

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110625\
 Data File : BP026087.D
 Acq On : 06 Nov 2025 21:52
 Operator : RC/JU
 Sample : Q3552-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-8

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % 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\8270E-BP102925.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP026087.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.231	46	50	60	rVB	175284	234648	1.20%	0.236%
2	4.596	274	282	287	rBV	158015	272029	1.39%	0.274%
3	5.296	394	401	411	rVB	3811215	5707569	29.08%	5.745%
4	6.855	659	666	677	rBV	3093372	5192238	26.46%	5.227%
5	7.678	799	806	814	rBV	1110840	1770479	9.02%	1.782%
6	8.425	927	933	944	rVB	1508470	2387128	12.16%	2.403%
7	8.813	992	999	1006	rVB	3613149	5827283	29.69%	5.866%
8	9.272	1071	1077	1084	rVB	174262	268946	1.37%	0.271%
9	10.443	1268	1276	1283	rBV	1432881	2400771	12.23%	2.417%
10	10.972	1360	1366	1374	rBV	207142	402227	2.05%	0.405%
11	12.919	1690	1697	1704	rBV	8609377	12940091	65.94%	13.026%
12	13.966	1869	1875	1882	rVB4	136194	221835	1.13%	0.223%
13	14.107	1894	1899	1908	rVB	284536	469151	2.39%	0.472%
14	14.307	1926	1933	1943	rVB	2050573	3169319	16.15%	3.190%
15	15.495	2131	2135	2142	rVB	195500	262999	1.34%	0.265%
16	15.689	2163	2168	2174	rVB	504648	726411	3.70%	0.731%
17	15.819	2183	2190	2199	rBV	6800753	10603622	54.03%	10.674%
18	16.519	2305	2309	2314	rBV	197499	274635	1.40%	0.276%
19	17.107	2403	2409	2416	rVB2	2318737	3684374	18.77%	3.709%
20	17.313	2439	2444	2451	rBV2	164128	274699	1.40%	0.277%
21	17.389	2452	2457	2462	rVV	204114	322546	1.64%	0.325%
22	17.825	2526	2531	2540	rVB2	319588	512705	2.61%	0.516%
23	17.930	2545	2549	2556	rVV	297283	450936	2.30%	0.454%
24	18.072	2568	2573	2582	rBV	1884193	3004638	15.31%	3.025%
25	18.166	2586	2589	2602	rVB2	200268	353267	1.80%	0.356%
26	18.289	2604	2610	2622	rBV	2621880	3937264	20.06%	3.963%
27	19.277	2775	2778	2786	rVB2	280398	435837	2.22%	0.439%
28	19.413	2796	2801	2813	rBV	467166	786558	4.01%	0.792%
29	19.860	2870	2877	2882	rBV	13603916	19623874	100.00%	19.754%
30	21.260	3111	3115	3124	rBV	882708	1210781	6.17%	1.219%
31	21.566	3162	3167	3176	rVB2	2709068	4492961	22.90%	4.523%
32	23.442	3480	3486	3494	rVB	803615	1722838	8.78%	1.734%
33	24.889	3723	3732	3748	rVB	1818701	5395737	27.50%	5.432%

