

Data Path : Z:\HPCHEM1\BNA N\DATA\BN042318\
 Data File : BN000824.D
 Acq On : 23 Apr 2018 12:35
 Operator : JU/SJ
 Sample : J2467-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3XT8

Integration Parameters: events.e
 Integrator: ChemStation
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 1 % 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:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041918-PAH.M
 Title : ??? S

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.799	642	648	657	rVB	127843	200462	8.24%	0.787%
2	6.969	670	677	689	rBV	470433	708328	29.11%	2.781%
3	7.157	703	709	717	rBV	463769	698035	28.68%	2.741%
4	7.622	781	788	795	rVB	464332	701647	28.83%	2.755%
5	8.322	900	907	917	rVB	358869	554280	22.78%	2.177%
6	8.763	975	982	993	rVB	513512	785095	32.26%	3.083%
7	9.210	1053	1058	1064	rVB	29911	41886	1.72%	0.164%
8	9.481	1097	1104	1112	rVB	393897	622842	25.59%	2.446%
9	10.010	1187	1194	1206	rBV	613270	958375	39.38%	3.763%
10	10.387	1251	1258	1265	rBV	632890	1014626	41.69%	3.984%
11	10.528	1275	1282	1289	rBV	18386	29311	1.20%	0.115%
12	13.669	1810	1816	1821	rBV	1055991	1361148	55.93%	5.345%
