

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF030220\
 Data File : BF119179.D
 Acq On : 2 Mar 2020 20:21
 Operator : CG/JU
 Sample : L1735-03 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-SF-01-009-B

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

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-BF022020.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.275	369	372	390	rBV	52173	81707	7.80%	0.740%
2	4.916	478	481	491	rBV	97197	112090	10.70%	1.015%
3	5.375	556	559	567	rBV	40721	64139	6.12%	0.581%
4	6.387	728	731	749	rBV	159346	223188	21.30%	2.020%
5	6.734	787	790	806	rBV	807555	707568	67.54%	6.405%
6	6.893	813	817	825	rBV	330709	287901	27.48%	2.606%
7	7.298	883	886	897	rBV	171211	177510	16.94%	1.607%
8	8.016	1003	1008	1018	rBV	1053657	942808	89.99%	8.535%
9	8.834	1143	1147	1152	rVB2	16720	19776	1.89%	0.179%
10	9.092	1187	1191	1196	rBV	446261	415981	39.71%	3.766%
11	9.175	1202	1205	1207	rBV	33232	30646	2.93%	0.277%
12	9.210	1207	1211	1215	rVB2	75937	73882	7.05%	0.669%
13	9.363	1232	1237	1240	rBV2	18463	23956	2.29%	0.217%
14	9.428	1245	1248	1251	rVV	30028	32438	3.10%	0.294%
15	9.486	1255	1258	1262	rVV2	30680	40583	3.87%	0.367%
16	9.551	1266	1269	1272	rBV3	17171	18922	1.81%	0.171%
