

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015556.D
 Acq On : 15 Jun 2023 15:19
 Operator : MA/JU
 Sample : 03088-13
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFR9

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_P\Methods\SFAM-EPA-BP060723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015556.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.417	204	209	213	rBV	54333	67454	0.16%	0.083%
2	7.246	683	690	704	rBV	196335	395337	0.96%	0.486%
3	7.410	712	718	726	rBV	178199	286114	0.69%	0.352%
4	7.610	744	752	762	rBV	230046	398150	0.97%	0.489%
5	8.087	822	833	848	rBV	530738	858644	2.08%	1.055%
6	8.805	949	955	967	rBV	244969	440797	1.07%	0.542%
7	8.922	969	975	986	rVB	146823	251624	0.61%	0.309%
8	9.263	1025	1033	1048	rBV	228853	414044	1.00%	0.509%
9	9.993	1150	1157	1169	rBV	205218	360070	0.87%	0.442%
10	10.540	1244	1250	1271	rBV	253084	519990	1.26%	0.639%
11	10.916	1307	1314	1320	rBV	777732	1307693	3.17%	1.607%
12	10.963	1320	1322	1332	rVB2	27888	42690	0.10%	0.052%
13	11.063	1332	1339	1354	rBV	118488	266771	0.65%	0.328%
14	14.134	1854	1861	1867	rBV	634352	877825	2.13%	1.079%
15	14.428	1903	1911	1927	rBV	585762	976542	2.37%	1.200%
16	14.728	1951	1962	1970	rBV	1212438	1849409	4.49%	2.273%
17	14.957	1995	2001	2018	rBV2	73510	228240	0.55%	0.280%
18	15.134	2026	2031	2039	rVB3	29922	52644	0.13%	0.065%
19	15.722	2124	2131	2138	rBV	858922	1274395	3.09%	1.566%
20	15.851	2147	2153	2163	rBV	204183	323977	0.79%	0.398%
21	16.086	2187	2193	2203	rVB2	84485	131471	0.32%	0.162%
22	16.651	2285	2289	2300	rVB4	26133	51515	0.12%	0.063%
23	17.357	2405	2409	2417	rBV	57768	90827	0.22%	0.112%
24	17.428	2417	2421	2425	rBV4	35730	46747	0.11%	0.057%
25	17.486	2425	2431	2436	rVV	1602843	2456498	5.96%	3.018%
26	17.528	2436	2438	2443	rVV	230374	343353	0.83%	0.422%
27	17.586	2443	2448	2461	rVV2	794192	1289233	3.13%	1.584%
28	17.828	2485	2489	2493	rBV3	36334	47331	0.11%	0.058%
29	18.280	2561	2566	2572	rBV2	214166	339506	0.82%	0.417%
30	18.345	2572	2577	2583	rBV	1076761	1351614	3.28%	1.661%
31	18.533	2605	2609	2613	rBV5	47230	78358	0.19%	0.096%
32	18.628	2620	2625	2632	rBV4	46135	110477	0.27%	0.136%
33	18.828	2655	2659	2661	rBV2	32210	50420	0.12%	0.062%
34	18.875	2665	2667	2670	rVV3	44644	56026	0.14%	0.069%
35	18.910	2670	2673	2677	rVB	47986	62935	0.15%	0.077%
36	18.980	2683	2685	2690	rBV5	25420	45482	0.11%	0.056%

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37	19.227	2723	2727	2731	rBV2	60891	83244	0.20%	0.102%
38	19.433	2758	2762	2764	rBV	244517	310433	0.75%	0.381%
39	19.480	2764	2770	2778	rVB	691320	1115414	2.71%	1.371%
40	19.569	2781	2785	2792	rBV	394073	471662	1.14%	0.580%
41	19.810	2819	2826	2829	rBV	537372	813501	1.97%	1.000%
42	20.304	2904	2910	2918	rVB2	132258	280747	0.68%	0.345%
43	20.416	2926	2929	2932	rBV2	47930	54668	0.13%	0.067%
44	20.616	2960	2963	2969	rBV3	99889	167303	0.41%	0.206%
45	20.780	2988	2991	2995	rBV	97362	119748	0.29%	0.147%
46	21.051	3034	3037	3040	rBV	69560	85425	0.21%	0.105%
47	21.245	3066	3070	3076	rVB2	481870	707495	1.72%	0.869%
48	21.569	3118	3125	3129	rBV2	1956813	3169163	7.69%	3.894%
49	21.610	3129	3132	3137	rVB	260287	345413	0.84%	0.424%
50	22.169	3224	3227	3230	rBV	216728	264501	0.64%	0.325%
51	22.204	3230	3233	3243	rVB3	251641	500246	1.21%	0.615%
52	23.286	3412	3417	3421	rBV	490077	703966	1.71%	0.865%
53	23.333	3421	3425	3426	rBV	185296	230374	0.56%	0.283%
54	23.957	3526	3531	3537	rVB2	269979	532194	1.29%	0.654%
55	24.115	3552	3558	3569	rVB	1037902	2159504	5.24%	2.654%
56	24.739	3658	3664	3676	rVB	790315	1731264	4.20%	2.127%
57	26.204	3909	3913	3927	rVB3	383232	973456	2.36%	1.196%
58	26.733	3995	4003	4027	rVB3	984749	3814397	9.26%	4.687%
59	27.809	4176	4186	4207	rVB4	785319	3789917	9.20%	4.657%
60	28.903	4351	4372	4388	rVB3	8540058	41213493	100.00%	50.642%

