

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
 Data File : BF133049.D
 Acq On : 26 Apr 2023 02:35
 Operator : CG\JU
 Sample : 02484-01 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CRCA-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_F\Methods\8270-BF042123.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF133049.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.710	272	276	282	rVB	18786	27056	1.55%	0.158%
2	4.269	365	371	376	rBV	181502	206575	11.85%	1.205%
3	4.910	476	480	485	rBV	311041	305327	17.52%	1.781%
4	5.340	550	553	561	rVB	160896	147298	8.45%	0.859%
5	6.369	725	728	739	rBV	633467	576674	33.09%	3.363%
6	6.587	759	765	775	rVB2	74200	77336	4.44%	0.451%
7	6.740	787	791	794	rBV	1717869	1284340	73.71%	7.490%
8	7.028	836	840	842	rVB3	20451	22659	1.30%	0.132%
9	7.057	842	845	848	rBV2	25640	25831	1.48%	0.151%
10	7.140	856	859	863	rBV2	34388	31871	1.83%	0.186%
11	7.298	879	886	889	rBV	582869	464617	26.66%	2.709%
12	7.345	892	894	897	rVV	121732	97208	5.58%	0.567%
13	7.392	897	902	905	rVB6	19089	29857	1.71%	0.174%
14	7.551	924	929	933	rVB3	34502	46356	2.66%	0.270%
15	7.663	939	948	953	rBV6	22244	62894	3.61%	0.367%
16	7.716	955	957	960	rVB2	26187	23569	1.35%	0.137%
17	7.751	960	963	964	rBV2	30288	26840	1.54%	0.157%
18	7.787	967	969	973	rVB2	29880	32696	1.88%	0.191%
19	7.863	978	982	985	rVB4	23924	33889	1.94%	0.198%
20	7.934	990	994	1001	rBV	163577	266446	15.29%	1.554%
21	8.022	1005	1009	1012	rBV	1945706	1742537	100.00%	10.162%
22	8.098	1020	1022	1024	rVB	39144	28916	1.66%	0.169%
23	8.345	1055	1064	1067	rBV8	55094	129253	7.42%	0.754%
24	8.381	1067	1070	1072	rVB	40709	38456	2.21%	0.224%
25	8.410	1072	1075	1078	rVB3	51915	39993	2.30%	0.233%
26	8.451	1080	1082	1084	rBV	39113	33814	1.94%	0.197%
27	8.528	1092	1095	1098	rBV2	64960	53407	3.06%	0.311%
28	8.622	1107	1111	1114	rBV	185050	157291	9.03%	0.917%
29	8.687	1119	1122	1124	rVB3	33701	29008	1.66%	0.169%
30	8.739	1129	1131	1134	rVB2	35947	31128	1.79%	0.182%
31	8.845	1143	1149	1151	rBV4	56123	71834	4.12%	0.419%
32	8.904	1157	1159	1161	rBV3	25760	24455	1.40%	0.143%
33	8.986	1171	1173	1176	rVB	45302	35298	2.03%	0.206%
34	9.051	1182	1184	1188	rVB2	48996	53899	3.09%	0.314%
35	9.098	1188	1192	1195	rBV	1172460	865985	49.70%	5.050%
36	9.186	1204	1207	1211	rBV	228280	184988	10.62%	1.079%

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37	9.251	1213	1218	1220	rVB2	34348	36469	2.09%	0.213%
38	9.357	1233	1236	1237	rBV3	25155	22334	1.28%	0.130%
39	9.416	1243	1246	1247	rBV3	26796	27083	1.55%	0.158%
40	9.439	1247	1250	1252	rVV2	47325	63449	3.64%	0.370%
41	9.463	1252	1254	1256	rVV2	48332	42608	2.45%	0.248%
42	9.510	1256	1262	1263	rVV3	76651	108959	6.25%	0.635%
43	9.528	1263	1265	1268	rVV2	59069	59111	3.39%	0.345%
44	9.563	1268	1271	1274	rVB3	39157	47823	2.74%	0.279%
45	9.716	1295	1297	1300	rBV	238937	168158	9.65%	0.981%
46	9.775	1304	1307	1310	rVB	2251303	1734094	99.52%	10.113%
47	9.886	1323	1326	1329	rVB4	18047	25503	1.46%	0.149%
48	9.957	1333	1338	1339	rBV4	24963	38648	2.22%	0.225%
49	9.981	1339	1342	1344	rVB2	38074	38835	2.23%	0.226%
50	10.039	1350	1352	1356	rVB4	44493	39731	2.28%	0.232%
51	10.216	1379	1382	1385	rVB	226807	162645	9.33%	0.948%
52	10.269	1389	1391	1396	rVB5	20265	25122	1.44%	0.147%
53	10.433	1416	1419	1425	rBV	65575	76478	4.39%	0.446%
54	10.486	1425	1428	1431	rVB	224852	194535	11.16%	1.134%
55	10.516	1431	1433	1437	rBV3	23331	28541	1.64%	0.166%
56	10.557	1437	1440	1445	rVB6	42642	46501	2.67%	0.271%
57	10.686	1459	1462	1470	rBV2	203301	259955	14.92%	1.516%
58	10.751	1470	1473	1476	rBV	51937	56044	3.22%	0.327%
59	10.792	1476	1480	1482	rVV2	36309	45682	2.62%	0.266%
60	10.822	1482	1485	1487	rVV2	27059	26568	1.52%	0.155%
61	10.880	1492	1495	1496	rBV3	28941	27708	1.59%	0.162%
62	10.969	1508	1510	1515	rBV2	44453	58550	3.36%	0.341%
63	11.010	1515	1517	1522	rVB2	55557	63177	3.63%	0.368%
64	11.133	1535	1538	1541	rBV	174231	149801	8.60%	0.874%
65	11.169	1541	1544	1549	rVB	68476	74656	4.28%	0.435%
66	11.257	1555	1559	1561	rBV	1874994	1481896	85.04%	8.642%
67	11.280	1561	1563	1565	rVV	127144	96099	5.51%	0.560%
68	11.322	1567	1570	1574	rVB2	33641	46485	2.67%	0.271%
69	11.563	1608	1611	1614	rBV	141375	115708	6.64%	0.675%
70	11.663	1625	1628	1632	rVB	122993	108866	6.25%	0.635%
71	11.786	1647	1649	1653	rVB2	48114	45317	2.60%	0.264%
72	11.863	1659	1662	1665	rBV2	36288	46480	2.67%	0.271%
73	11.969	1677	1680	1683	rBV	118426	93415	5.36%	0.545%
74	12.069	1694	1697	1703	rVB6	46263	58551	3.36%	0.341%
75	12.310	1735	1738	1740	rBV2	32413	30664	1.76%	0.179%
76	12.345	1741	1744	1746	rBV	370682	346458	19.88%	2.020%