17	9.633	1279	1283	1287	rBV	26727	27139	2.59%	0.246%
18	9.704	1293	1295	1299	rVB	26341	19975	1.91%	0.181%
19	9.769	1302	1306	1310	rVV	1183437	990704	94.57%	8.969%
20	9.798	1310	1311	1316	rVB	43368	42117	4.02%	0.381%
21	9.875	1322	1324	1328	rVB3	14058	15117	1.44%	0.137%
22	9.975	1337	1341	1345	rBV	38718	37103	3.54%	0.336%
23	10.016	1345	1348	1354	rVB3	17325	23211	2.22%	0.210%
24	10.086	1357	1360	1363	rBV2	15980	18895	1.80%	0.171%
25	10.175	1372	1375	1377	rBV2	14382	17522	1.67%	0.159%
26	10.204	1377	1380	1382	rVB2	43999	37278	3.56%	0.337%
27	10.316	1395	1399	1404	rBV2	54185	62078	5.93%	0.562%
28	10.428	1415	1418	1423	rVB2	20188	22918	2.19%	0.207%
29	10.475	1423	1426	1429	rVB2	21381	21223	2.03%	0.192%
30	10.675	1457	1460	1467	rVB3	57787	84668	8.08%	0.766%
31	10.863	1488	1492	1494	rBV4	24615	26608	2.54%	0.241%
32	10.892	1494	1497	1500	rVB3	15415	18460	1.76%	0.167%
33	11.063	1523	1526	1530	rVV	25054	28408	2.71%	0.257%
34	11.122	1530	1536	1539	rVV3	58222	69516	6.64%	0.629%

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35	11.151	1539	1541	1546	rVB	63219	58540	5.59%	0.530%
36	11.251	1555	1558	1561	rBV	1032804	969408	92.53%	8.776%
37	11.275	1561	1562	1566	rVV	420042	333972	31.88%	3.023%
38	11.328	1569	1571	1575	rVV	67851	66153	6.31%	0.599%
