

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100621\
 Data File : BP007361.D
 Acq On : 06 Oct 2021 19:09
 Operator : CG/JU
 Sample : M4039-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-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\8270E-BP092121.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP007361.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.440	393	400	418	rBV	2128096	3246759	53.04%	9.010%
2	7.040	664	672	691	rBV	1840544	3153462	51.52%	8.751%
3	7.857	801	811	819	rBV	604300	969383	15.84%	2.690%
4	9.022	1002	1009	1017	rBV	1157990	2036538	33.27%	5.651%
5	10.451	1244	1252	1267	rBV	206819	525557	8.59%	1.458%
6	10.657	1276	1287	1302	rBV	745289	1380079	22.55%	3.830%
7	12.075	1522	1528	1534	rBV	46338	75659	1.24%	0.210%
8	13.116	1696	1705	1725	rBV	2918845	4589993	74.99%	12.737%
9	13.316	1735	1739	1746	rVB	62700	94220	1.54%	0.261%
10	14.386	1915	1921	1928	rBV2	84138	130814	2.14%	0.363%
11	14.492	1930	1939	1952	rBV	1073467	1695469	27.70%	4.705%
12	15.339	2076	2083	2088	rVB2	85367	111780	1.83%	0.310%
13	15.745	2141	2152	2163	rBV	38055	96540	1.58%	0.268%
14	15.986	2186	2193	2215	rBV	2067055	3446416	56.31%	9.564%
15	16.204	2226	2230	2233	rBV	104110	151269	2.47%	0.420%
16	16.998	2361	2365	2369	rBV	92357	115683	1.89%	0.321%
17	17.051	2371	2374	2386	rVB2	46724	86958	1.42%	0.241%
18	17.245	2401	2407	2413	rBV	1321456	1942948	31.74%	5.392%
19	17.739	2487	2491	2495	rBV	83993	104886	1.71%	0.291%
20	18.015	2533	2538	2544	rBV3	44446	107238	1.75%	0.298%
21	18.133	2554	2558	2562	rBV	136384	180161	2.94%	0.500%
22	18.415	2602	2606	2610	rBV	90091	110770	1.81%	0.307%
23	19.021	2706	2709	2718	rVB	77791	107816	1.76%	0.299%
24	19.257	2745	2749	2755	rVB	148616	194124	3.17%	0.539%
25	19.580	2801	2804	2806	rBV2	93464	107854	1.76%	0.299%
26	19.610	2806	2809	2818	rVB	164726	257235	4.20%	0.714%
27	19.804	2837	2842	2858	rVB	5058190	6120789	100.00%	16.985%
28	20.586	2972	2975	2978	rBV	117043	113785	1.86%	0.316%
29	21.045	3049	3053	3061	rBV2	268732	405479	6.62%	1.125%
30	21.333	3097	3102	3107	rBV	1753556	2254774	36.84%	6.257%
31	23.739	3506	3511	3522	rVB2	1079205	2122534	34.68%	5.890%