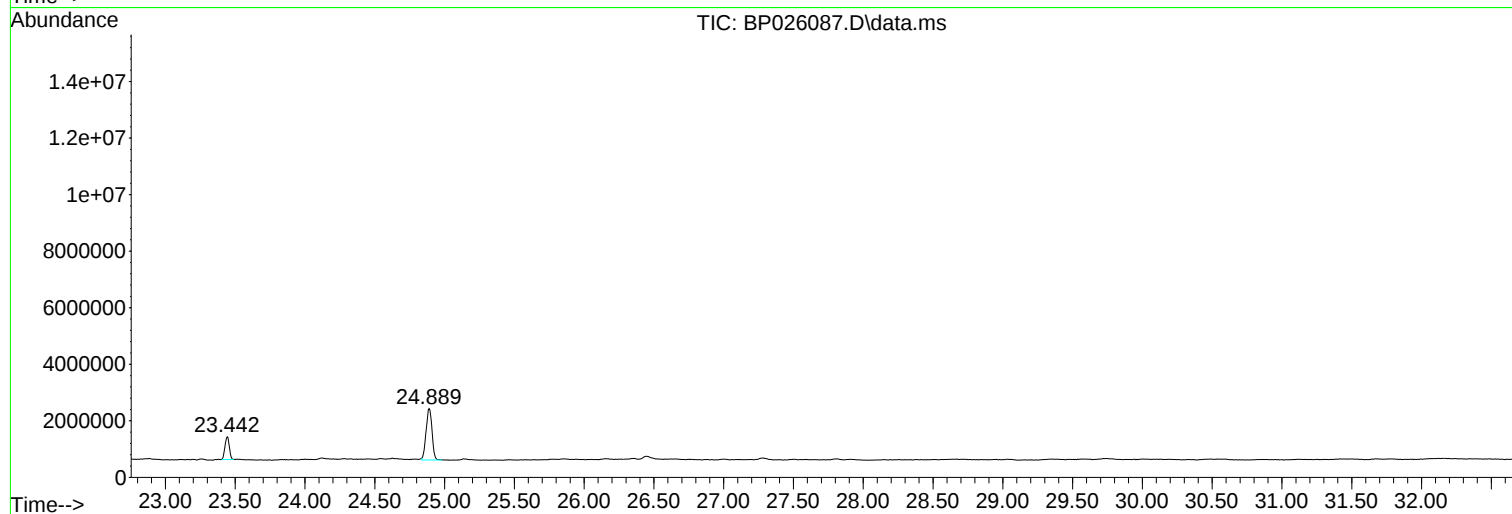
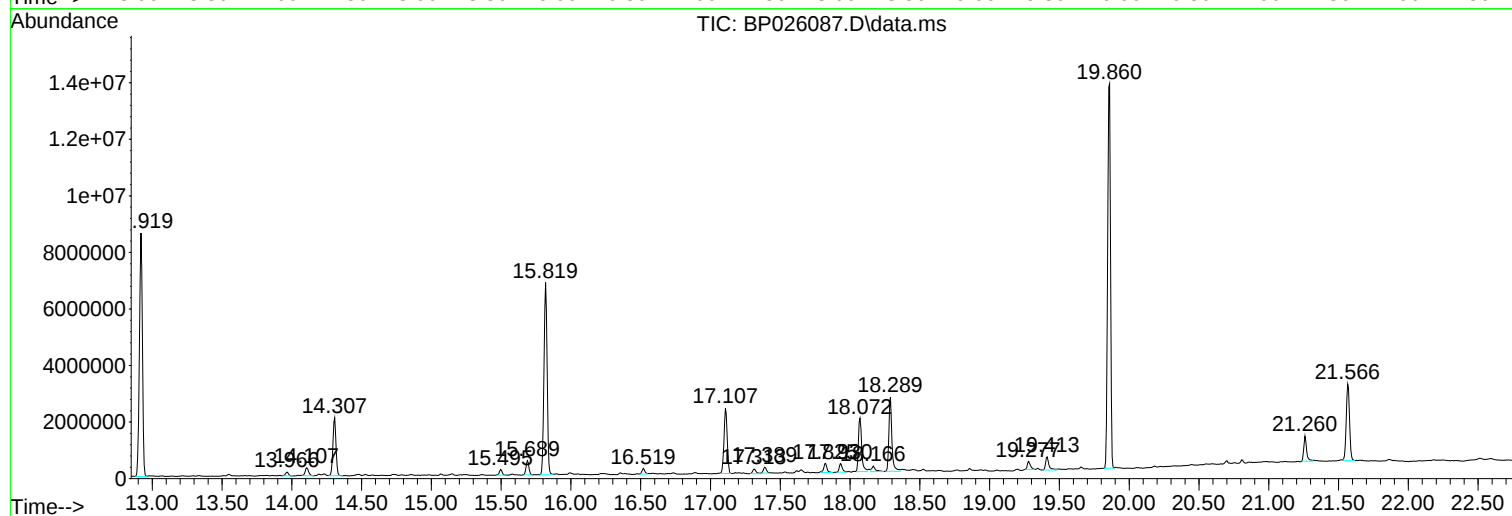
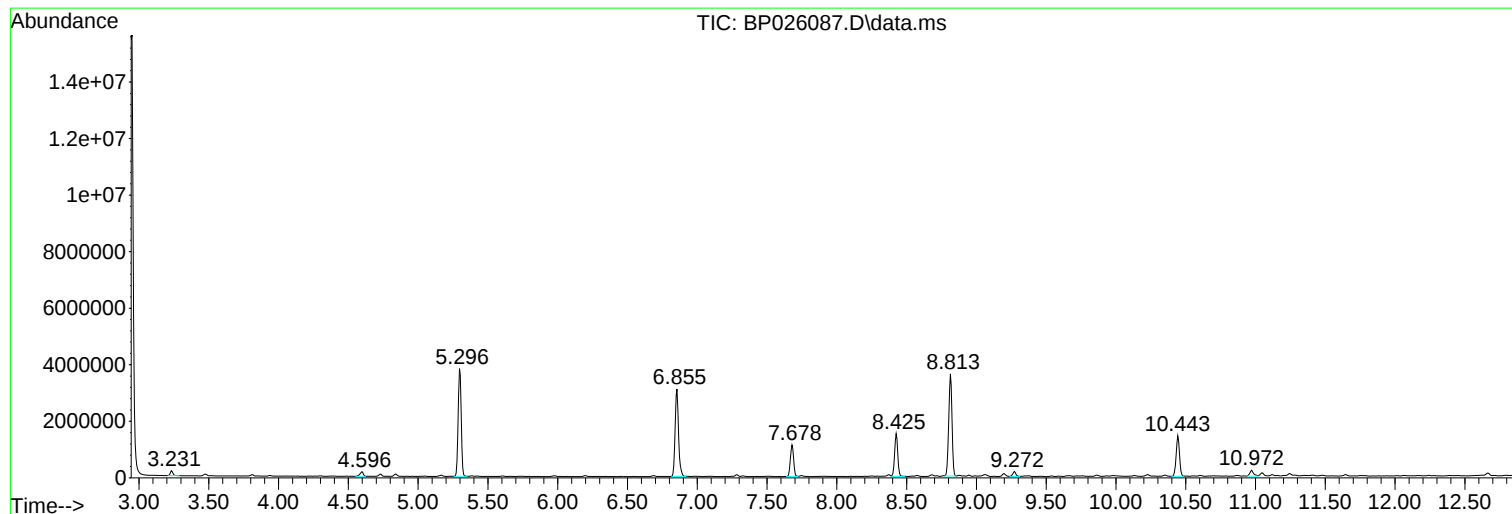
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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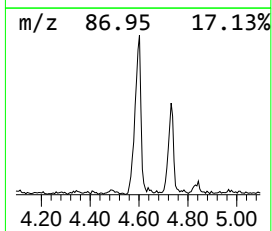
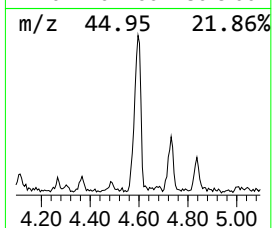
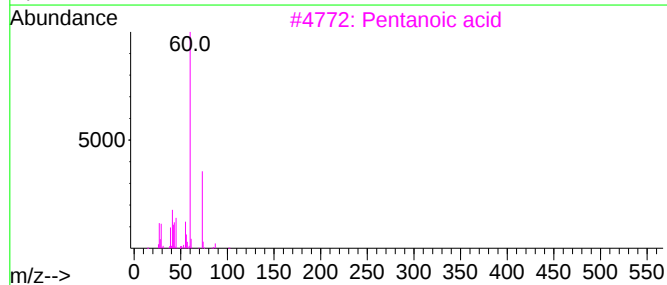
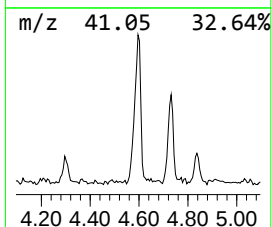
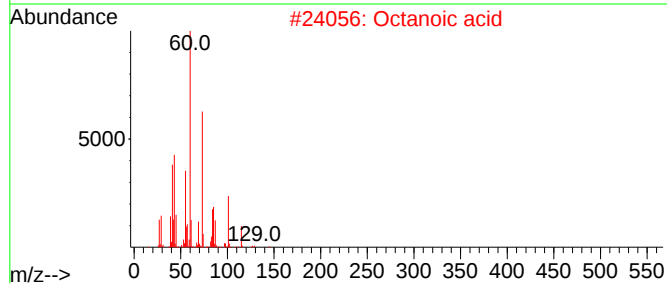
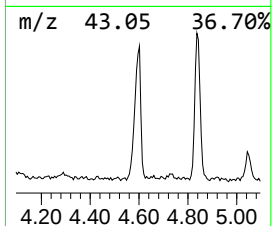
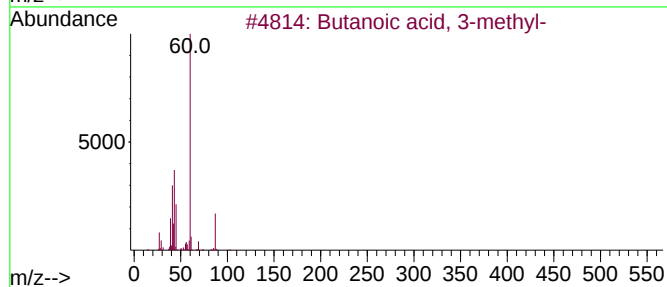
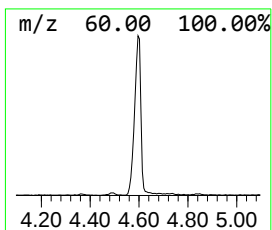
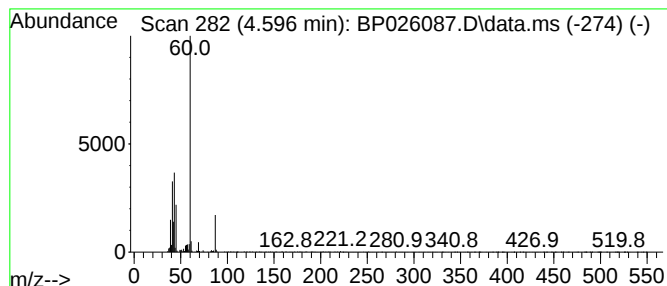
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butanoic acid, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.596	3.07 ng	272029	1,4-Dichlorobenzene-d4	7.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	90
2			Octanoic acid	144	C8H16O2	000124-07-2	38
3			Pentanoic acid	102	C5H10O2	000109-52-4	38
4			Heptanoic acid	130	C7H14O2	000111-14-8	36
5			Urea, N-methyl-N-nitroso-	103	C2H5N3O2	000684-93-5	36