13	13.716	1821	1824	1829	rVB	99070	123259	5.07%	0.484%
14	13.939	1856	1862	1875	rBV	1212682	1755303	72.13%	6.893%
15	14.180	1898	1903	1909	rVB	31604	39840	1.64%	0.156%
16	14.251	1909	1915	1926	rVB2	819704	1236643	50.82%	4.856%
17	14.451	1943	1949	1958	rBV	57220	88589	3.64%	0.348%
18	15.133	2060	2065	2070	rVB	40559	47970	1.97%	0.188%
19	15.251	2078	2085	2094	rVB	1472416	2096718	86.16%	8.233%
20	15.363	2099	2104	2112	rBV	404804	538833	22.14%	2.116%
21	15.998	2207	2212	2221	rBV	31040	52226	2.15%	0.205%
22	16.998	2376	2382	2389	rBV2	997561	1370881	56.33%	5.383%
23	17.098	2392	2399	2412	rVB2	1500819	2128670	87.47%	8.359%
24	17.916	2534	2538	2546	rBV	47858	63580	2.61%	0.250%
25	19.204	2753	2757	2767	rBV	26343	33025	1.36%	0.130%
26	19.398	2784	2790	2799	rBV	1787485	2392408	98.31%	9.394%
27	20.945	3050	3053	3057	rBV3	20887	30246	1.24%	0.119%
28	21.204	3091	3097	3104	rBV	1247041	1520507	62.48%	5.971%
29	22.274	3276	3279	3285	rVB4	21003	26341	1.08%	0.103%
30	22.768	3360	3363	3369	rVB	36737	48089	1.98%	0.189%
31	23.256	3438	3446	3453	rBV2	1342725	2433474	100.00%	9.556%
32	23.315	3454	3456	3461	rVV	48283	71690	2.95%	0.282%