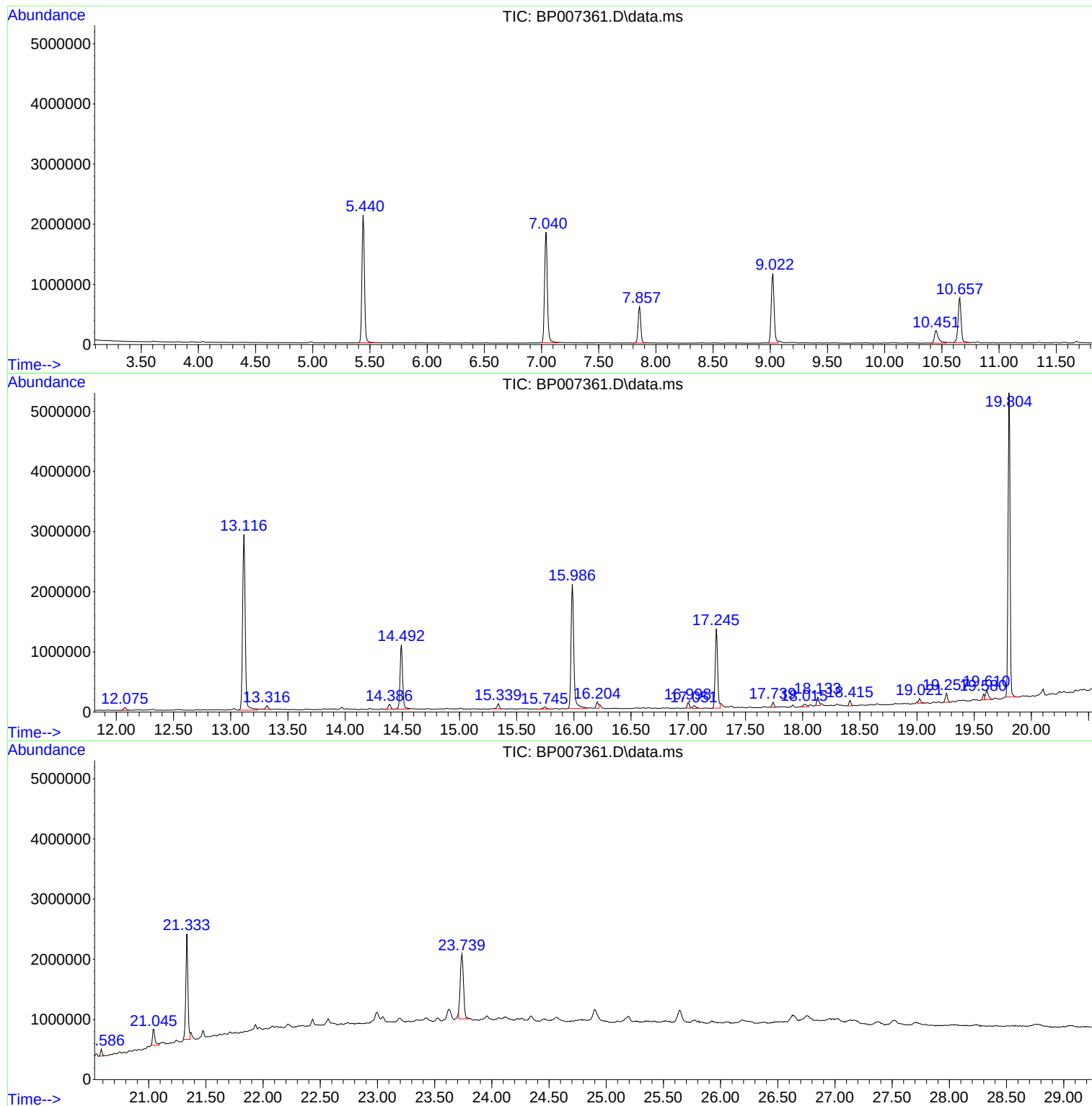
Sum of corrected areas: 36036972

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ClientSampled :
TP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.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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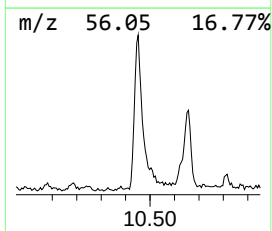
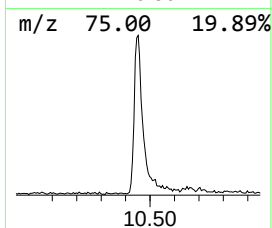
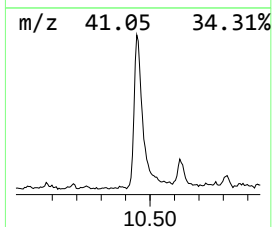
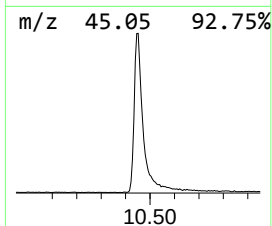
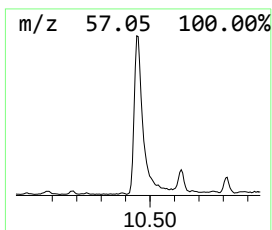
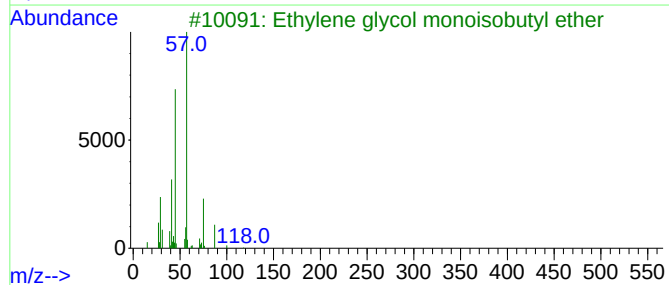
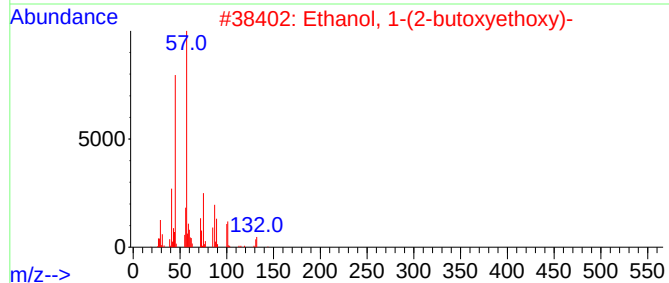
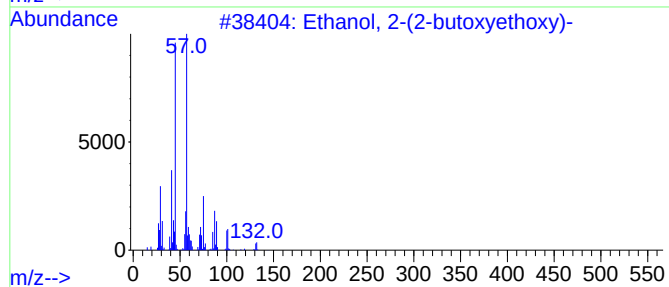
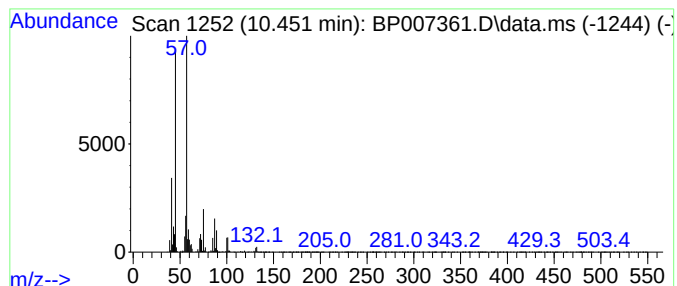
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.451	7.62 ng	525557	Naphthalene-d8	10.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	91
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
4			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	56
5			3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	50



