

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080524\
 Data File : BF138794.D
 Acq On : 05 Aug 2024 19:51
 Operator : RC/JU
 Sample : P3448-03 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP2

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-BF073024.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF138794.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.052	498	503	511	rBB	7371	10529	3.28%	0.362%
2	5.469	570	574	589	rBB	122676	172331	53.68%	5.924%
3	6.481	742	746	764	rBV	118984	175752	54.74%	6.042%
4	6.846	803	808	813	rBV	226493	277841	86.54%	9.551%
5	7.404	898	903	917	rVB	79986	103881	32.36%	3.571%
6	8.128	1020	1026	1031	rBV	251827	321040	100.00%	11.036%
7	8.204	1036	1039	1042	rVB2	3460	4099	1.28%	0.141%
8	8.416	1072	1075	1079	rVV2	5721	7112	2.22%	0.244%
9	8.534	1092	1095	1098	rBV3	2667	4041	1.26%	0.139%
10	8.581	1098	1103	1107	rVB3	7425	12660	3.94%	0.435%
11	8.657	1110	1116	1122	rBV8	5873	13979	4.35%	0.481%
12	8.769	1129	1135	1138	rBV3	7629	14451	4.50%	0.497%
13	8.804	1138	1141	1145	rVB3	7574	10233	3.19%	0.352%
14	8.863	1146	1151	1154	rBV5	5854	9940	3.10%	0.342%
15	8.898	1154	1157	1162	rVV5	3906	6537	2.04%	0.225%
16	8.998	1171	1174	1181	rBV7	5808	11961	3.73%	0.411%
17	9.081	1185	1188	1190	rBV3	4710	5762	1.79%	0.198%
18	9.110	1190	1193	1196	rBV3	9988	13410	4.18%	0.461%
19	9.198	1203	1208	1216	rBV	129488	178289	55.53%	6.129%
20	9.269	1216	1220	1226	rVV4	30551	45290	14.11%	1.557%
21	9.581	1270	1273	1278	rBV2	16281	19933	6.21%	0.685%
22	9.639	1280	1283	1287	rBV4	12080	15957	4.97%	0.549%
23	9.739	1296	1300	1302	rBV2	22282	27329	8.51%	0.939%
24	9.781	1305	1307	1311	rVB3	25414	30658	9.55%	1.054%
25	9.822	1311	1314	1316	rBV3	11341	14947	4.66%	0.514%
26	9.881	1317	1324	1328	rVV	220580	292229	91.03%	10.046%
27	10.281	1389	1392	1396	rBV	27942	30946	9.64%	1.064%
