

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120822\
 Data File : BP012999.D
 Acq On : 08 Dec 2022 23:56
 Operator : CG/JU
 Sample : PB149468BL
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK468

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012999.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.352	24	28	38	rVB	140233	207803	1.23%	0.125%
2	3.781	97	101	123	rBV	1768388	2846751	16.87%	1.713%
3	4.110	154	157	163	rVB	38749	58616	0.35%	0.035%
4	7.128	663	670	681	rBV	2155595	3325206	19.70%	2.001%
5	7.298	692	699	712	rVB	1905497	2922250	17.31%	1.759%
6	7.504	727	734	742	rBV	2222881	3520895	20.86%	2.119%
7	7.975	807	814	824	rBV	2147606	3363047	19.93%	2.024%
8	8.681	927	934	947	rBV	2805791	4351971	25.79%	2.619%
9	9.139	1005	1012	1023	rBV	2091762	3438266	20.37%	2.069%
10	9.863	1128	1135	1150	rBV	1735215	2873576	17.03%	1.729%
11	10.404	1220	1227	1241	rBV	2850013	4640721	27.50%	2.793%
12	10.792	1285	1293	1303	rVB	3031252	5133217	30.41%	3.089%
13	10.922	1308	1315	1327	rBV	2680643	4786979	28.36%	2.881%
14	12.369	1555	1561	1567	rBV	53350	94794	0.56%	0.057%
15	13.586	1764	1768	1777	rVB6	9595	18005	0.11%	0.011%
16	14.022	1835	1842	1852	rBV	5694904	7891706	46.76%	4.750%
17	14.310	1880	1891	1909	rBV	6240904	9179019	54.39%	5.524%
18	14.616	1934	1943	1955	rBV	5179575	7658537	45.38%	4.609%
19	14.804	1967	1975	1994	rBV	2282425	3458392	20.49%	2.081%
20	15.022	2009	2012	2018	rVB8	22090	36356	0.22%	0.022%
21	15.610	2104	2112	2123	rBV	8146986	11815962	70.01%	7.111%
22	15.722	2125	2131	2147	rVV	2736950	3625497	21.48%	2.182%
23	15.845	2149	2152	2157	rVB2	41835	58263	0.35%	0.035%
24	17.045	2352	2356	2360	rVB5	10900	17961	0.11%	0.011%
25	17.157	2371	2375	2383	rVB3	13501	23441	0.14%	0.014%
26	17.368	2402	2411	2421	rBV	6817514	9582607	56.78%	5.767%
27	17.468	2421	2428	2442	rBV2	9185417	13115090	77.71%	7.893%
