

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111820\
 Data File : BP004002.D
 Acq On : 18 Nov 2020 19:58
 Operator : CG/JU
 Sample : L4770-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FJ-BH-01-0-1

Integration Parameters: rteint.p

Integrator: RTE

Smoother: 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\8270-BP110320.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP004002.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.922	3	6	25	rVB	3568848	5186289	63.43%	10.709%
2	3.081	27	33	45	rVV	3897366	8175788	100.00%	16.881%
3	3.175	45	49	58	rVB	299774	486563	5.95%	1.005%
4	3.387	80	85	97	rVB	59094	112246	1.37%	0.232%
5	5.281	401	407	414	rVB	139552	209746	2.57%	0.433%
6	5.811	489	497	515	rBV	2219656	3398413	41.57%	7.017%
7	7.463	770	778	790	rBV	2061451	3402389	41.62%	7.025%
8	7.852	839	844	851	rBV	83687	126779	1.55%	0.262%
9	8.287	911	918	925	rBV	575130	938580	11.48%	1.938%
10	9.452	1103	1116	1125	rBV	1291354	2228544	27.26%	4.601%
11	11.116	1389	1399	1405	rBV	736570	1295745	15.85%	2.675%
12	13.528	1802	1809	1822	rBV	3054165	4606484	56.34%	9.511%
13	13.728	1838	1843	1851	rBV5	58289	136853	1.67%	0.283%
14	14.334	1941	1946	1951	rBV	476120	701570	8.58%	1.449%
15	14.381	1951	1954	1958	rVB	67543	84582	1.03%	0.175%
16	14.634	1992	1997	2002	rBV	69977	114275	1.40%	0.236%
17	14.910	2038	2044	2051	rVB2	1009076	1505356	18.41%	3.108%
18	15.728	2180	2183	2188	rVB2	66997	84665	1.04%	0.175%
19	16.139	2246	2253	2261	rVB2	76387	172393	2.11%	0.356%
20	16.392	2291	2296	2310	rBV	2007602	2859877	34.98%	5.905%
21	16.616	2330	2334	2340	rVB2	175885	256075	3.13%	0.529%
22	17.151	2421	2425	2430	rBV2	139475	228361	2.79%	0.472%
23	17.433	2469	2473	2477	rBV	111550	154114	1.89%	0.318%
24	17.645	2503	2509	2513	rBV	1171167	1686506	20.63%	3.482%
25	17.686	2513	2516	2524	rVV	222812	345185	4.22%	0.713%
26	17.775	2528	2531	2537	rVB	95858	138470	1.69%	0.286%
27	18.486	2648	2652	2656	rBV2	131554	216078	2.64%	0.446%
28	18.580	2665	2668	2671	rBV3	66941	91162	1.12%	0.188%
29	18.680	2681	2685	2688	rBV4	92306	155146	1.90%	0.320%
30	19.227	2776	2778	2785	rBV2	135074	216944	2.65%	0.448%
31	19.622	2841	2845	2850	rVB	436785	574468	7.03%	1.186%
32	19.969	2901	2904	2911	rVB	494630	637137	7.79%	1.316%
33	20.139	2928	2933	2939	rVB	4364057	5411003	66.18%	11.173%
34	21.327	3132	3135	3146	rVB	814946	1130831	13.83%	2.335%
35	21.680	3192	3195	3199	rVB	1096004	1362651	16.67%	2.814%

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ClientSampleId :
FJ-BH-01-0-1

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\8270-BP110320.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