28	10.669	1454	1458	1463	rBV	55085	73955	23.04%	2.542%
29	11.369	1572	1577	1586	rBV	212051	292488	91.11%	10.055%
30	12.580	1779	1783	1788	rVB	52550	65529	20.41%	2.253%
31	12.810	1818	1822	1829	rVB	48246	68052	21.20%	2.339%
32	12.945	1841	1845	1852	rVB	114433	155997	48.59%	5.363%
33	14.010	2021	2026	2036	rVB2	164670	263327	82.02%	9.052%
34	15.474	2270	2275	2283	rVB	99026	158483	49.37%	5.448%

Sum of corrected areas: 2908968

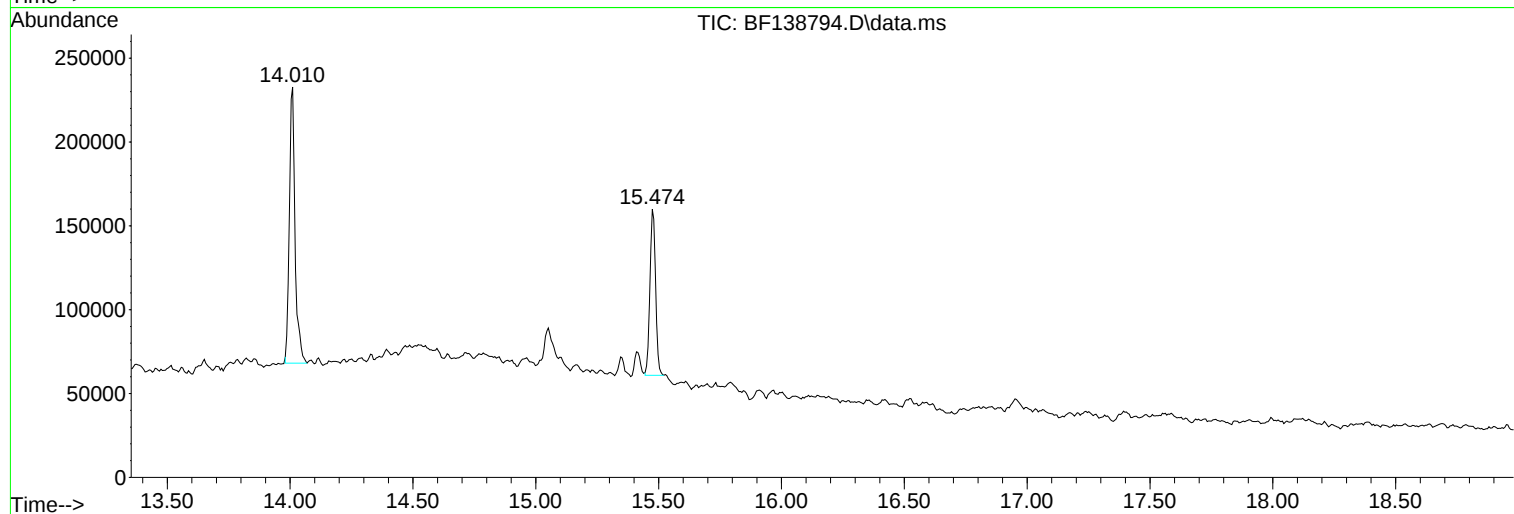
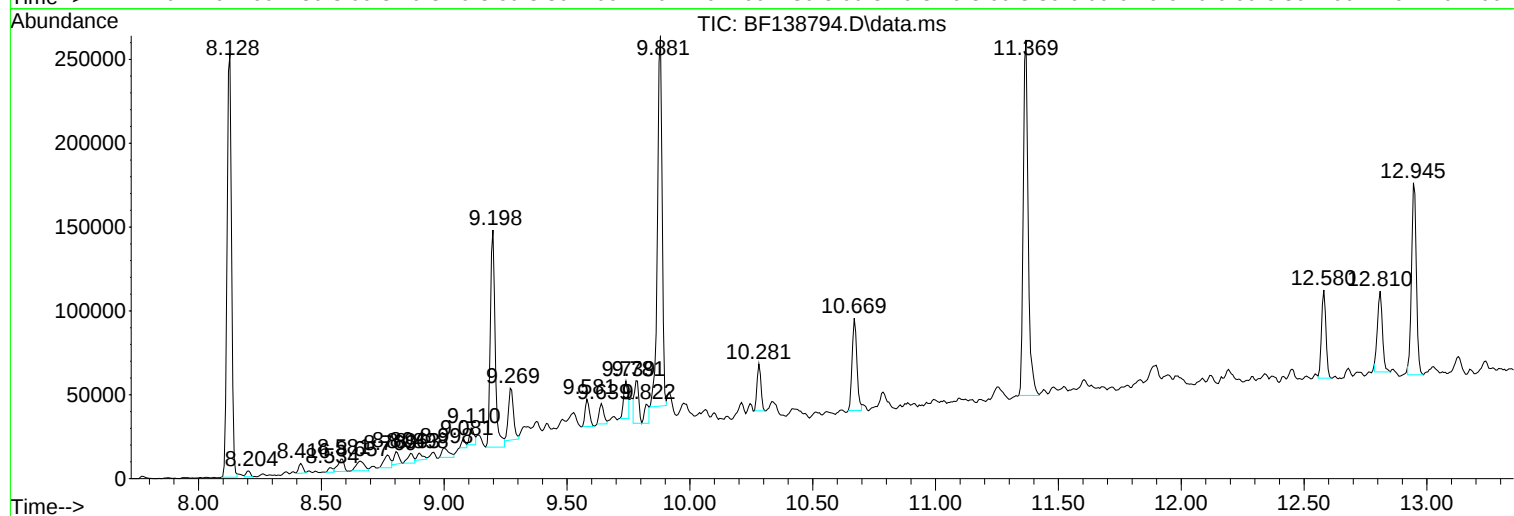
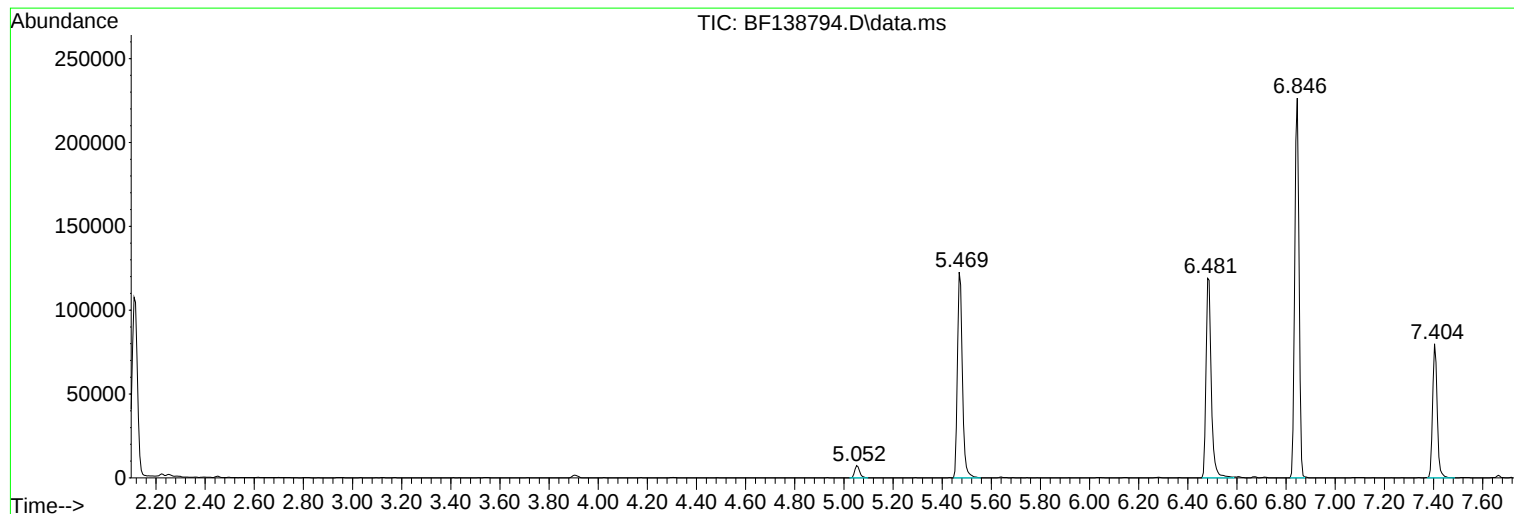
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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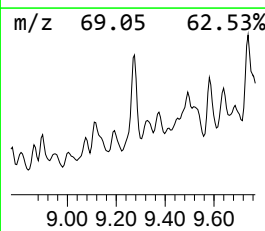
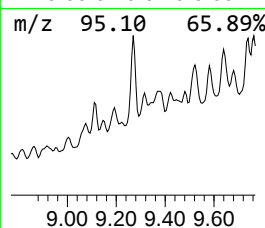
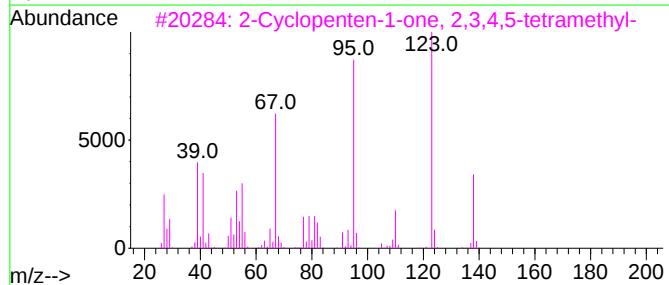
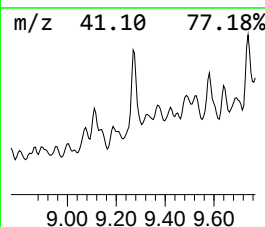
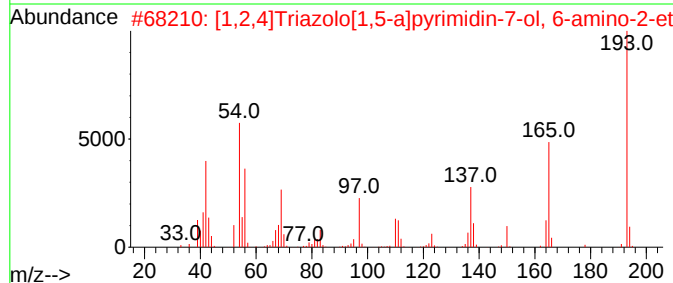
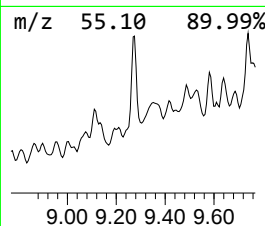
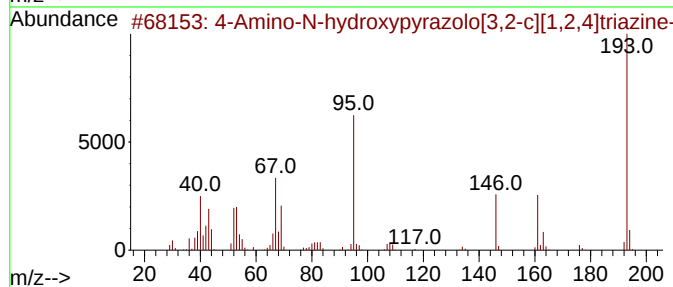
