

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012420\
 Data File : BM024617.D
 Acq On : 24 Jan 2020 12:31
 Operator : CG/JU
 Sample : PB125860BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB125860BL

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_M\METHODS\8270-BM012320.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	5.087	367	374	382	rBV	4174799	5860538	30.29%	7.373%
2	6.640	631	638	658	rBV	4055303	6321671	32.67%	7.954%
3	7.439	768	774	784	rBV	800277	1235777	6.39%	1.555%
4	7.757	821	828	838	rBV	3378696	5159376	26.66%	6.491%
5	8.581	961	968	979	rBV	2645210	4155563	21.47%	5.228%
6	10.198	1236	1243	1256	rVB	1040251	1691042	8.74%	2.128%
7	12.704	1662	1669	1680	rBV	7370138	10453220	54.02%	13.152%
8	14.092	1898	1905	1916	rVB	1720360	2539680	13.12%	3.195%
9	15.586	2150	2159	2169	rBV	6611234	9244114	47.77%	11.631%
10	16.839	2365	2372	2381	rBV	2313928	3229197	16.69%	4.063%
11	17.780	2527	2532	2538	rBV2	142654	200089	1.03%	0.252%
12	19.498	2818	2824	2834	rBV	15238445	19351250	100.00%	24.347%
13	20.809	3042	3047	3052	rBV3	339730	547955	2.83%	0.689%
14	21.044	3082	3087	3101	rBV	3158002	4085940	21.11%	5.141%
15	21.674	3191	3194	3201	rVB2	185510	227222	1.17%	0.286%
16	22.115	3265	3269	3276	rBV	211823	259152	1.34%	0.326%
17	22.591	3346	3350	3354	rBV	228642	286603	1.48%	0.361%
18	23.121	3436	3440	3441	rBV	212119	260477	1.35%	0.328%
19	23.162	3441	3447	3455	rVB	2231154	4083406	21.10%	5.138%
20	23.721	3538	3542	3548	rVB	172793	289180	1.49%	0.364%