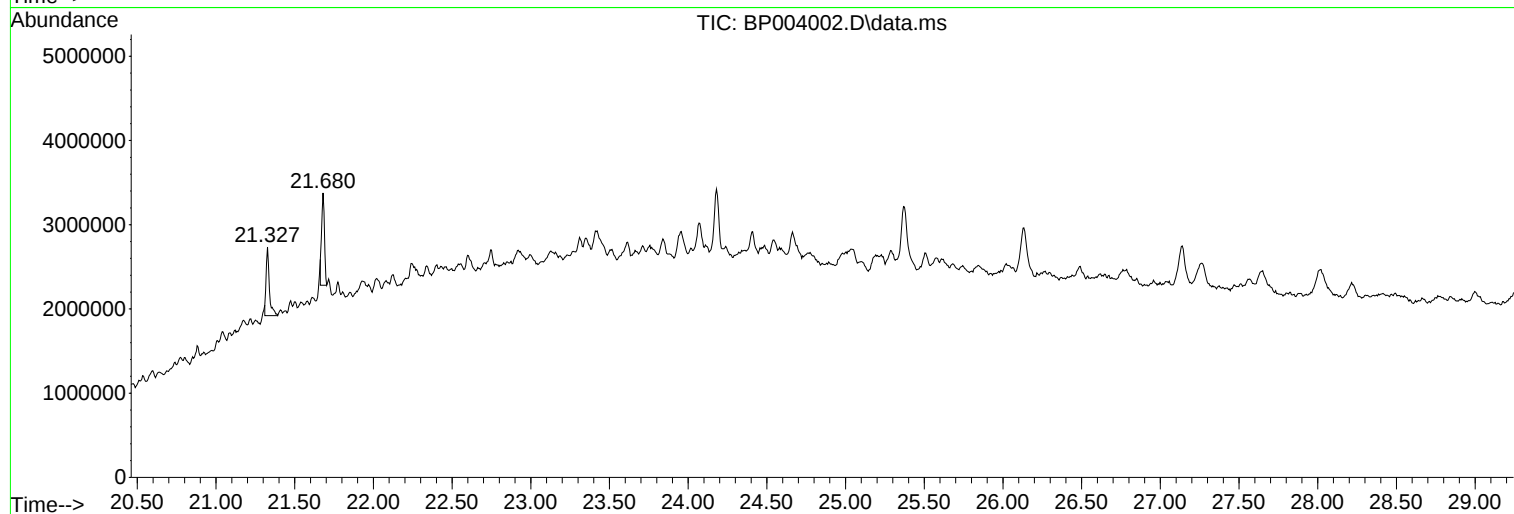
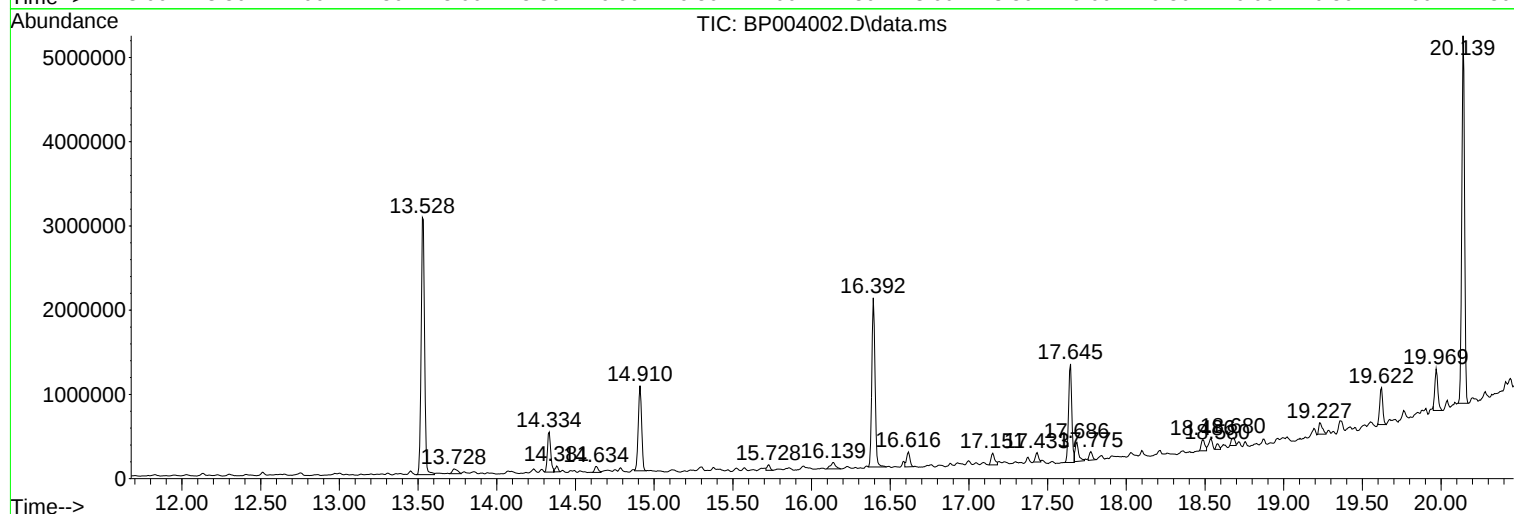
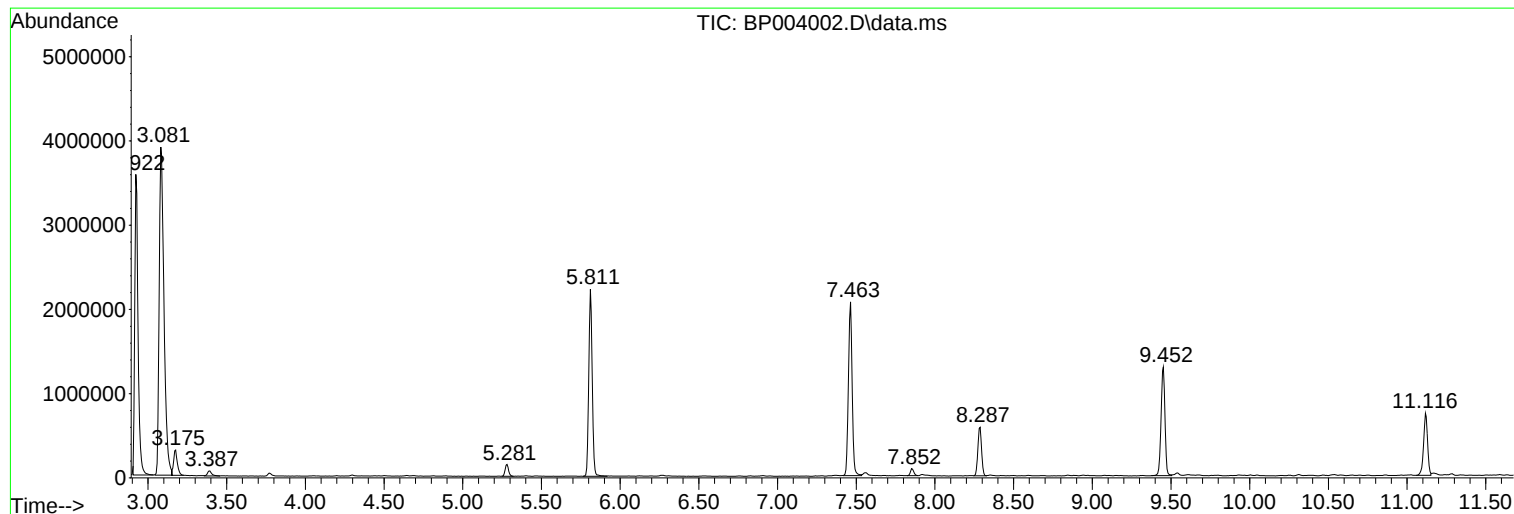
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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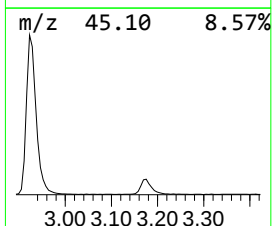
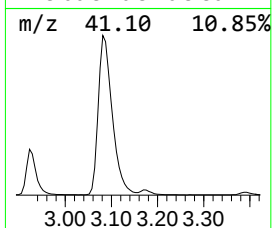
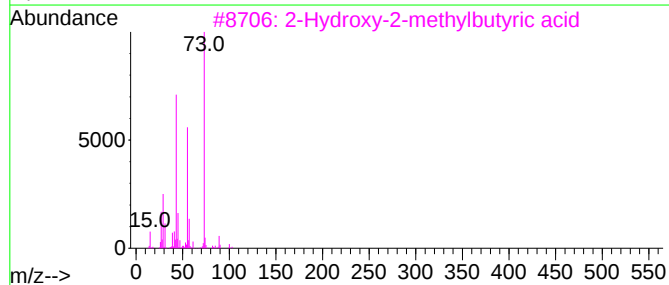
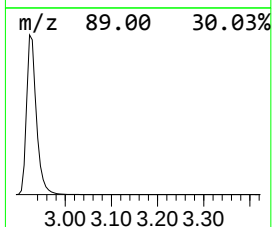
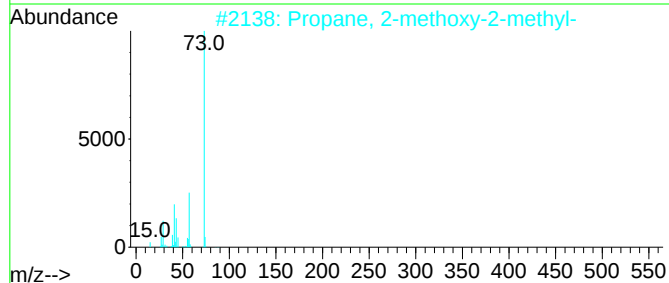
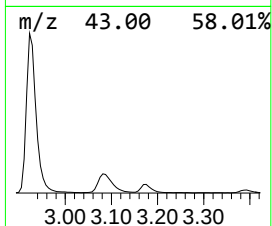
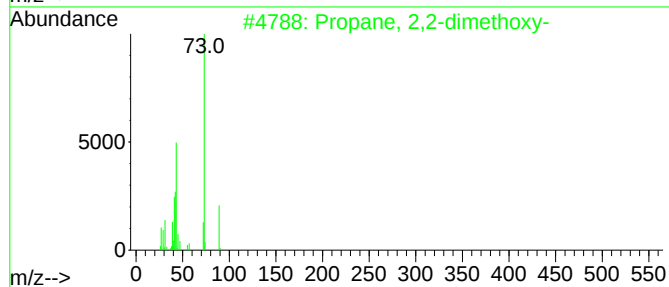
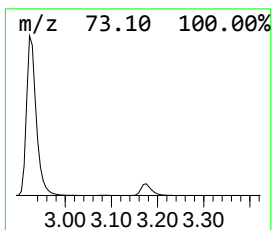
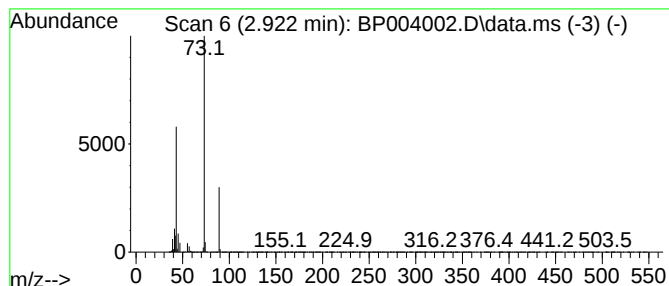
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.922	110.51 ng	5186290	1,4-Dichlorobenzene-d4	8.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	72
2			Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
3			2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
4			1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9
5			2-Butanone, 3-methoxy-3-methyl-	116	C6H12O2	036687-98-6	9



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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
FJ-BH-01-0-1

