LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
Data File : BP017285.D
Acq On : 22 Sep 2023 14:03
Operator : MA/JU
Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BBBK3

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-BP092023.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP017285.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.258	24	29	34	rBV	579008	750234	14.57%	1.522%
2	3.305	34	37	41	rVB	40207	51838	1.01%	0.105%
3	3.740	108	111	122	rBV	211842	328663	6.38%	0.667%
4	3.928	139	143	149	rVB	35823	49054	0.95%	0.100%
5	4.040	160	162	168	rVB3	9381	12013	0.23%	0.024%
6	4.187	184	187	194	rVB7	16679	24902	0.48%	0.051%
7	4.434	227	229	235	rVB5	12291	16461	0.32%	0.033%
8	4.569	246	252	264	rVB	663432	927408	18.01%	1.882%
9	4.987	319	323	326	rVB3	6840	8692	0.17%	0.018%
10	5.569	419	422	425	rVB5	5210	6835	0.13%	0.014%
11	5.734	445	450	453	rBV7	3452	6430	0.12%	0.013%
12	6.340	549	553	557	rBV3	11637	16336	0.32%	0.033%
13	7.093	675	681	690	rBV	186768	293044	5.69%	0.595%
14	7.275	706	712	720	rBV	631367	957245	18.59%	1.942%
15	7.457	736	743	755	rBV	666731	1032932	20.06%	2.096%
16	7.546	755	758	760	rBV4	4412	5675	0.11%	0.012%
17	7.752	788	793	796	rBV7	3403	5665	0.11%	0.011%
18	7.934	816	824	835	rBV	582857	905798	17.60%	1.838%
19	8.652	939	946	954	rBV	495958	807070	15.68%	1.637%
20	9.134	1020	1028	1036	rBV	722932	1146078	22.26%	2.325%
21	9.334	1059	1062	1068	rVB5	8256	14533	0.28%	0.029%
22	9.846	1142	1149	1160	rBV	579242	941816	18.29%	1.911%
23	10.034	1171	1181	1187	rBV4	19412	36628	0.71%	0.074%
24	10.304	1220	1227	1232	rBV2	15271	30142	0.59%	0.061%
25	10.375	1232	1239	1256	rVV	842773	1409046	27.37%	2.859%
26	10.757	1295	1304	1312	rBV	755558	1238314	24.05%	2.512%
27	10.922	1323	1332	1346	rBV	569869	1014111	19.70%	2.058%
28	11.040	1349	1352	1357	rVB7	3439	6464	0.13%	0.013%
29	11.451	1418	1422	1426	rBV7	3611	6937	0.13%	0.014%
30	11.522	1429	1434	1437	rBV7	2989	5389	0.10%	0.011%
31	11.693	1460	1463	1468	rVB7	3574	5910	0.11%	0.012%
32	11.787	1474	1479	1482	rVV7	3700	6987	0.14%	0.014%
33	11.828	1484	1486	1492	rVB7	4913	6830	0.13%	0.014%
34	12.240	1549	1556	1559	rBV7	6571	11505	0.22%	0.023%
35	12.340	1568	1573	1578	rVB	14877	23967	0.47%	0.049%
36	12.840	1652	1658	1664	rVB4	14712	25153	0.49%	0.051%

LSC Area Percent Report

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Data File : BP017285.D
Acq On : 22 Sep 2023 14:03
Operator : MA/JU
Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBBK3

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-BP092023.MA.M
Title : SVOA CALIBRATION