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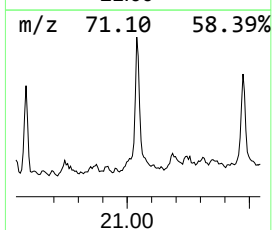
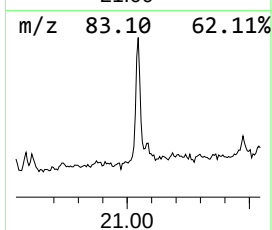
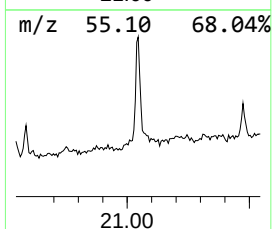
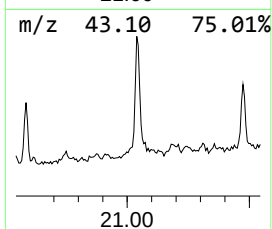
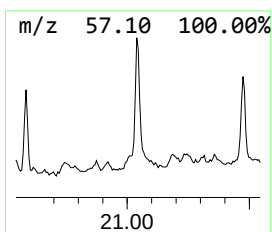
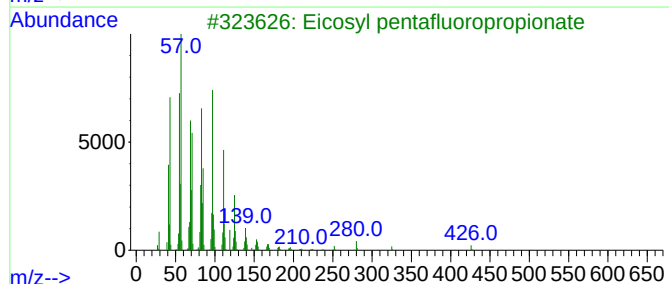
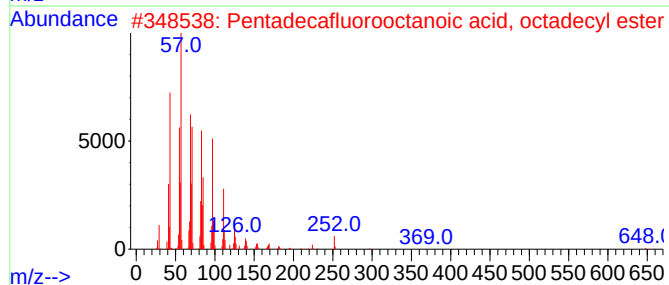
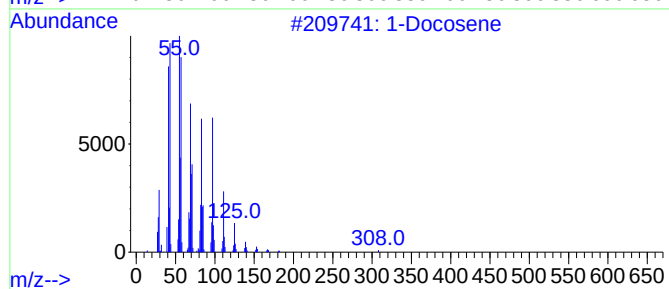
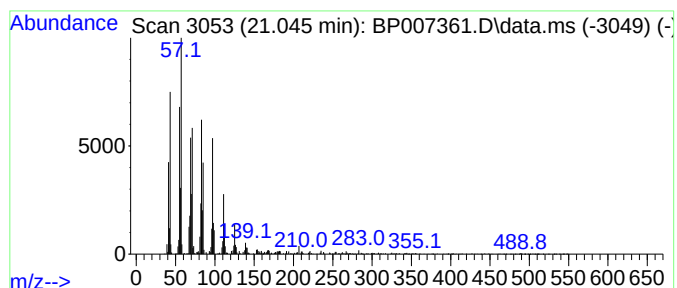
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Peak Number 3 1-Docosene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.045	3.60 ng	405479	Chrysene-d12	21.333

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene			308	C22H44	001599-67-3	95
2	Pentadecafluorooctanoic acid, oc...			666	C26H37F15O2	1000406-04-8	93
3	Eicosyl pentafluoropropionate			444	C23H41F5O2	1000351-80-8	91
4	1-Triacontanol			438	C30H62O	000593-50-0	91
5	Docosyl pentafluoropropionate			472	C25H45F5O2	1000351-80-9	91



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-b...	10.451	7.6	ng	525557	2	10.657	1380080	20.0
1-Docosene	21.045	3.6	ng	405479	5	21.333	2254770	20.0