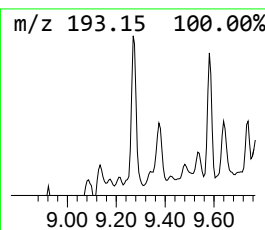
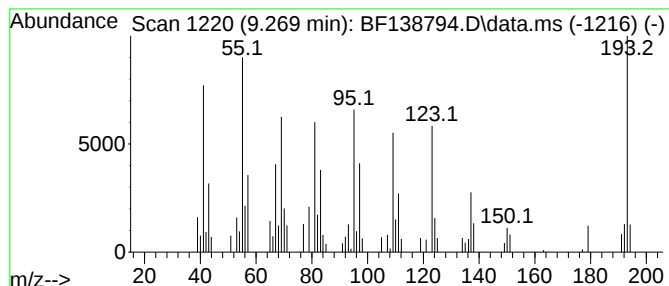
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown9.269 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.269	3.10 ng	45290	Acenaphthene-d10	9.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Amino-N-hydroxypyrazolo[3,2-c]...	193	C6H7N7O	1000443-50-2	38
2			[1,2,4]Triazolo[1,5-a]pyrimidin-...	193	C8H11N5O	1000351-50-1	38
3			2-Cyclopenten-1-one, 2,3,4,5-tet...	138	C9H14O	054458-61-6	30
4			2-Propanone, 1-(1-cyclohexen-1-yl)-	138	C9H14O	000768-50-3	25
5			2,4,6-Trimethoxybenzonitrile	193	C10H11NO3	002571-54-2	18



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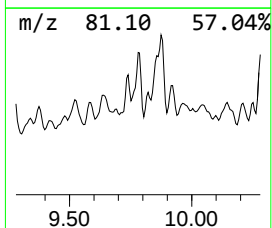
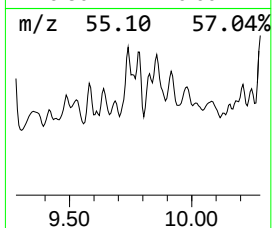
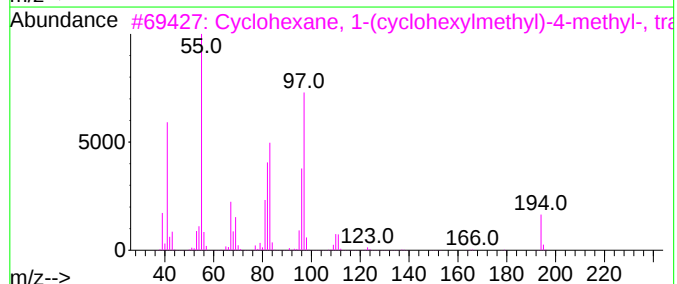
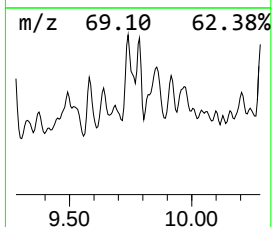
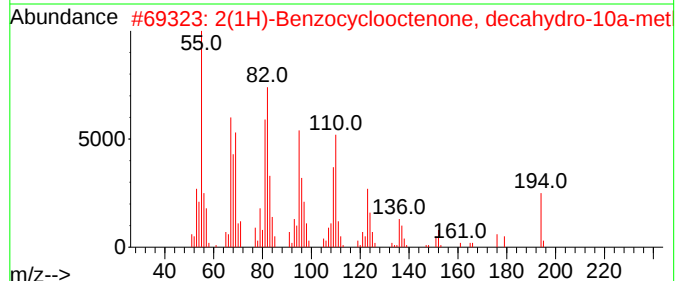
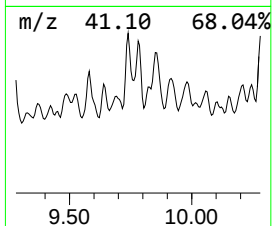