37	13.175	1710	1715	1724	rVB6	11214	20766	0.40%	0.042%
38	13.345	1740	1744	1748	rVB7	4275	6971	0.14%	0.014%
39	13.387	1748	1751	1754	rBV5	4085	5209	0.10%	0.011%
40	13.440	1754	1760	1766	rVB	42075	69472	1.35%	0.141%
41	13.504	1766	1771	1772	rBV5	4028	5650	0.11%	0.011%
42	13.551	1776	1779	1783	rBV6	6063	8120	0.16%	0.016%
43	14.010	1850	1857	1870	rVV	1696331	2234242	43.40%	4.533%
44	14.292	1898	1905	1916	rBV	1805580	2647042	51.42%	5.371%
45	14.522	1940	1944	1947	rVB6	5708	6629	0.13%	0.013%
46	14.598	1947	1957	1965	rBV	1034883	1540239	29.92%	3.125%
47	14.781	1985	1988	1992	rBV3	19356	25415	0.49%	0.052%
48	14.834	1992	1997	2010	rVV	113804	217239	4.22%	0.441%
49	14.975	2017	2021	2028	rVV5	19561	32110	0.62%	0.065%
50	15.039	2028	2032	2038	rVB9	5493	9843	0.19%	0.020%
51	15.216	2057	2062	2067	rBV8	6379	12248	0.24%	0.025%
52	15.381	2086	2090	2093	rBV6	7390	9387	0.18%	0.019%
53	15.592	2118	2126	2139	rBV	2483563	3448277	66.98%	6.996%
54	15.722	2142	2148	2154	rBV	650235	866487	16.83%	1.758%
55	15.828	2163	2166	2170	rVV5	10427	14151	0.27%	0.029%
56	15.869	2170	2173	2179	rVB8	9325	16689	0.32%	0.034%
57	16.128	2214	2217	2223	rVB6	8606	11455	0.22%	0.023%
58	16.257	2234	2239	2245	rBV2	28804	46459	0.90%	0.094%
59	16.516	2279	2283	2284	rBV4	4371	5966	0.12%	0.012%
60	16.716	2314	2317	2326	rVB4	13488	21420	0.42%	0.043%
61	17.092	2377	2381	2385	rVB5	14528	23560	0.46%	0.048%
62	17.369	2421	2428	2435	rBV	1353872	1907789	37.06%	3.871%
63	17.469	2438	2445	2453	rBV	2875085	4027154	78.23%	8.171%
64	17.992	2530	2534	2538	rBV7	9256	14673	0.29%	0.030%
65	18.210	2566	2571	2583	rBV	446853	705449	13.70%	1.431%
66	19.080	2717	2719	2722	rVB2	19278	19518	0.38%	0.040%
67	19.280	2750	2753	2755	rBV2	35372	44332	0.86%	0.090%
68	19.310	2756	2758	2760	rVV3	41839	46021	0.89%	0.093%
69	19.345	2760	2764	2769	rVB2	61351	115089	2.24%	0.234%
70	19.445	2777	2781	2794	rBV	177216	276918	5.38%	0.562%
71	19.645	2813	2815	2820	rVB	48799	51299	1.00%	0.104%
72	19.710	2820	2826	2832	rBV	4062148	5128908	99.63%	10.406%
73	20.169	2901	2904	2910	rVB	173028	191799	3.73%	0.389%
74	20.657	2984	2987	2990	rBV	261936	273144	5.31%	0.554%
75	21.116	3062	3065	3067	rBV	327453	369182	7.17%	0.749%
76	21.145	3067	3070	3078	rVB	454954	685214	13.31%	1.390%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
Data File : BP017285.D
Acq On : 22 Sep 2023 14:03
Operator : MA/JU
Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBBK3

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

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Title : SVOA CALIBRATION

77	21.474	3121	3126	3131	rVB	1673910	2156226	41.88%	4.375%
78	21.557	3137	3140	3144	rVB	303800	328213	6.38%	0.666%
79	22.033	3217	3221	3226	rBV	314336	399317	7.76%	0.810%
80	22.545	3305	3308	3312	rVB	265417	338609	6.58%	0.687%
81	23.121	3403	3406	3414	rVB	347068	517565	10.05%	1.050%
82	23.827	3520	3526	3535	rVB2	2573998	5148011	100.00%	10.445%
83	23.992	3548	3554	3563	rVB	1197357	2485869	48.29%	5.044%
84	24.539	3644	3647	3655	rVB	332547	615361	11.95%	1.249%