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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	12.369	1746	1748	1750	rVB2	351485	283529	16.27%	1.653%
78	12.463	1760	1764	1768	rBV2	143041	167269	9.60%	0.975%
79	12.692	1800	1803	1807	rBV	106313	85736	4.92%	0.500%
80	12.733	1807	1810	1812	rBV2	84282	71088	4.08%	0.415%
81	12.839	1825	1828	1832	rBV	790300	680158	39.03%	3.966%
82	13.086	1867	1870	1874	rBV	82463	82093	4.71%	0.479%
83	13.427	1926	1928	1931	rBV	54850	48595	2.79%	0.283%
84	13.892	2003	2007	2010	rBV	1400935	1259804	72.30%	7.347%
85	15.315	2245	2249	2253	rVB	845853	941329	54.02%	5.489%

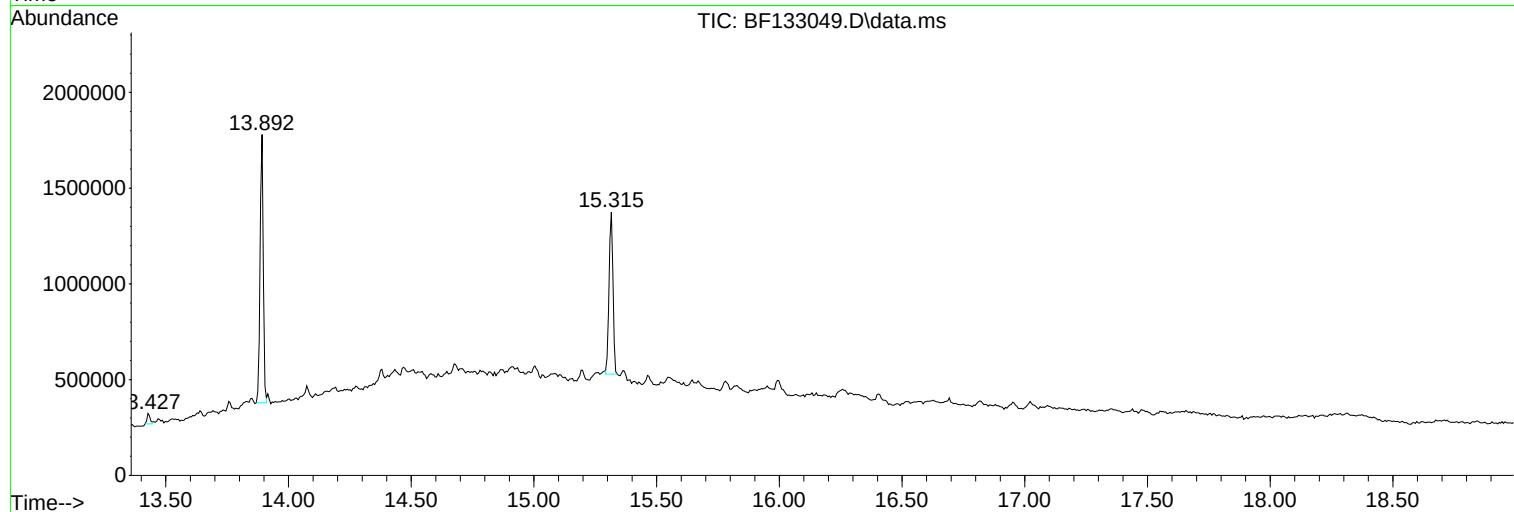
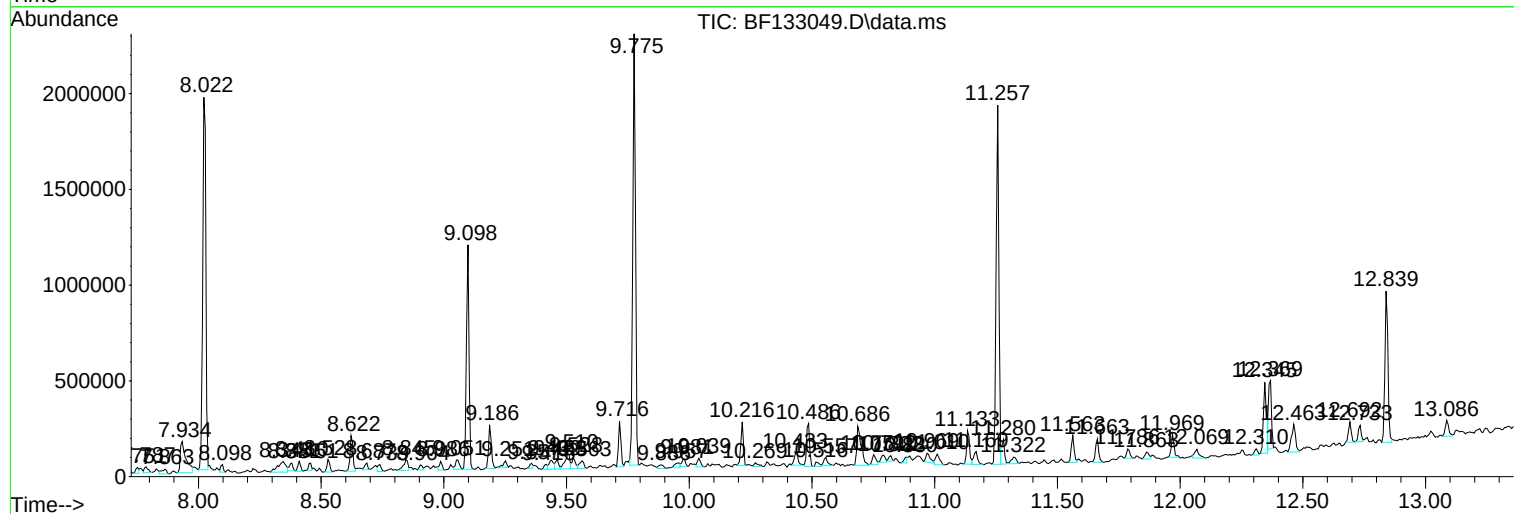
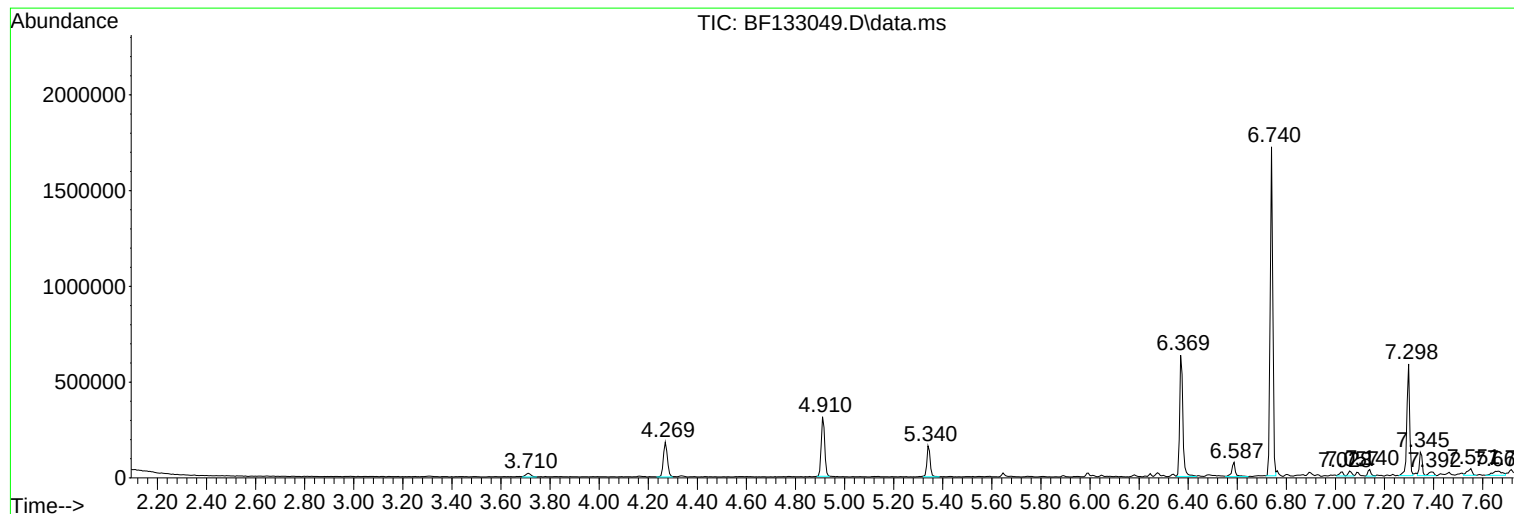