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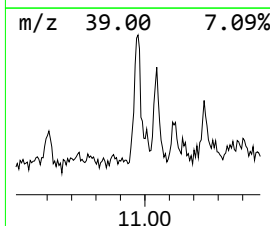
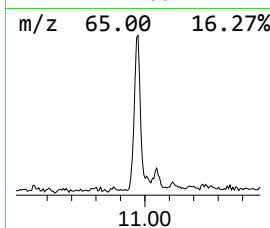
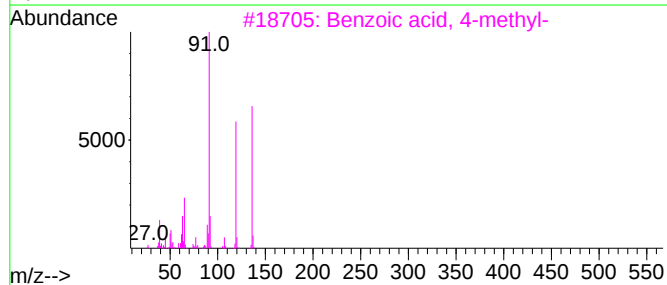
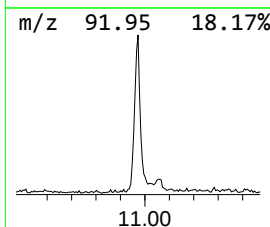
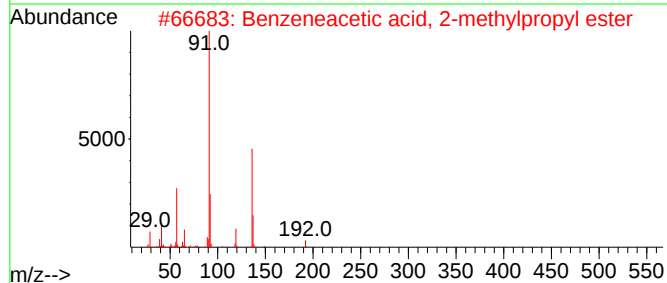
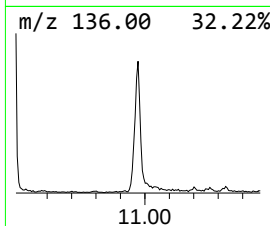
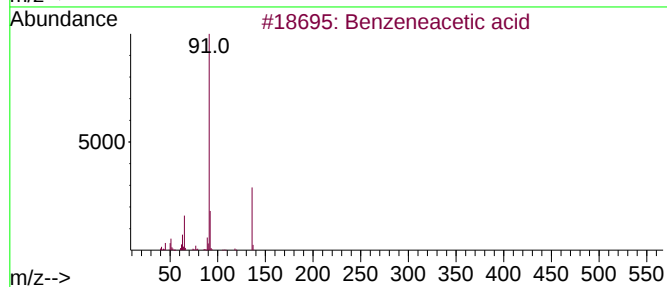
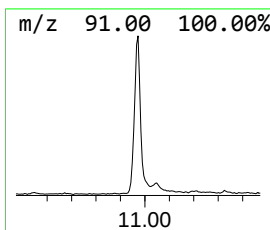
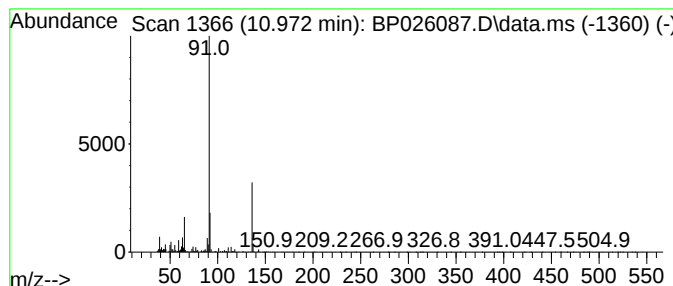
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Benzeneacetic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.972	3.35 ng	402227	Naphthalene-d8	10.443

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneacetic acid	136	C8H8O2	000103-82-2	93
2			Benzeneacetic acid, 2-methylprop...	192	C12H16O2	000102-13-6	59
3			Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	53
4			Benzoic acid, 3-methyl-	136	C8H8O2	000099-04-7	49
5			Benzeneacetaldehyde	120	C8H8O	000122-78-1	47



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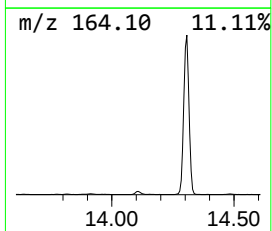
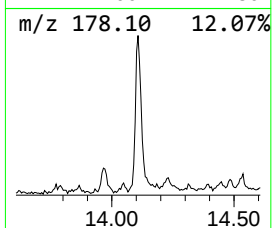
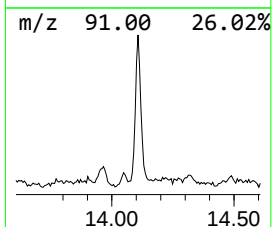
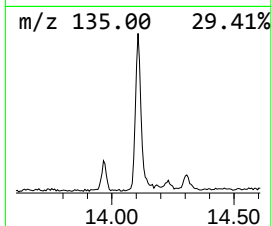
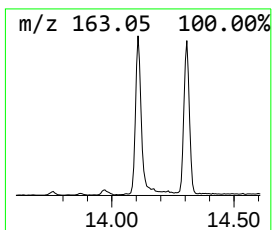
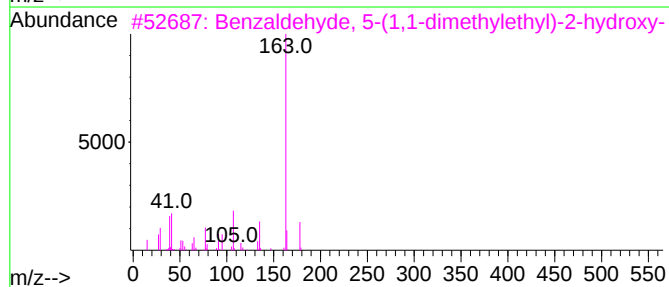
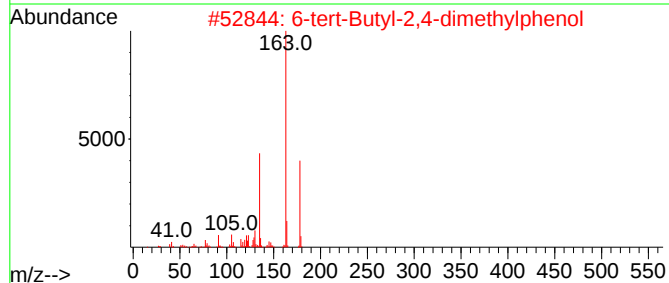
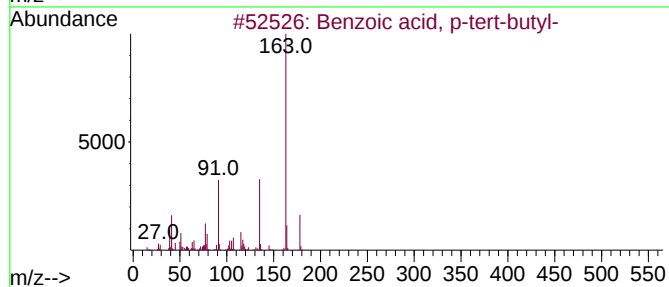
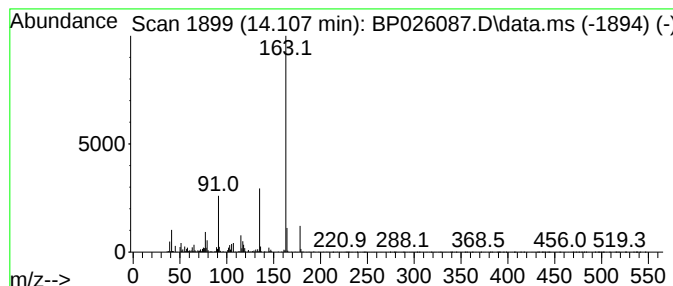
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 Benzoic acid, p-tert-butyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.107	2.96 ng	469151	Acenaphthene-d10	14.307

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	93
2			6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	80
3			Benzaldehyde, 5-(1,1-dimethylethyl)-	178	C11H14O2	002725-53-3	80
4			Silane, (4-ethylphenyl)trimethyl-	178	C11H18Si	017988-50-0	64
5			3-tert-Butylbenzoic acid	178	C11H14O2	007498-54-6	60