Sum of corrected areas: 49286811

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
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Acq On : 22 Sep 2023 14:03
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Sample : 04410-05
Misc :
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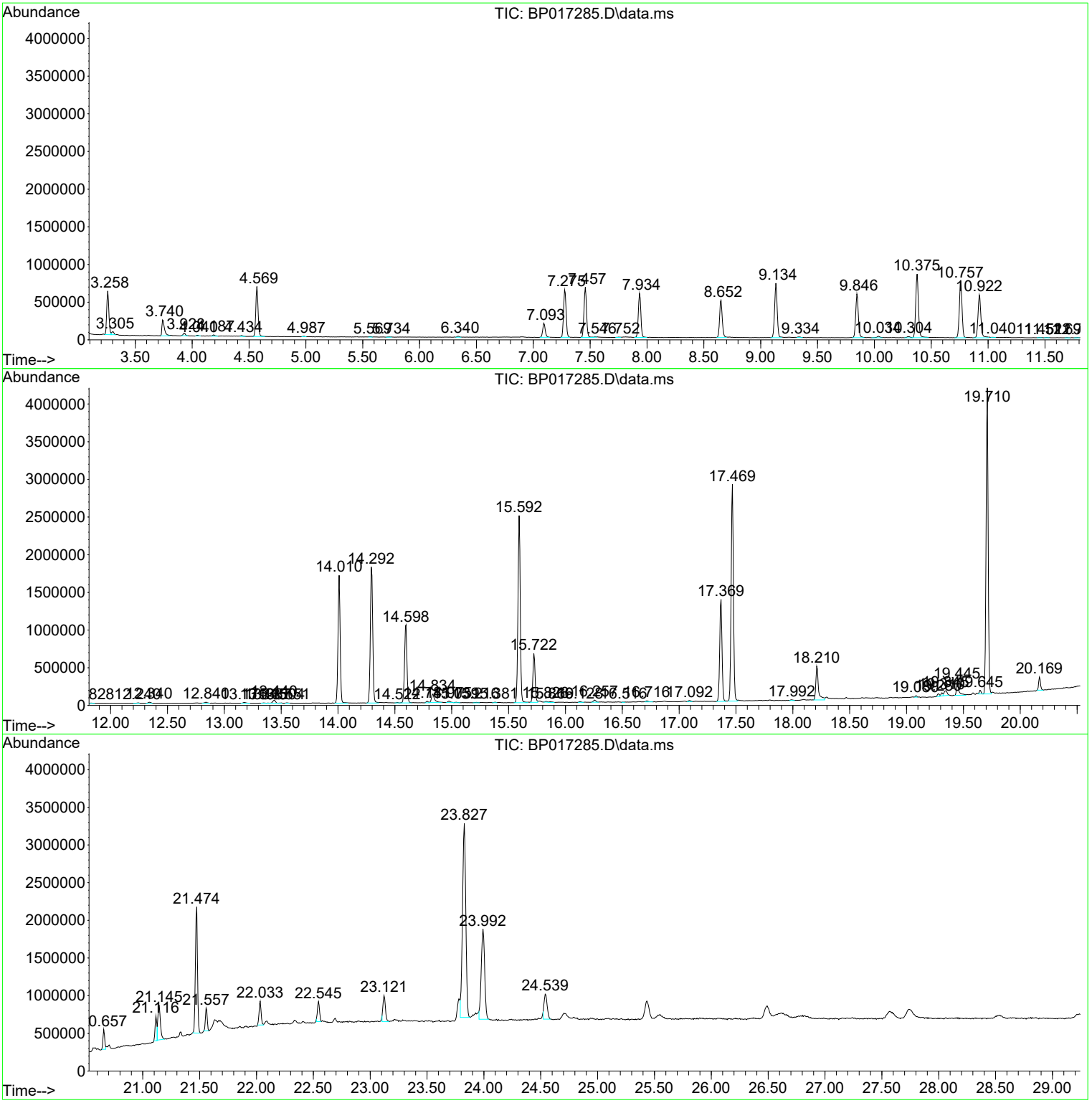
Instrument :