33	23.392	3462	3469	3479	rVB2	844277	1545097	63.49%	6.067%
34	23.945	3560	3563	3569	rVB3	41575	66492	2.73%	0.261%

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Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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35	24.668	3683	3686	3693	rvB3	42205	80390	3.30%	0.316%
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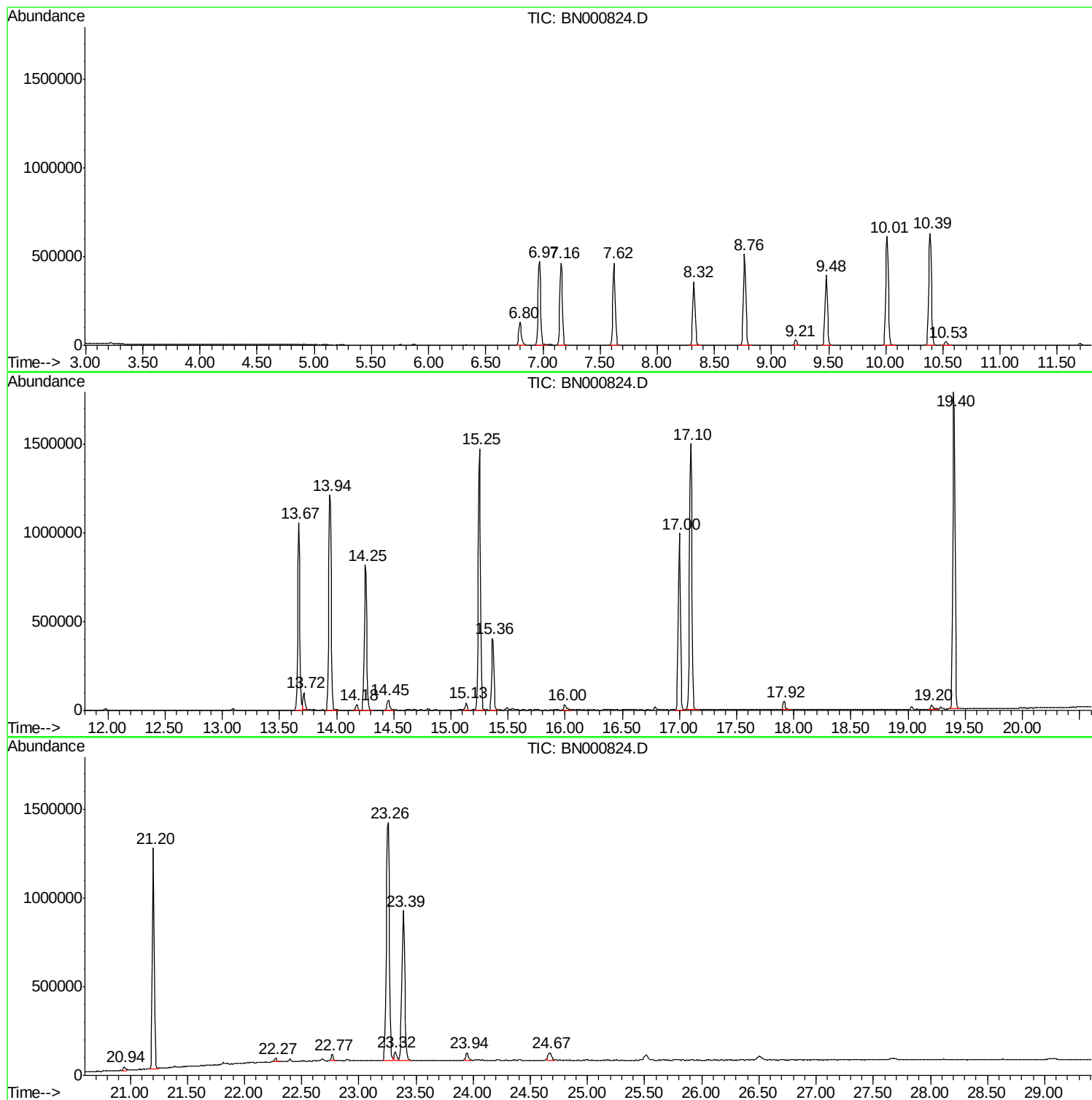
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Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041918-PAH.M
Quant Title : SVOA CALIBRATION

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



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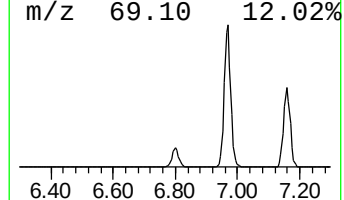
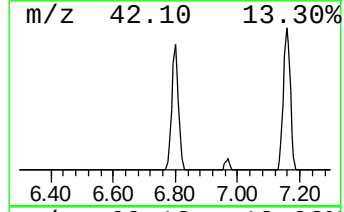
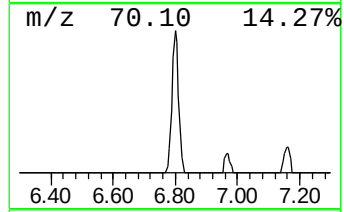
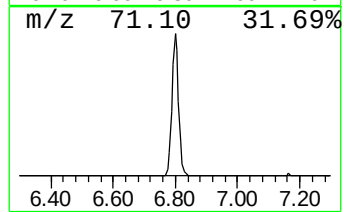
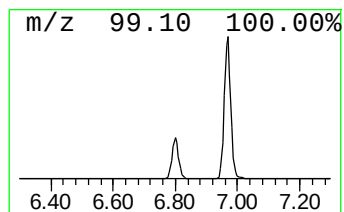
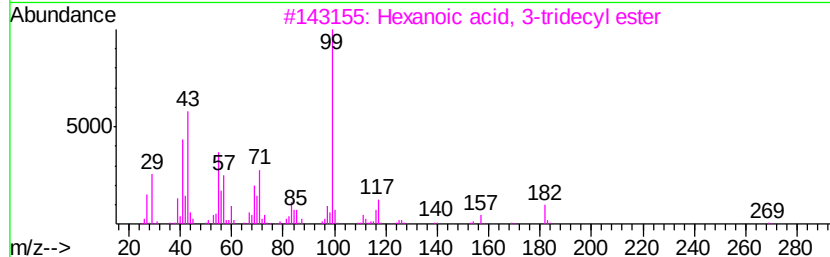
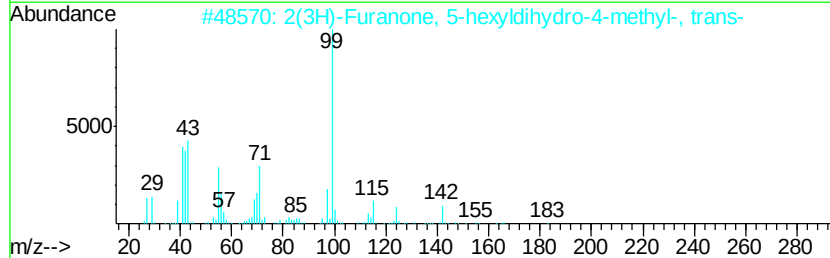
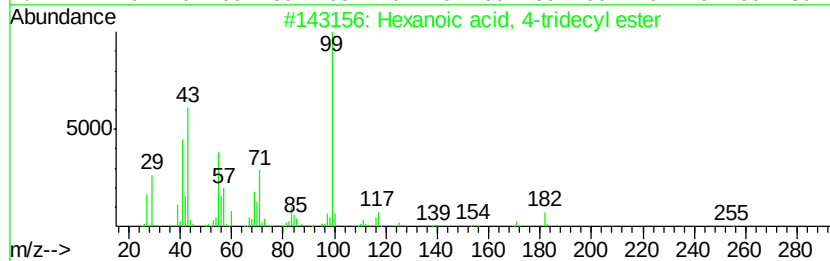
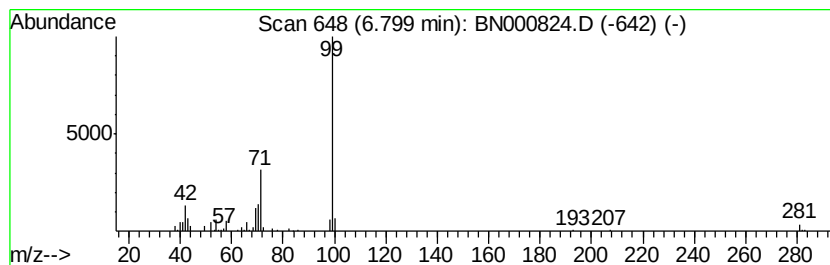