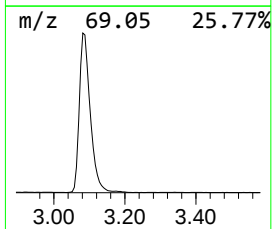
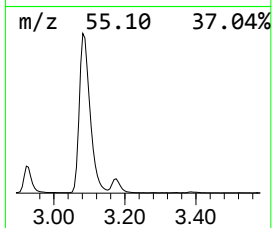
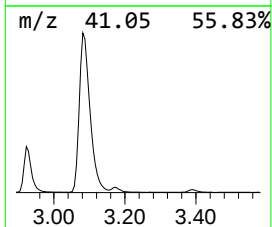
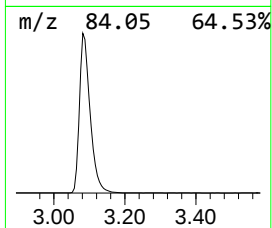
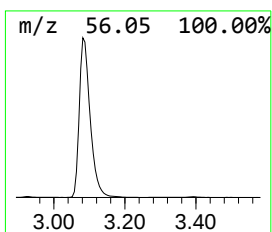
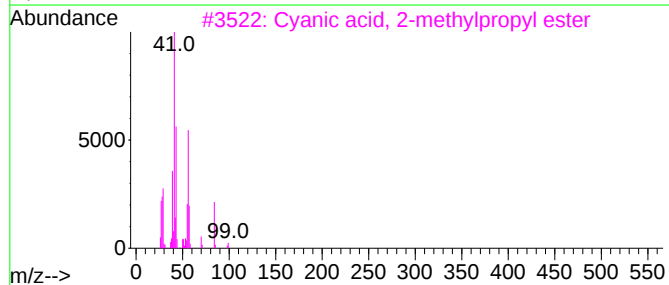
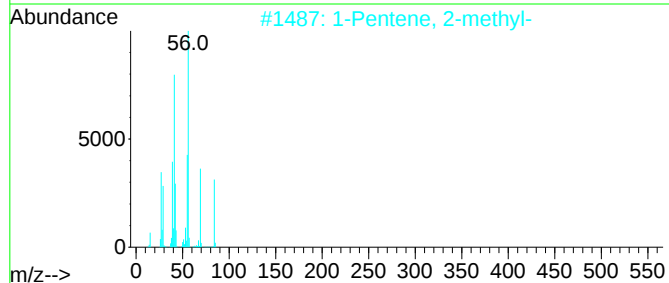
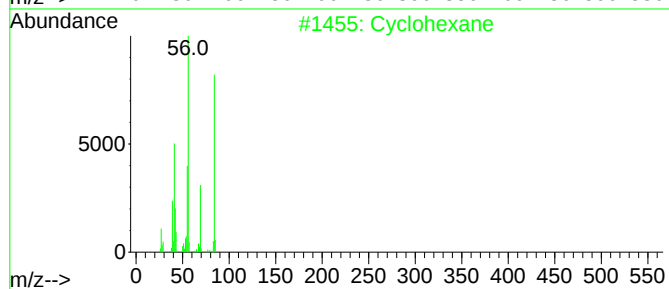
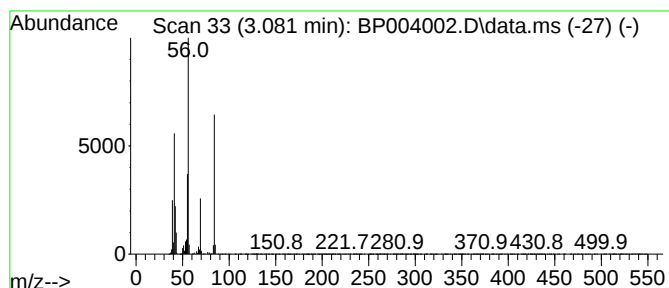
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Cyclohexane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.081	174.22 ng	8175790	1,4-Dichlorobenzene-d4	8.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane	84	C6H12	000110-82-7	95
2			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72
3			Cyanic acid, 2-methylpropyl ester	99	C5H9NO	001768-25-8	64
4			1-Hexene	84	C6H12	000592-41-6	53
5			Cyclopentane, methyl-	84	C6H12	000096-37-7	50



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Instrument :
BNA_P
ClientSampleId :
FJ-BH-01-0-1

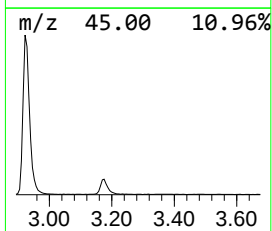
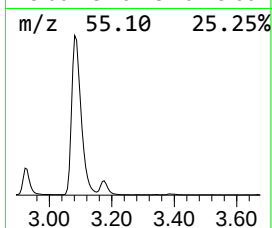
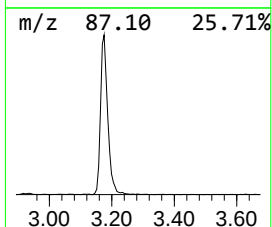
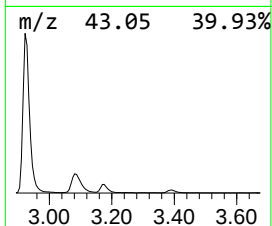
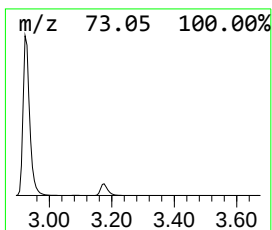
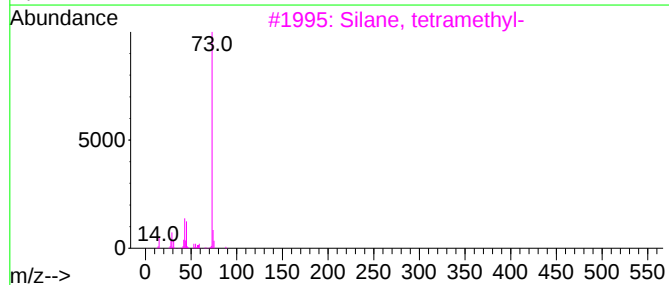
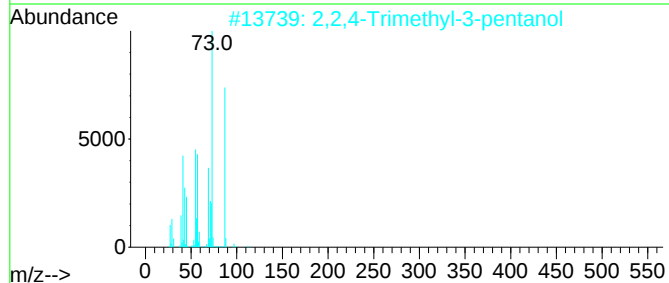
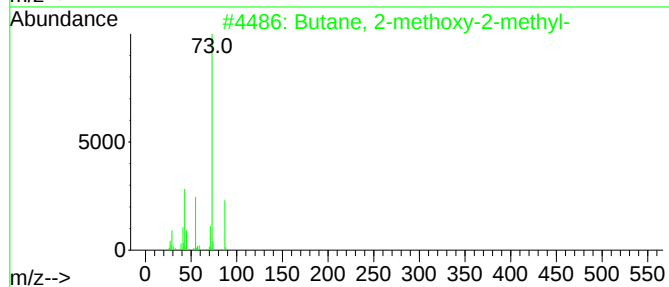
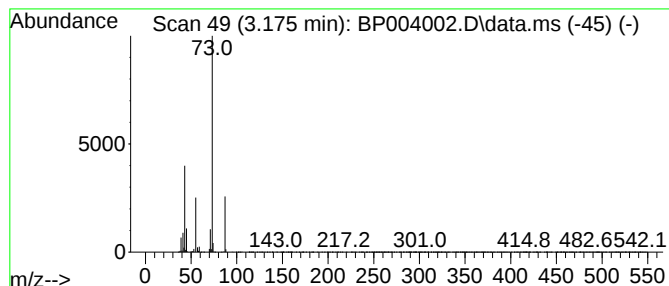
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.175	10.37 ng	486563	1,4-Dichlorobenzene-d4	8.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	23
4			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
FJ-BH-01-0-1