39	11.369	1575	1578	1581	rVV2	23440	22043	2.10%	0.200%
40	11.492	1595	1599	1603	rBV	48580	66250	6.32%	0.600%
41	11.545	1606	1608	1612	rVV	59118	52783	5.04%	0.478%
42	11.586	1612	1615	1620	rVB2	28257	38916	3.71%	0.352%
43	11.710	1634	1636	1640	rBV4	14233	22225	2.12%	0.201%
44	11.757	1641	1644	1646	rVV	73679	75806	7.24%	0.686%
45	11.780	1646	1648	1652	rVV	94008	104105	9.94%	0.942%
46	11.833	1654	1657	1659	rVV	21936	26581	2.54%	0.241%
47	11.869	1659	1663	1665	rVV2	88414	106622	10.18%	0.965%
48	11.886	1665	1666	1670	rVB	50759	43071	4.11%	0.390%
49	11.951	1675	1677	1681	rVV	54071	54899	5.24%	0.497%
50	11.992	1681	1684	1686	rVB	20092	18469	1.76%	0.167%
51	12.051	1691	1694	1696	rBV	39259	41284	3.94%	0.374%
52	12.080	1697	1699	1702	rVB	42978	41448	3.96%	0.375%
53	12.180	1713	1716	1718	rBV	29991	38100	3.64%	0.345%
54	12.233	1722	1725	1727	rBV3	38342	37759	3.60%	0.342%
55	12.304	1734	1737	1739	rBV	78444	82986	7.92%	0.751%
56	12.345	1739	1744	1749	rVB3	114859	171250	16.35%	1.550%
57	12.398	1749	1753	1755	rBV3	39486	50565	4.83%	0.458%
58	12.469	1761	1765	1768	rBV	437142	429674	41.01%	3.890%
59	12.622	1788	1791	1793	rBV3	36554	39910	3.81%	0.361%
60	12.692	1798	1803	1810	rBV	429414	532192	50.80%	4.818%
61	12.751	1811	1813	1817	rVB2	35248	38181	3.64%	0.346%