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Hexanoic acid, 4-tridecyl e... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.80	5.71 ng/ul	200462	1,4-Dichlorobenzene-d4	7.62	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 4-tridecyl ester	298	C19H38O2	1000279-29-3	64
2	2(3H)-Furanone, 5-hexyldihydro-4...	184	C11H20O2	147254-33-9	56
3	Hexanoic acid, 3-tridecyl ester	298	C19H38O2	1000279-29-2	56
4	Hexanoic acid, 5-tridecyl ester	298	C19H38O2	1000279-29-4	56
5	Hexanoic acid, 4-tetradecyl ester	312	C20H40O2	1000279-29-8	56



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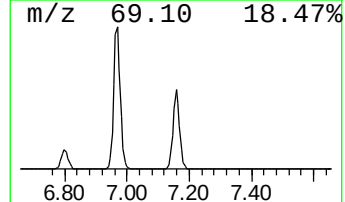
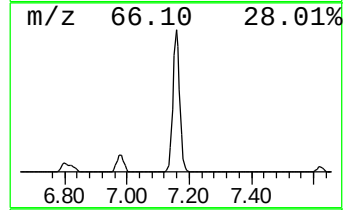
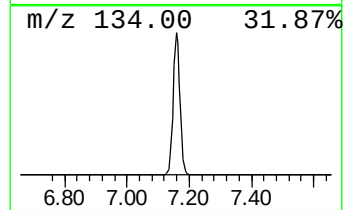
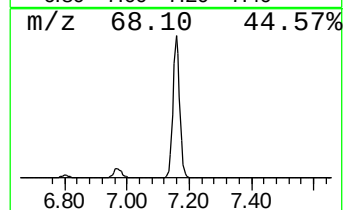
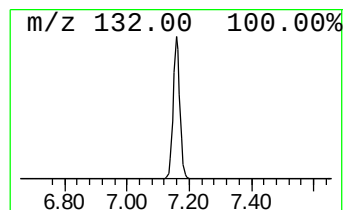
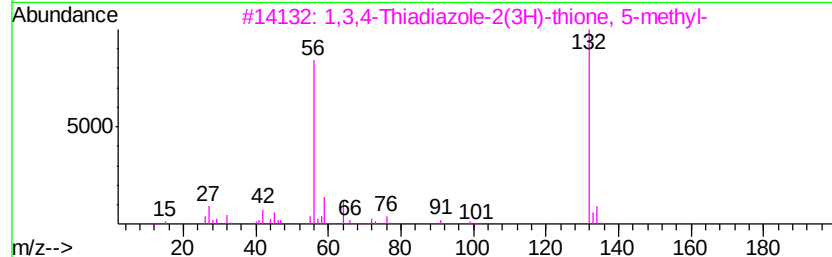
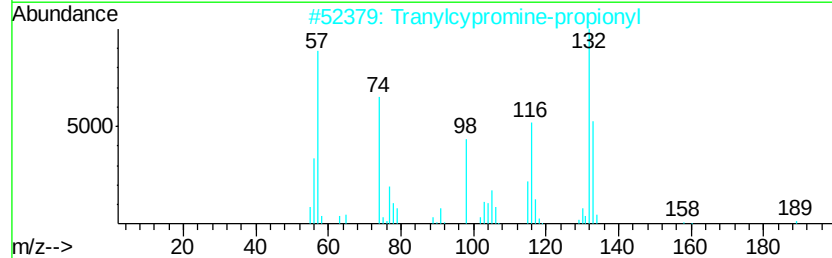
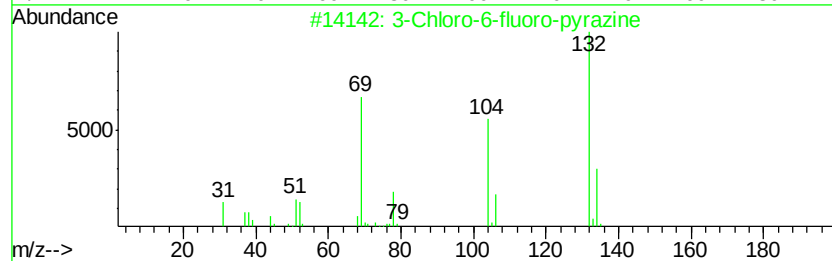
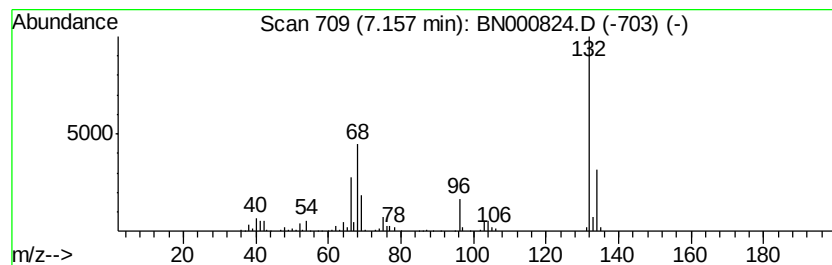
Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN041918-PAH.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.16	19.90 ng/ul	698035	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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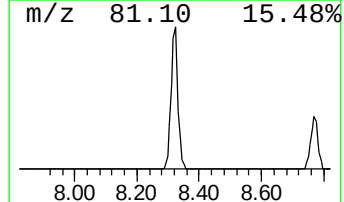