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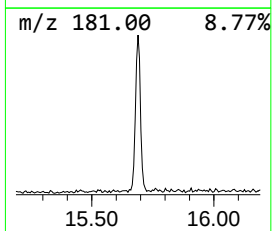
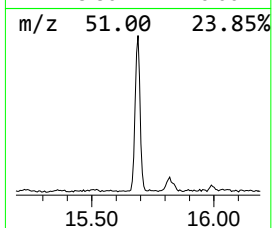
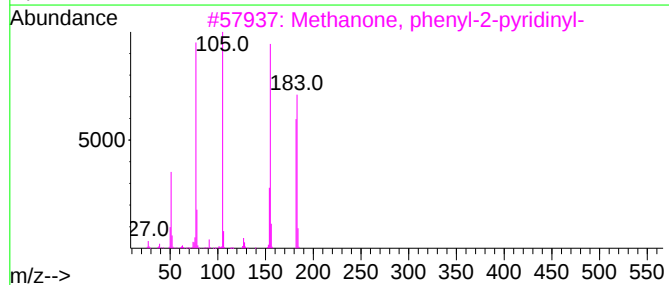
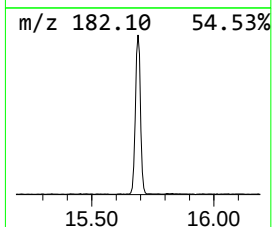
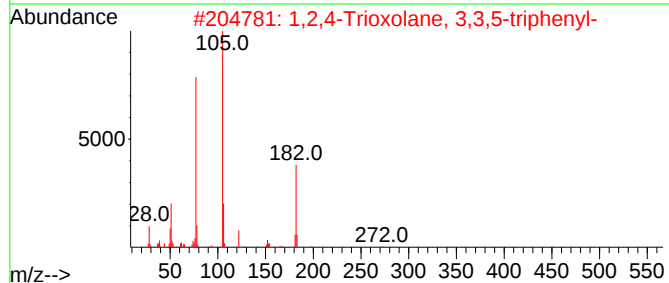
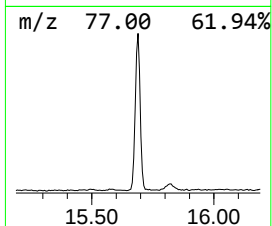
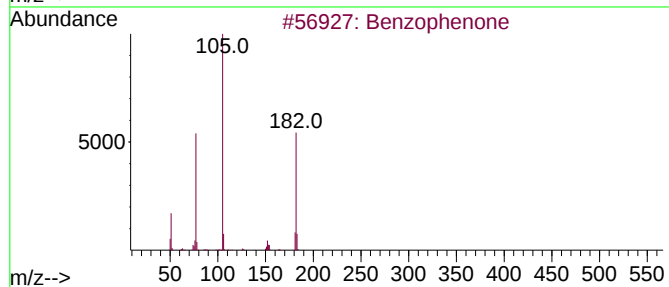
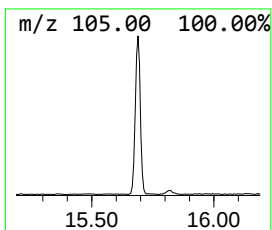
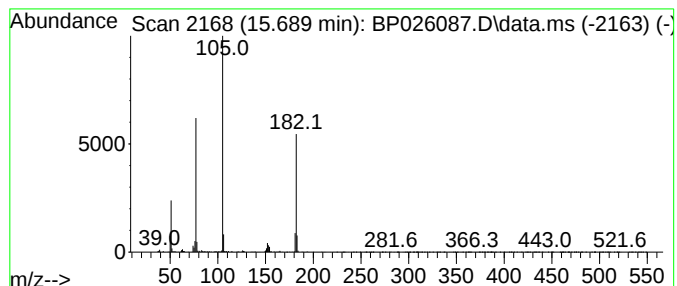
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Peak Number 5 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.689	4.58 ng	726411	Acenaphthene-d10	14.307

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			1,2,4-Trioxolane, 3,3,5-triphenyl-	304	C20H16O3	023246-12-0	50
3			Methanone, phenyl-2-pyridinyl-	183	C12H9NO	000091-02-1	50
4			Benzenamine, N-(3-pyridinylmethy...	182	C12H10N2	029722-97-2	45
5			Azobenzene	182	C12H10N2	000103-33-3	45



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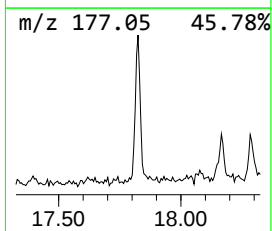
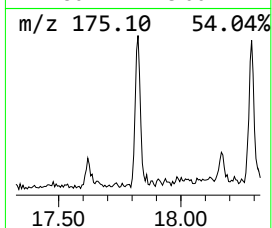
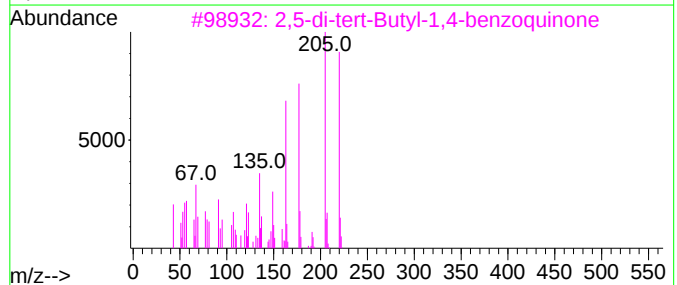
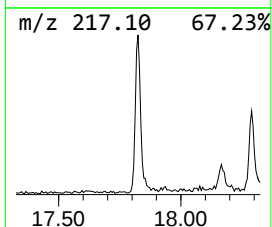
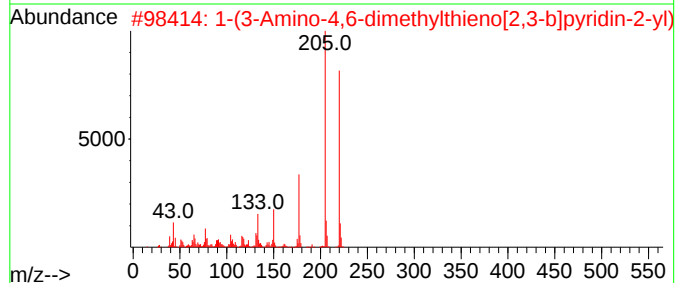
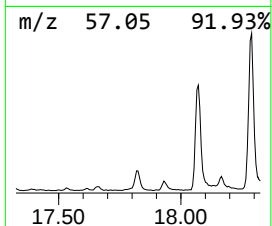
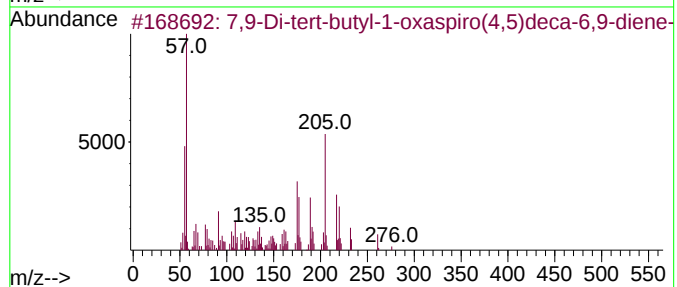
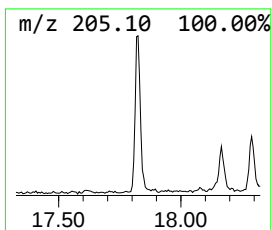
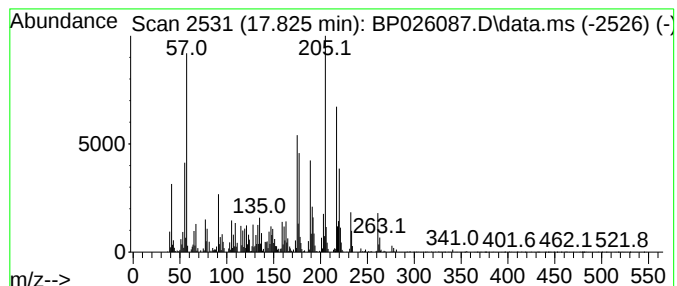
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TIC Library : C:\Database\NIST20.L
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Peak Number 6 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.825	2.78 ng	512705	Phenanthrene-d10	17.107