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Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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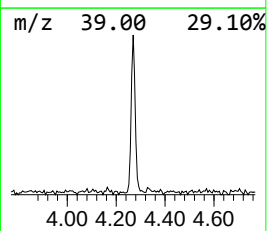
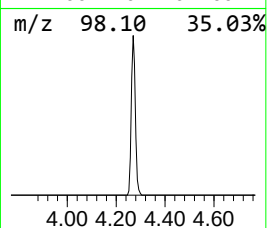
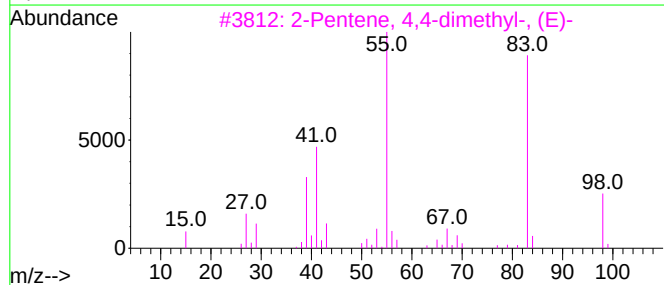
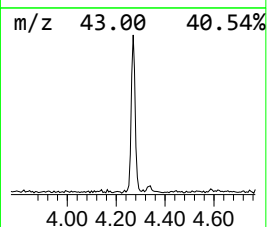
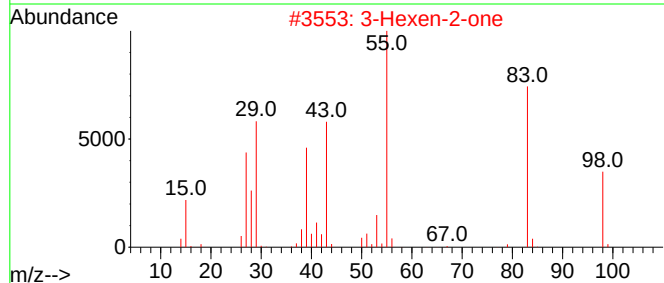
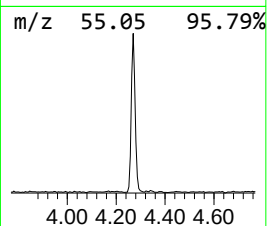
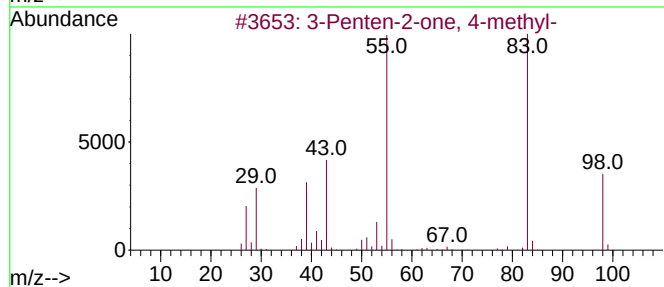
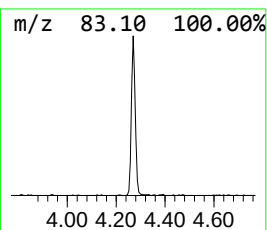
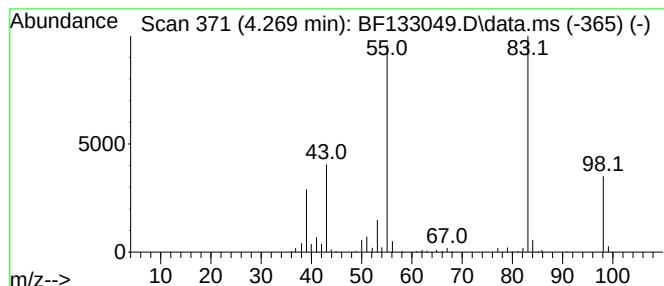
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.269	3.22 ng	206575	1,4-Dichlorobenzene-d4	6.740

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	86
4		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	86