62	12.833	1823	1827	1830	rBV	445188	413453	39.47%	3.743%
63	13.069	1864	1867	1869	rBV	81984	79489	7.59%	0.720%
64	13.410	1922	1925	1927	rBV	79011	70666	6.75%	0.640%
65	13.892	2003	2007	2010	rBV2	981455	1047639	100.00%	9.484%
66	14.051	2032	2034	2038	rVB4	113767	96568	9.22%	0.874%
67	14.651	2134	2136	2139	rVB	165676	146932	14.03%	1.330%
68	15.321	2247	2250	2257	rBV2	675302	812502	77.56%	7.355%

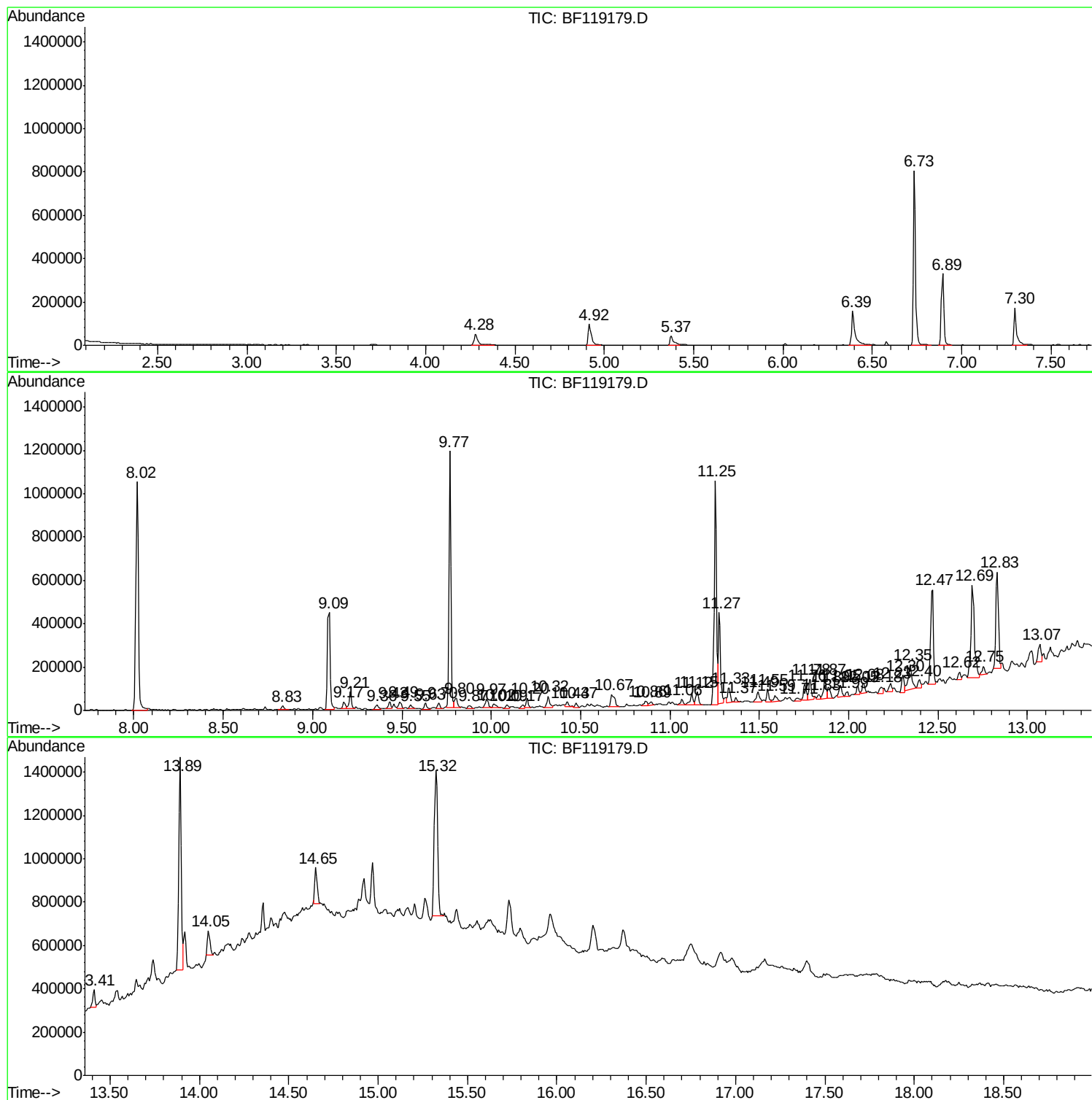
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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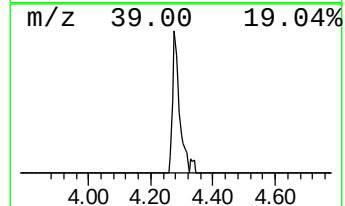
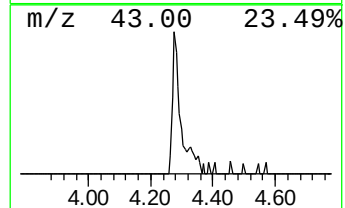
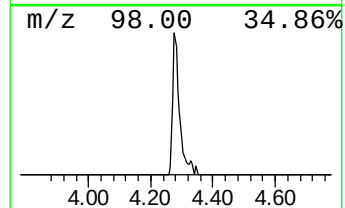
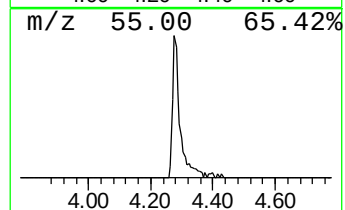
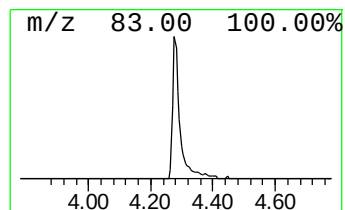
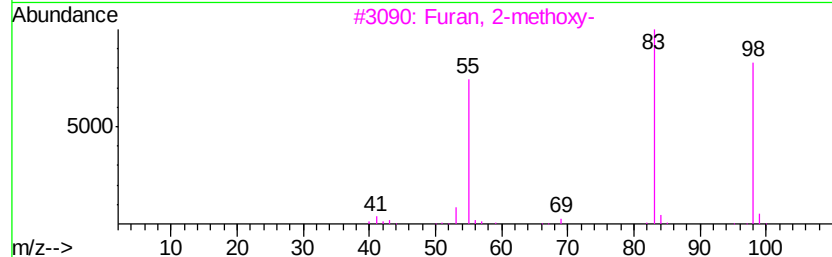
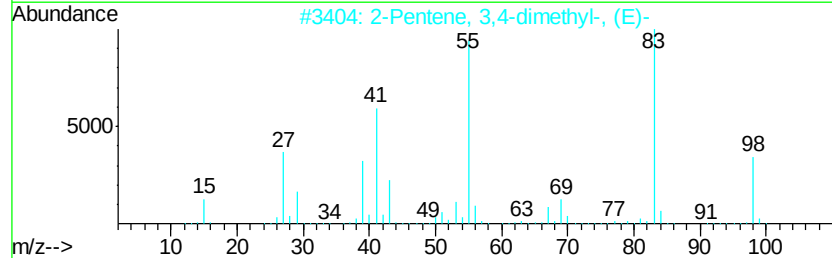
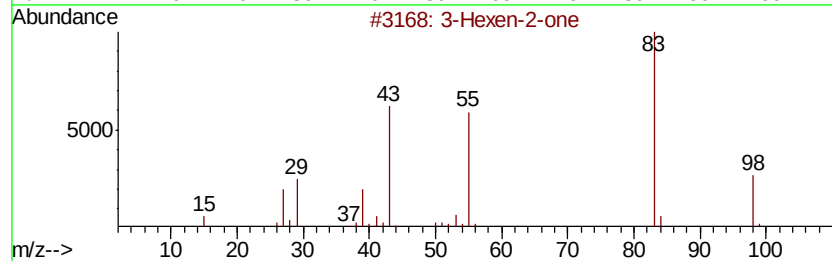
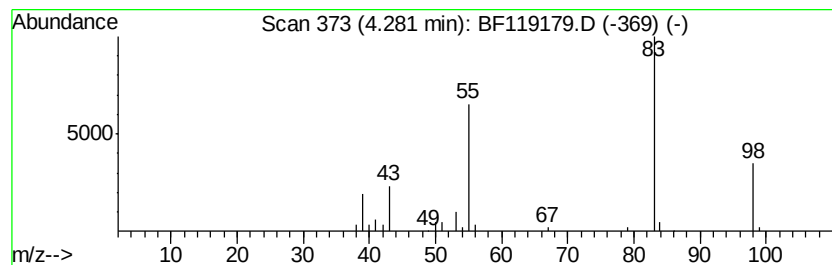
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 3-Hexen-2-one Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.28	2.31 ng	81707	1,4-Dichlorobenzene-d4	6.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Hexen-2-one	98	C6H10O	000763-93-9	90
2		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
3		Furan, 2-methoxy-	98	C5H6O2	025414-22-6	86
