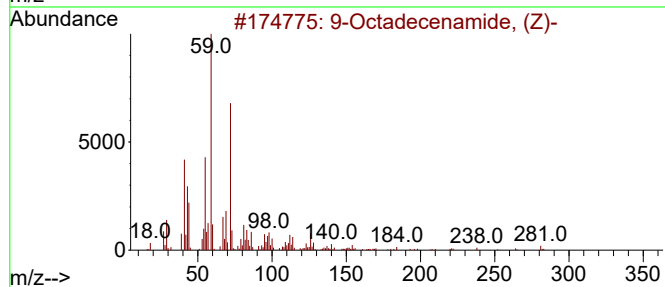
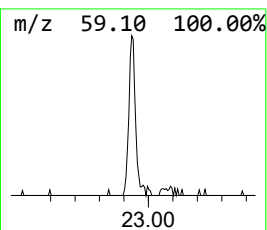
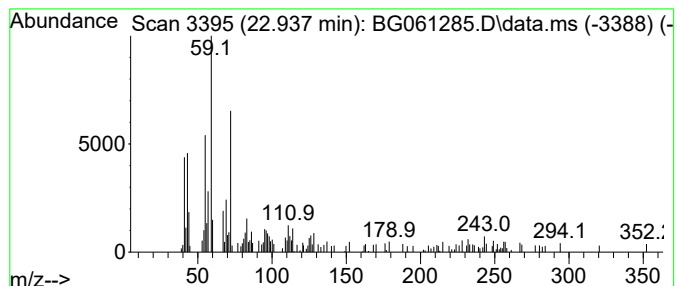
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9-Octadecenamide, (Z)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.937	3.25 ng/ul	129807	Chrysene-d12	21.515

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	68
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	64
3			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	64
4			Pentanal, oxime	101	C5H11NO	000628-79-5	47
5			1,1,2-Trimethyl-1-silacyclobutane	114	C6H14Si	030681-90-4	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061285.D
Acq On : 24 May 2024 13:09
Operator : MA/JU
Sample : P2548-10
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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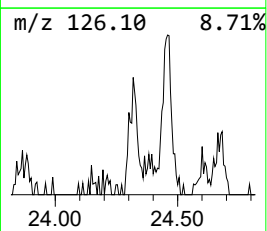
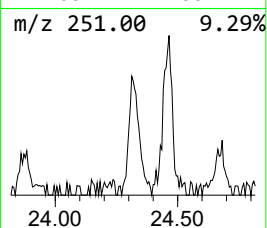
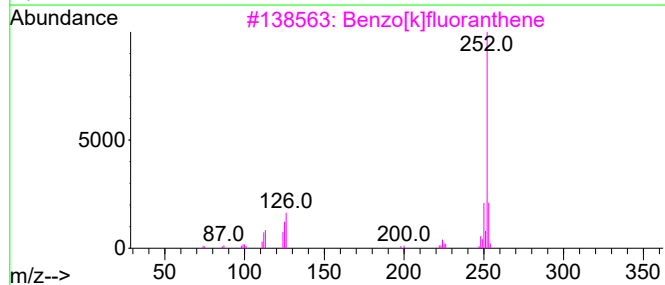
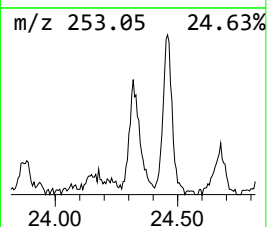
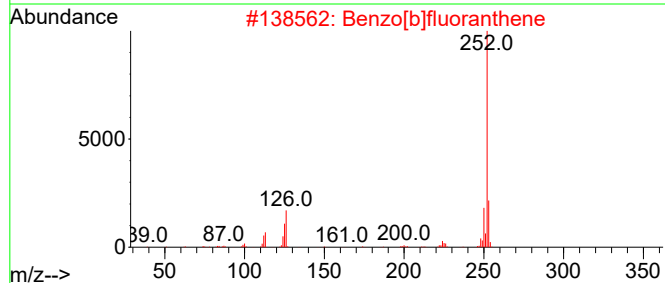
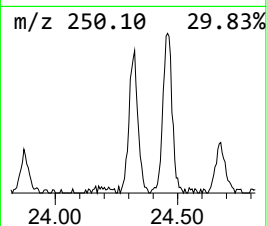
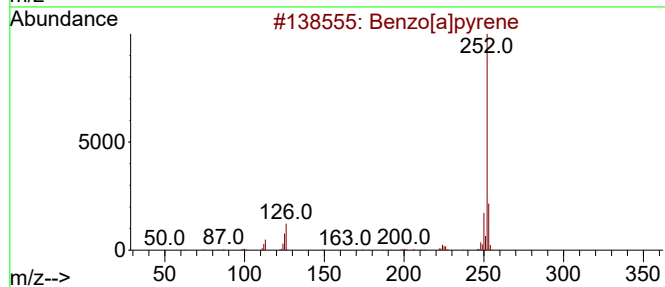
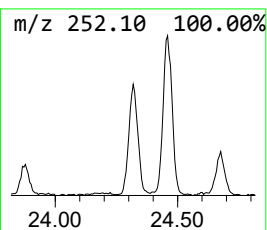