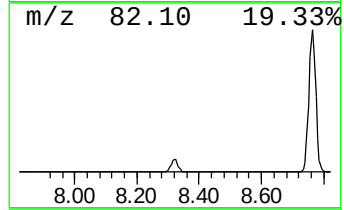
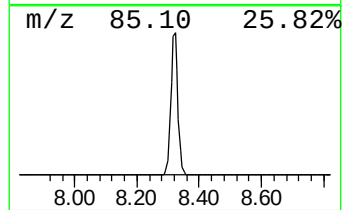
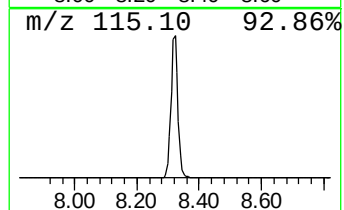
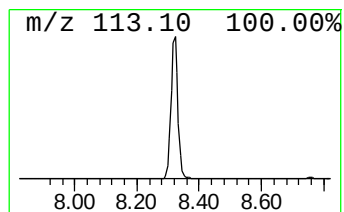
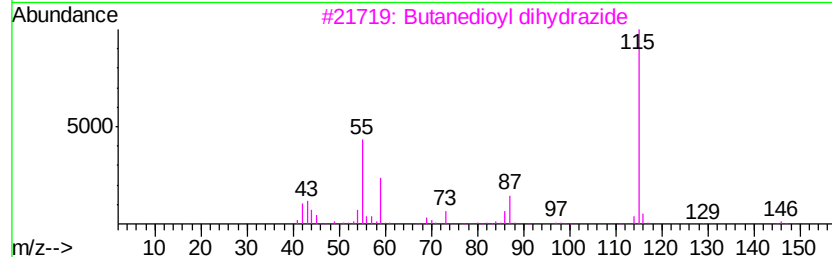
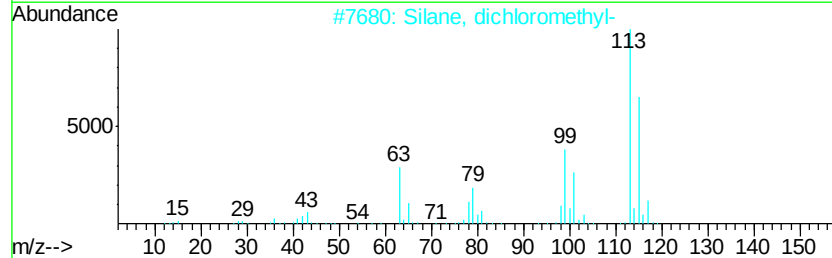
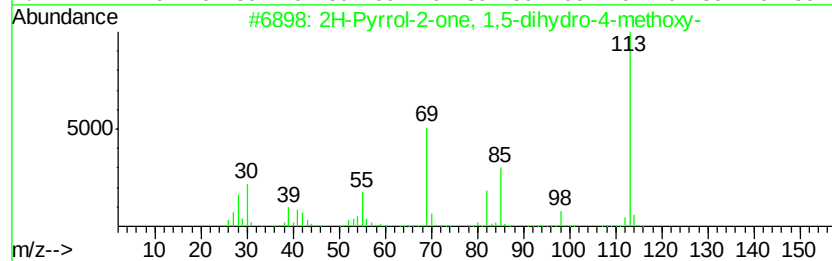
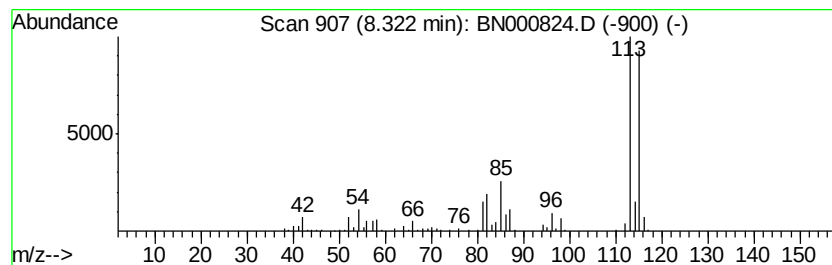
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.32	15.80 ng/ul	554280	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-Pyrrol-2-one, 1,5-dihydro-4-m...	113	C5H7NO2	069778-83-2	37
2		Silane, dichloromethyl-	114	CH4Cl2Si	000075-54-7	36
3		Butanedioyl dihydrazide	146	C4H10N4O2	004146-43-4	27
4		Heptanoic acid, anhydride	242	C14H26O3	000626-27-7	25
5		2(3H)-Oxazolethione, 4-methyl-	115	C4H5NOS	013016-17-6	22



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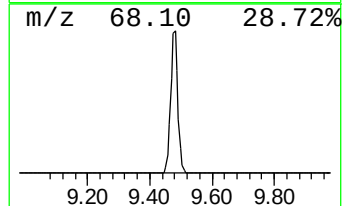
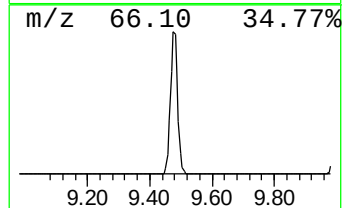
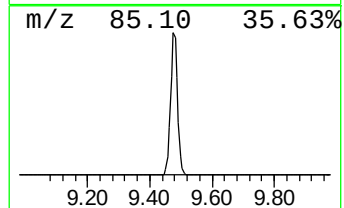
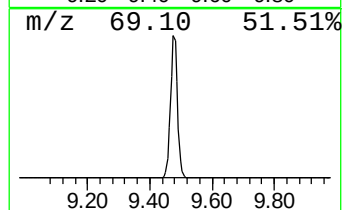
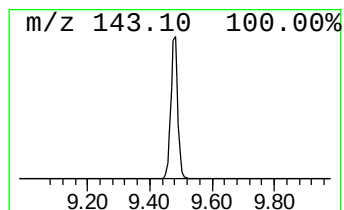
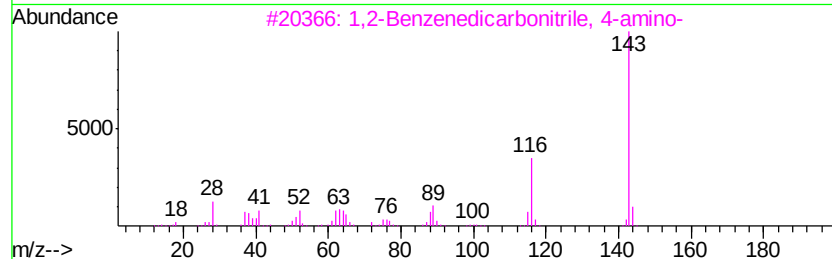