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	95
2			1-(3-Amino-4,6-dimethylthieno[2,...	220	C11H12N2OS	052505-42-7	55
3			2,5-di-tert-Butyl-1,4-benzoquinone	220	C14H20O2	002460-77-7	50
4			Ethanone, 1-(9-anthracenyl)-	220	C16H12O	000784-04-3	45
5			(1R,3aR,5aR,9aS)-1,4,4,7-Tetrame...	220	C15H24O	104188-25-2	42



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Misc :
ALS Vial : 13 Sample Multiplier: 1

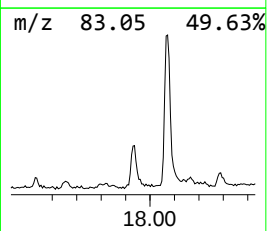
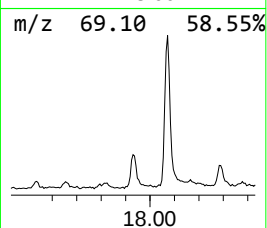
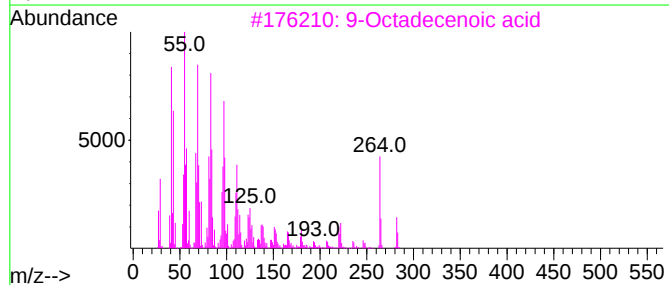
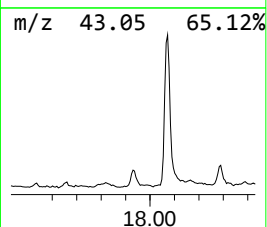
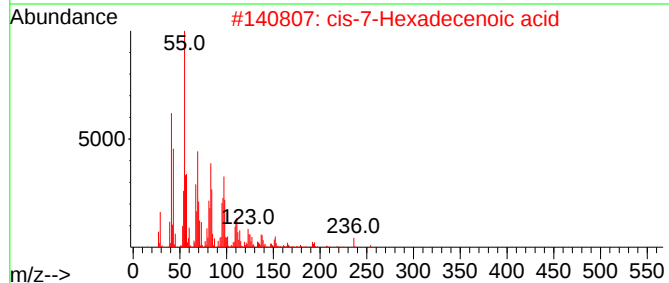
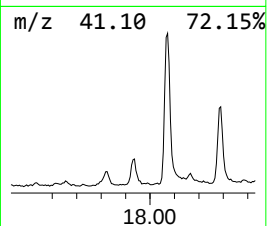
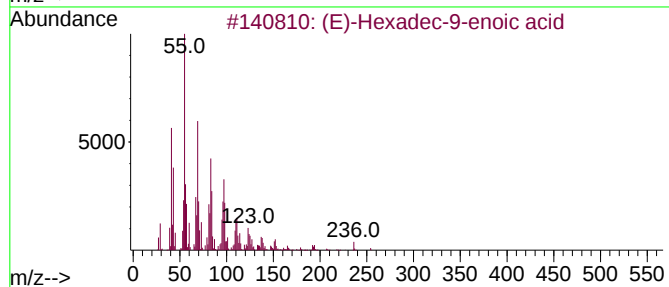
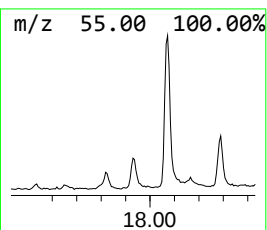
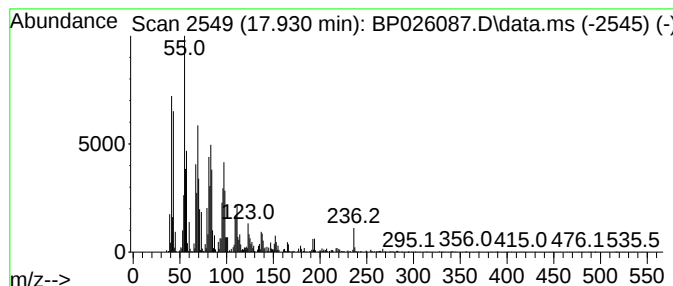
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 (E)-Hexadec-9-enoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.930	2.45 ng	450936	Phenanthrene-d10	17.107	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	99
2	cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	97
3	9-Octadecenoic acid	282	C18H34O2	002027-47-6	94
4	Palmitoleic acid	254	C16H30O2	000373-49-9	91
5	Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110625\
Data File : BP026087.D
Acq On : 06 Nov 2025 21:52
Operator : RC/JU
Sample : Q3552-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

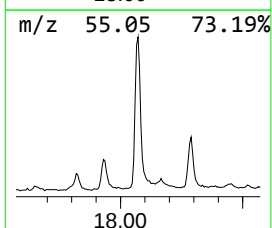
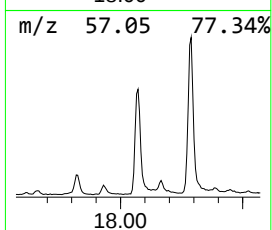
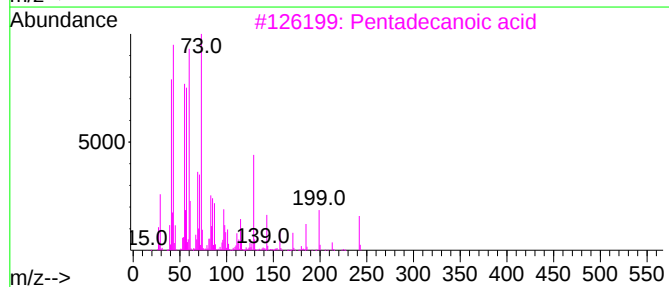
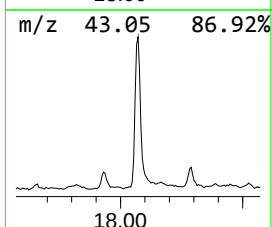
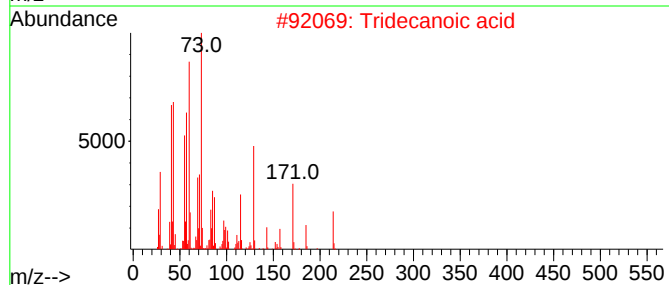
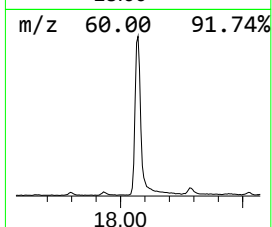
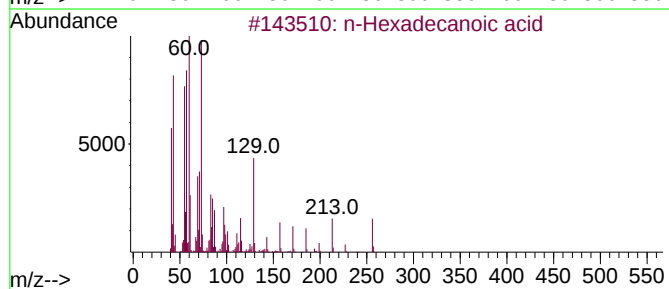
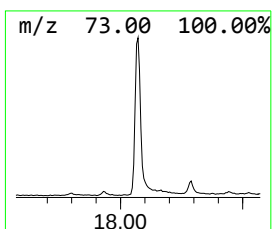
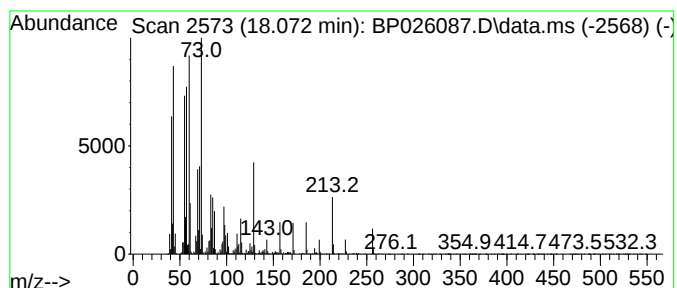
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.072	16.31 ng	3004640	Phenanthrene-d10	17.107	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	95
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4	Undecanoic acid	186	C11H22O2	000112-37-8	80
5	n-Decanoic acid	172	C10H20O2	000334-48-5	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110625\
Data File : BP026087.D
Acq On : 06 Nov 2025 21:52
Operator : RC/JU
Sample : Q3552-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