28	17.751	2472	2476	2479	rBV6	10648	18663	0.11%	0.011%
29	19.274	2731	2735	2741	rVB	122442	157262	0.93%	0.095%
30	19.339	2742	2746	2751	rBV	107236	152238	0.90%	0.092%
31	19.698	2801	2807	2816	rBV	12464718	16064062	95.18%	9.668%
32	21.451	3100	3105	3112	rBV	7859699	10263920	60.81%	6.177%
33	23.156	3392	3395	3405	rVB	495099	809684	4.80%	0.487%
34	23.780	3494	3501	3518	rVB2	7418712	16877632	100.00%	10.158%
35	23.945	3522	3529	3544	rVB2	5110917	10426735	61.78%	6.275%
36	24.574	3631	3636	3651	rVB	791685	1645962	9.75%	0.991%

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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Title : SVOA CALIBRATION

37	25.456	3781	3786	3798	rVB2	696259	1695664	10.05%	1.021%
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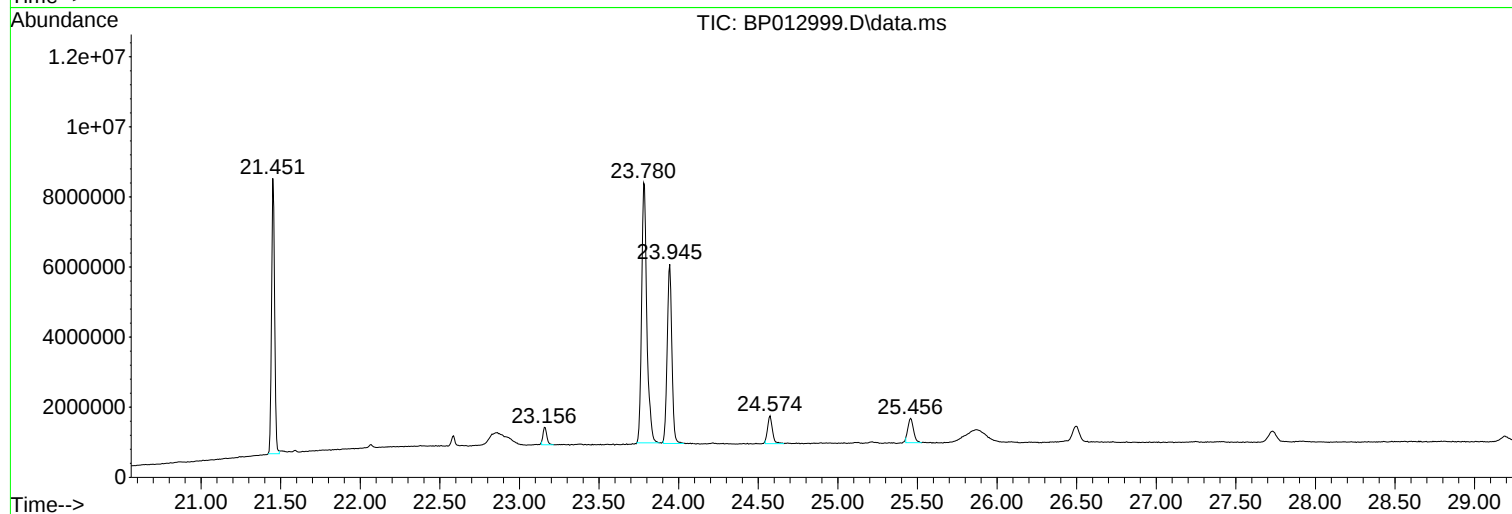
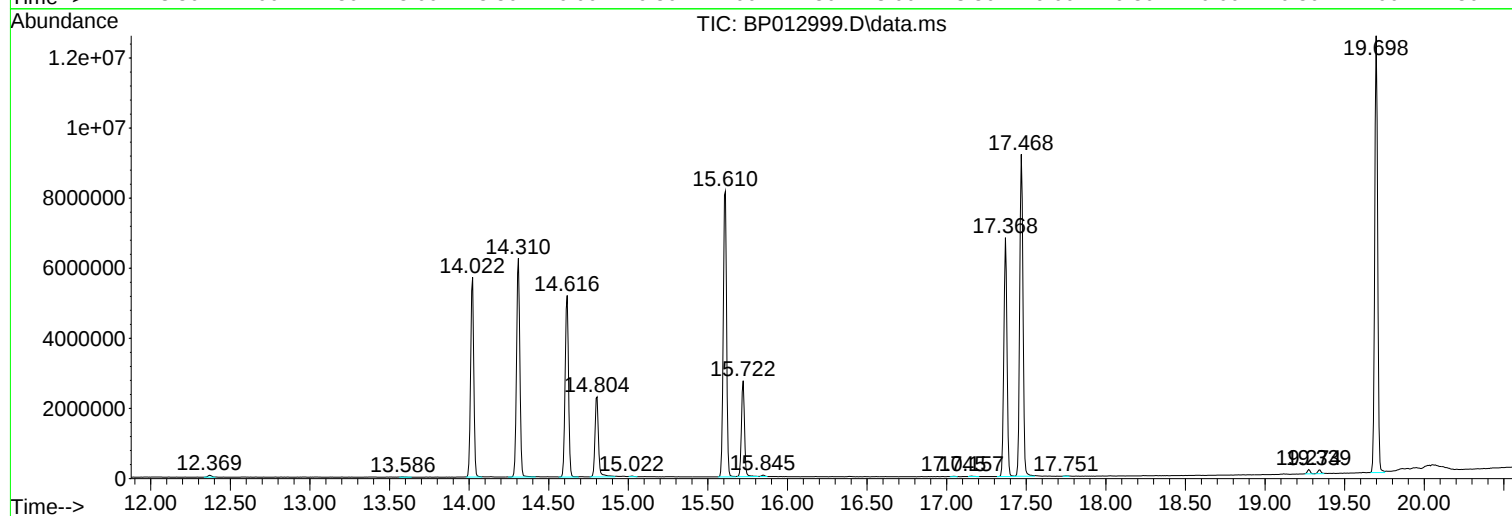
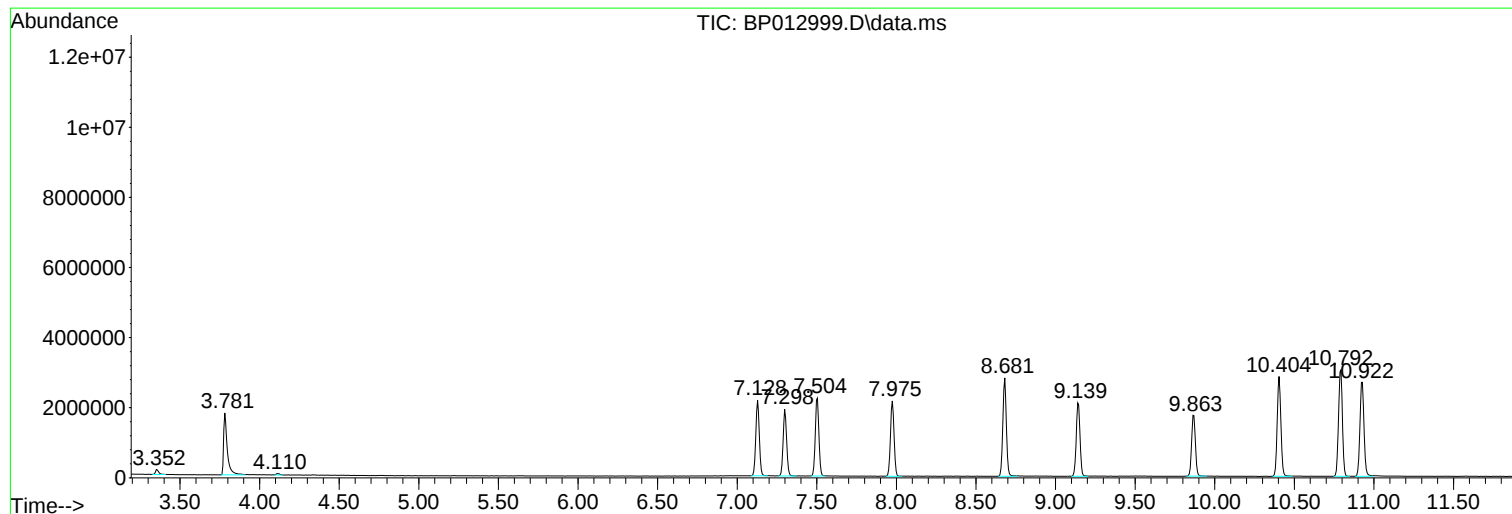
Sum of corrected areas: 166156750