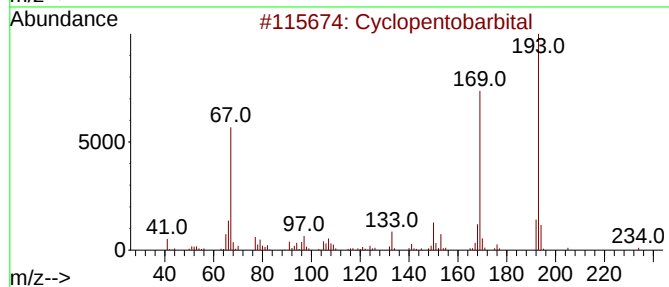
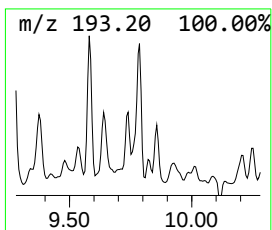
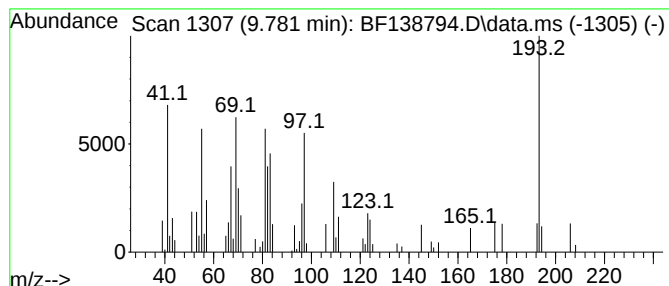
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Peak Number 2 unknown9.781 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.781	2.10 ng	30658	Acenaphthene-d10	9.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentobarbital	234	C12H14N2O3	000076-68-6	35
2			2(1H)-Benzocyclooctenone, decahy...	194	C13H22O	055103-68-9	30
3			Cyclohexane, 1-(cyclohexylmethyl)...	194	C14H26	054823-98-2	30
4			Cyclohexane, 1-(cyclohexylmethyl)...	194	C14H26	054823-97-1	25
5			6-Methylphenanthridine	193	C14H11N	003955-65-5	25



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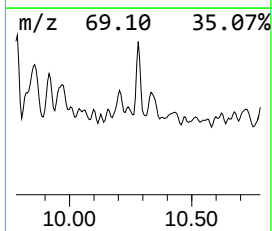
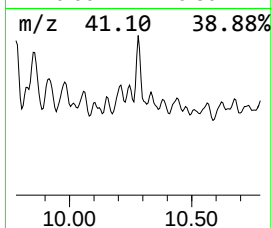
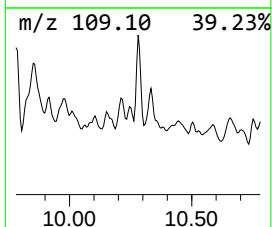
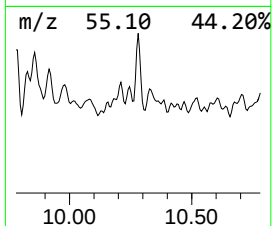
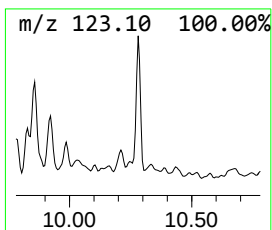