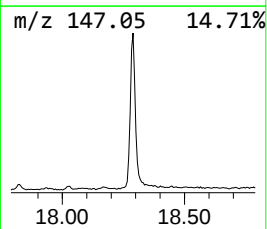
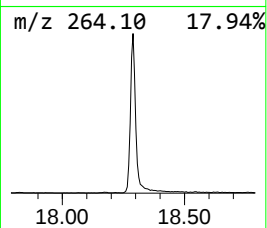
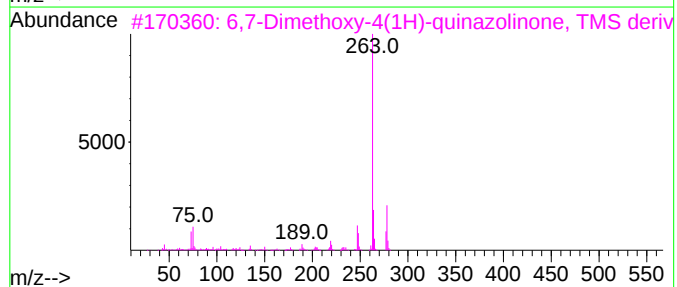
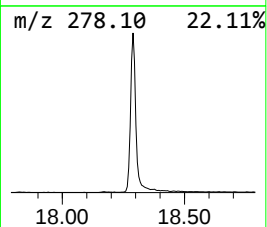
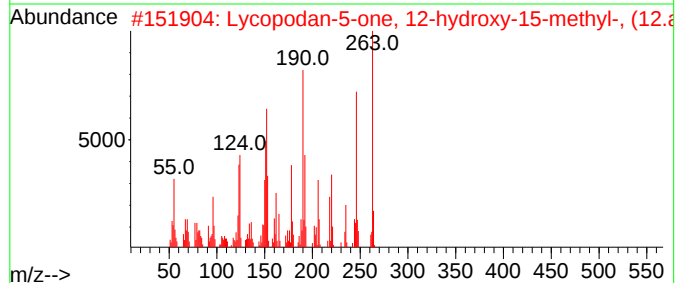
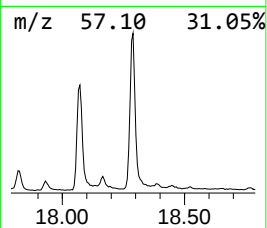
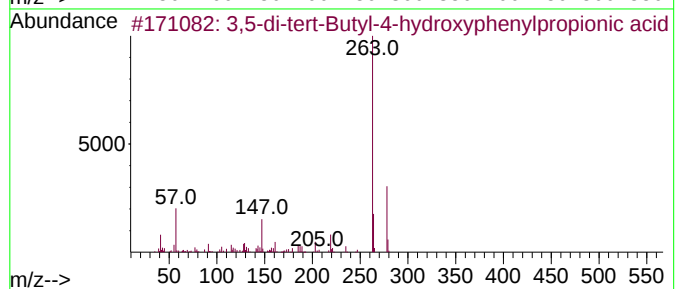
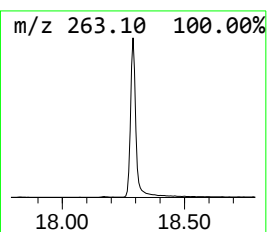
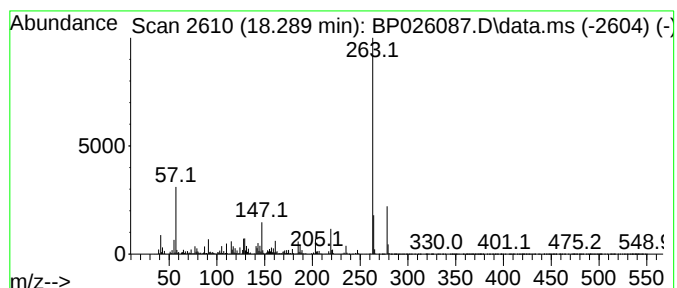
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.289	21.37 ng	3937260	Phenanthrene-d10		17.107	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	3,5-di-tert-Butyl-4-hydroxypheny...		278	C17H26O3	020170-32-5	94
2	Lycopodan-5-one, 12-hydroxy-15-m...		263	C16H25NO2	021061-90-5	91
3	6,7-Dimethoxy-4(1H)-quinazolinon...		278	C13H18N2O3Si	1000479-65-5	72
4	5-(2,3-Difluorophenyl)-2-pyridin...		278	C14H16F2N2Si	1000504-49-4	64
5	Cyanic acid, (1-methylethylidene...		278	C17H14N2O2	001156-51-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110625\
Data File : BP026087.D
Acq On : 06 Nov 2025 21:52
Operator : RC/JU
Sample : Q3552-02
Misc :
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Instrument :
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ClientSampleId :
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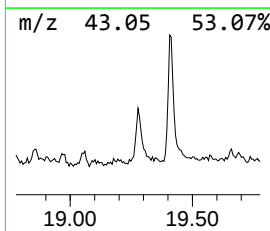
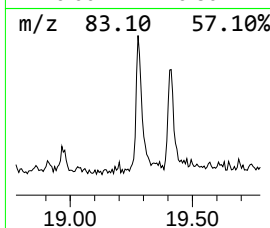
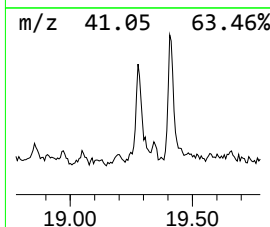
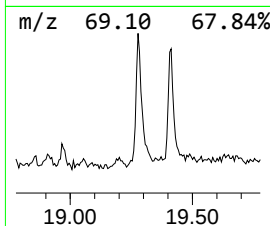
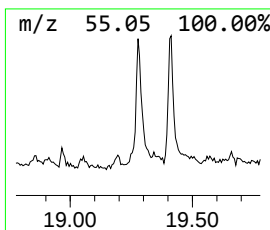
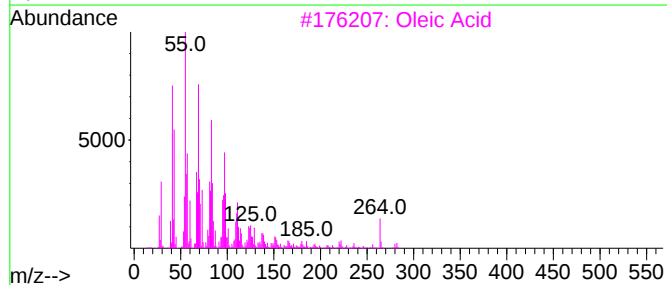
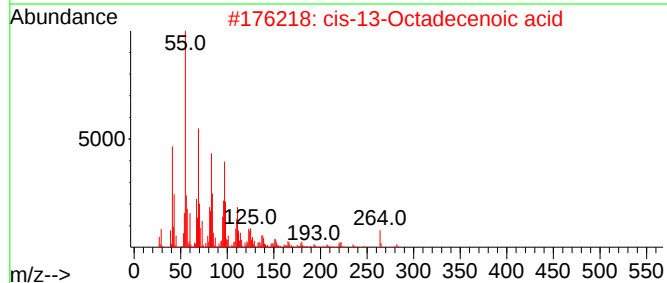
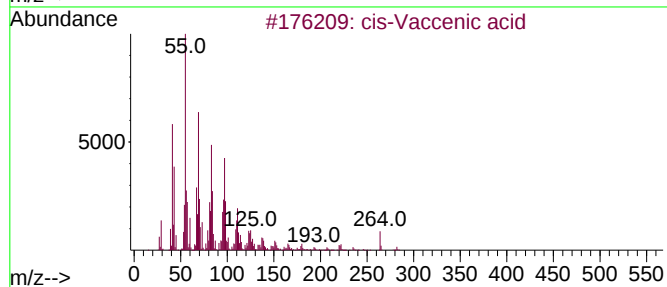
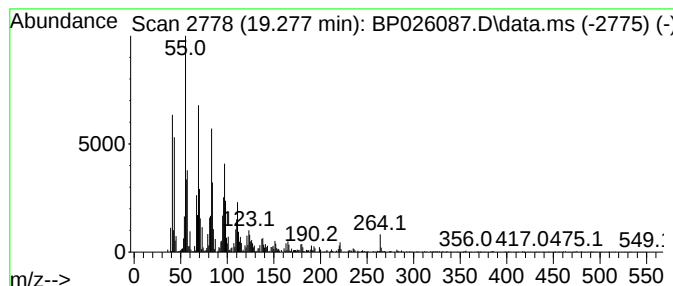
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 cis-Vaccenic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.277	2.37 ng	435837	Phenanthrene-d10	17.107