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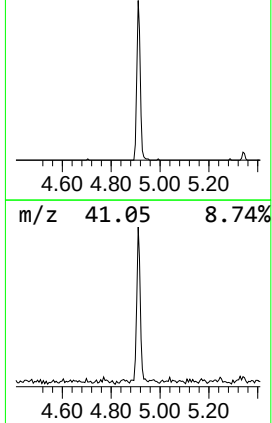
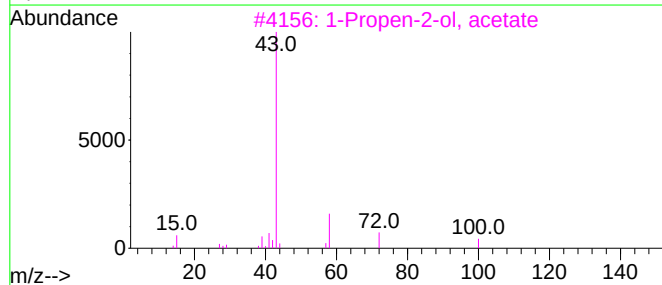
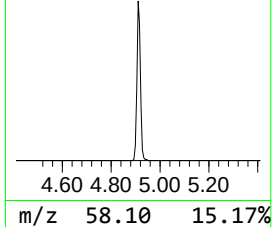
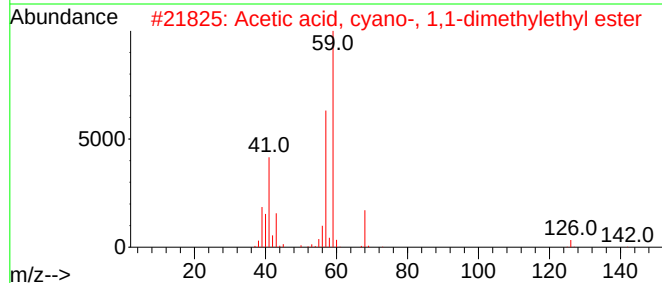
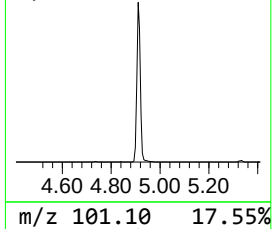
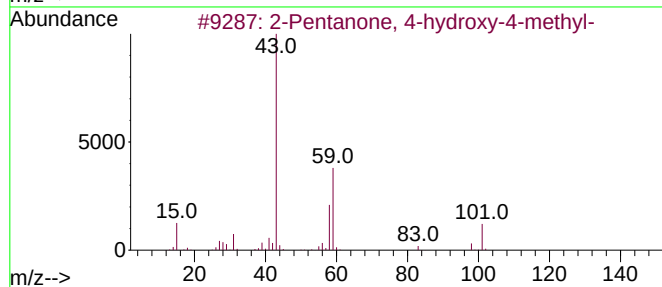
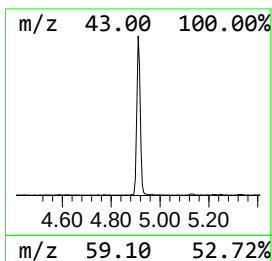
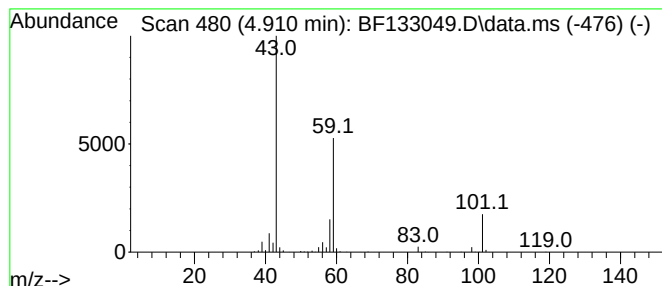
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
4.910	4.75 ng	305327	1,4-Dichlorobenzene-d4			6.740
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59	
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25	
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12	
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9	
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9	



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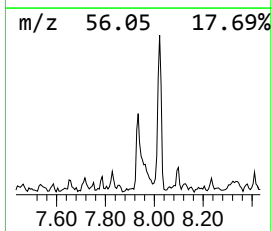
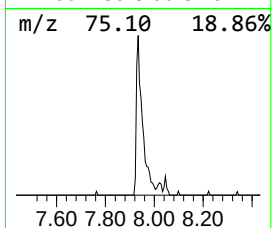
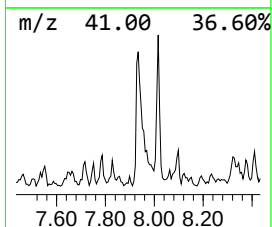
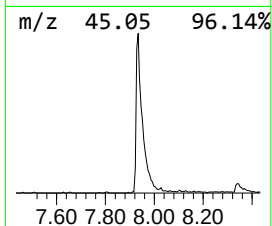
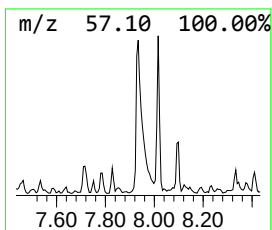
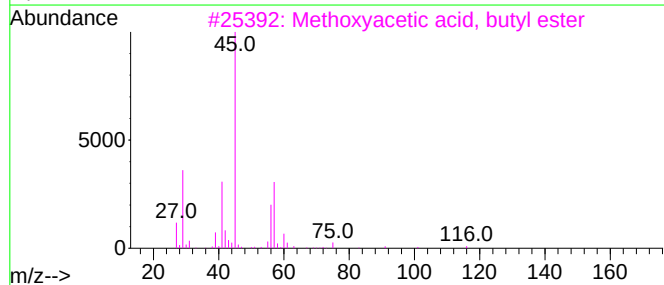
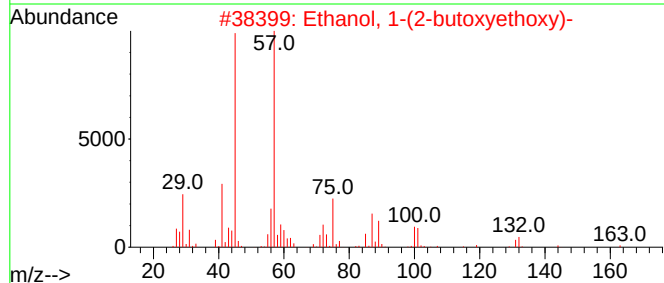
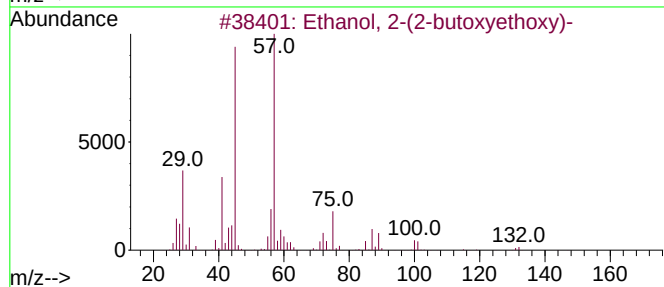
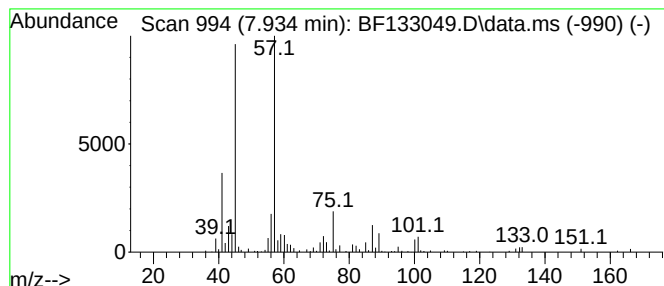
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.934	3.06 ng	266446	Naphthalene-d8	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3			Methoxyacetic acid, butyl ester	146	C7H14O3	017640-22-1	53
4			Acetaldehyde, di-sec-butyl acetal	174	C10H22O2	005314-41-0	53
5			3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	53