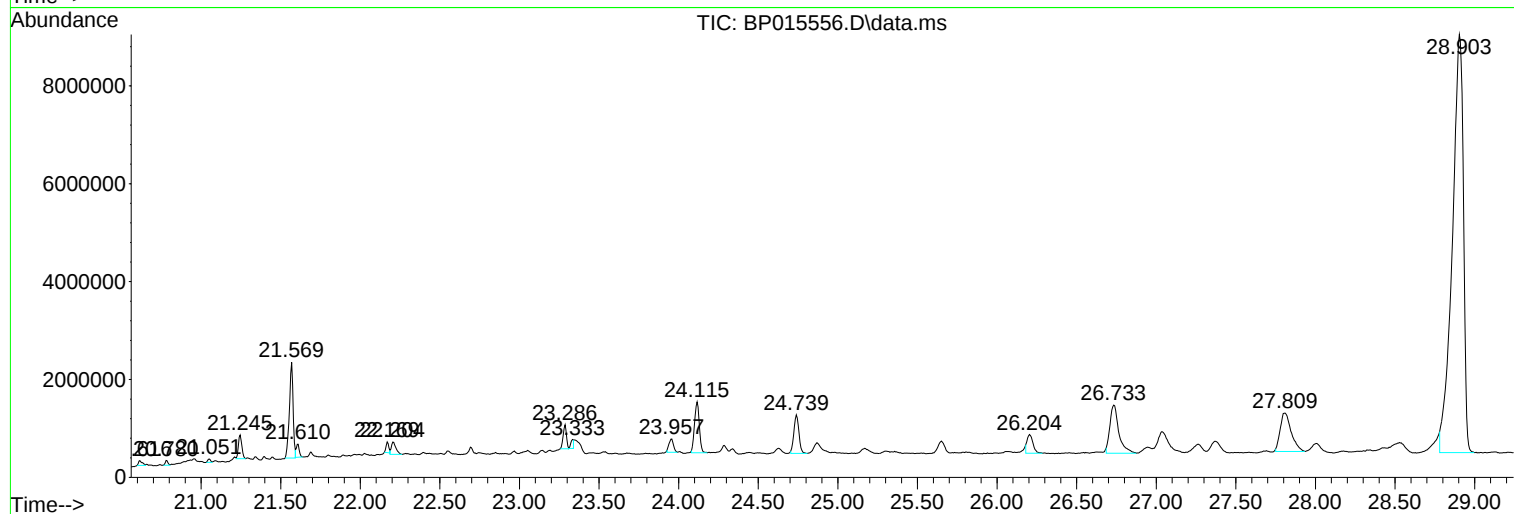
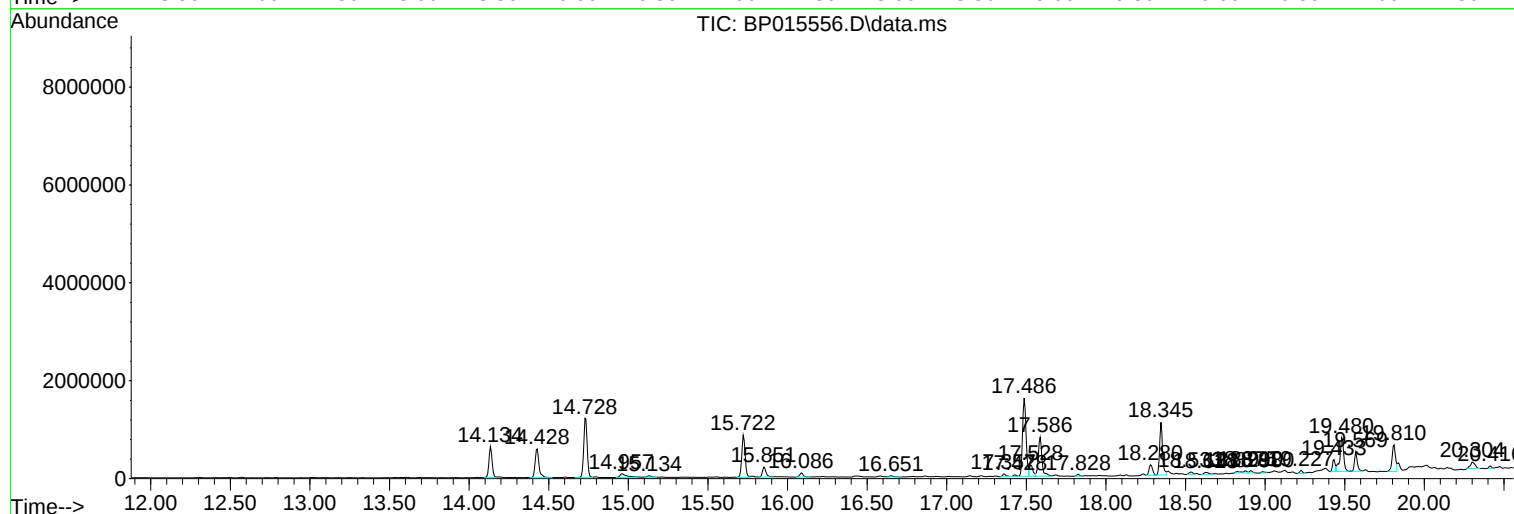
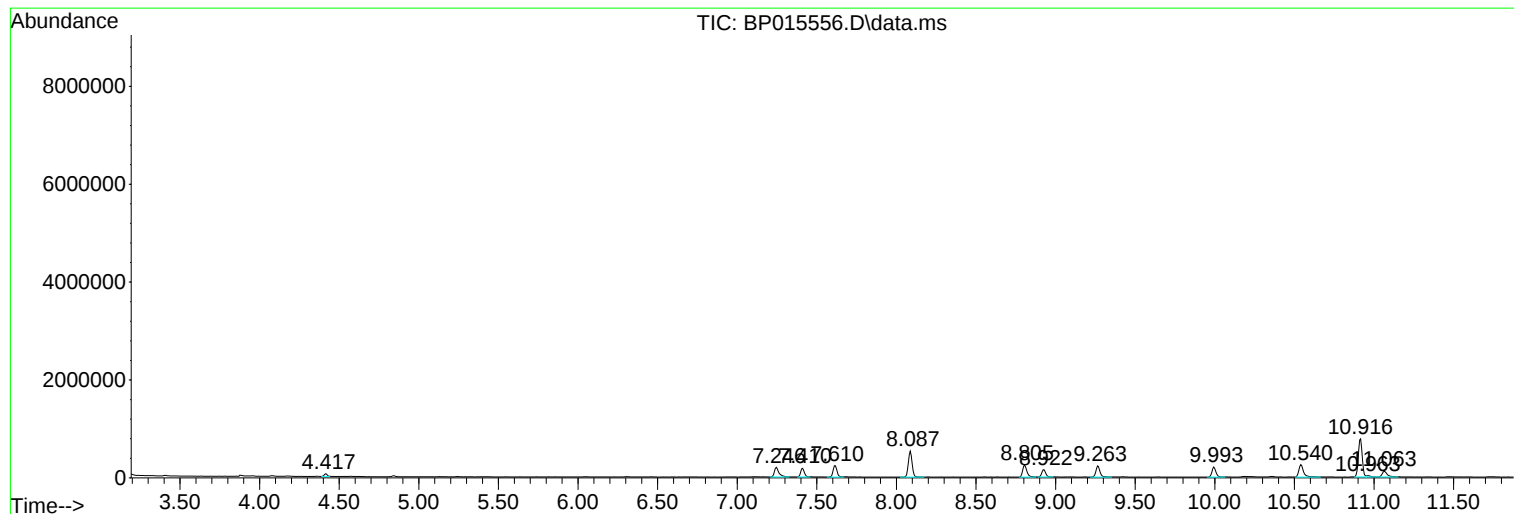
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Quant Title : SVOA CALIBRATION

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



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ALS Vial : 50 Sample Multiplier: 1

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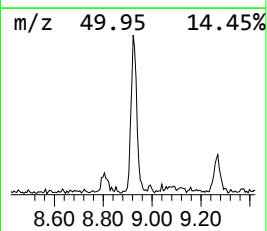
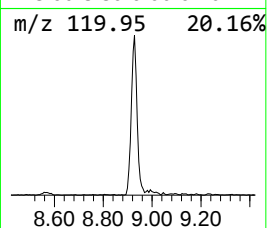
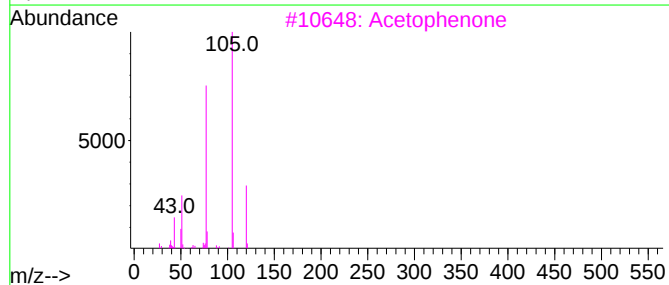
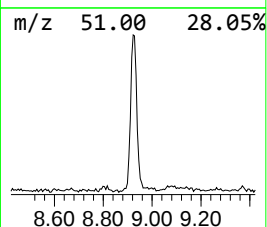
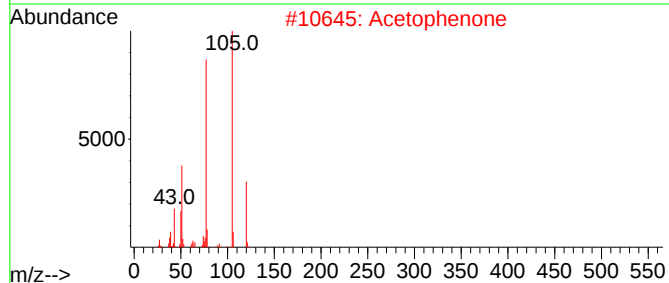
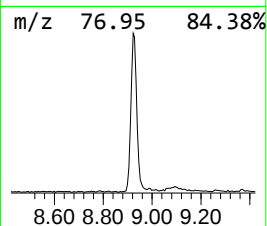
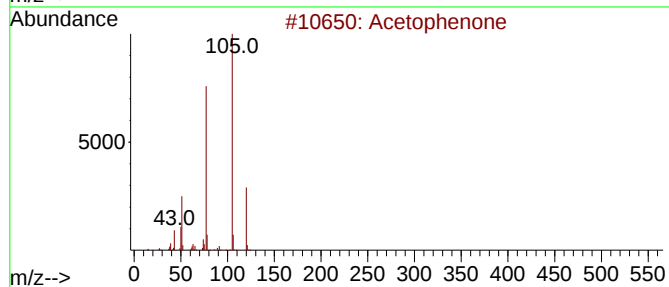
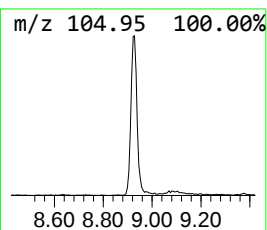
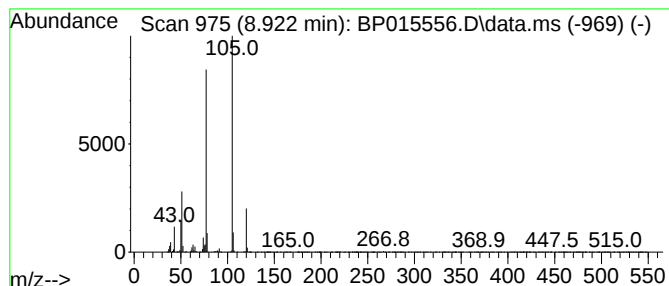
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Aceto phenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.922	5.86 ng/ul	251624	1,4-Dichlorobenzene-d4	8.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone			120	C8H8O	000098-86-2	94
2	Acetophenone			120	C8H8O	000098-86-2	91
3	Acetophenone			120	C8H8O	000098-86-2	91
4	Acetophenone			120	C8H8O	000098-86-2	91
5	Acetophenone			120	C8H8O	000098-86-2	91