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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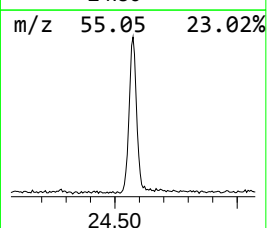
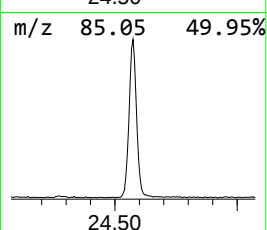
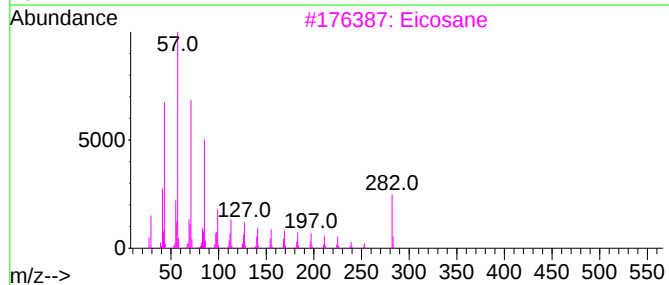
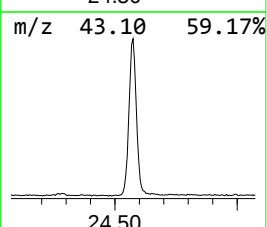
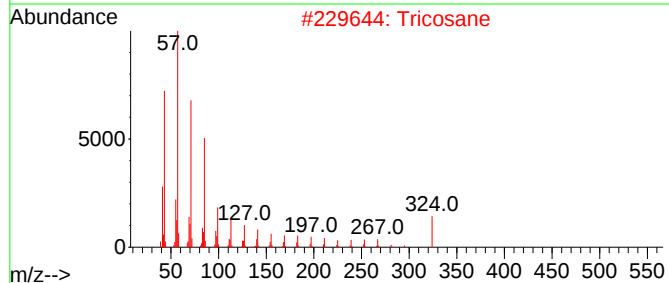
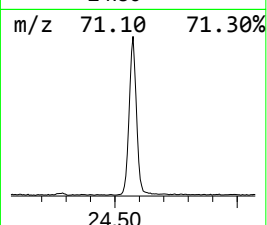
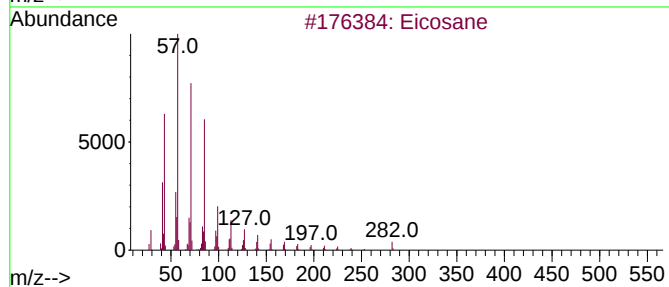
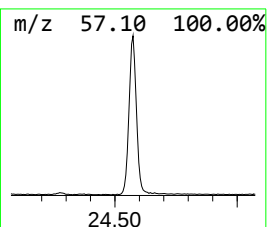
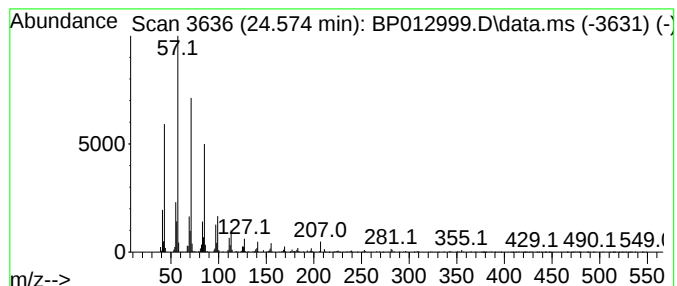
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.574	3.16 ng/ul	1645960	Perylene-d12	23.945

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	95
2			Tricosane	324	C23H48	000638-67-5	94
3			Eicosane	282	C20H42	000112-95-8	93
4			Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
5			Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	93



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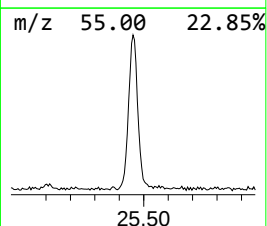