4		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	83



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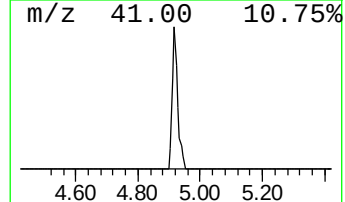
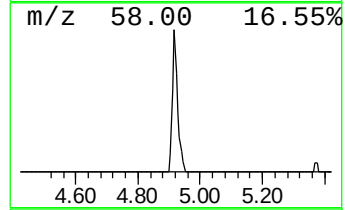
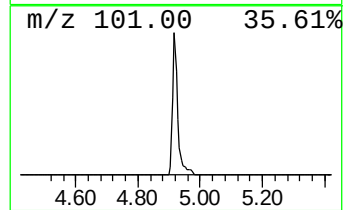
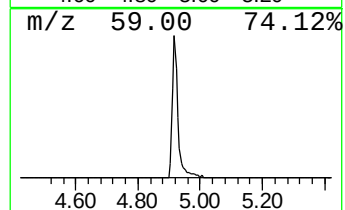
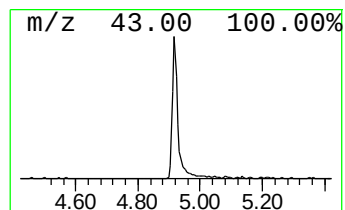
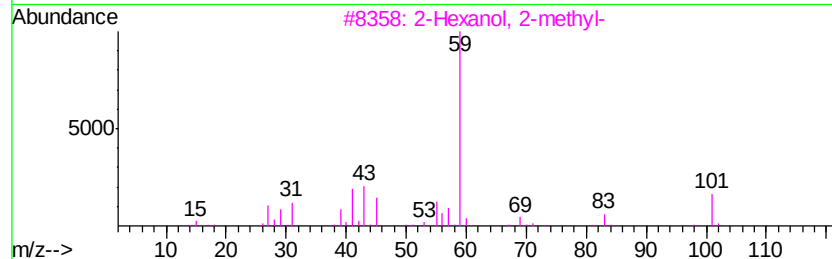
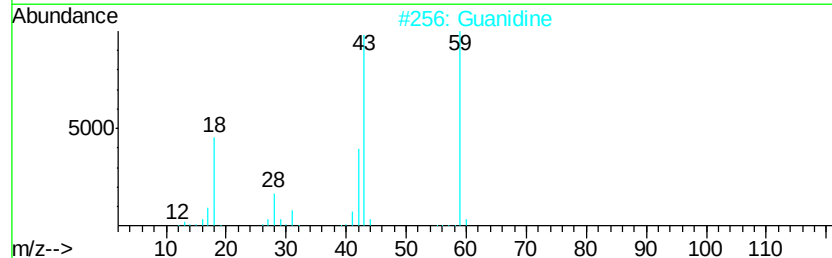
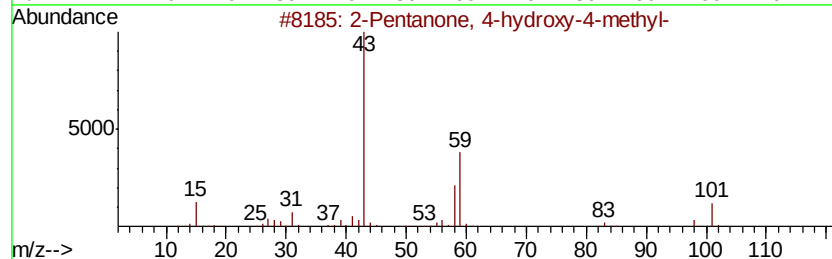
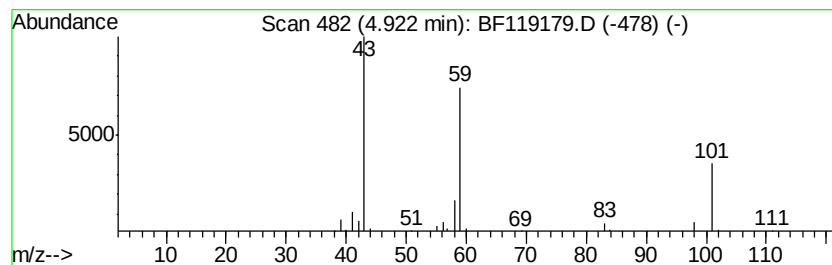
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	3.17 ng	112090	1,4-Dichlorobenzene-d4	6.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Guanidine	59	CH5N3	000113-00-8	43
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	40
4		5-Ethyl-3-methylhept-1-en-4-ol	156	C10H20O	286424-80-4	28
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25



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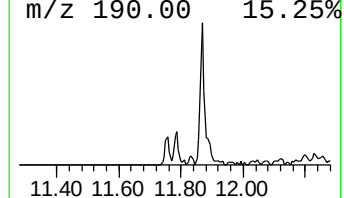
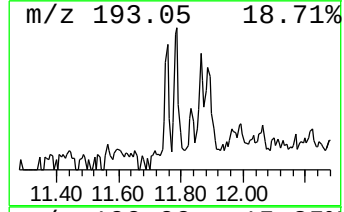
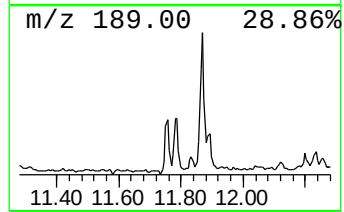
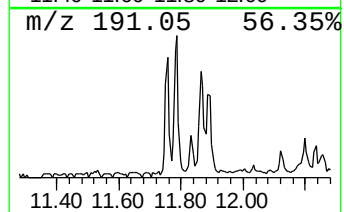
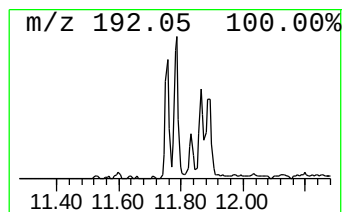
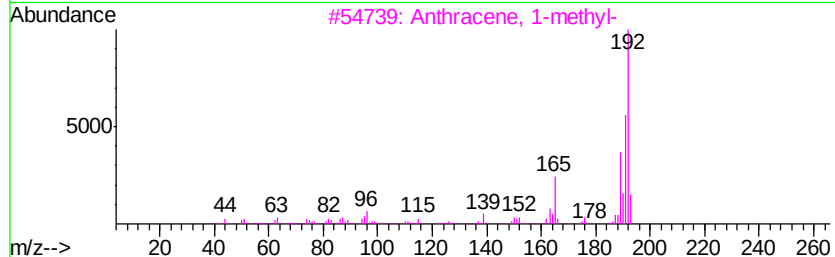
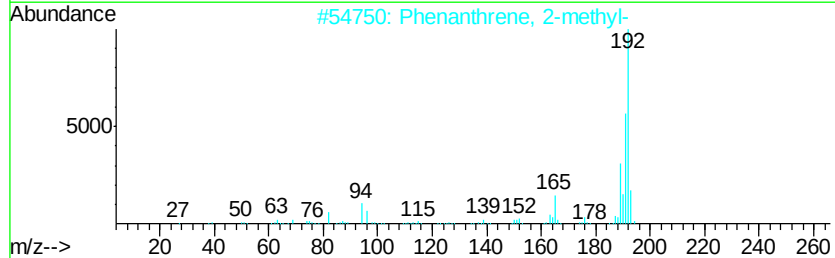
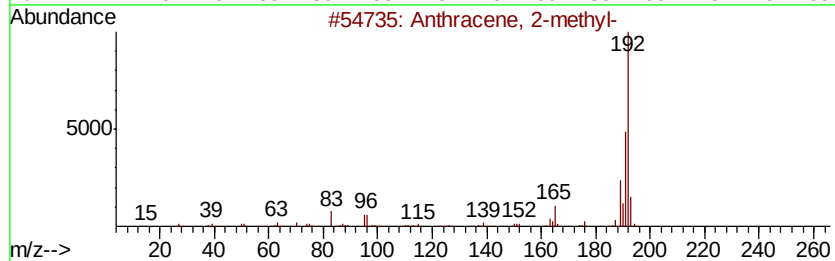
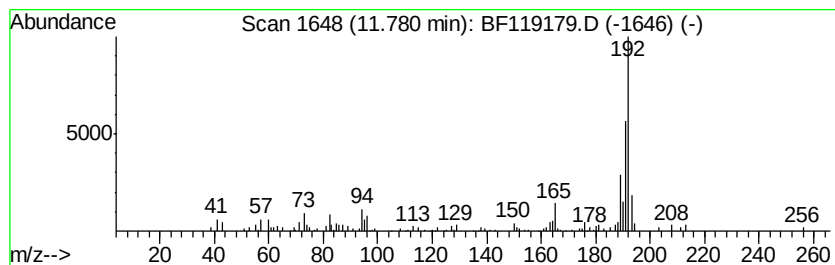