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ALS Vial : 50 Sample Multiplier: 1

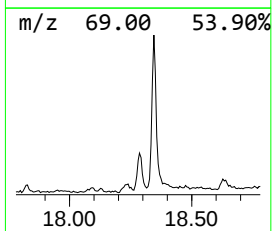
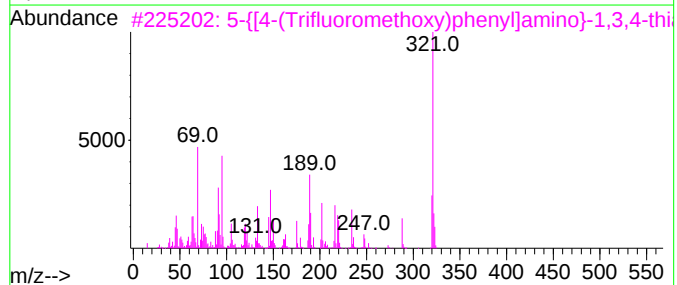
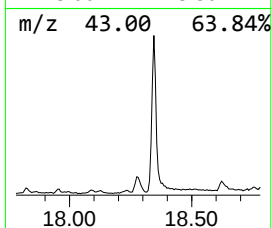
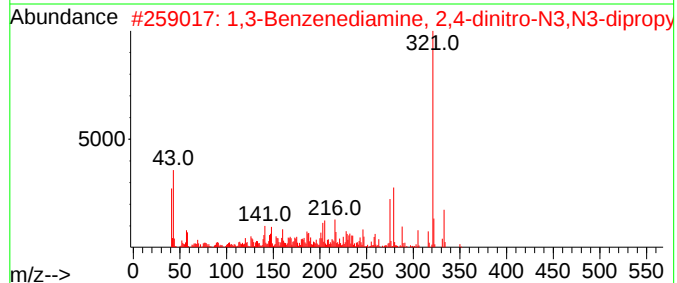
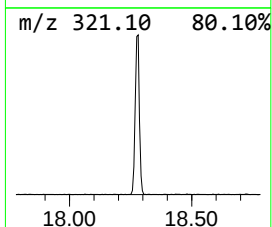
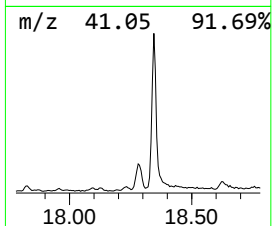
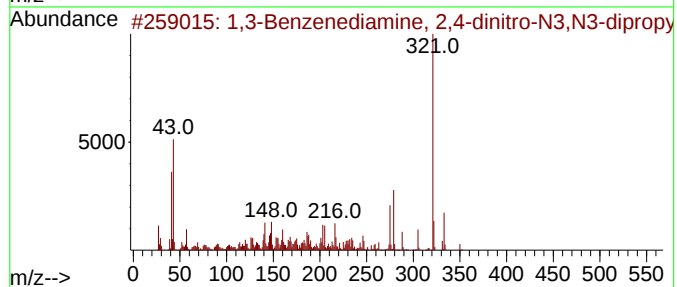
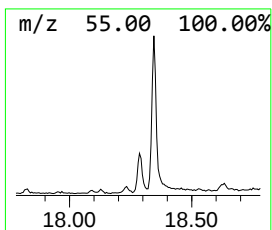
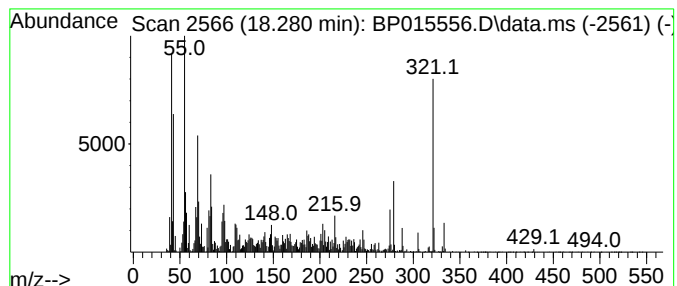
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1,3-Benzenediamine, 2,4-din... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.281	2.76 ng/ul	339506	Phenanthrene-d10	17.486	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenediamine, 2,4-dinitro-...	350	C13H17F3N4O4	029091-21-2	87
2	1,3-Benzenediamine, 2,4-dinitro-...	350	C13H17F3N4O4	029091-21-2	81
3	5-{{4-(Trifluoromethoxy)phenyl}a...	321	C11H10F3N3O52	1000507-70-1	49
4	Fumaric acid, monoamide, N-(2,5-...	321	C17H23NO5	1000345-40-2	46
5	Cyclohexanone, 2-(2-nitro-1(5H)-...	321	C18H15N3O3	021589-32-2	41



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ALS Vial : 50 Sample Multiplier: 1

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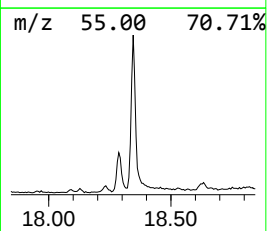
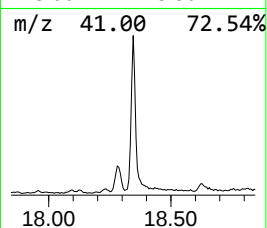
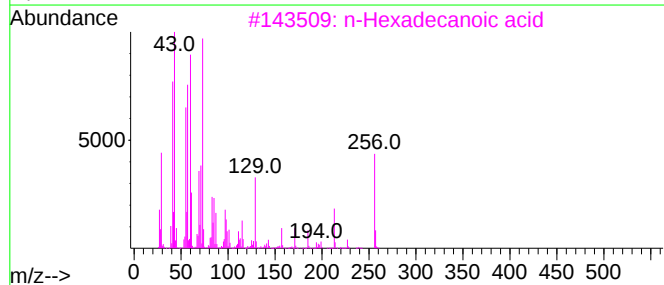
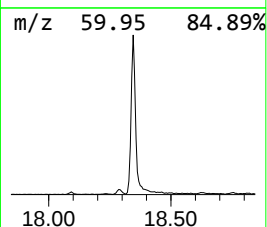
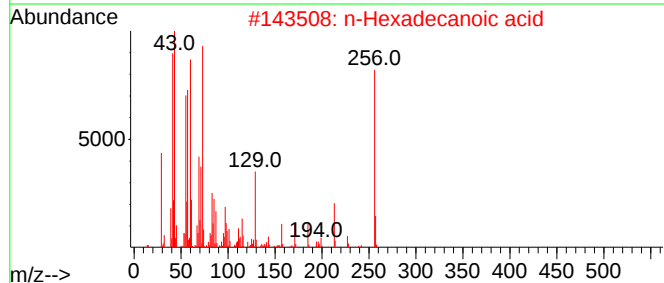
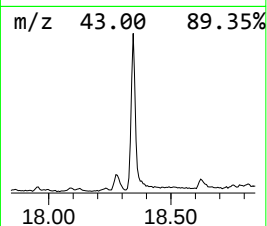
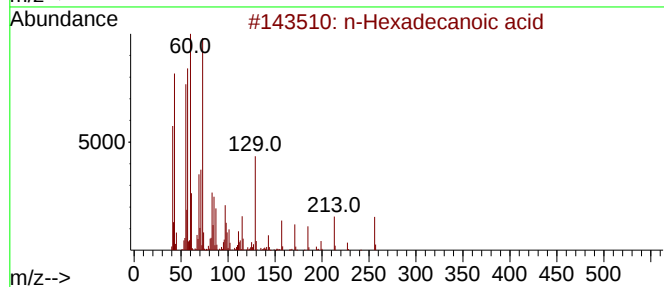
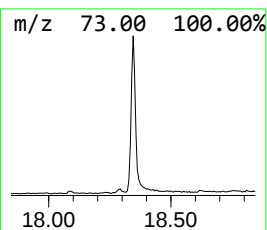
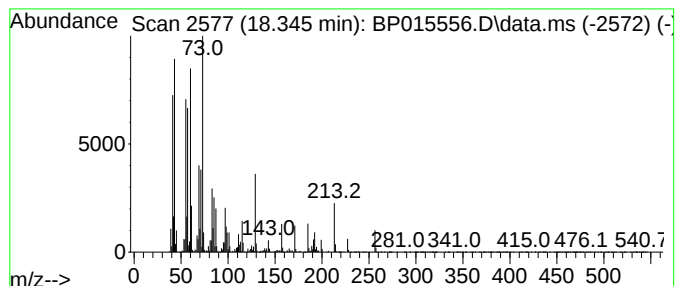
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.345	11.00 ng/ul	1351610	Phenanthrene-d10	17.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Tridecanoic acid	214	C13H26O2	000638-53-9	95