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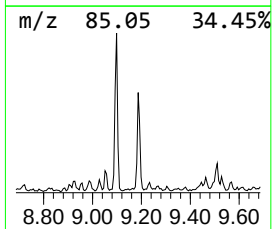
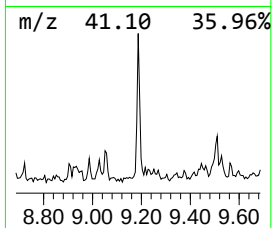
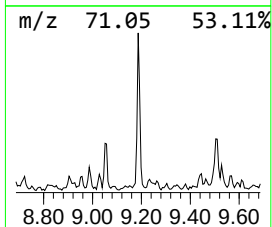
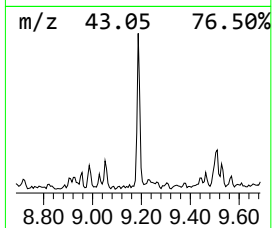
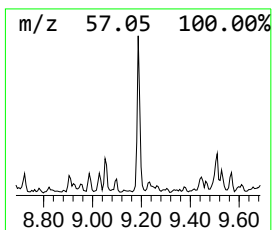
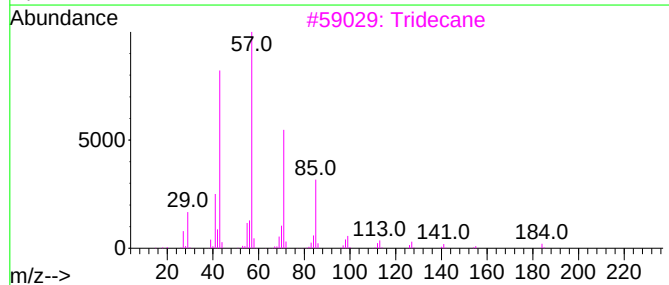
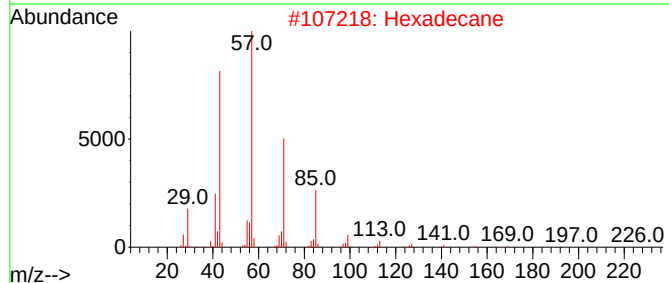
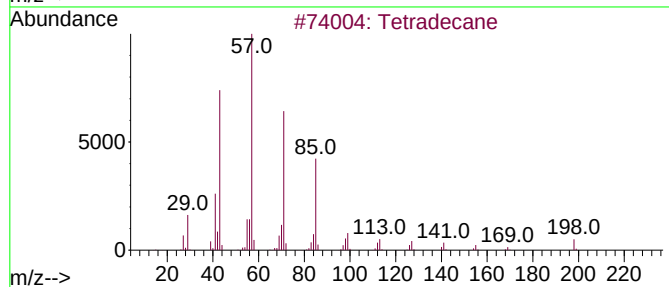
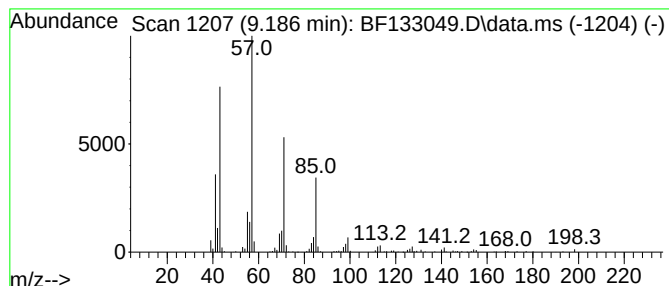
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Tetradecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.186	2.13 ng	184988	Acenaphthene-d10	9.775

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	95
2		Hexadecane	226	C16H34	000544-76-3	91
3		Tridecane	184	C13H28	000629-50-5	90
4		Pentadecane	212	C15H32	000629-62-9	90
5		Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133049.D
Acq On : 26 Apr 2023 02:35
Operator : CG\JU
Sample : 02484-01 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CRCA-1