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Anthracene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.78	2.15 ng	104105	Phenanthrene-d10		11.25
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	86
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	81
5	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	81



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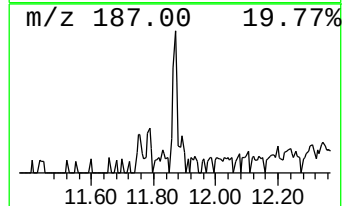
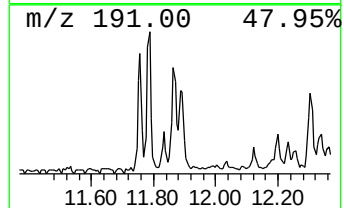
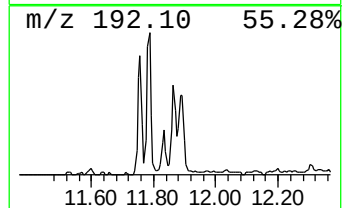
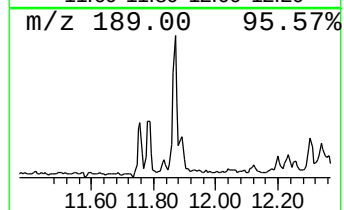
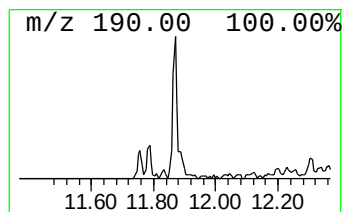
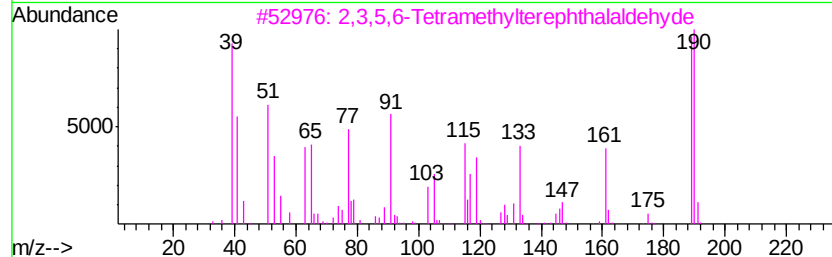
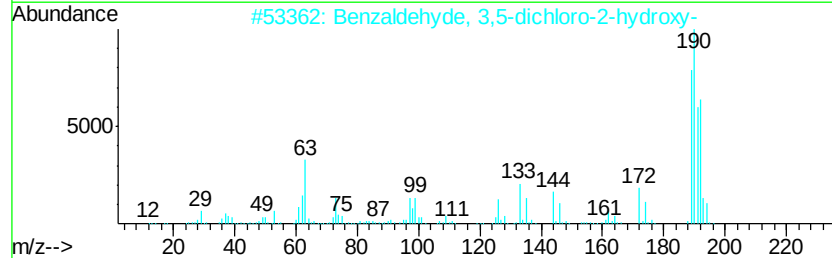
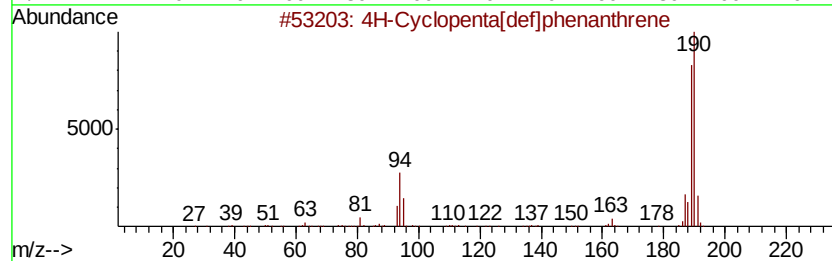
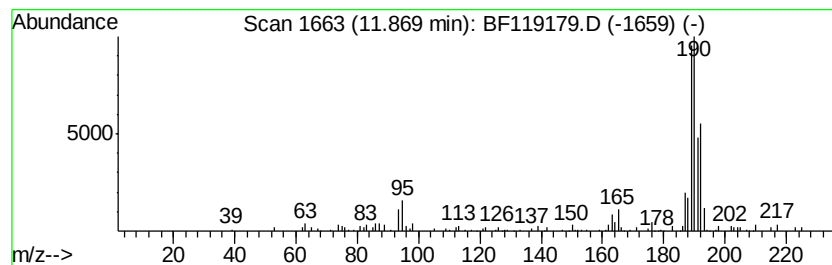
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Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.87	2.20 ng	106622	Phenanthrene-d10		11.25	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	50
3		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	43
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	30
5		Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	25



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Sample : L1735-03 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-SF-01-009-B