BNA_P
ClientSampleId :
BBBK3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
Data File : BP017285.D
Acq On : 22 Sep 2023 14:03
Operator : MA/JU
Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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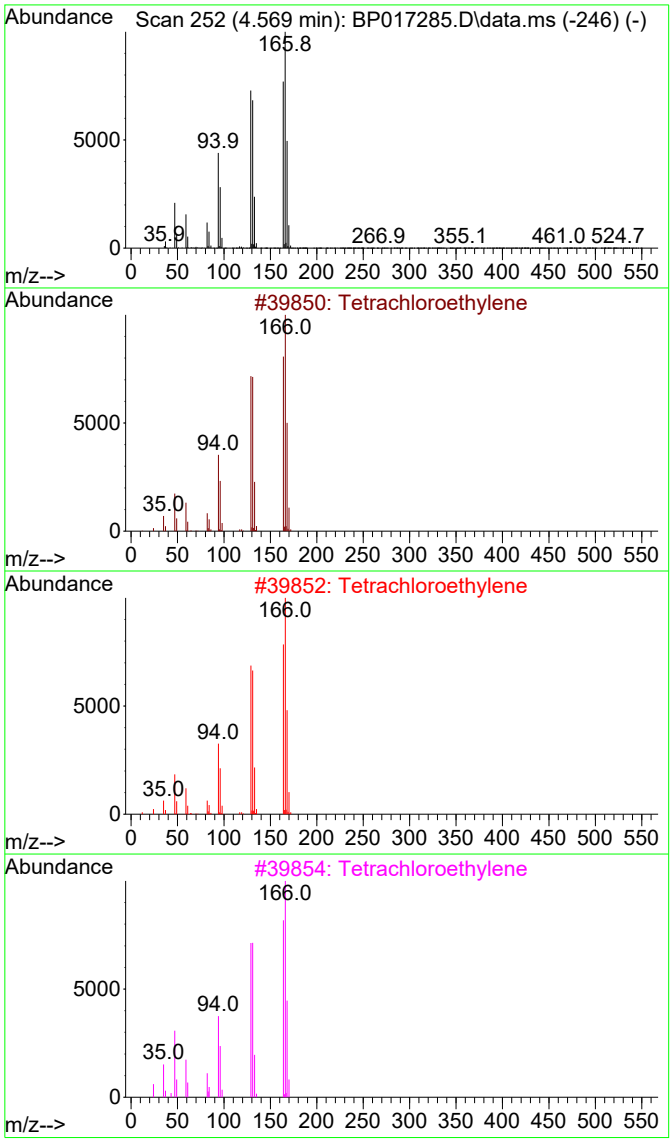
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