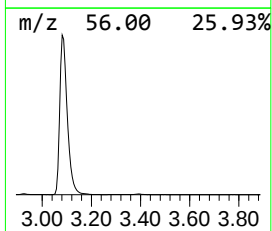
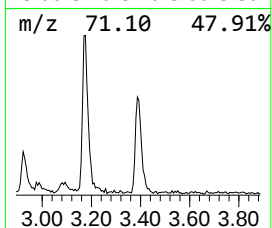
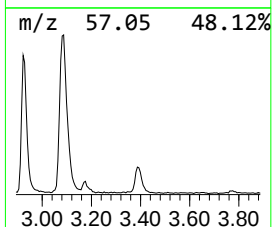
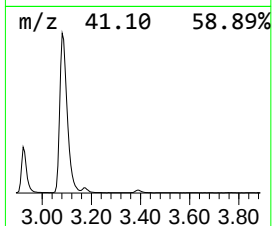
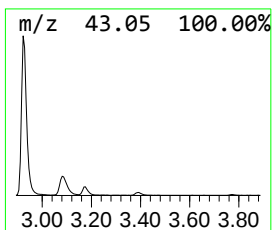
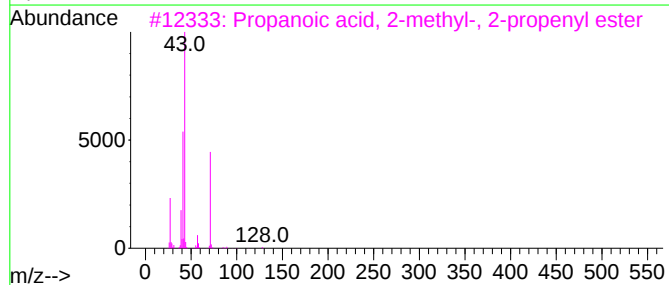
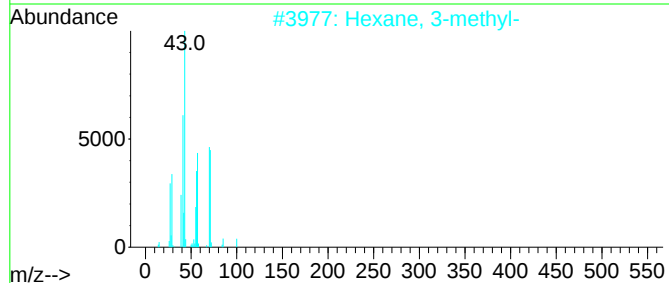
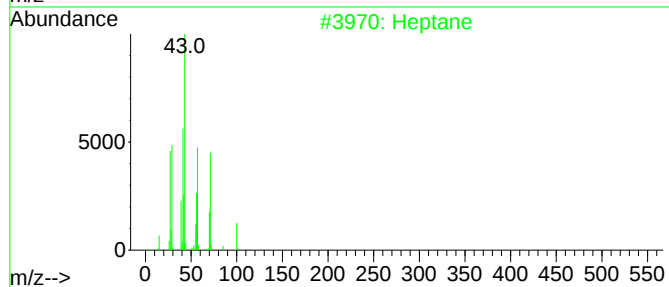
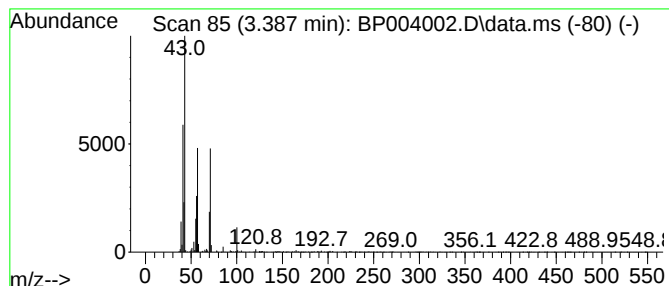
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Heptane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.387	2.39 ng	112246	1,4-Dichlorobenzene-d4	8.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	91
2			Hexane, 3-methyl-	100	C7H16	000589-34-4	40
3			Propanoic acid, 2-methyl-, 2-pro...	128	C7H12O2	015727-77-2	38
4			Octane, 3-ethyl-	142	C10H22	005881-17-4	37
5			Ether, 6-methylheptyl vinyl	156	C10H20O	010573-35-0	33



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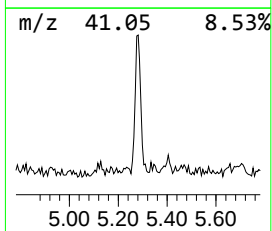
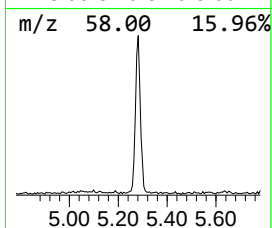
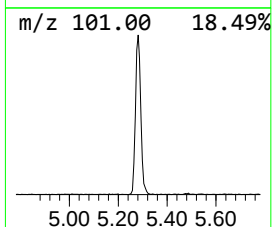
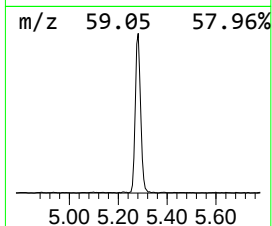
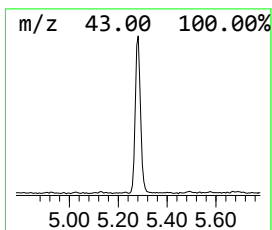
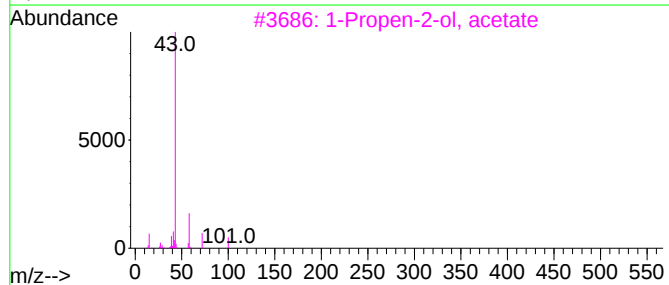
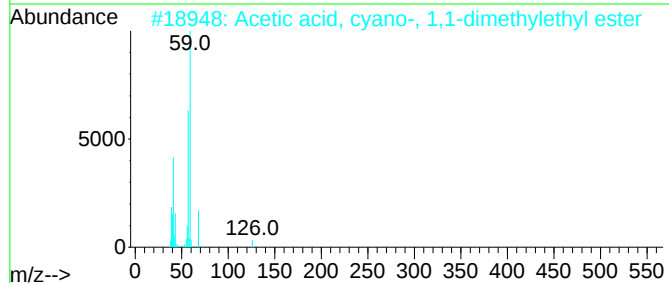
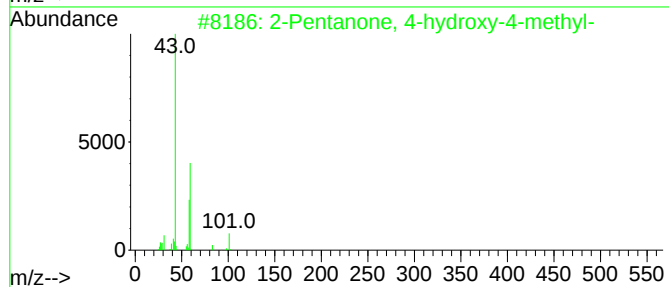
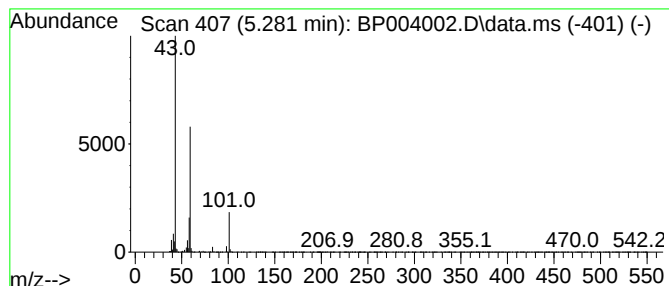
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.281	4.47 ng	209746	1,4-Dichlorobenzene-d4	8.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111820\
Data File : BP004002.D
Acq On : 18 Nov 2020 19:58
Operator : CG/JU
Sample : L4770-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