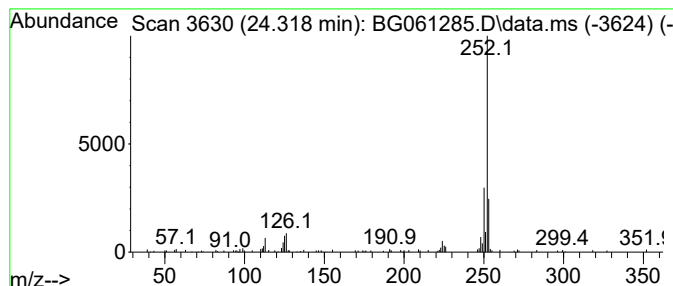
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.318	3.92 ng/ul	113994	Perylene-d12	24.606

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[a]pyrene	252	C20H12	000050-32-8	90
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	90
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	89
4			Benzo[a]pyrene	252	C20H12	000050-32-8	89
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061285.D
Acq On : 24 May 2024 13:09
Operator : MA/JU
Sample : P2548-10
Misc :
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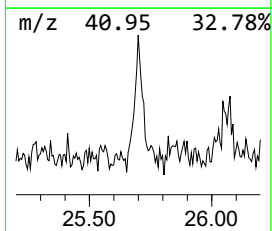
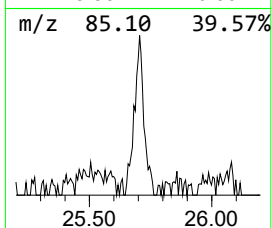
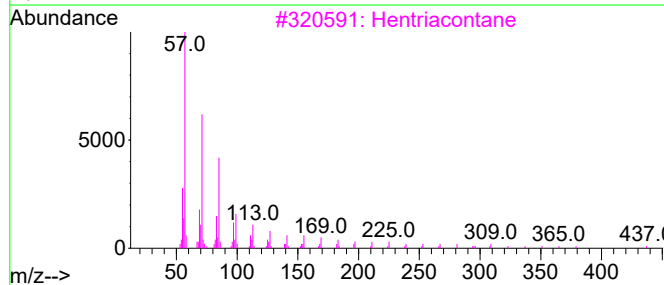
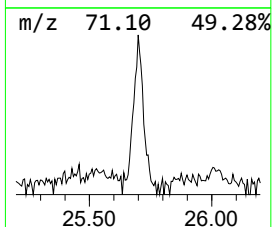
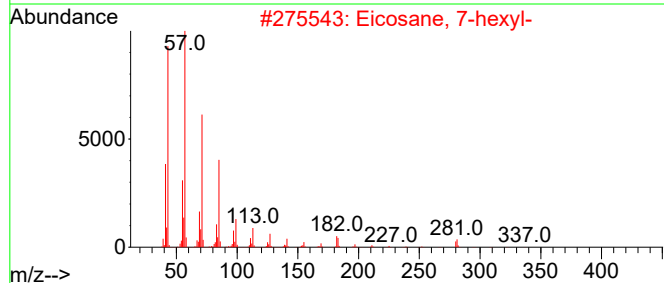
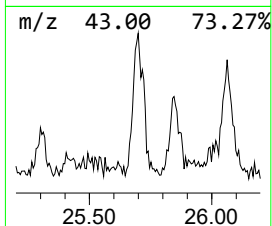
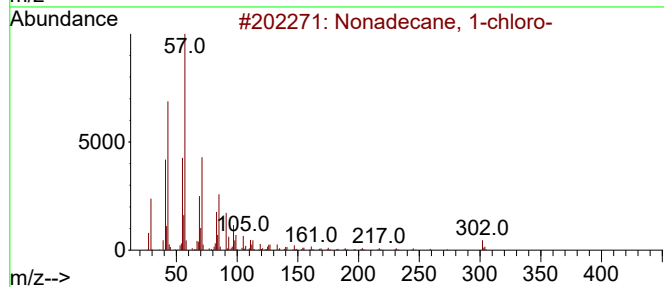
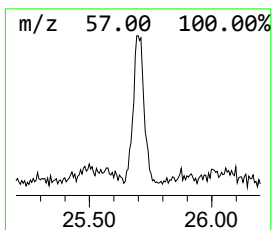
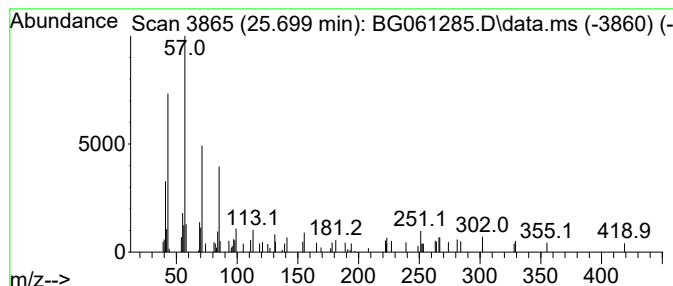
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 Nonadecane, 1-chloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