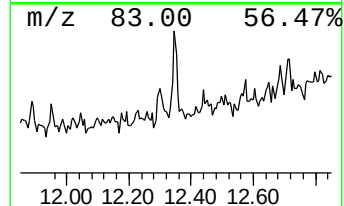
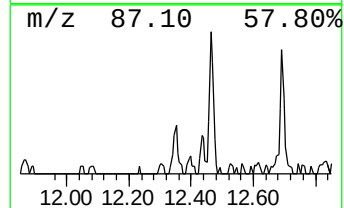
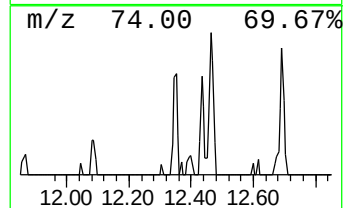
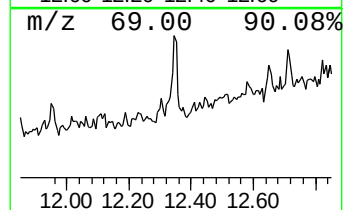
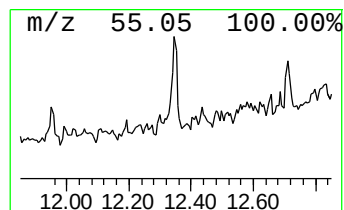
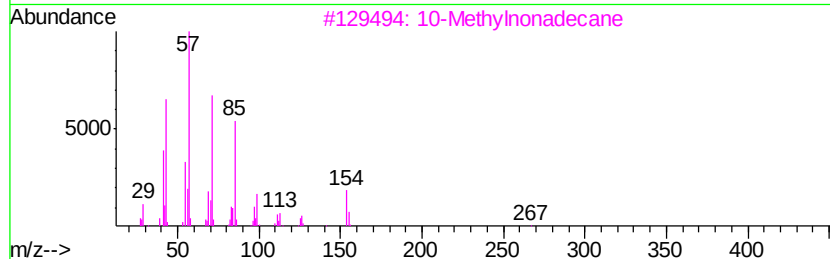
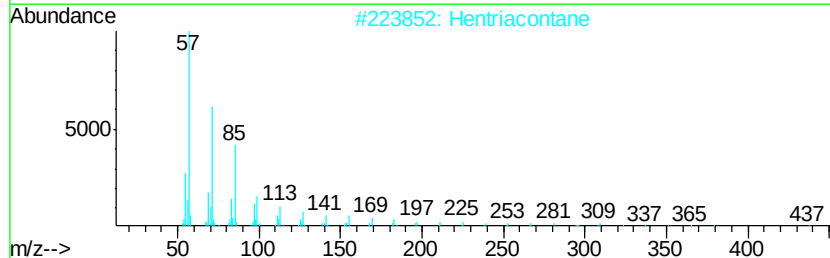
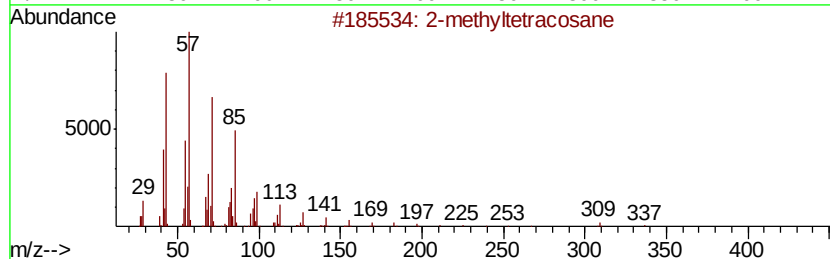
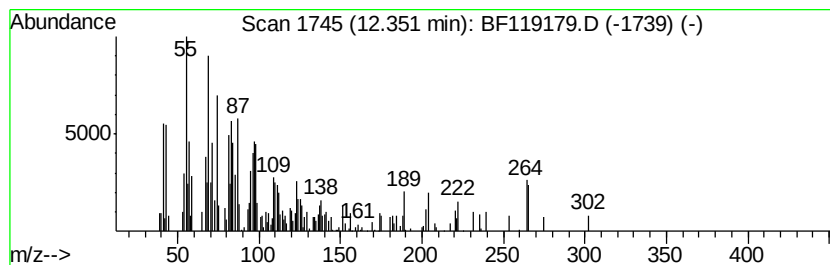
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown12.35 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	3.53 ng	171250	Phenanthrene-d10	11.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyltetracosane	352	C25H52	1000376-72-6	46
2		Hentriacontane	437	C31H64	000630-04-6	46
3		10-Methylnonadecane	282	C20H42	056862-62-5	43
4		Pentadecane	212	C15H32	000629-62-9	43
5		2-Bromotetradecane	276	C14H29Br	074036-95-6	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030220\
Data File : BF119179.D
Acq On : 2 Mar 2020 20:21
Operator : CG/JU
Sample : L1735-03 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

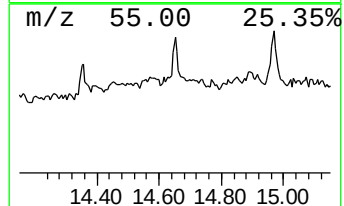
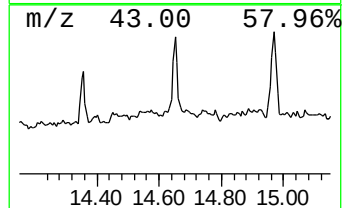
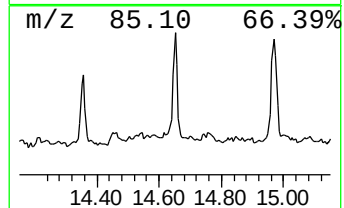
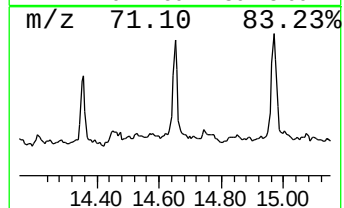
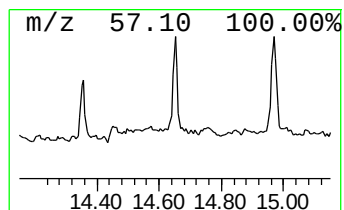