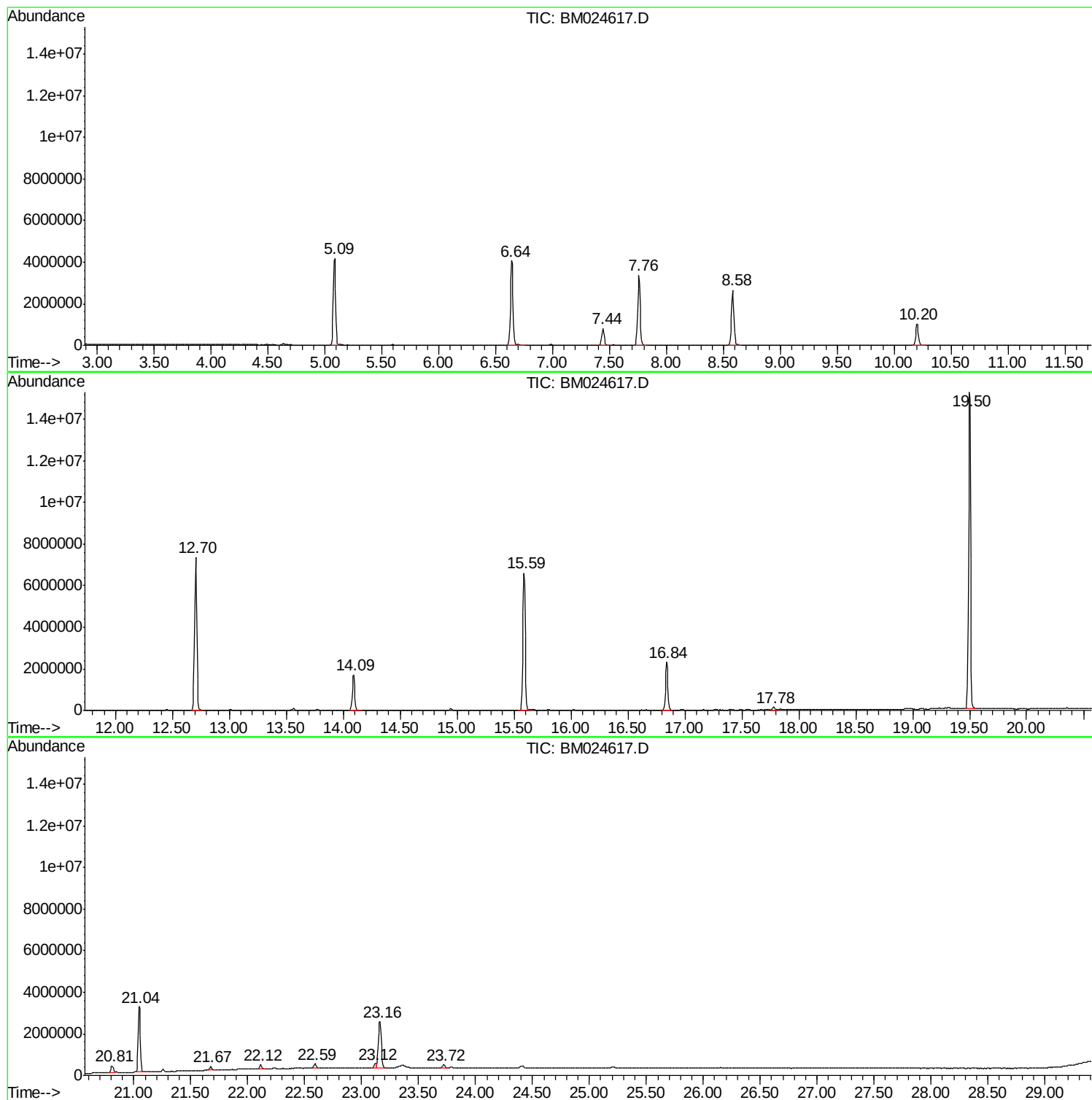
Sum of corrected areas: 79481452

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM012320.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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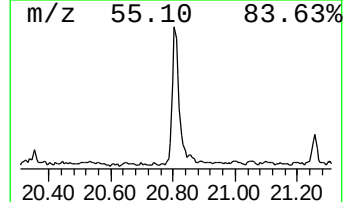
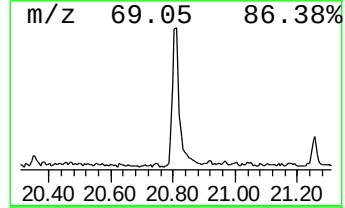
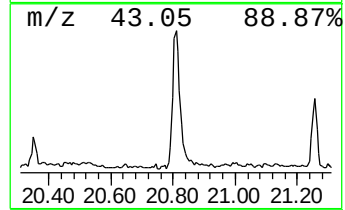
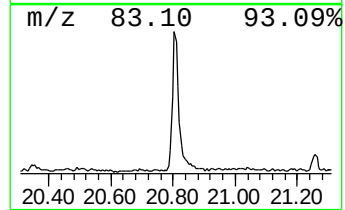
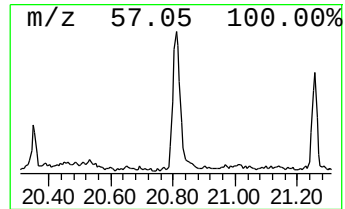
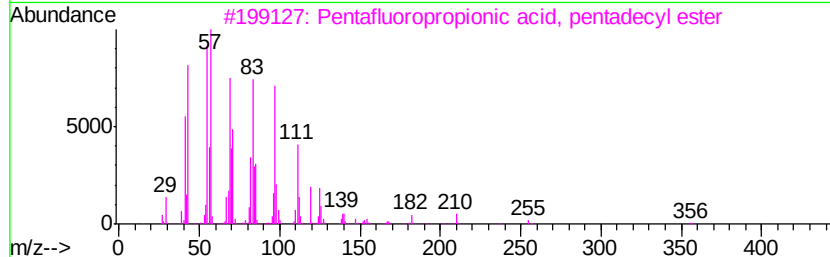
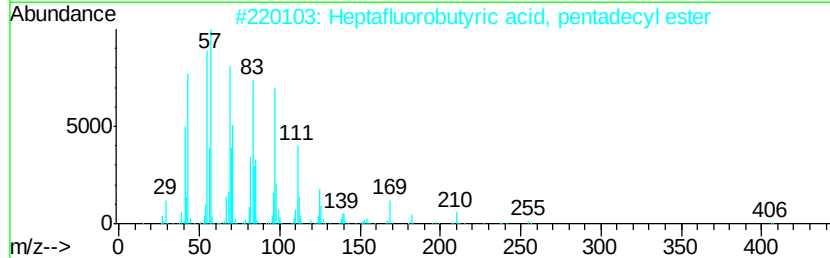
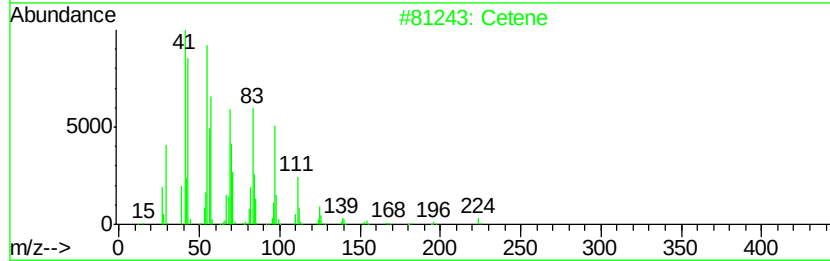
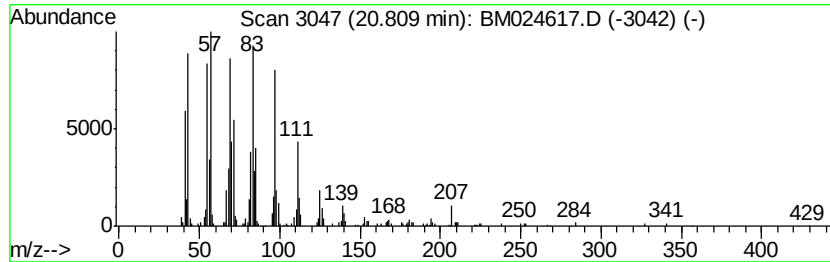
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Peak Number 2 Cetene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.81	2.68 ng	547955	Chrysene-d12	21.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cetene	224	C16H32	000629-73-2	96
2		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
3		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
4		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
5		Behenic alcohol	326	C22H46O	000661-19-8	93



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cetene	20.81	2.7	ng	547955	5	21.04	4085940	20.0