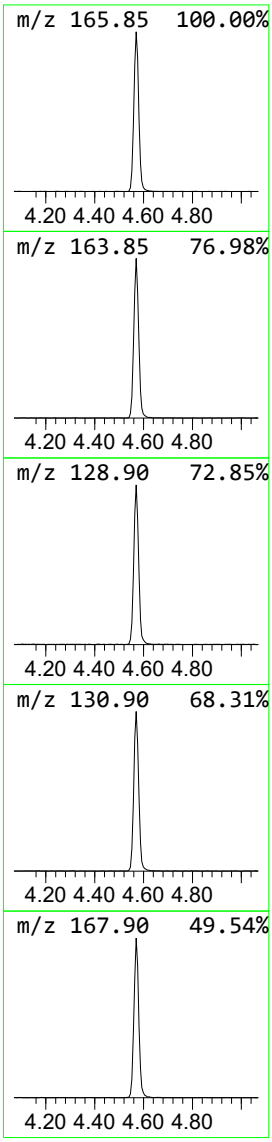
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Tetrachloroethylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.569	20.48 ng/ul	927408	1,4-Dichlorobenzene-d4	7.934

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetrachloroethylene	164	C2Cl4	000127-18-4	98
2	Tetrachloroethylene	164	C2Cl4	000127-18-4	98
3	Tetrachloroethylene	164	C2Cl4	000127-18-4	97
4	Tetrachloroethylene	164	C2Cl4	000127-18-4	96
5	Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HCl2FN2	002927-71-1	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
Data File : BP017285.D
Acq On : 22 Sep 2023 14:03
Operator : MA/JU
Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBBK3

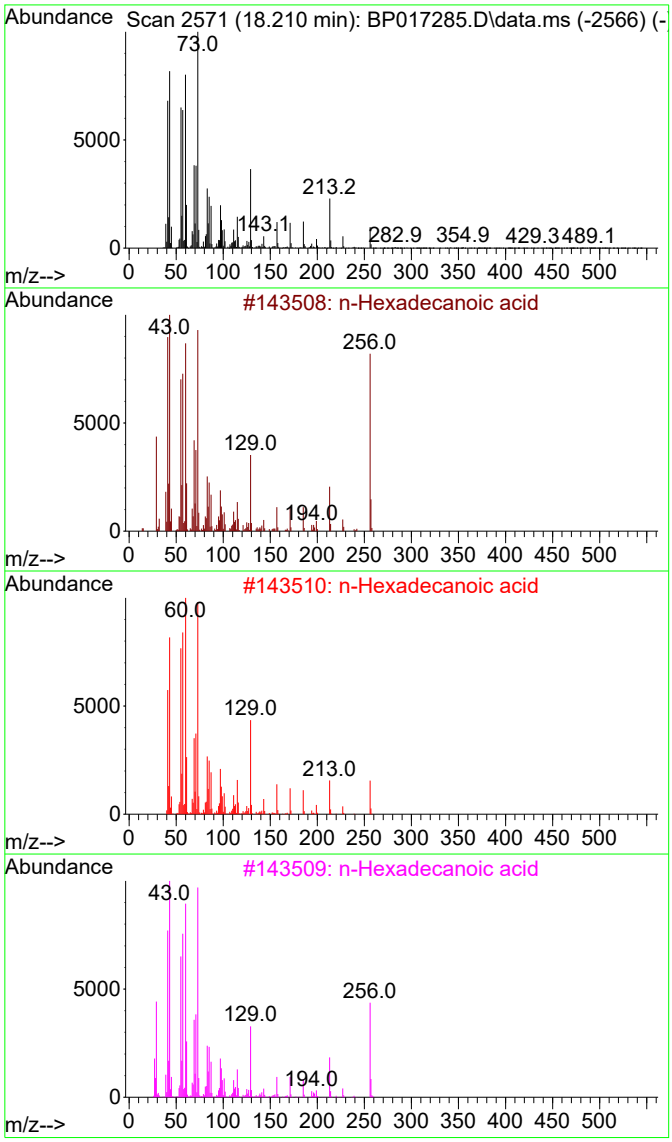
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