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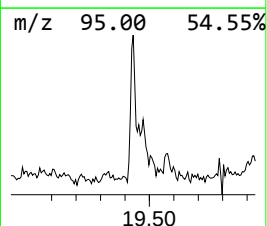
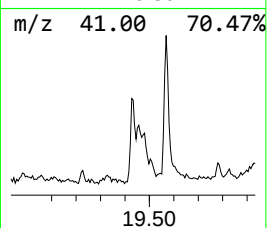
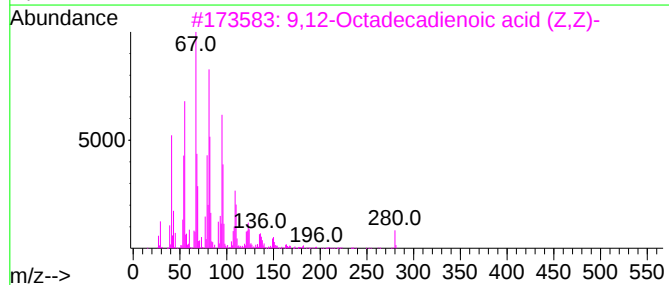
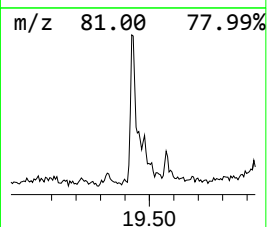
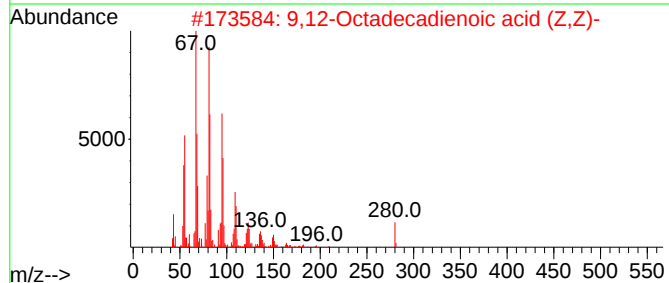
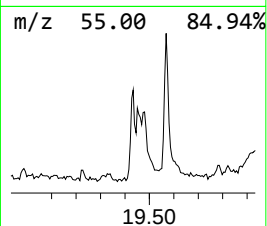
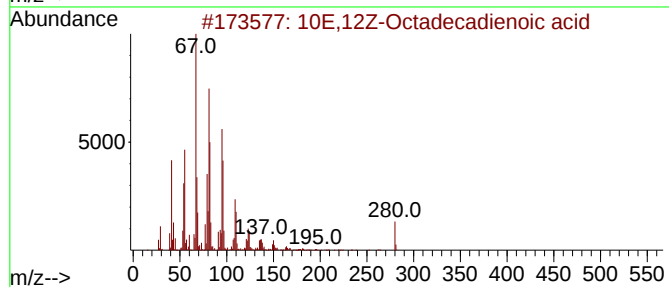
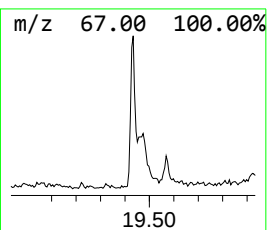
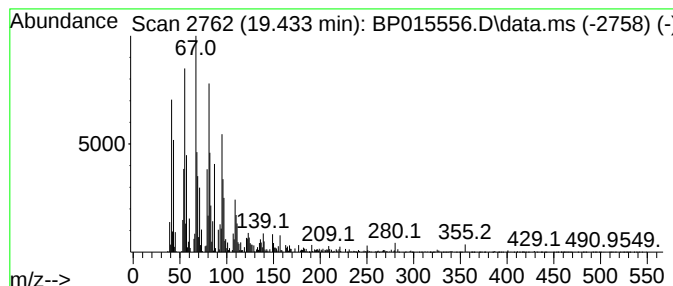
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TIC Library : C:\DATABASE\NIST20.L
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Peak Number 4 10E,12Z-Octadecadienoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.433	2.53 ng/ul	310433	Phenanthrene-d10	17.486	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10E,12Z-Octadecadienoic acid	280	C18H32O2	002420-56-6	97
2	9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	97
3	9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	97
4	Linolelaidic acid	280	C18H32O2	000506-21-8	93
5	(9E,11E)-Octadecadienoic acid	280	C18H32O2	000544-71-8	91