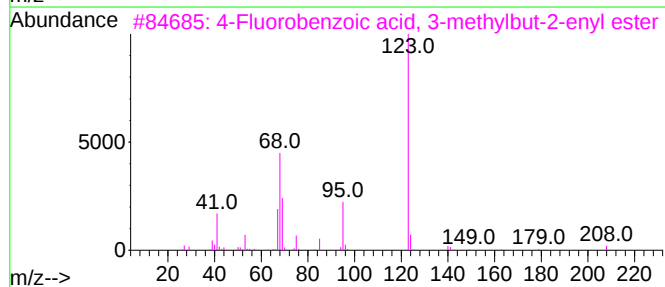
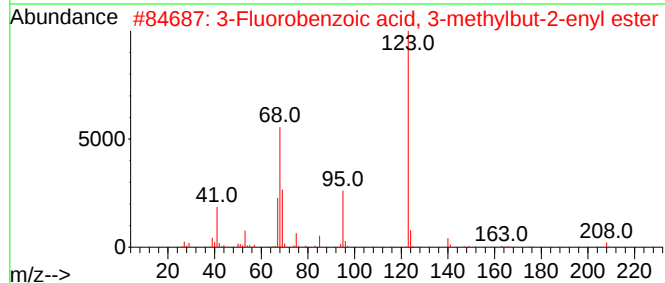
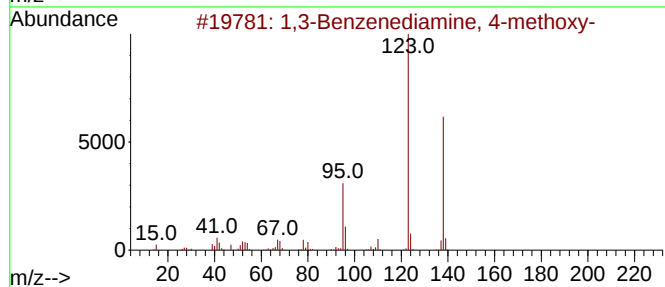
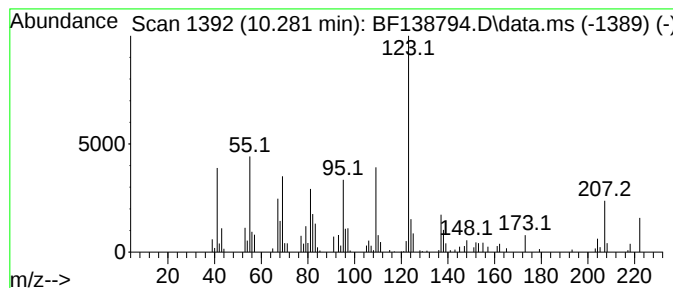
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown10.281 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.281	2.12 ng	30946	Acenaphthene-d10	9.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Benzenediamine, 4-methoxy-	138	C7H10N2O	000615-05-4	46
2			3-Fluorobenzoic acid, 3-methylbu...	208	C12H13F02	154559-38-3	43
3			4-Fluorobenzoic acid, 3-methylbu...	208	C12H13F02	1000299-15-1	43
4			4-Fluorobenzoic acid, 3-pentadec...	350	C22H35F02	1000280-68-4	43
5			2-Pyrimidinamine, 4,6-dimethyl-	123	C6H9N3	000767-15-7	43



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown9.269	9.269	3.1	ng	45290	3	9.881	292229	20.0
unknown9.781	9.781	2.1	ng	30658	3	9.881	292229	20.0
unknown10.281	10.281	2.1	ng	30946	3	9.881	292229	20.0