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-Vaccenic acid	282	C18H34O2	000506-17-2	94
2			cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	94
3			Oleic Acid	282	C18H34O2	000112-80-1	94
4			trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	93
5			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110625\
Data File : BP026087.D
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Sample : Q3552-02
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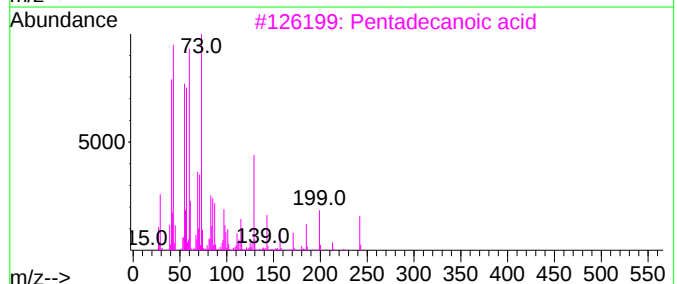
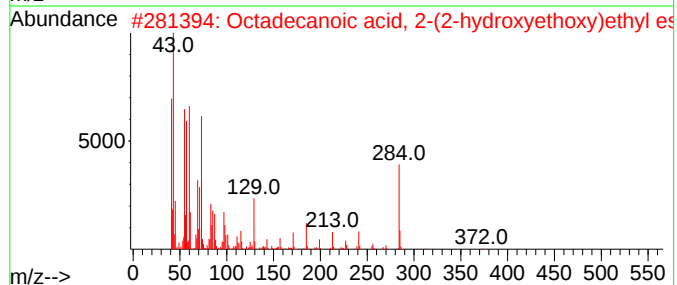
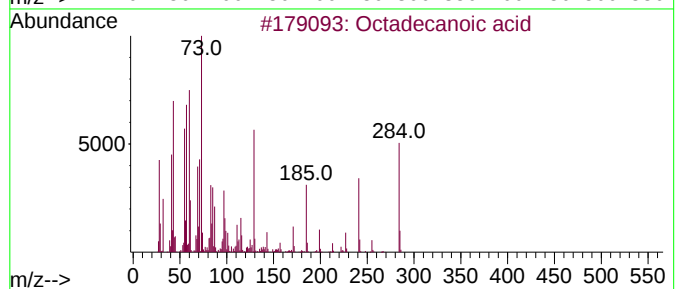
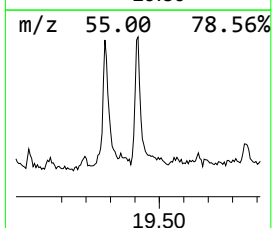
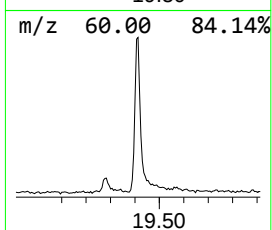
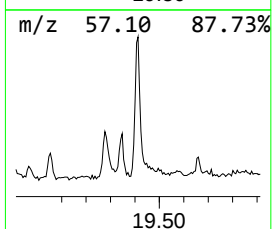
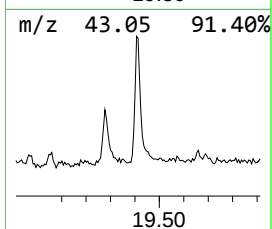
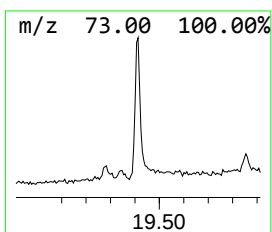
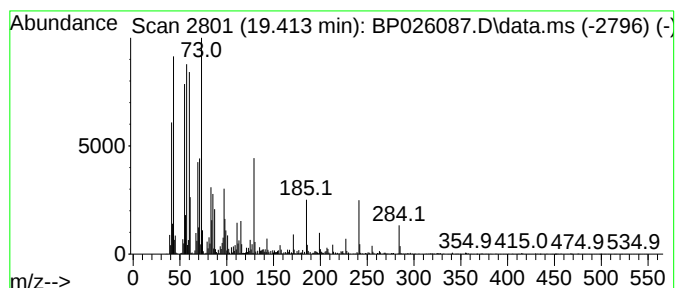
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.413	3.50 ng	786558	Chrysene-d12	21.566	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	93
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5	Tridecanoic acid	214	C13H26O2	000638-53-9	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110625\
Data File : BP026087.D
Acq On : 06 Nov 2025 21:52
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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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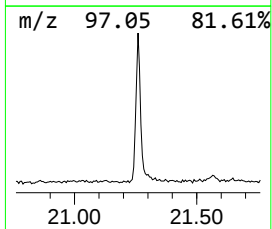
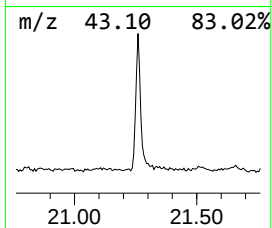
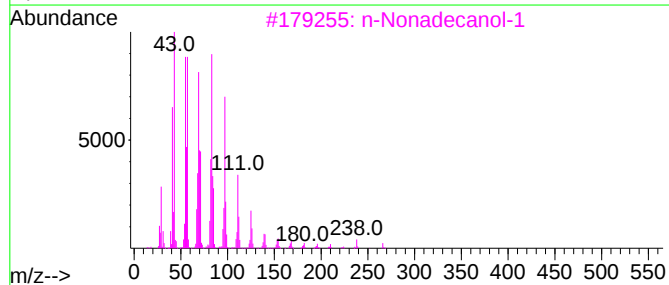
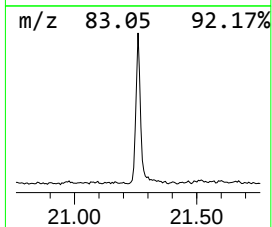
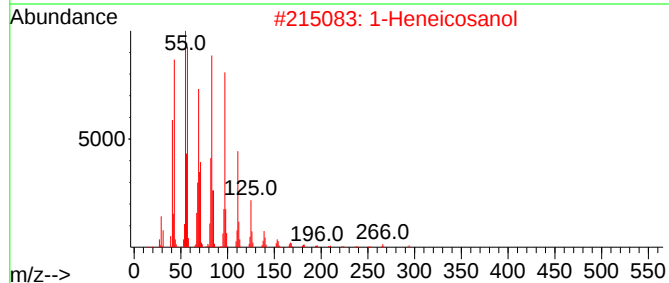
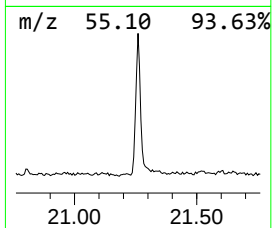
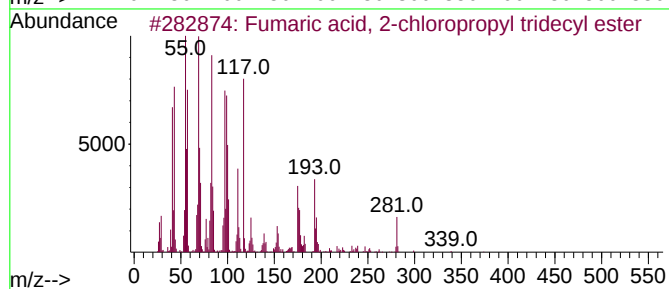
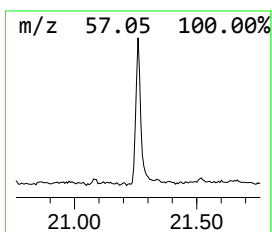
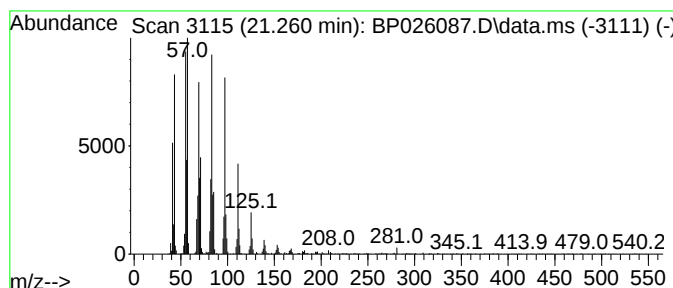
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Fumaric acid, 2-chloropropyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.260	5.39 ng	1210780	Chrysene-d12	21.566