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ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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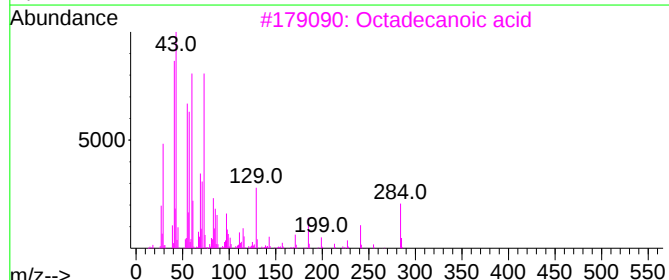
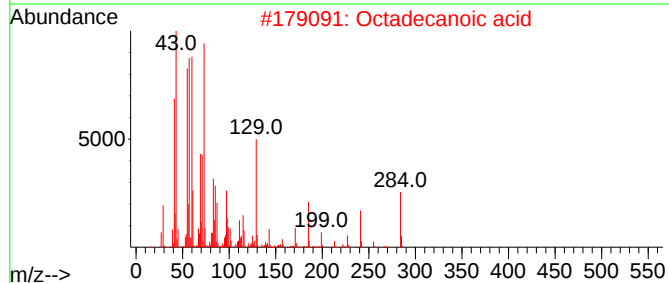
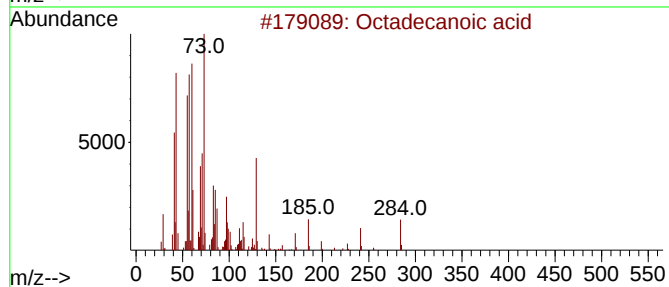
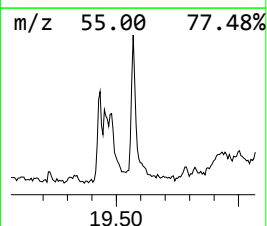
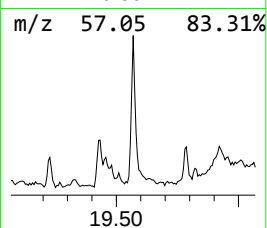
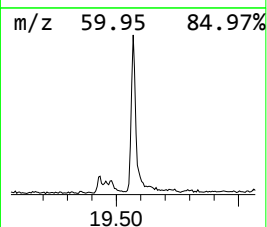
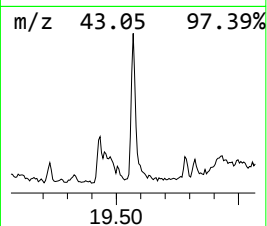
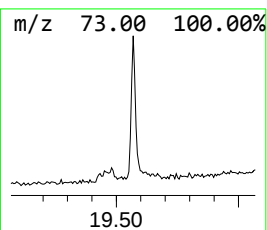
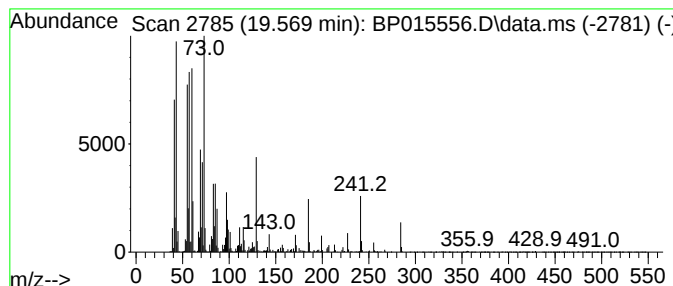
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.569	2.98 ng/ul	471662	Chrysene-d12	21.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	97
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015556.D
Acq On : 15 Jun 2023 15:19
Operator : MA/JU
Sample : 03088-13
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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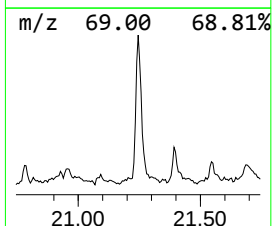
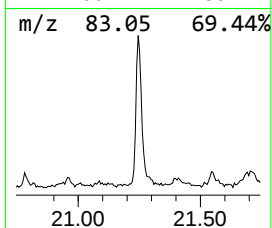
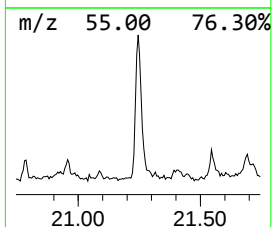
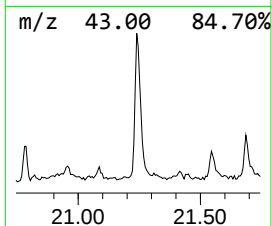
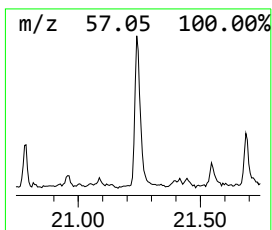
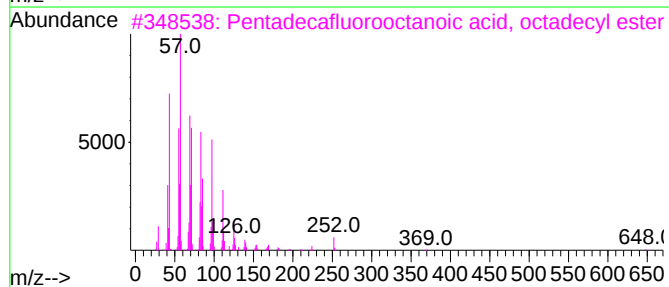
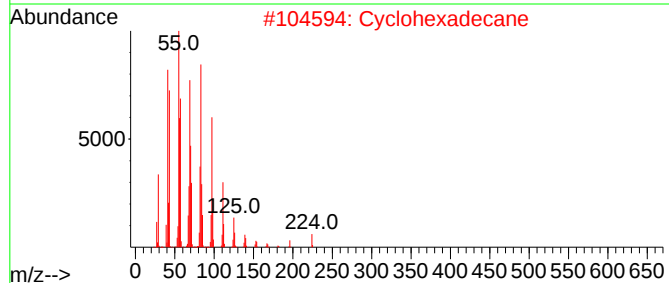
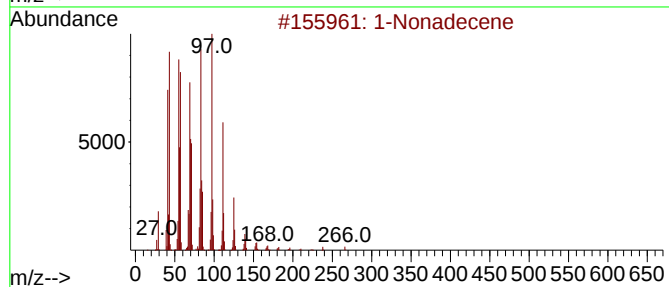
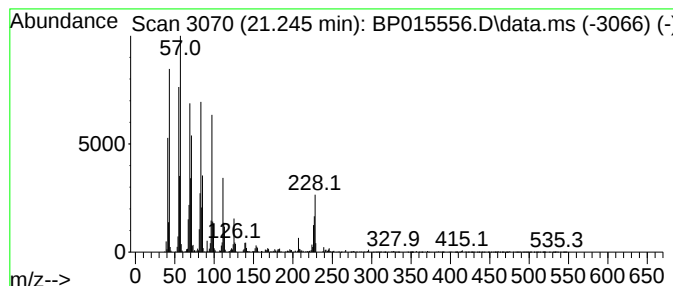
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.245	4.46 ng/ul	707495	Chrysene-d12	21.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene			266	C19H38	018435-45-5	93
2	Cyclohexadecane			224	C16H32	000295-65-8	92
3	Pentadecafluorooctanoic acid, oc...			666	C26H37F15O2	1000406-04-8	90
4	Pentacos-1-ene			350	C25H50	016980-85-1	90
5	1-Hexacosanol			382	C26H54O	000506-52-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015556.D
Acq On : 15 Jun 2023 15:19
Operator : MA/JU
Sample : 03088-13
Misc :
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Instrument :
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ClientSampleId :
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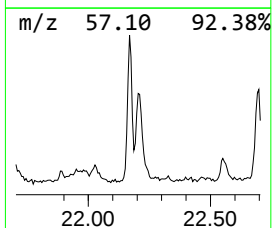
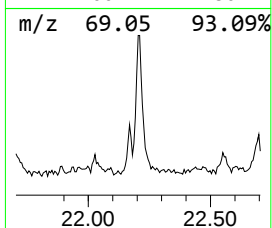
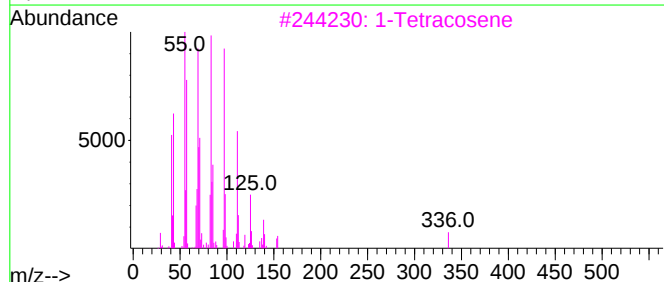
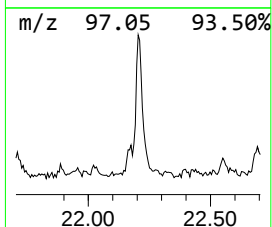
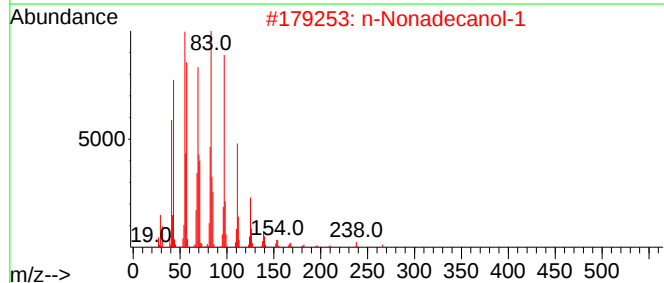
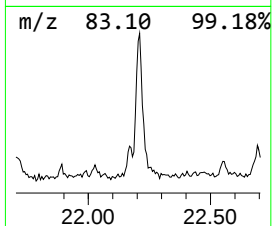
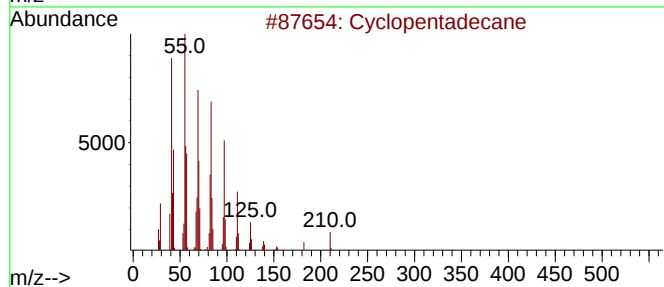
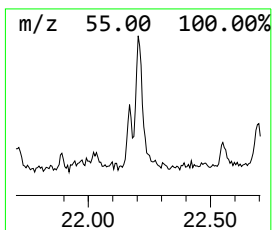
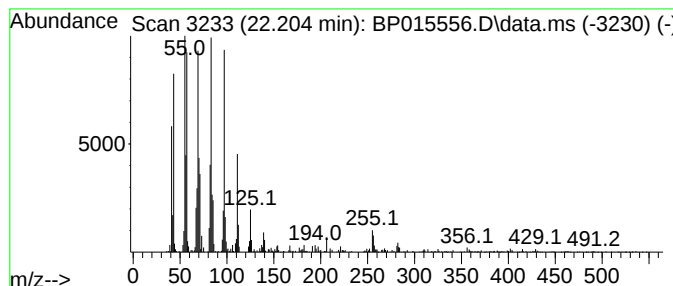
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Cyclic22.204 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.204	3.16 ng/ul	500246	Chrysene-d12	21.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentadecane	210	C15H30	000295-48-7	96
2			n-Nonadecanol-1	284	C19H40O	001454-84-8	95
3			1-Tetracosene	336	C24H48	010192-32-2	95
4			1-Heneicosanol	312	C21H44O	015594-90-8	94
5			Behenic alcohol	326	C22H46O	000661-19-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015556.D
Acq On : 15 Jun 2023 15:19
Operator : MA/JU
Sample : 03088-13
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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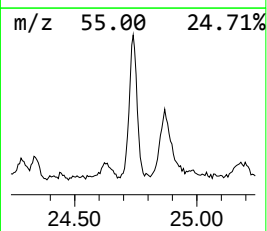
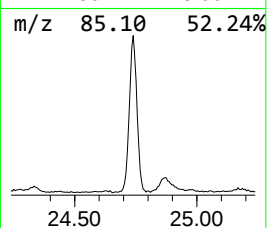
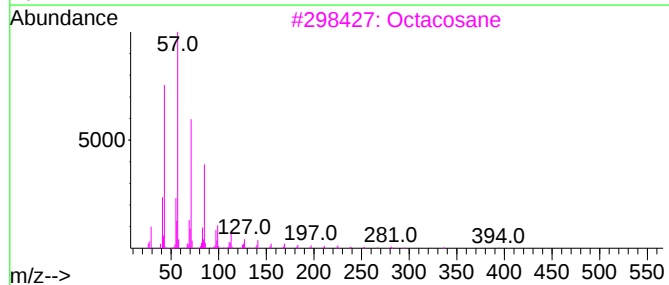
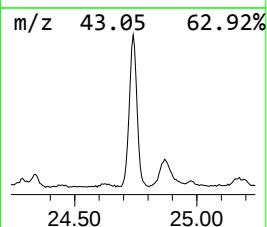
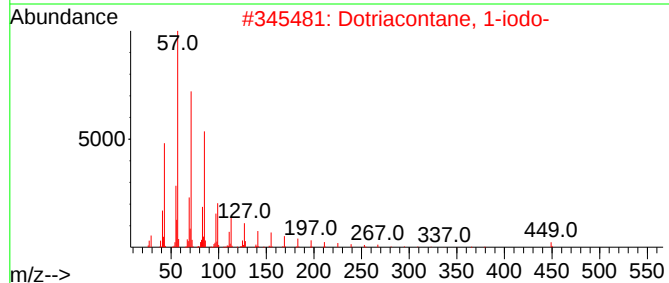
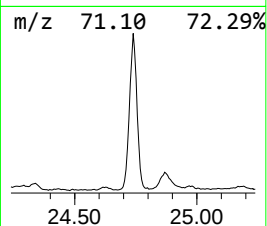
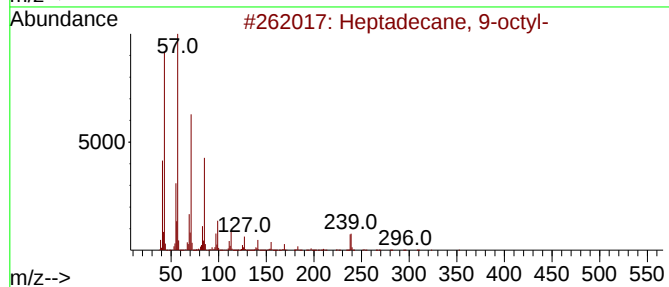
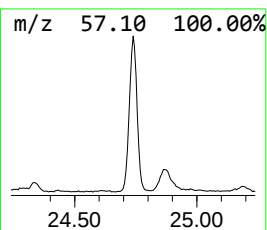
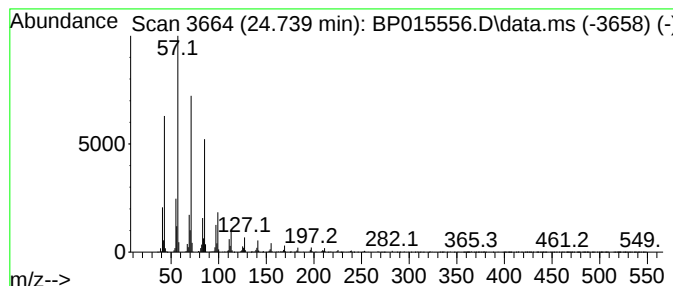
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.739	16.03 ng/ul	1731260	Perylene-d12	24.116