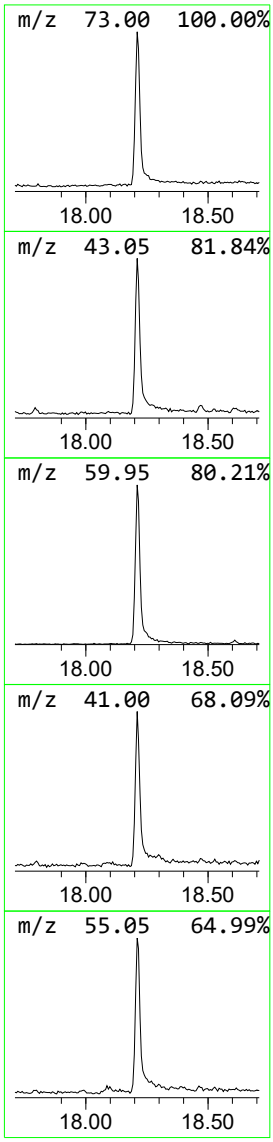
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.210	7.40 ng/ul	705449	Phenanthrene-d10	17.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
Data File : BP017285.D
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Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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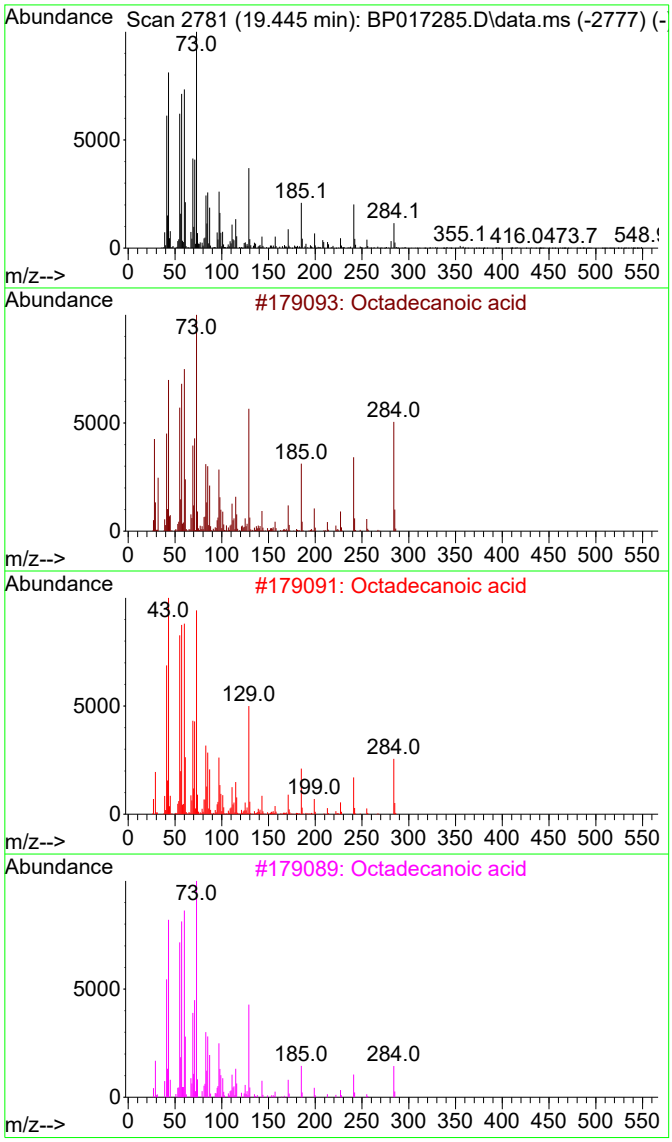
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

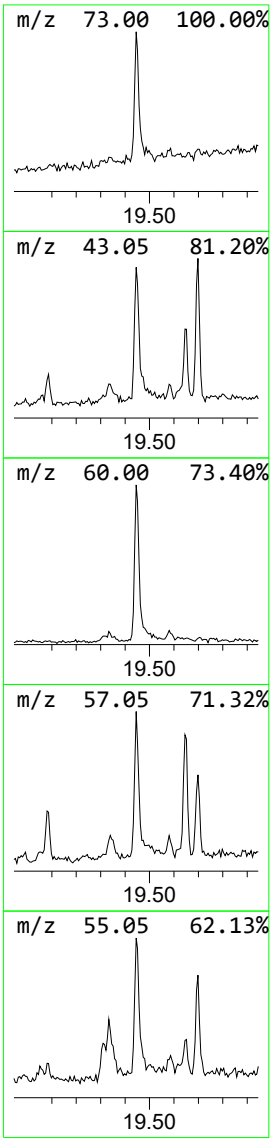
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.445	2.57 ng/ul	276918	Chrysene-d12	21.474