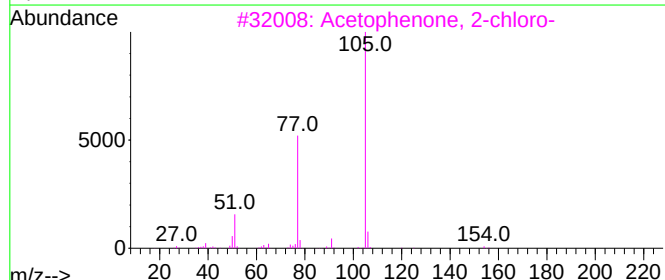
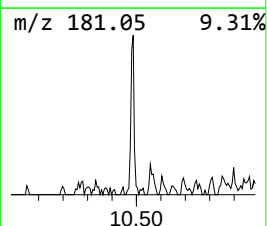
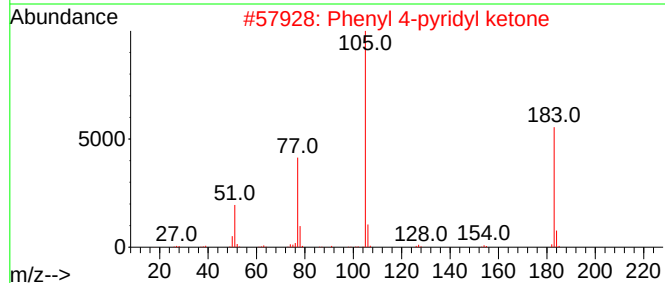
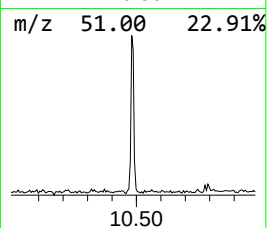
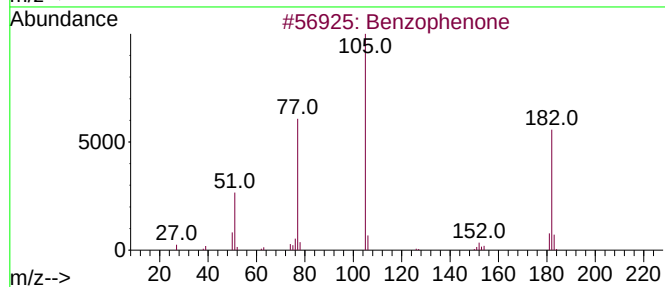
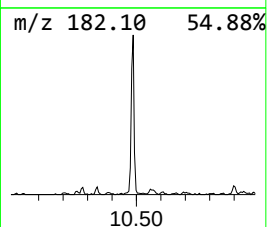
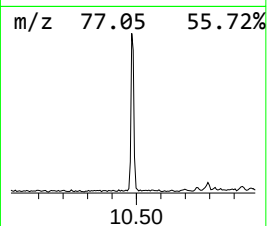
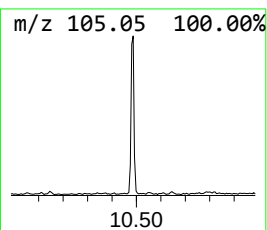
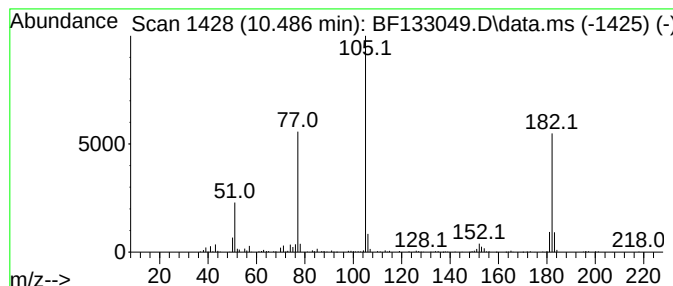
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.486	2.24 ng	194535	Acenaphthene-d10	9.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	58
3			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49
4			Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	46
5			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133049.D
Acq On : 26 Apr 2023 02:35
Operator : CG\JU
Sample : 02484-01 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CRCA-1

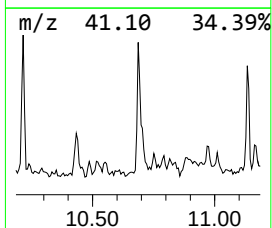
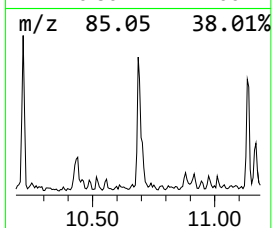
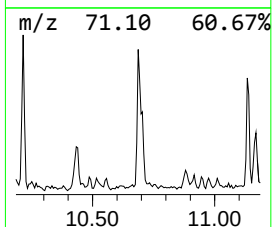
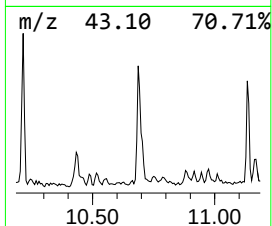
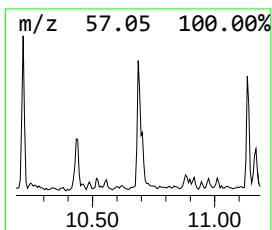
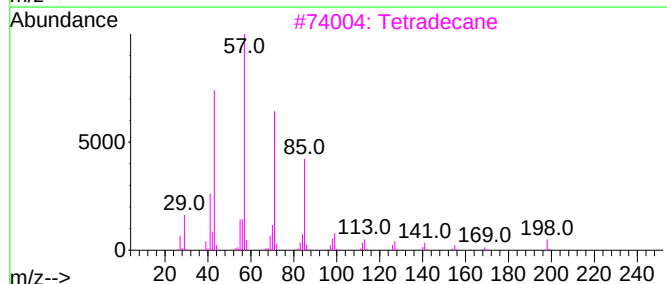
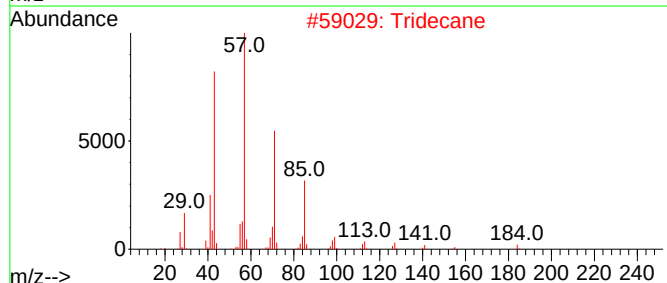
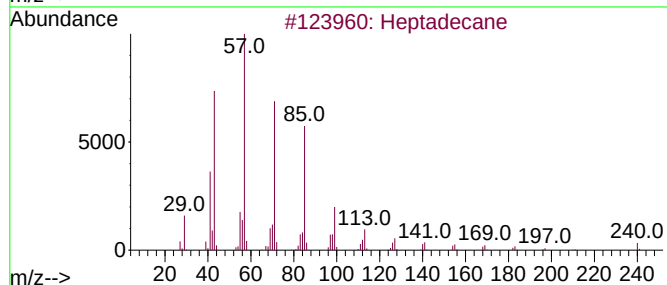
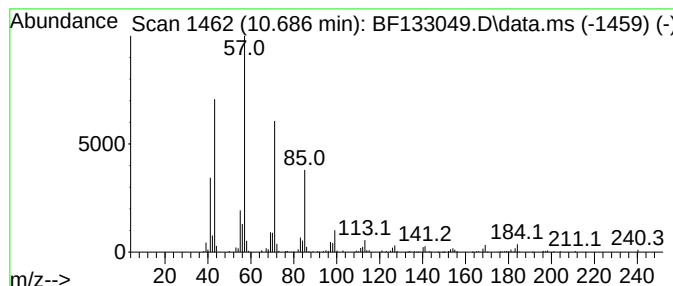
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heptadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.686	3.51 ng	259955	Phenanthrene-d10	11.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	97
2			Tridecane	184	C13H28	000629-50-5	96
3			Tetradecane	198	C14H30	000629-59-4	93
4			Eicosane	282	C20H42	000112-95-8	87
5			Tridecane, 6-methyl-	198	C14H30	013287-21-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133049.D
Acq On : 26 Apr 2023 02:35
Operator : CG\JU
Sample : 02484-01 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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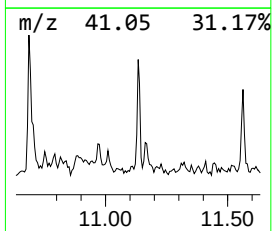
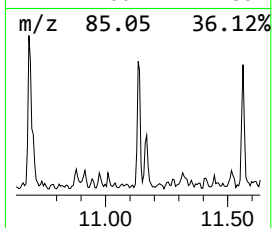
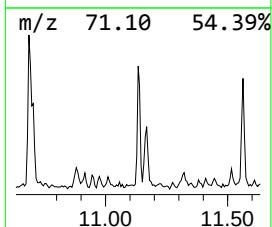
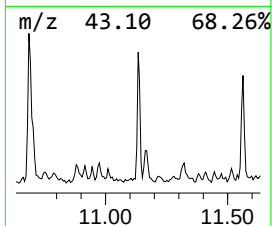
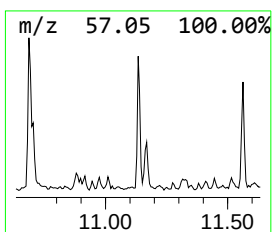
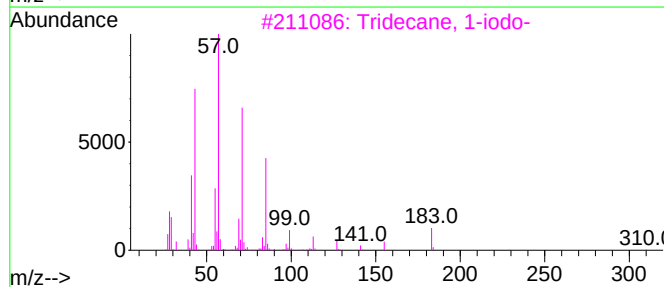
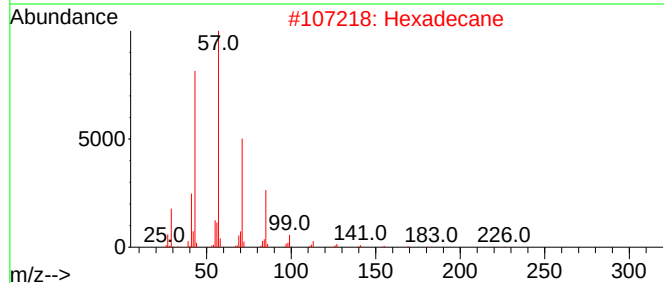
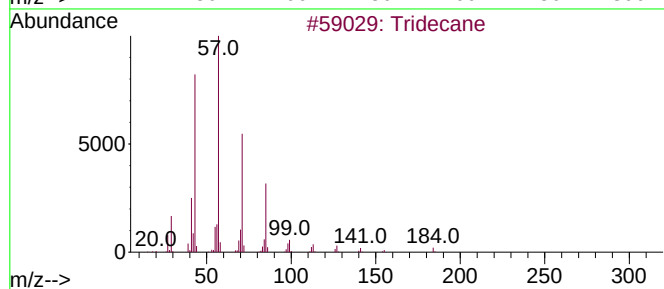
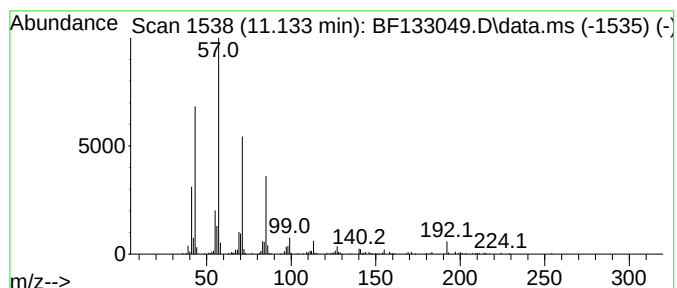
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tridecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.133	2.02 ng	149801	Phenanthrene-d10	11.257