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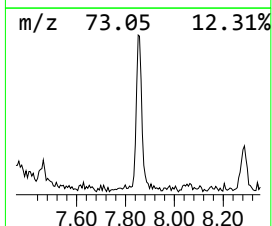
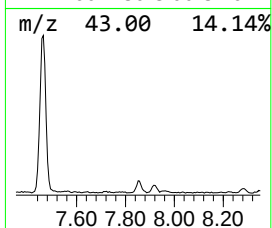
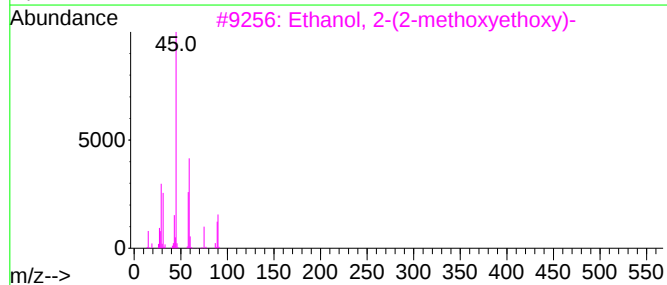
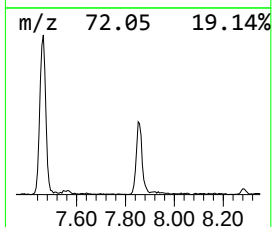
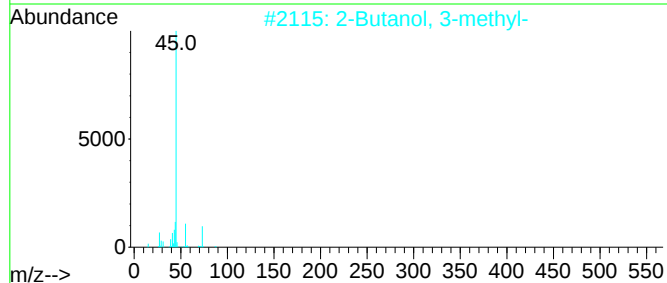
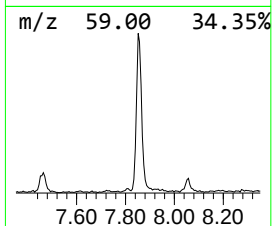
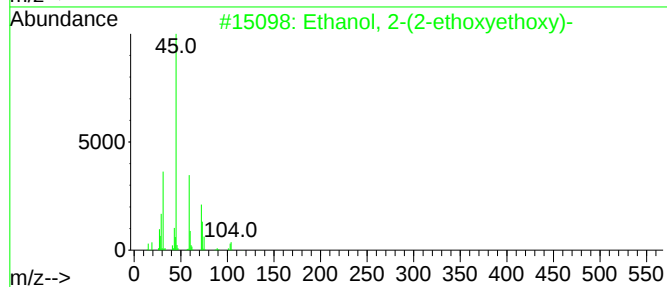
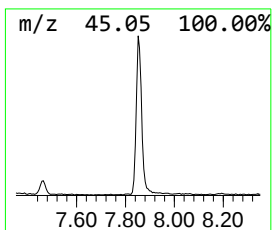
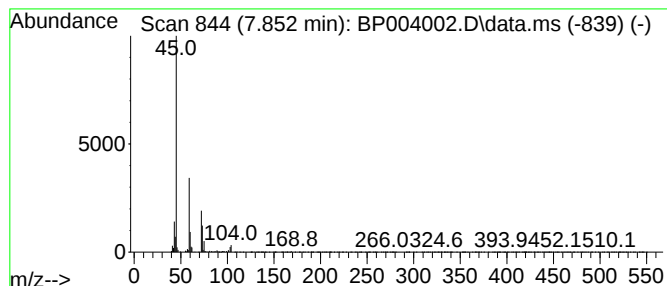
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.852	2.70 ng	126779	1,4-Dichlorobenzene-d4	8.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2			2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	47
3			Ethanol, 2-(2-methoxyethoxy)-	120	C5H12O3	000111-77-3	45
4			2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	45
5			L-Lactic acid	90	C3H6O3	000079-33-4	37