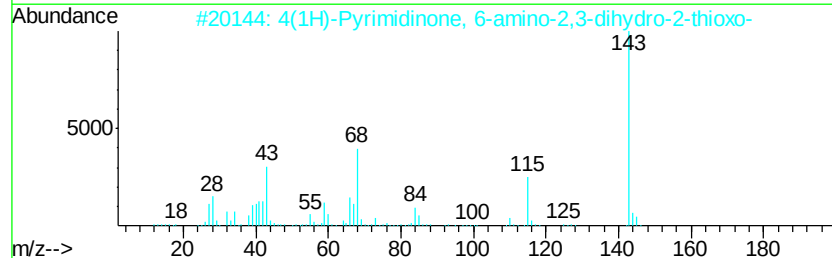
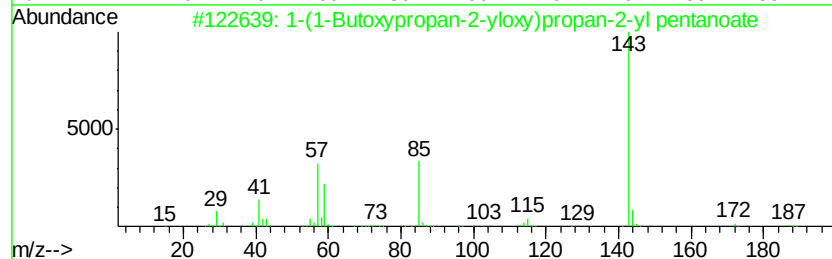
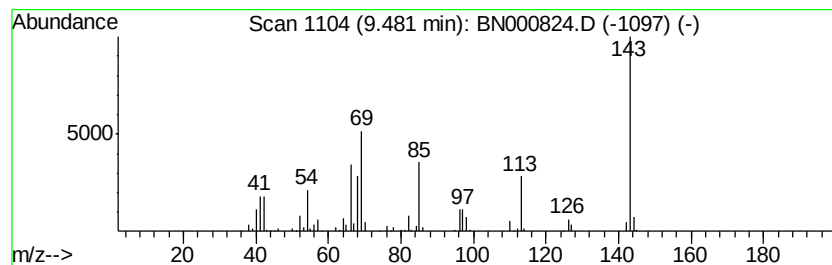
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 unknown-03 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.48	12.28 ng/ul	622842	Napthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-(1-Butoxypropan-2-yloxy)propan...	274	C15H30O4	1000367-13-0	22
2		4(1H)-Pyrimidinone, 6-amino-2,3-...	143	C4H5N3O	001004-40-6	17
3		1,2-Benzenedicarbonitrile, 4-amino-	143	C8H5N3	056765-79-8	14
4		1-(1-Butoxypropan-2-yloxy)propan...	274	C15H30O4	1000367-12-9	12
5		6-Chloro-N-methyl-3-pyridazinamine	143	C5H6ClN3	014959-32-1	12



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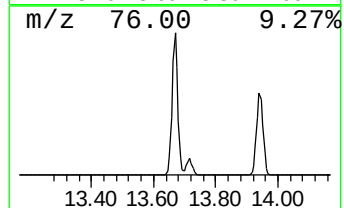
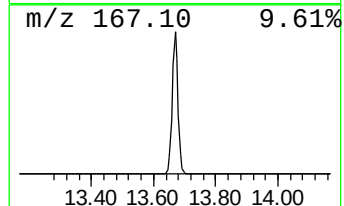
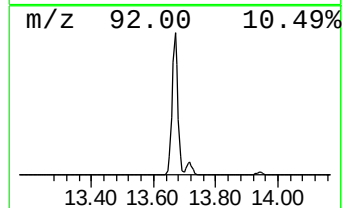
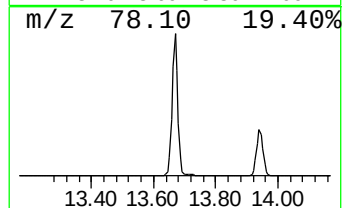
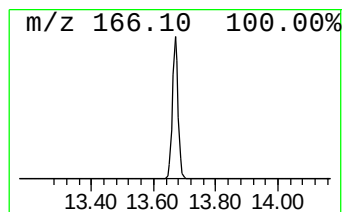
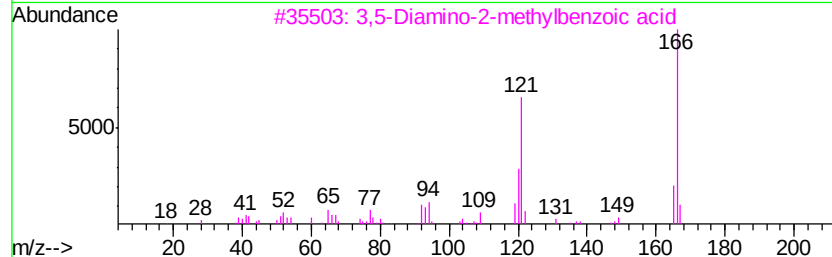
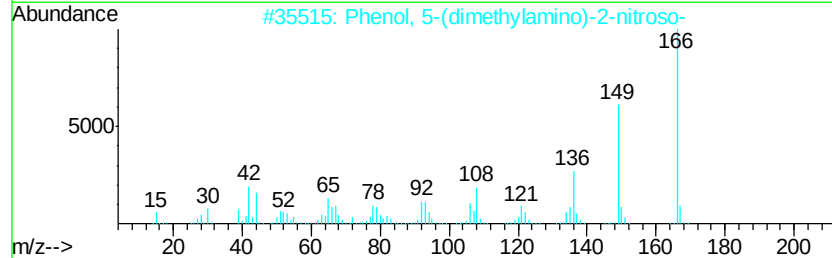
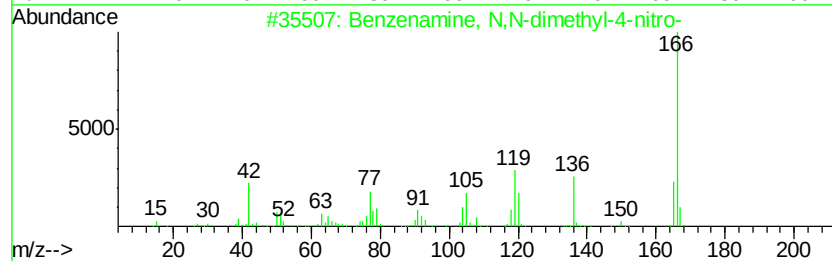
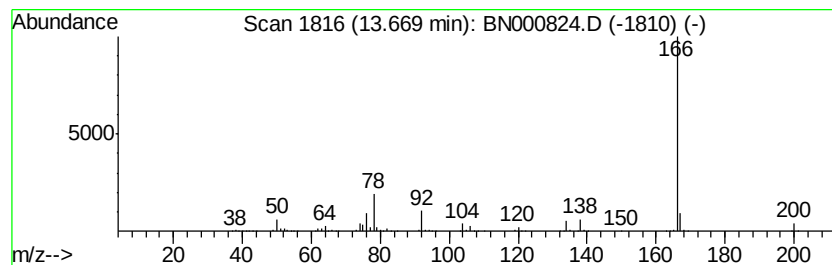