25.699	2.30 ng/ul	66959	Perylene-d12	24.606	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	64
2	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	52
3	Hentriacontane	437	C31H64	000630-04-6	52
4	Octadecane, 3-ethyl-5-(2-ethylbu...	366	C26H54	055282-12-7	52
5	10-Methylnonadecane	282	C20H42	056862-62-5	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061285.D
Acq On : 24 May 2024 13:09
Operator : MA/JU
Sample : P2548-10
Misc :
ALS Vial : 9 Sample Multiplier: 1

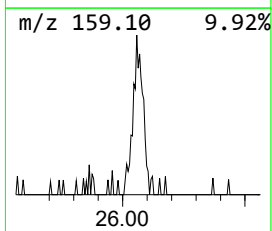
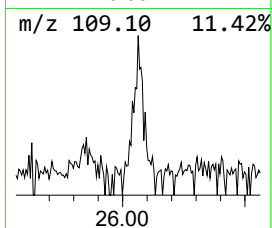
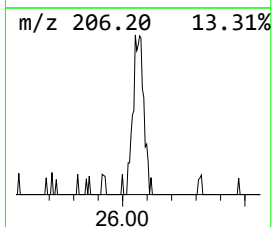
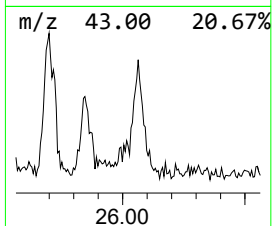
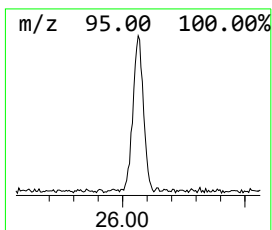
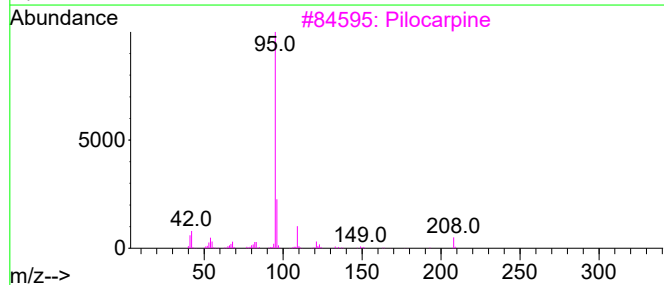
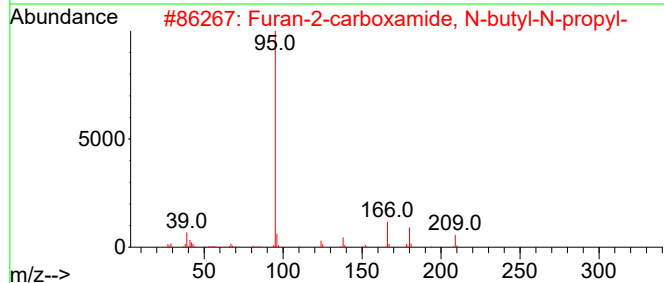
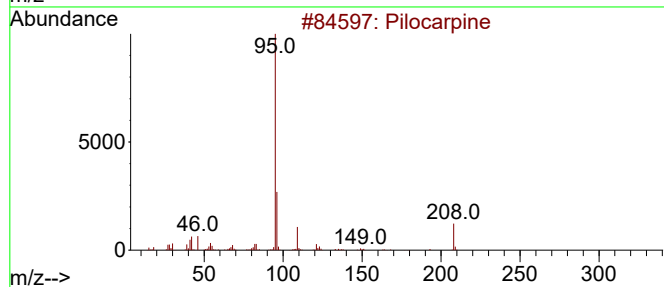
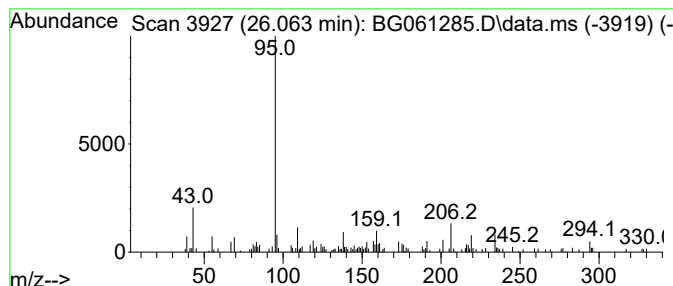
Instrument :
BNA_G
ClientSampleId :
BH5Y7

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 10 Pilocarpine Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.063	5.77 ng/ul	167717	Perylene-d12	24.606	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pilocarpine	208	C11H16N2O2	000092-13-7	50