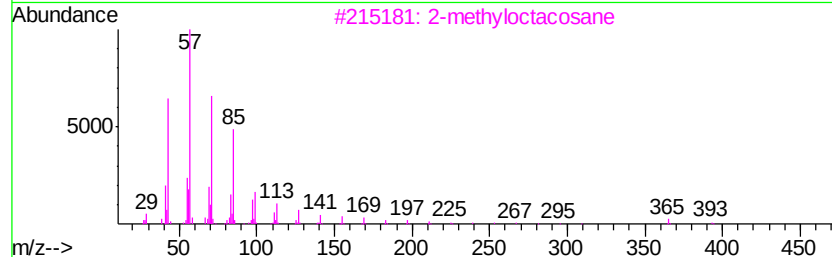
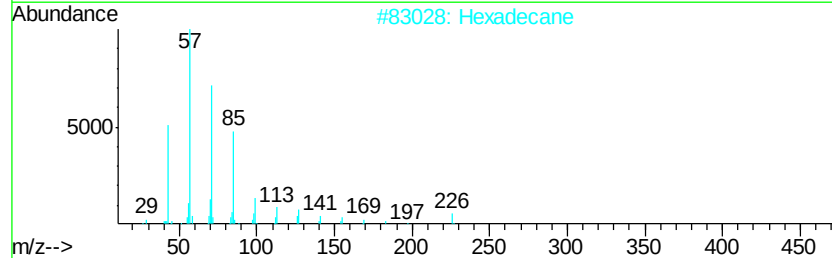
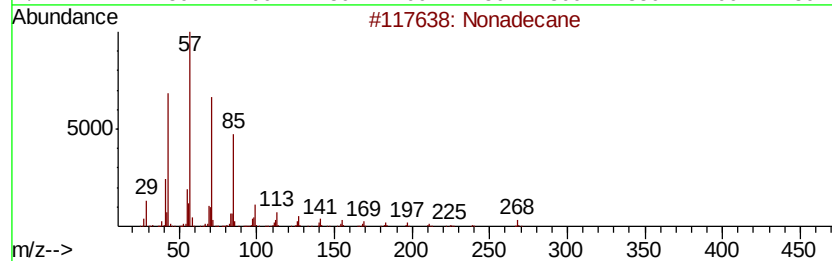
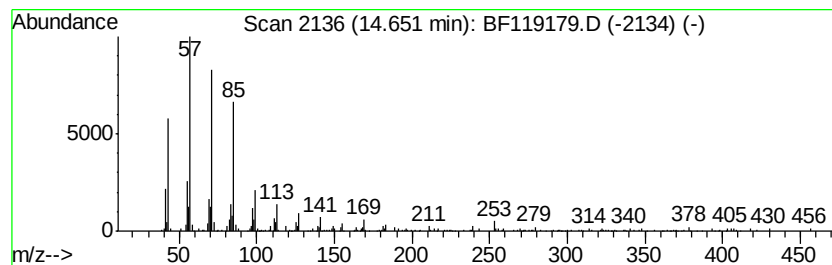
Instrument :
BNA_F
ClientSampleId :
FK-SF-01-009-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Nonadecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.65	3.62 ng	146932	Perylene-d12	15.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Nonadecane	268 C19H40	000629-92-5	93
2	Hexadecane	226 C16H34	000544-76-3	90
3	2-methyloctacosane	408 C29H60	1000376-72-8	87
4	Hexacosane	366 C26H54	000630-01-3	87
5	Heneicosane	296 C21H44	000629-94-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF030220\
Data File : BF119179.D
Acq On : 2 Mar 2020 20:21
Operator : CG/JU
Sample : L1735-03 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-SF-01-009-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.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
3-Hexen-2-one	4.28	2.3	ng	81707	1	6.73	707568	20.0
2-Pentanone, 4-hy...	4.92	3.2	ng	112090	1	6.73	707568	20.0
Anthracene, 2-met...	11.78	2.1	ng	104105	4	11.25	969408	20.0
4H-Cyclopenta[def...	11.87	2.2	ng	106622	4	11.25	969408	20.0
unknown12.35	12.35	3.5	ng	171250	4	11.25	969408	20.0
Nonadecane	14.65	3.6	ng	146932	6	15.32	812502	20.0