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	86
2		Hexadecane	226	C16H34	000544-76-3	86
3		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86
4		Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	86
5		Tetradecane	198	C14H30	000629-59-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133049.D
Acq On : 26 Apr 2023 02:35
Operator : CG\JU
Sample : 02484-01 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

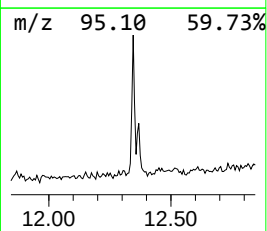
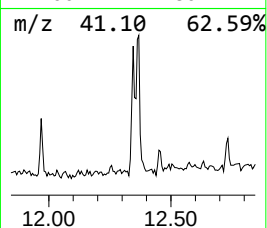
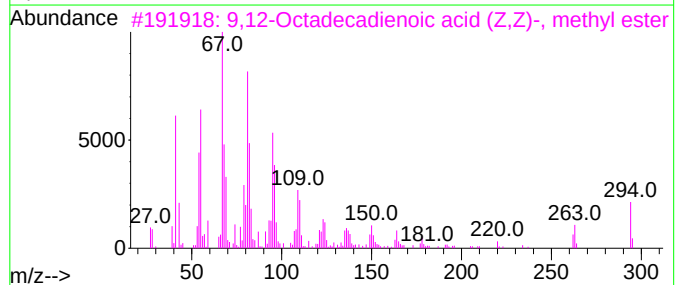
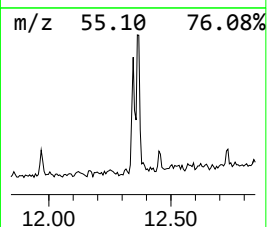
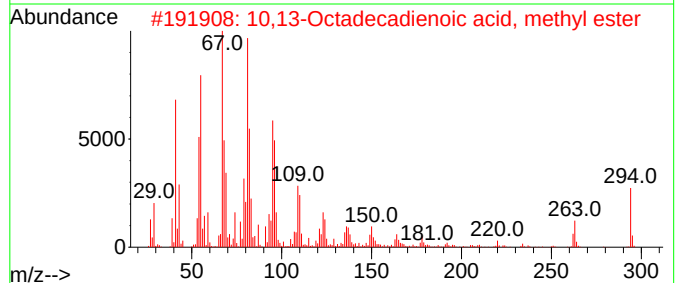
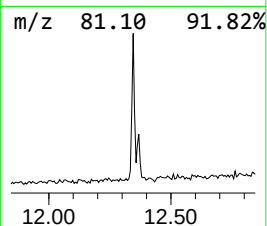
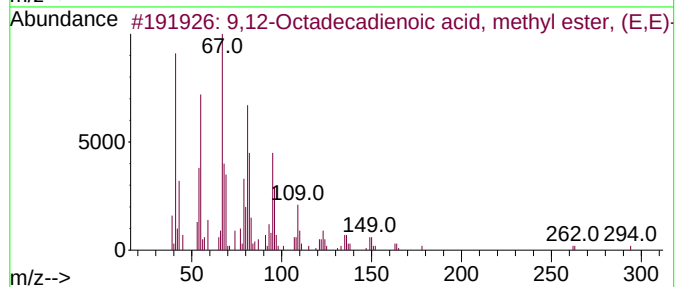
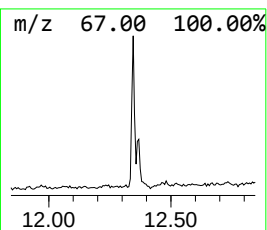
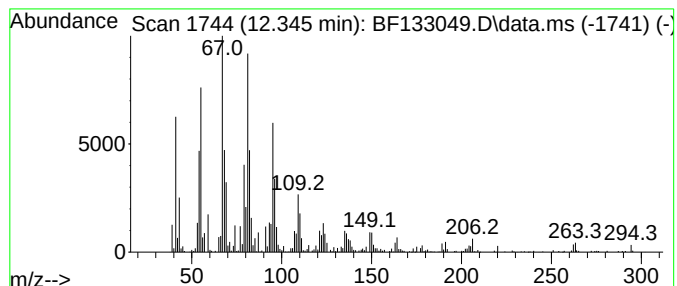
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9,12-Octadecadienoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.345	4.68 ng	346458	Phenanthrene-d10	11.257		
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99	
2	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99	
3	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99	
4	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99	
5	Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	99	