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fumaric acid, 2-chloropropyl tri...	374	C20H35ClO4	1010348-57-1	97
2			1-Heneicosanol	312	C21H44O	015594-90-8	95
3			n-Nonadecanol-1	284	C19H40O	001454-84-8	95
4			Octacosyl acetate	452	C30H60O2	018206-97-8	95
5			1-Heneicosyl formate	340	C22H44O2	077899-03-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110625\
Data File : BP026087.D
Acq On : 06 Nov 2025 21:52
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Sample : Q3552-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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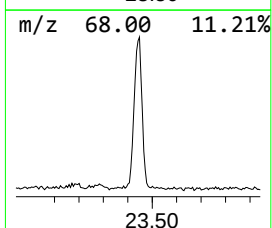
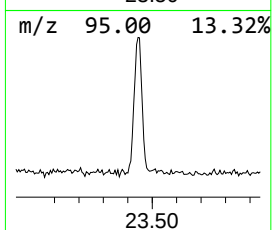
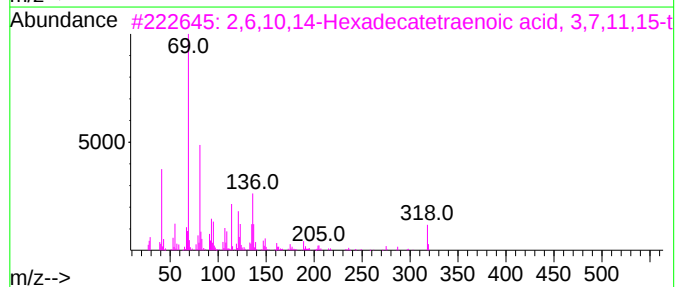
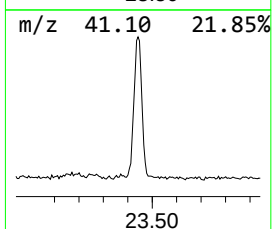
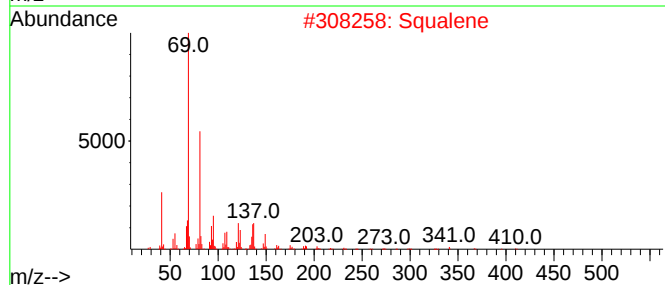
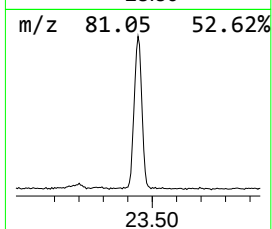
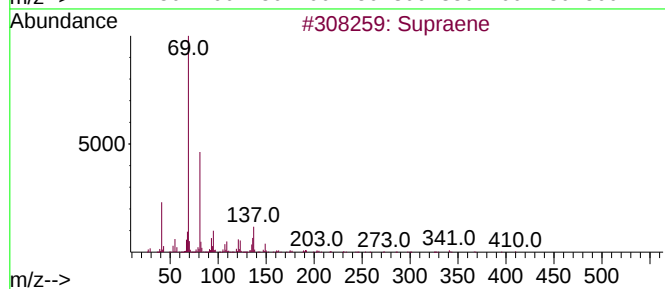
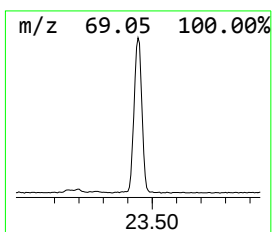
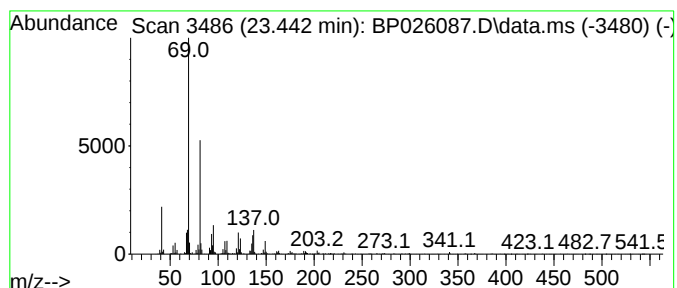
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Supraene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.442	6.39 ng	1722840	Perylene-d12	24.889

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	91
2			Squalene	410	C30H50	000111-02-4	91
3			2,6,10,14-Hexadecatetraenoic aci...	318	C21H34O2	042207-88-5	78
4			Farnesol isomer a	222	C15H26O	1000108-92-4	64
5			2,3,5-Trimethyl-2,3,5-hexanetric...	203	C12H17N3	003974-79-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110625\
Data File : BP026087.D
Acq On : 06 Nov 2025 21:52
Operator : RC/JU
Sample : Q3552-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Butanoic acid, ...	4.596	3.1	ng	272029	1	7.678	1770480	20.0
Benzeneacetic acid	10.972	3.4	ng	402227	2	10.443	2400770	20.0
Benzoic acid, p...	14.107	3.0	ng	469151	3	14.307	3169320	20.0
Benzophenone	15.689	4.6	ng	726411	3	14.307	3169320	20.0
7,9-Di-tert-but...	17.825	2.8	ng	512705	4	17.107	3684370	20.0
(E)-Hexadec-9-e...	17.930	2.5	ng	450936	4	17.107	3684370	20.0
n-Hexadecanoic ...	18.072	16.3	ng	3004640	4	17.107	3684370	20.0
3,5-di-tert-But...	18.289	21.4	ng	3937260	4	17.107	3684370	20.0
cis-Vaccenic acid	19.277	2.4	ng	435837	4	17.107	3684370	20.0
Octadecanoic acid	19.413	3.5	ng	786558	5	21.566	4492960	20.0
Fumaric acid, 2...	21.260	5.4	ng	1210780	5	21.566	4492960	20.0
Supraene	23.442	6.4	ng	1722840	6	24.889	5395740	20.0