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2		Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015556.D
Acq On : 15 Jun 2023 15:19
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Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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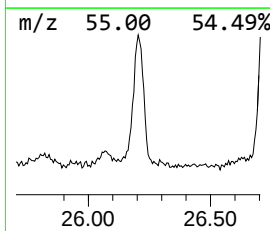
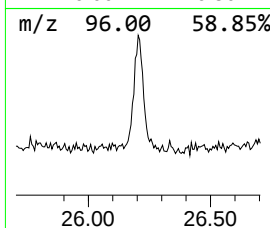
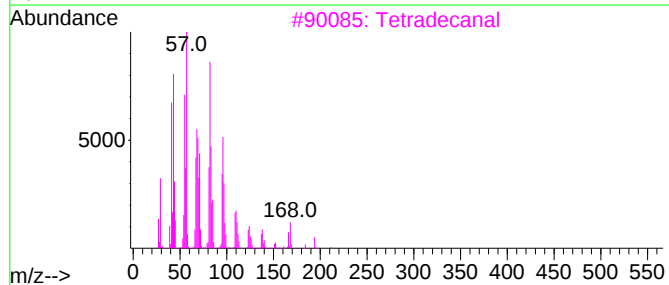
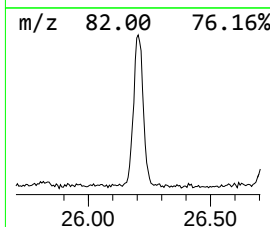
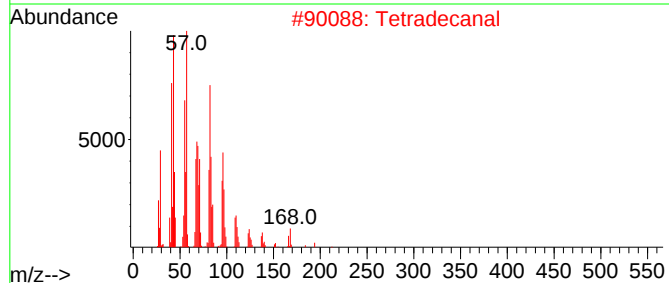
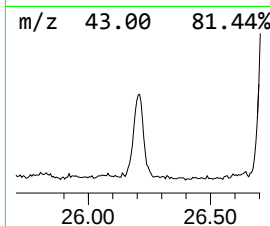
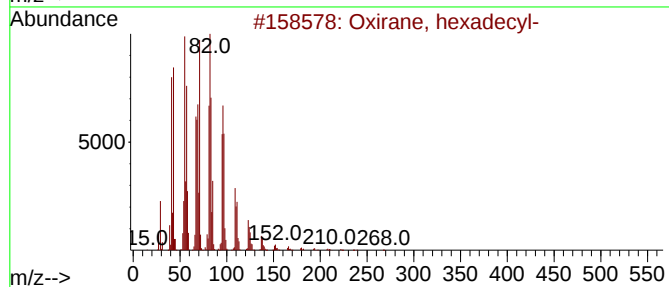
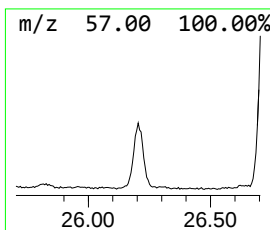
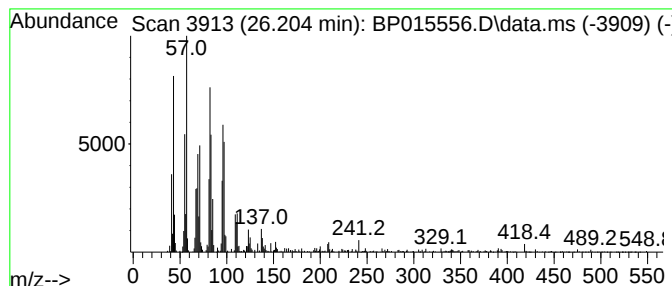
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Oxirane, hexadecyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.204	9.02 ng/ul	973456	Perylene-d12	24.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	90
2			Tetradecanal	212	C14H28O	000124-25-4	90
3			Tetradecanal	212	C14H28O	000124-25-4	87
4			Cyclohexane, (1-hexadecylheptade...	547	C39H78	055517-75-4	86
5			13-Methyltetradecanal	226	C15H30O	075853-51-9	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
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Sample : 03088-13
Misc :
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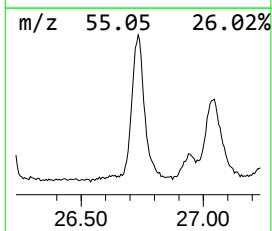
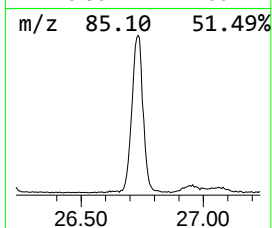
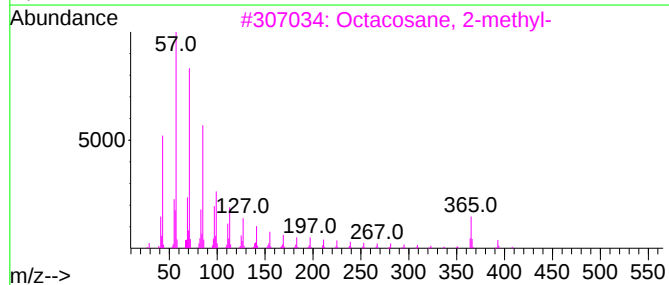
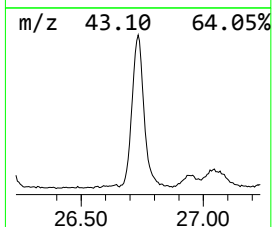
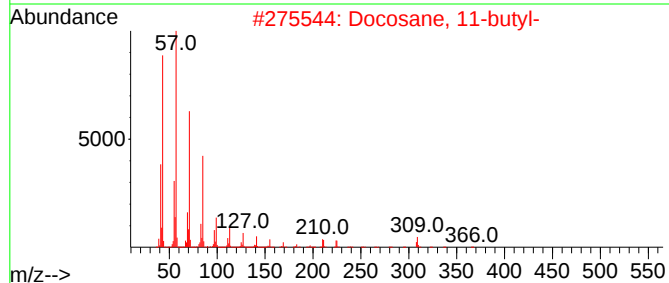
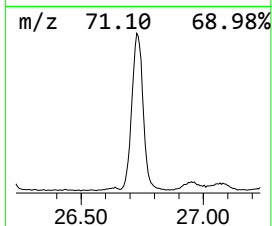
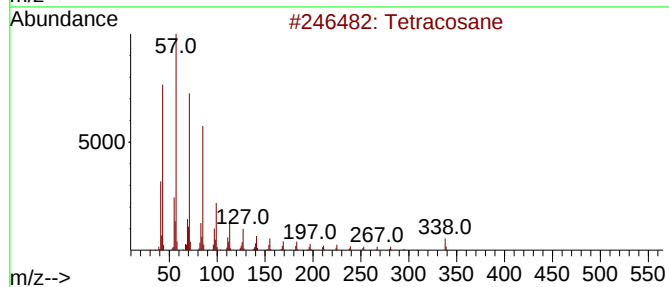
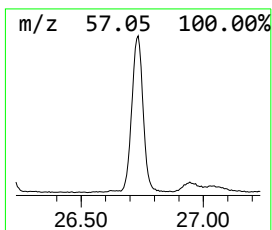
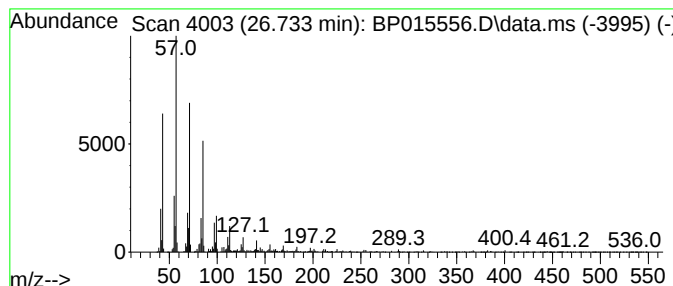
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
26.733	35.33 ng/ul	3814400	Perylene-d12		24.116	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetracosane		338	C24H50	000646-31-1	97
2	Docosane, 11-butyl-		366	C26H54	013475-76-8	94
3	Octacosane, 2-methyl-		408	C29H60	001560-98-1	93
4	Docosane, 7-hexyl-		394	C28H58	055373-86-9	93
5	Heneicosane, 11-pentyl-		366	C26H54	014739-72-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015556.D
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ALS Vial : 50 Sample Multiplier: 1