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Peak Number 7 unknown-04 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	22.01 ng/ul	1361150	Acenaphthene-d10	14.25
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzenamine, N,N-dimethyl-4-nitro-	166 C8H10N2O2	000100-23-2	36
2	Phenol, 5-(dimethylamino)-2-nitr...	166 C8H10N2O2	016761-04-9	33
3	3,5-Diamino-2-methylbenzoic acid	166 C8H10N2O2	090007-30-0	33
4	1H-Purine-2,6-dione, 3,7-dihydro...	166 C6H6N4O2	006136-37-4	9
5	p-Anisamidoxime	166 C8H10N2O2	005373-87-5	9



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Misc :
ALS Vial : 5 Sample Multiplier: 1

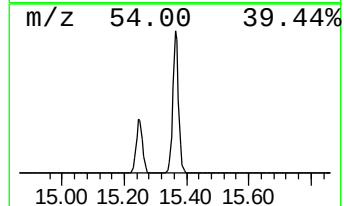
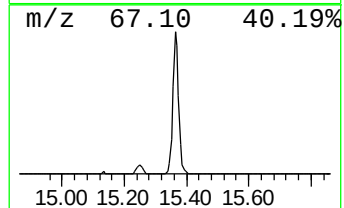
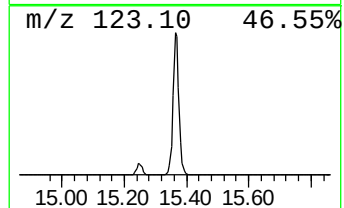
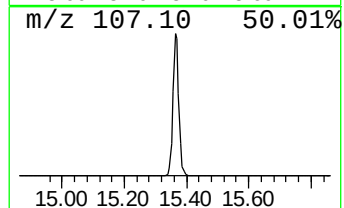
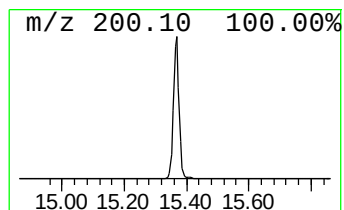
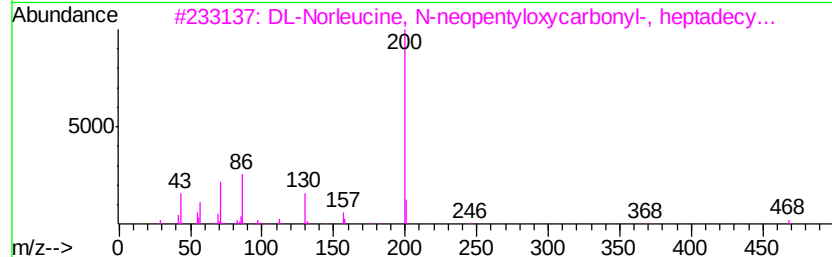
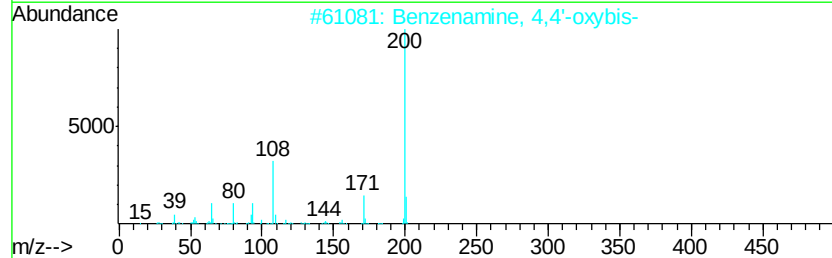
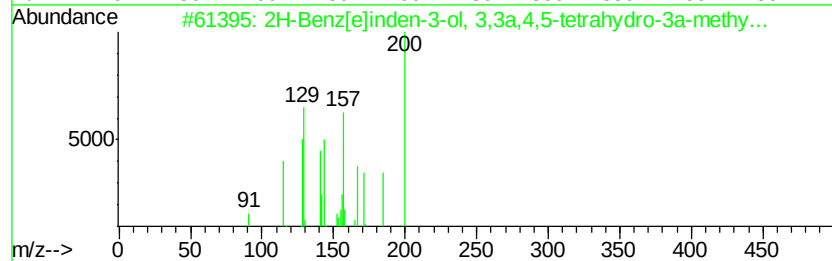
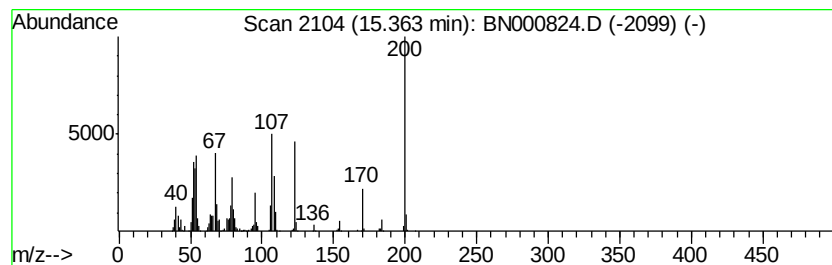
Instrument :
BNA_N
ClientSampleId :
A3XT8

Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041918-PAH.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-05 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.36	8.71 ng/ul	538833	Acenaphthene-d10	14.25