2	Furan-2-carboxamide, N-butyl-N-p...	209	C12H19NO2	1000415-64-1	50
3	Pilocarpine	208	C11H16N2O2	000092-13-7	50
4	Pilocarpine	208	C11H16N2O2	000092-13-7	50
5	4-Fluoro-N-(2-fluorophenyl)benze...	365	C14H8F5NO3S	1000487-94-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061285.D
Acq On : 24 May 2024 13:09
Operator : MA/JU
Sample : P2548-10
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH5Y7

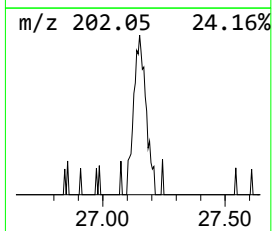
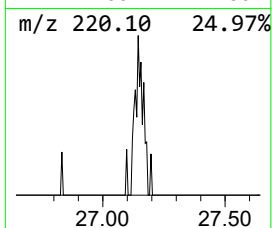
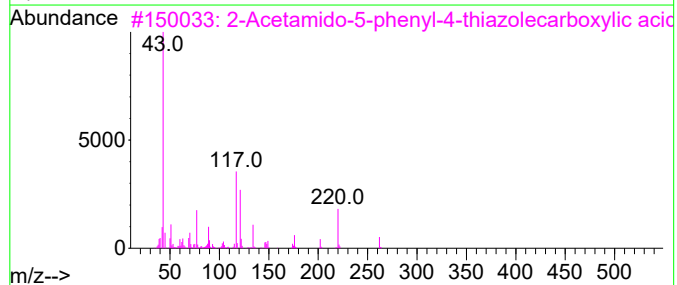
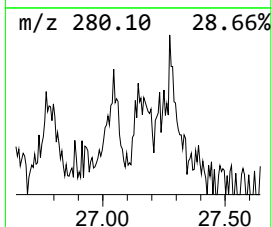
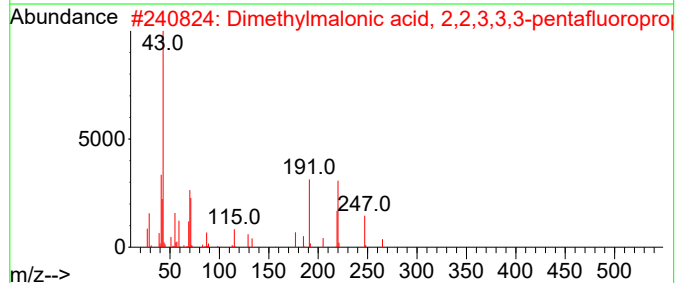
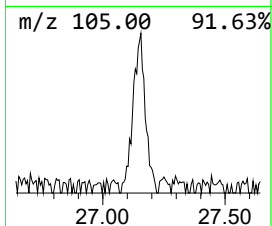
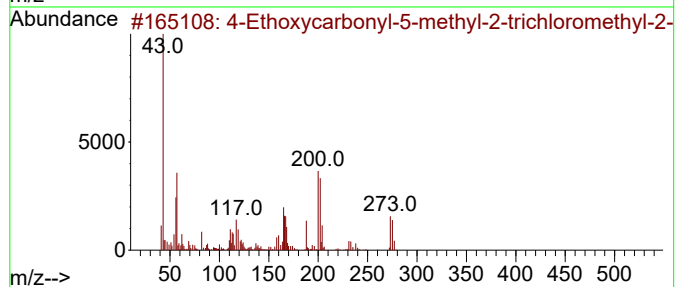
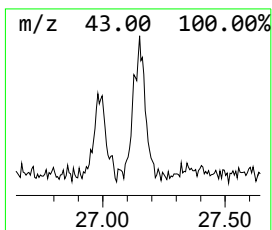
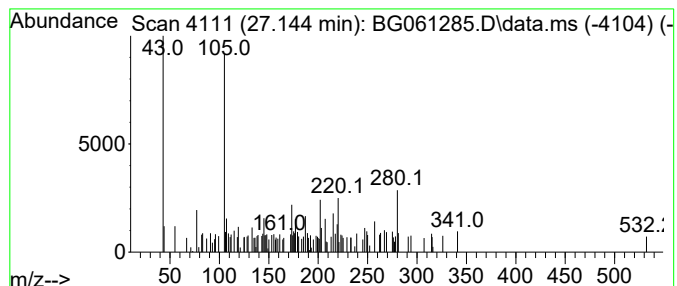
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-03 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.144	3.11 ng/ul	90515	Perylene-d12	24.606

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Ethoxycarbonyl-5-methyl-2-tric...	273	C8H10Cl3NO3	1000251-31-3	10		
2	Dimethylmalonic acid, 2,2,3,3,3-...	334	C13H19F5O4	1000361-93-8	10		