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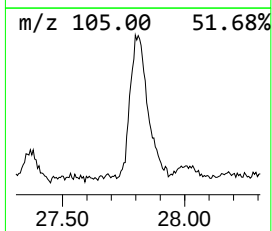
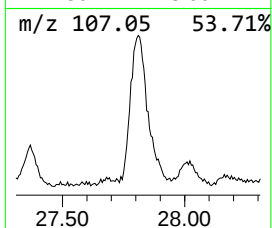
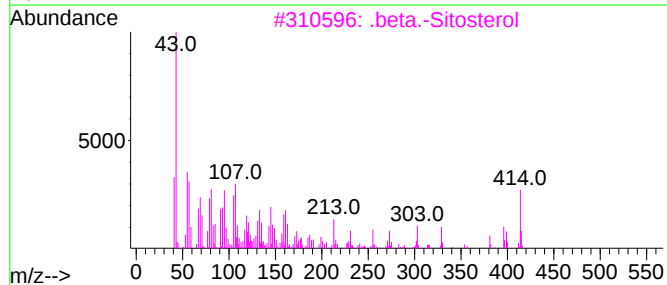
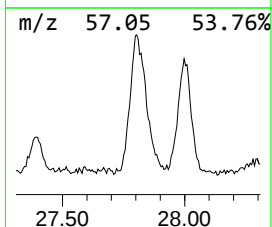
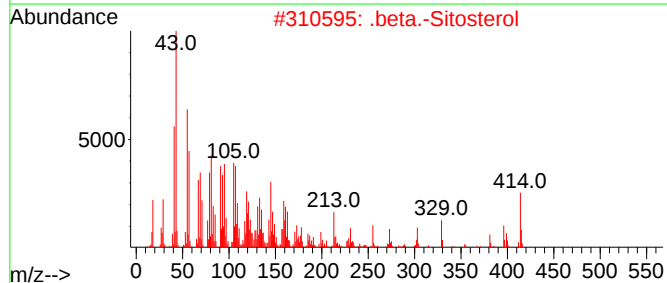
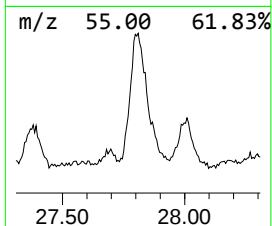
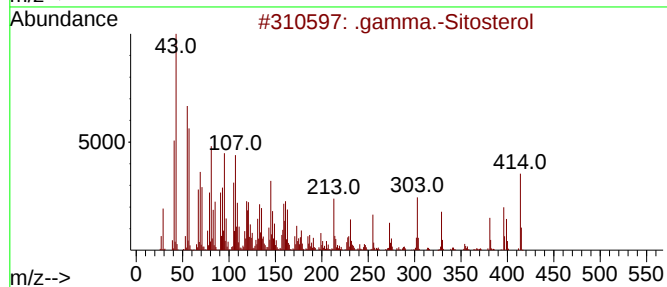
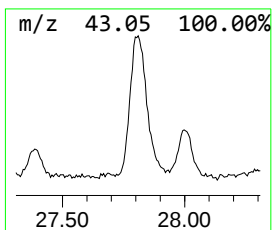
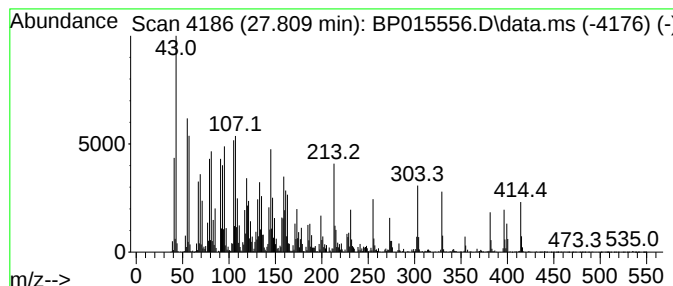
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 .gamma.-Sitosterol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.809	35.10 ng/ul	3789920	Perylene-d12	24.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.gamma.-Sitosterol	414	C29H50O	000083-47-6	99
2			.beta.-Sitosterol	414	C29H50O	000083-46-5	93
3			.beta.-Sitosterol	414	C29H50O	000083-46-5	93
4			(3S,8S,9S,10R,13R,14S,17R)-17-((... 428	C30H52O	020194-50-7	52	
5			(3S,8S,9S,10R,13R,14S,17R)-17-((... 414	C29H50O	029365-29-5	50	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015556.D
Acq On : 15 Jun 2023 15:19
Operator : MA/JU
Sample : 03088-13
Misc :
ALS Vial : 50 Sample Multiplier: 1