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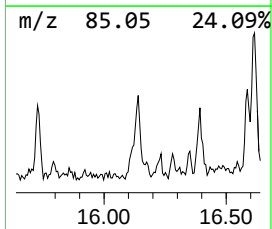
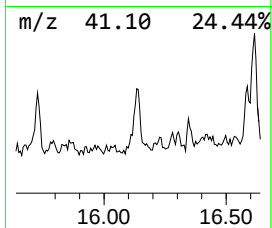
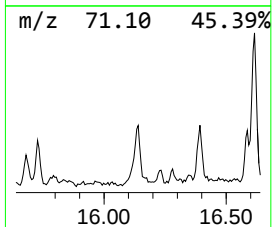
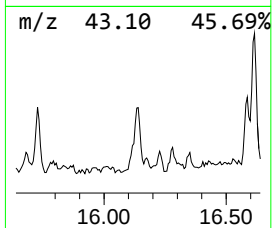
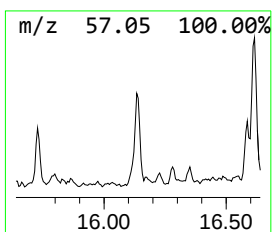
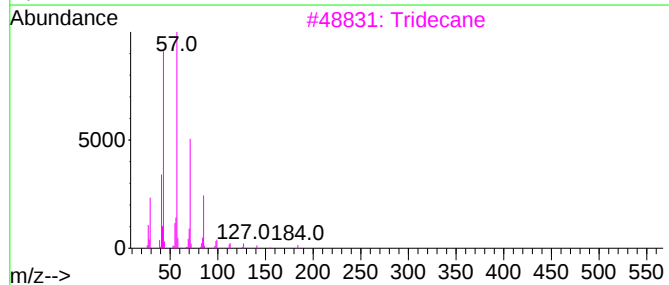
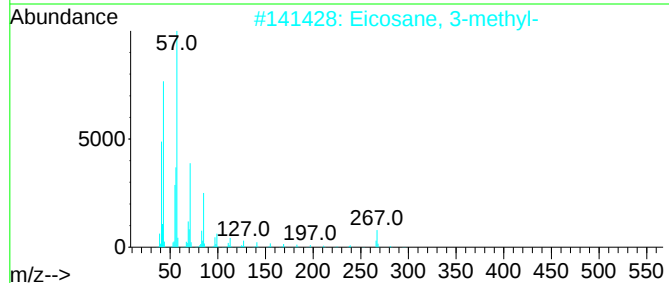
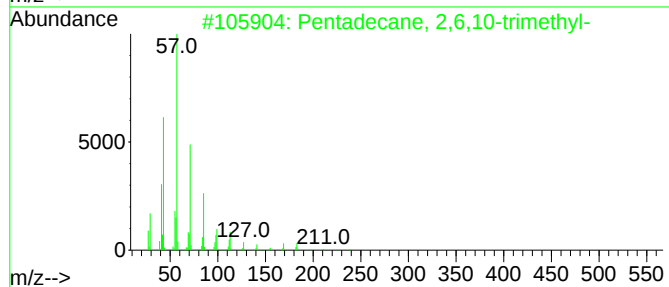
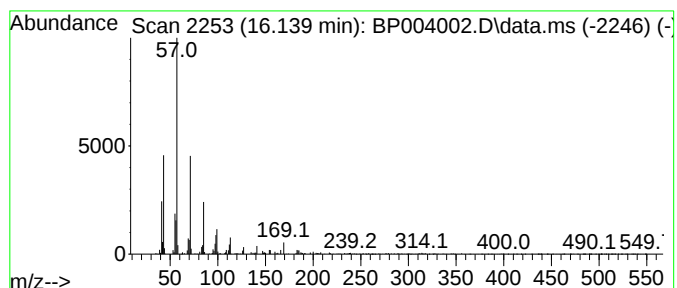
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pentadecane, 2,6,10-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.139	2.29 ng	172393	Acenaphthene-d10	14.910	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	90
2	Eicosane, 3-methyl-	296	C21H44	006418-46-8	86
3	Tridecane	184	C13H28	000629-50-5	72
4	Octacosane	394	C28H58	000630-02-4	72
5	Heneicosane	296	C21H44	000629-94-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111820\
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Sample : L4770-01
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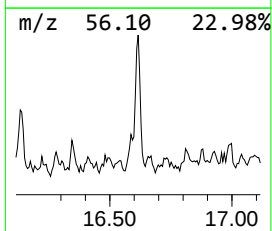
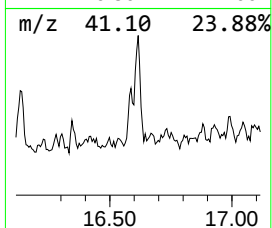
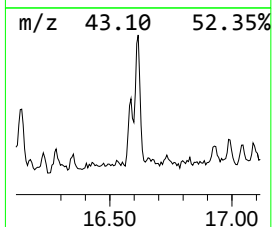
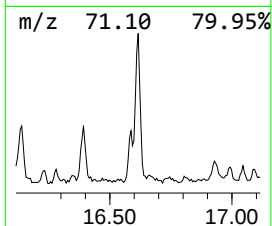
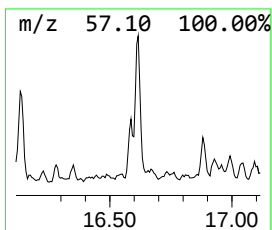
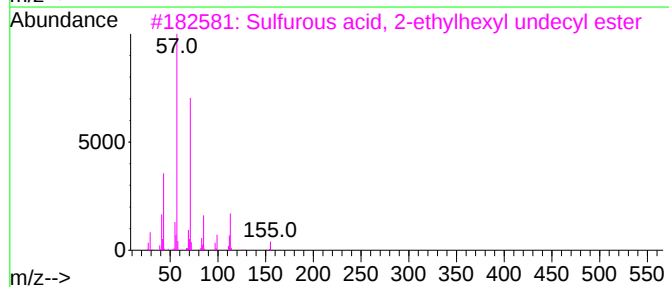
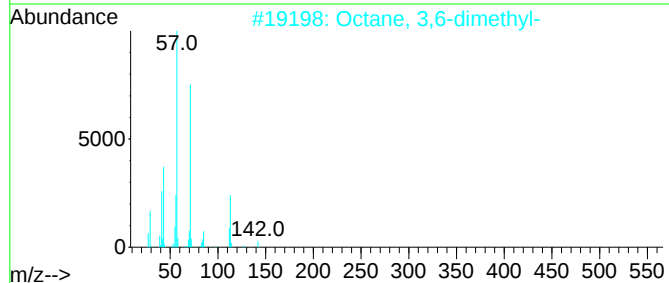
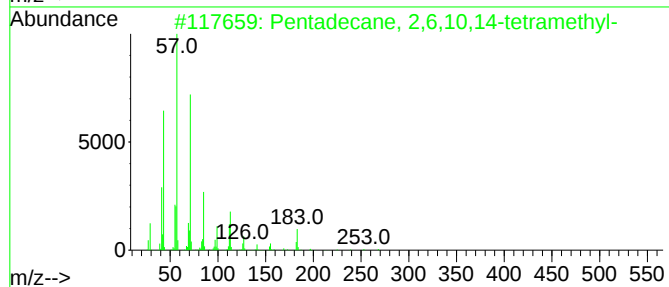
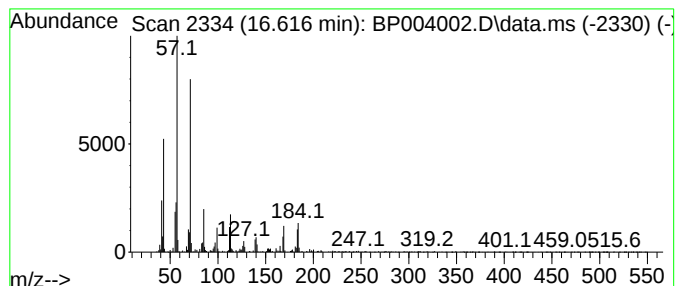
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pentadecane, 2,6,10,14-tetr... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.616	3.04 ng	256075	Phenanthrene-d10		17.645	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10,14-tetramethyl-		268	C19H40	001921-70-6	86
2	Octane, 3,6-dimethyl-		142	C10H22	015869-94-0	70
3	Sulfurous acid, 2-ethylhexyl und...		348	C19H40O3S	1000309-19-4	68
4	Sulfurous acid, decyl 2-ethylhex...		334	C18H38O3S	1000309-19-3	64
5	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111820\
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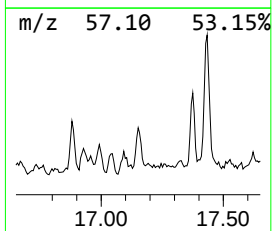
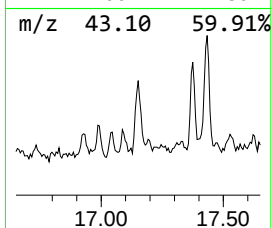
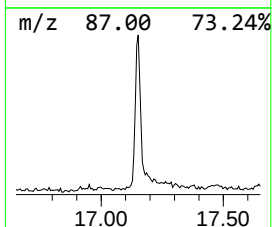
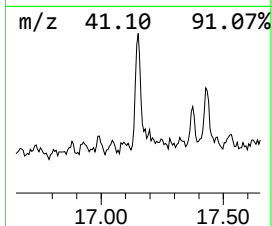
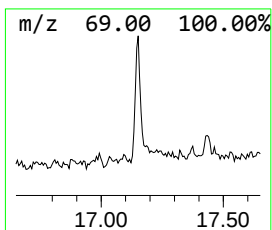
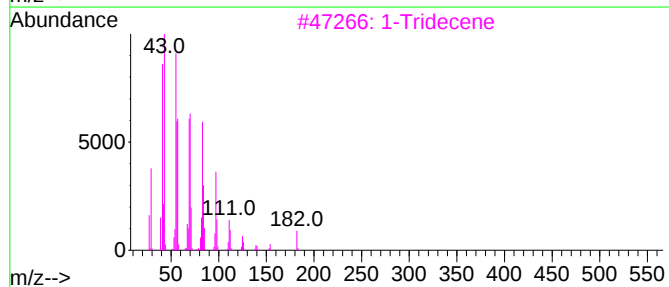
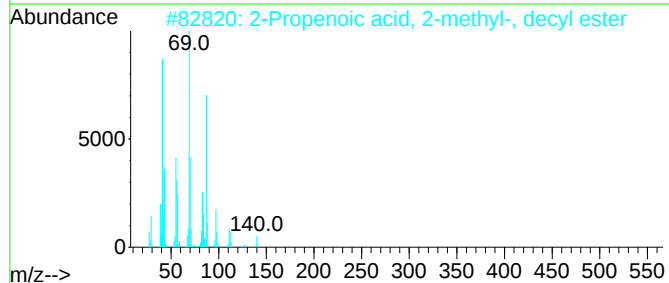
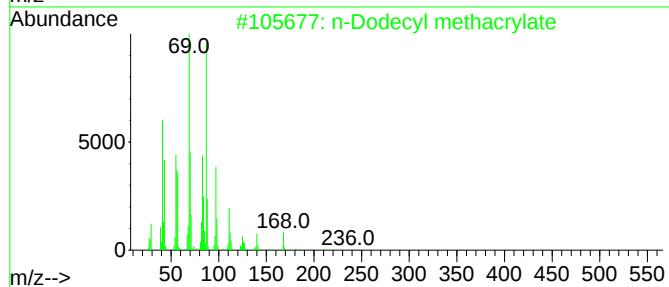
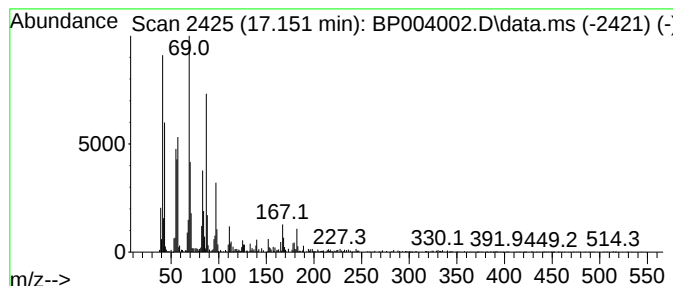
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Dodecyl methacrylate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.151	2.71 ng	228361	Phenanthrene-d10	17.645	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Dodecyl methacrylate	254	C16H30O2	000142-90-5	87
2	2-Propenoic acid, 2-methyl-, dec...	226	C14H26O2	003179-47-3	83
3	1-Tridecene	182	C13H26	002437-56-1	81
4	Methacrylic acid, tetradecyl ester	282	C18H34O2	1000340-29-0	64
5	Cyclohexane, 2-butyl-1,1,3-trime...	182	C13H26	054676-39-0	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111820\
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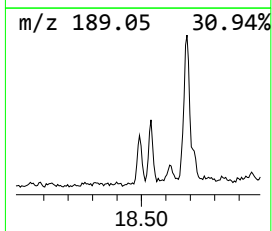
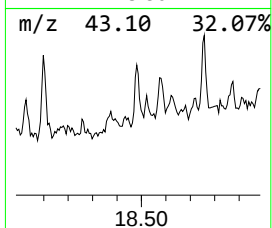
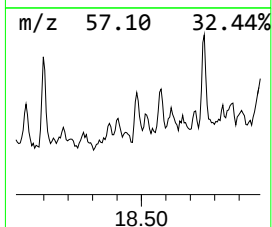
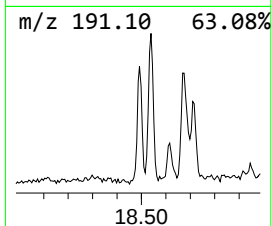
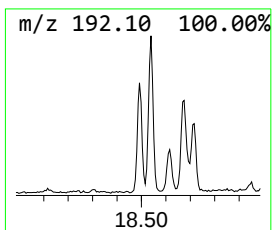
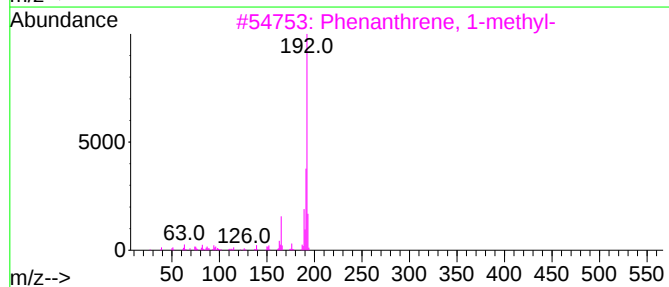
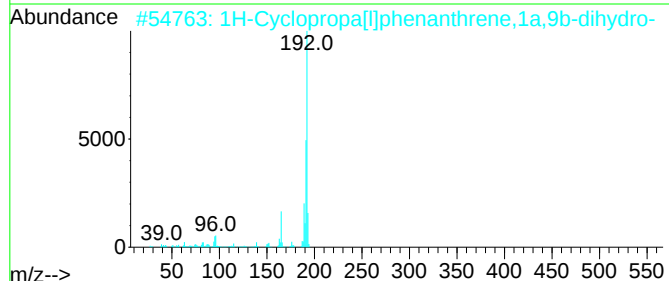
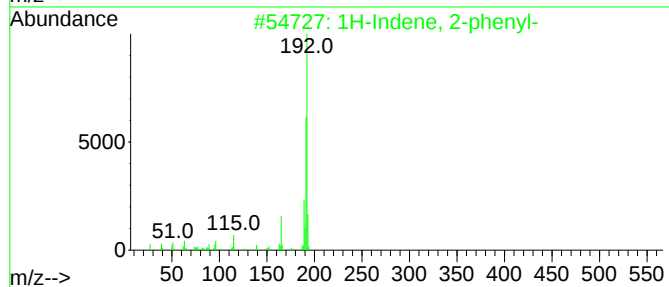
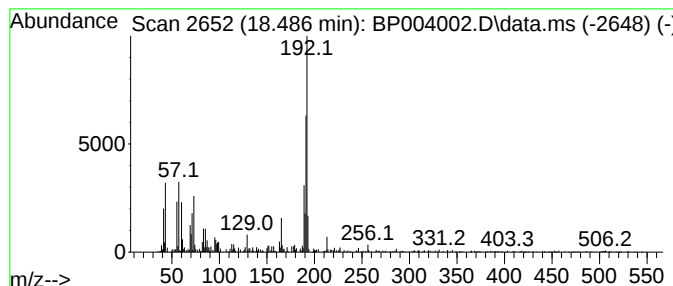
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1H-Indene, 2-phenyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.486	2.56 ng	216078	Phenanthrene-d10	17.645