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2H-Benz[e]indene-3-ol, 3,3a,4,5-t...	200 C14H16O	071805-91-9	22
2	Benzenamine, 4,4'-oxybis-	200 C12H12N2O	000101-80-4	10
3	DL-Norleucine, N-neopentylloxycar...	483 C29H57NO4	1000339-05-9	9
4	2,3-Dihydro-6-isopropyl-1,2,4-tr...	200 C6H8N4S2	014778-89-3	9
5	Benzene, 1-methoxy-4-phenoxy-	200 C13H12O2	001655-69-2	9



Data Path : Z:\HPCHEM1\BNA_N\DATA\BN042318\
Data File : BN000824.D
Acq On : 23 Apr 2018 12:35
Operator : JU/SJ
Sample : J2467-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A3XT8

Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041918-PAH.M
Quant Title : SVOA 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
Hexanoic acid, 4-...	6.80	5.7	ng/ul	200462	1	7.62	701647	20.0
unknown-01	7.16	19.9	ng/ul	698035	1	7.62	701647	20.0
unknown-02	8.32	15.8	ng/ul	554280	1	7.62	701647	20.0
unknown-03	9.48	12.3	ng/ul	622842	2	10.39	1014630	20.0
unknown-04	13.67	22.0	ng/ul	1361150	3	14.25	1236640	20.0
unknown-05	15.36	8.7	ng/ul	538833	3	14.25	1236640	20.0