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133049.D
Acq On : 26 Apr 2023 02:35
Operator : CG\JU
Sample : 02484-01 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CRCA-1

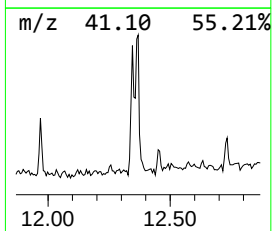
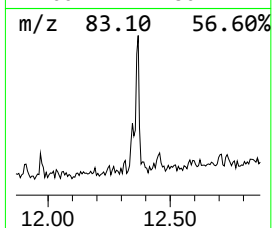
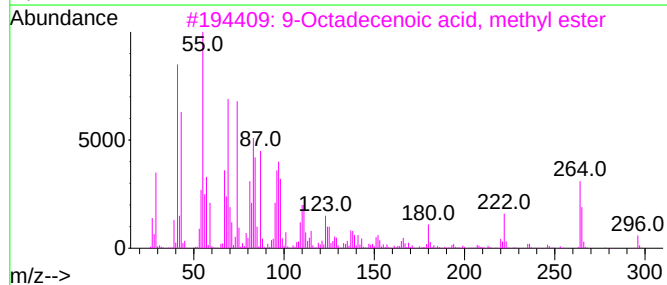
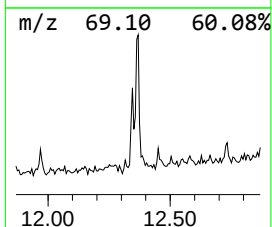
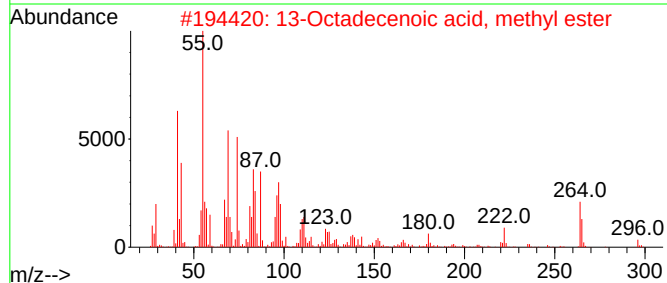
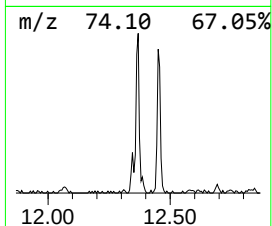
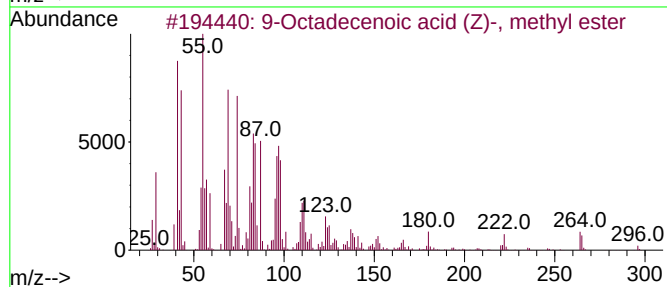
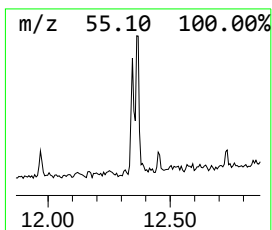
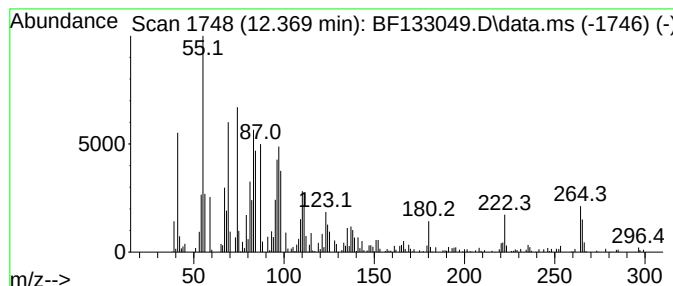
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9-Octadecenoic acid (Z)-, m... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.369	3.83 ng	283529	Phenanthrene-d10	11.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2			13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	99
3			9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	99
4			16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99
5			15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
Data File : BF133049.D
Acq On : 26 Apr 2023 02:35
Operator : CG\JU
Sample : 02484-01 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.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
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2-Pentanone, 4-...	4.910	4.8	ng	305327	1	6.740	1284340	20.0
Ethanol, 2-(2-b...	7.934	3.1	ng	266446	2	8.022	1742540	20.0
Tetradecane	9.186	2.1	ng	184988	3	9.775	1734090	20.0
Benzophenone	10.486	2.2	ng	194535	3	9.775	1734090	20.0
Heptadecane	10.686	3.5	ng	259955	4	11.257	1481900	20.0
Tridecane	11.133	2.0	ng	149801	4	11.257	1481900	20.0
9,12-Octadecadi...	12.345	4.7	ng	346458	4	11.257	1481900	20.0
9-Octadecenoic ...	12.369	3.8	ng	283529	4	11.257	1481900	20.0