3	2-Acetamido-5-phenyl-4-thiazolec...	262	C12H10N2O3S	1000224-27-1	10		
4	4-Acetyloxyimino-6,6-dimethyl-3-...	341	C15H19NO4S2	1000301-45-7	10		
5	9.beta.-Acetoxy-3.beta.-hydroxy-...	280	C17H28O3	1000140-72-2	10		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052424\
Data File : BG061285.D
Acq On : 24 May 2024 13:09
Operator : MA/JU
Sample : P2548-10
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH5Y7

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
Butane, 2-metho...	3.078	299.3	ng/ul	1950060	1	7.884	130293	20.0
(DEL) Alkane: S...	4.306	3.5	ng/ul	22882	1	7.884	130293	20.0
unknown-01	4.718	2.3	ng/ul	15258	1	7.884	130293	20.0
Phenanthrene, 4...	18.190	2.6	ng/ul	60370	4	17.268	463992	20.0
4H-Cyclopenta[d...	18.337	2.6	ng/ul	60642	4	17.268	463992	20.0
5(10H)-Pyrido[3...	21.192	3.7	ng/ul	146244	5	21.515	800022	20.0
9-Octadecenamid...	22.937	3.3	ng/ul	129807	5	21.515	800022	20.0
unknown-02	24.318	3.9	ng/ul	113994	6	24.606	581572	20.0
Nonadecane, 1-c...	25.699	2.3	ng/ul	66959	6	24.606	581572	20.0
Pilocarpine	26.063	5.8	ng/ul	167717	6	24.606	581572	20.0
unknown-03	27.144	3.1	ng/ul	90515	6	24.606	581572	20.0