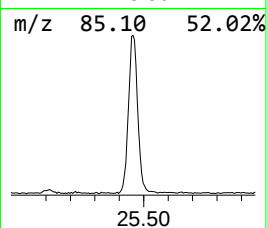
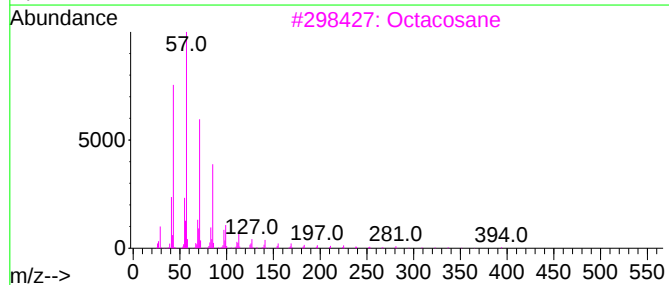
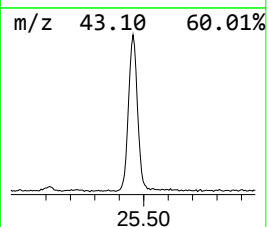
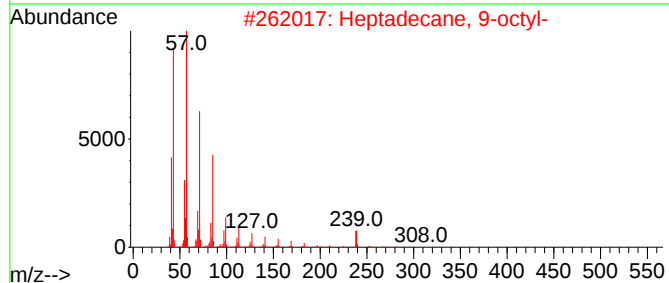
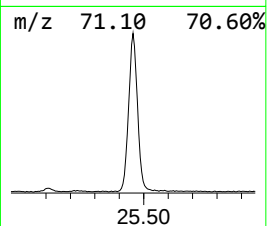
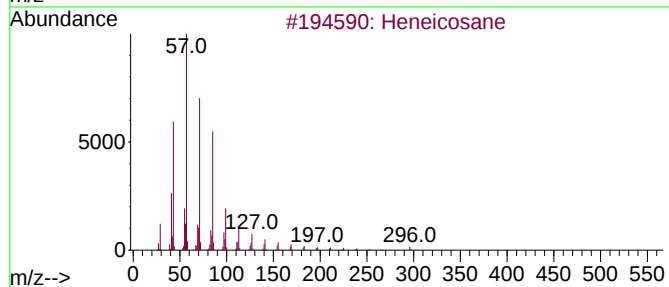
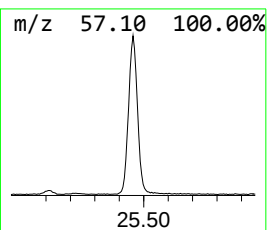
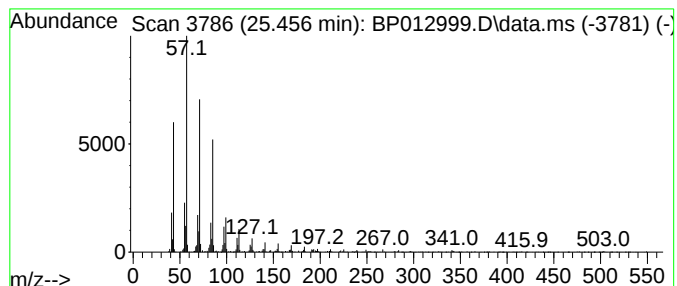
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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
25.456	3.25 ng/ul	1695660	Perylene-d12		23.945	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	97
2	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
3	Octacosane		394	C28H58	000630-02-4	91
4	Hexacosane, 1-iodo-		492	C26H53I	1000406-32-1	91
5	Triacontane, 1-iodo-		548	C30H61I	1000406-32-3	91



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	24.574	3.2	ng/u1	1645960	6	23.945	10426700	20.0
(DEL) Alkane: S...	25.456	3.3	ng/u1	1695660	6	23.945	10426700	20.0