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	91
2		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	90
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	86
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	78
5		Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111820\
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Acq On : 18 Nov 2020 19:58
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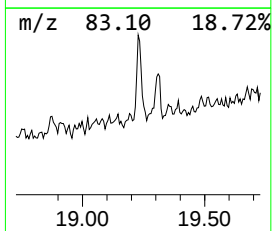
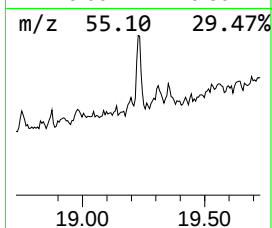
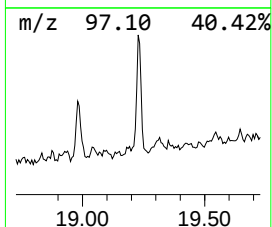
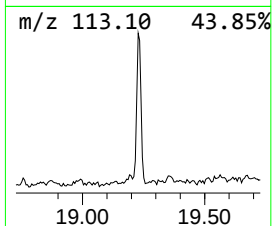
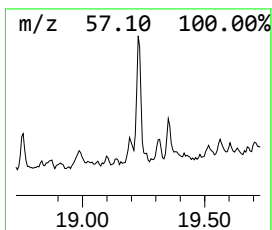
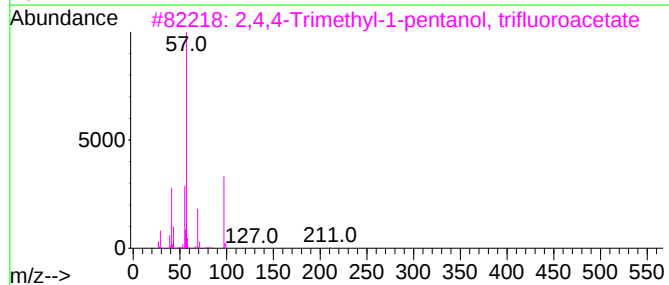
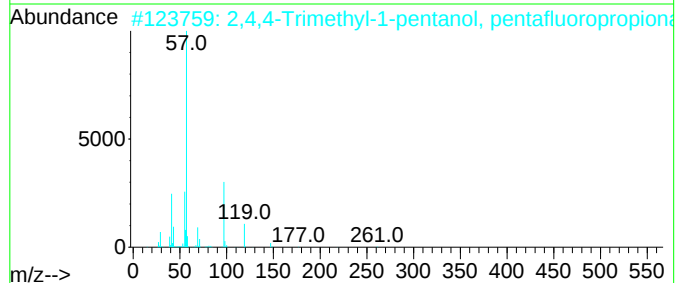
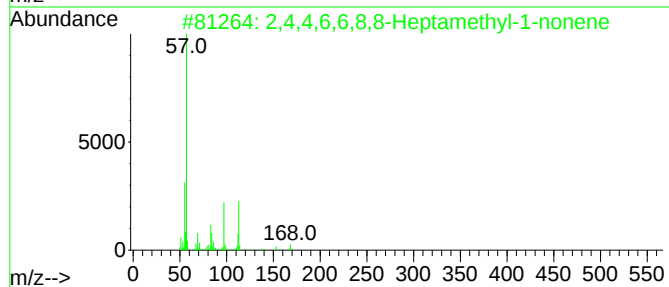
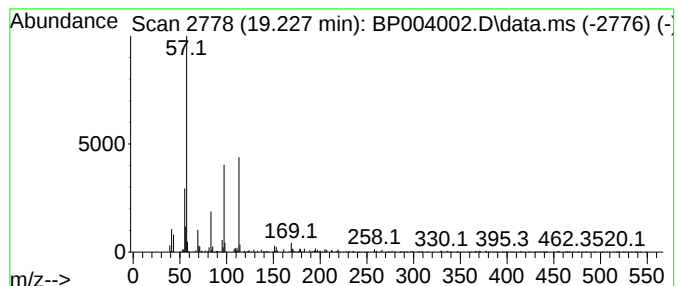
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown19.227 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.227	2.57 ng	216944	Phenanthrene-d10	17.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	43		
2	2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F502	1000365-51-0	38		
3	2,4,4-Trimethyl-1-pentanol, trif...	226	C10H17F302	1000365-19-5	38		
4	Octane, 2-bromo-	192	C8H17Br	000557-35-7	37		
5	2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F702	1000365-01-4	32		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111820\
Data File : BP004002.D
Acq On : 18 Nov 2020 19:58
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ClientSampleId :
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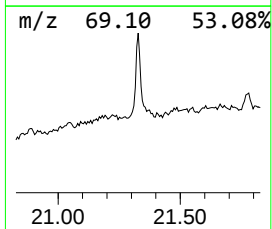
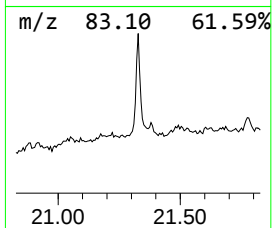
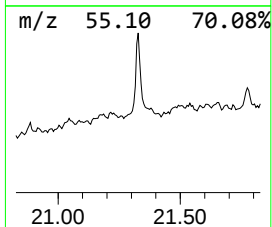
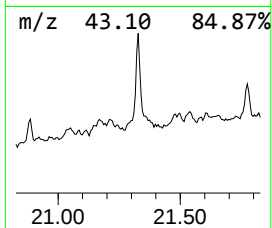
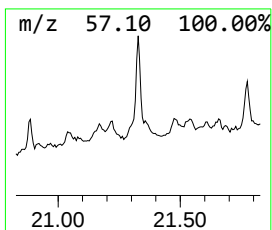
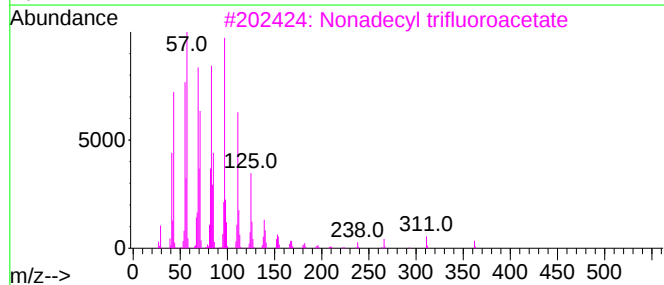
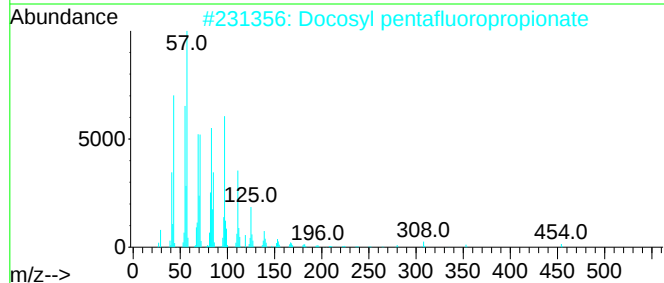
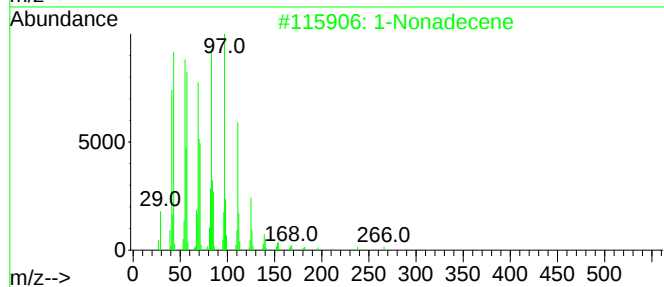
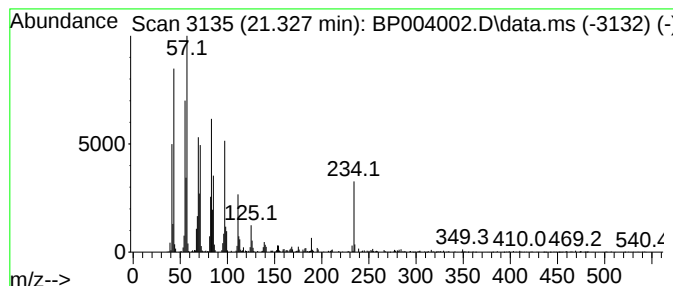
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Nonadecene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.327	16.60 ng	1130830	Chrysene-d12	21.680