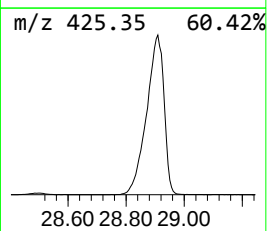
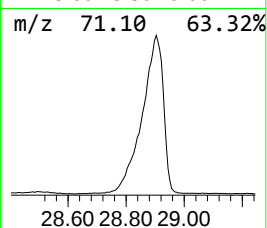
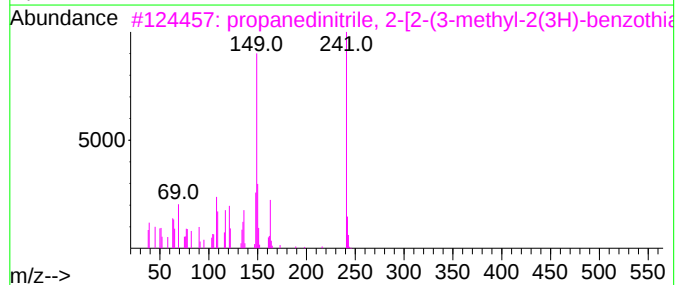
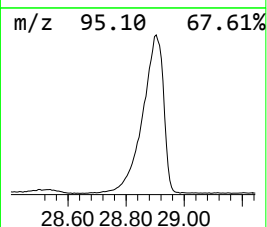
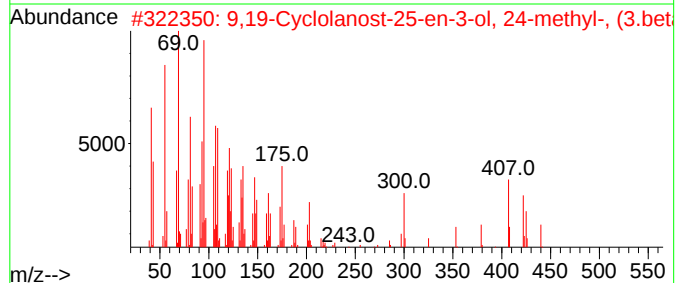
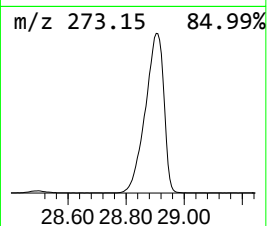
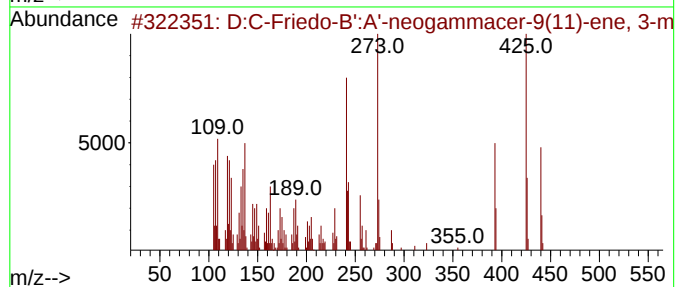
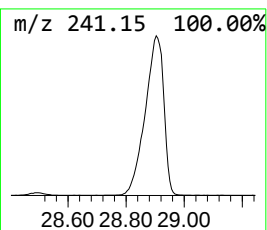
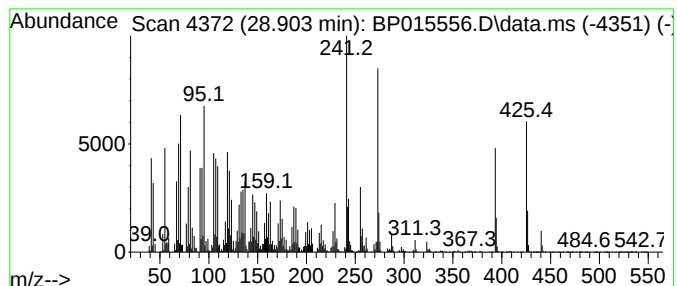
Instrument :
BNA_P
ClientSampleId :
DCFR9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 D:C-Friedo-B':A'-neogammace... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
28.904	381.69 ng/ul	41213500	Perylene-d12			24.116
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	D:C-Friedo-B':A'-neogammacer-9(1...		440	C31H52O	004555-56-0	50
2	9,19-Cyclolanost-25-en-3-ol, 24-...		440	C31H52O	000511-61-5	35
3	propanedinitrile, 2-[2-(3-methyl...		241	C11H7N5S	1000396-51-8	11
4	3-Bromo-4-hydroxy-5-methoxypheny...		241	C9H8BrNO2	081038-44-0	11
5	benzenamine, 2,6-dimethyl-N,N-di...		273	C20H19N	1010401-78-4	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015556.D
Acq On : 15 Jun 2023 15:19
Operator : MA/JU
Sample : 03088-13
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFR9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.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
Aceto phenone	8.922	5.9	ng/ul	251624	1	8.087	858644	20.0
1,3-Benzenediam...	18.281	2.8	ng/ul	339506	4	17.486	2456500	20.0
n-Hexadecanoic ...	18.345	11.0	ng/ul	1351610	4	17.486	2456500	20.0
10E,12Z-Octadec...	19.433	2.5	ng/ul	310433	4	17.486	2456500	20.0
Octadecanoic acid	19.569	3.0	ng/ul	471662	5	21.569	3169160	20.0
(DEL) Alkane: S...	21.245	4.5	ng/ul	707495	5	21.569	3169160	20.0
(DEL) Alkane: C...	22.204	3.2	ng/ul	500246	5	21.569	3169160	20.0
(DEL) Alkane: S...	24.739	16.0	ng/ul	1731260	6	24.116	2159500	20.0
Oxirane, hexade...	26.204	9.0	ng/ul	973456	6	24.116	2159500	20.0
(DEL) Alkane: S...	26.733	35.3	ng/ul	3814400	6	24.116	2159500	20.0
.gamma.-Sitosterol	27.809	35.1	ng/ul	3789920	6	24.116	2159500	20.0
D:C-Friedo-B':A...	28.904	381.7	ng/ul	41213500	6	24.116	2159500	20.0