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene			266	C19H38	018435-45-5	96
2	Docosyl pentafluoropropionate			472	C25H45F5O2	1000351-80-9	87
3	Nonadecyl trifluoroacetate			380	C21H39F3O2	1000351-76-3	87
4	Heneicosyl heptafluorobutyrate			508	C25H43F7O2	1000351-83-8	87
5	Heptafluorobutyric acid, n-octad...			466	C22H37F7O2	000400-57-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111820\
Data File : BP004002.D
Acq On : 18 Nov 2020 19:58
Operator : CG/JU
Sample : L4770-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
FJ-BH-01-0-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 2,2-di...	2.922	110.5	ng	5186290	1	8.287	938580	20.0
Cyclohexane	3.081	174.2	ng	8175790	1	8.287	938580	20.0
Butane, 2-metho...	3.175	10.4	ng	486563	1	8.287	938580	20.0
Heptane	3.387	2.4	ng	112246	1	8.287	938580	20.0
2-Pentanone, 4-...	5.281	4.5	ng	209746	1	8.287	938580	20.0
Ethanol, 2-(2-e...	7.852	2.7	ng	126779	1	8.287	938580	20.0
Pentadecane, 2,...	16.139	2.3	ng	172393	3	14.910	1505360	20.0
Pentadecane, 2,...	16.616	3.0	ng	256075	4	17.645	1686510	20.0
n-Dodecyl metha...	17.151	2.7	ng	228361	4	17.645	1686510	20.0
1H-Indene, 2-ph...	18.486	2.6	ng	216078	4	17.645	1686510	20.0
unknown19.227	19.227	2.6	ng	216944	4	17.645	1686510	20.0
1-Nonadecene	21.327	16.6	ng	1130830	5	21.680	1362650	20.0