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	98
4	Octadecanoic acid	284	C18H36O2	000057-11-4	95
5	Octadecanoic acid	284	C18H36O2	000057-11-4	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
Data File : BP017285.D
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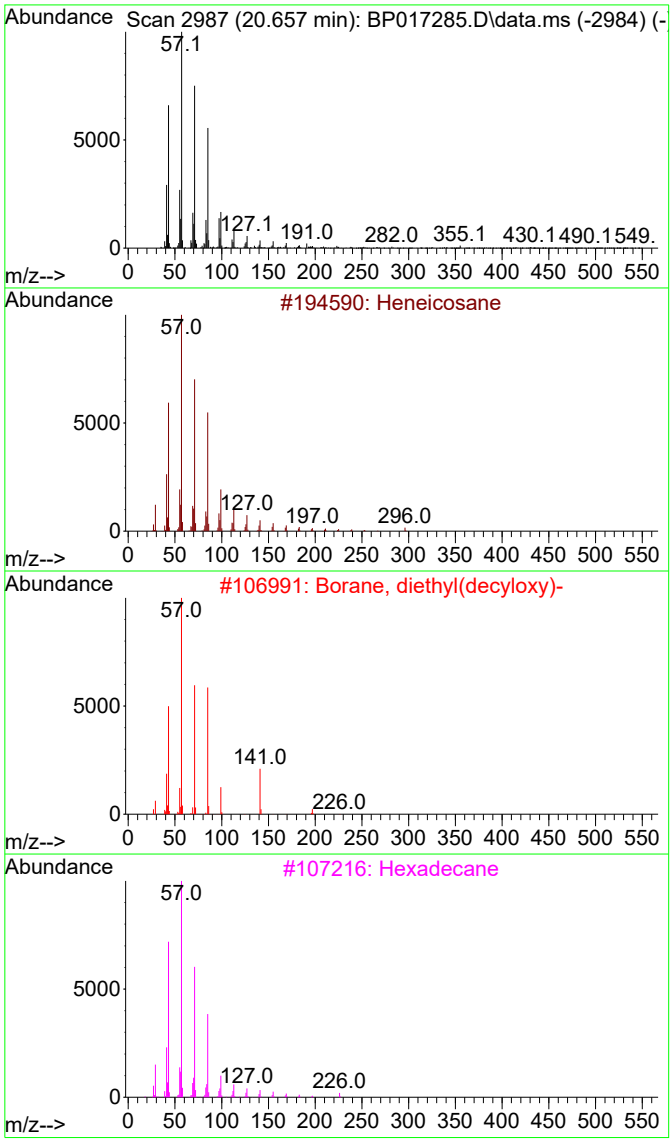
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

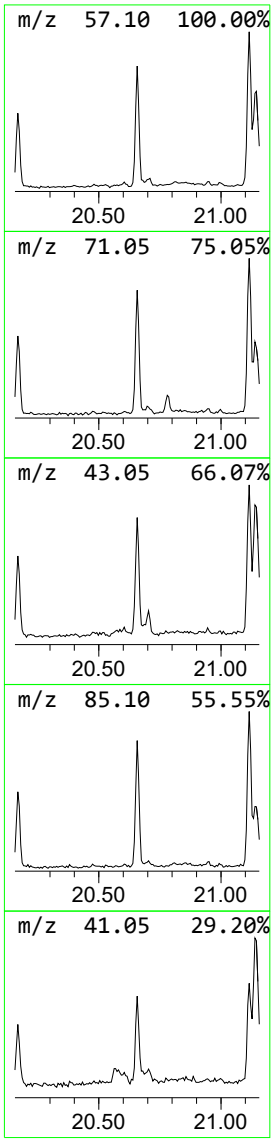
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.657	2.53 ng/ul	273144	Chrysene-d12	21.474

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	90
2	Borane, diethyl(decyloxy)-	226	C14H31BO	1000152-34-3	83
3	Hexadecane	226	C16H34	000544-76-3	80
4	Octacosane	394	C28H58	000630-02-4	74
5	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	74





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
Data File : BP017285.D
Acq On : 22 Sep 2023 14:03
Operator : MA/JU
Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBBK3

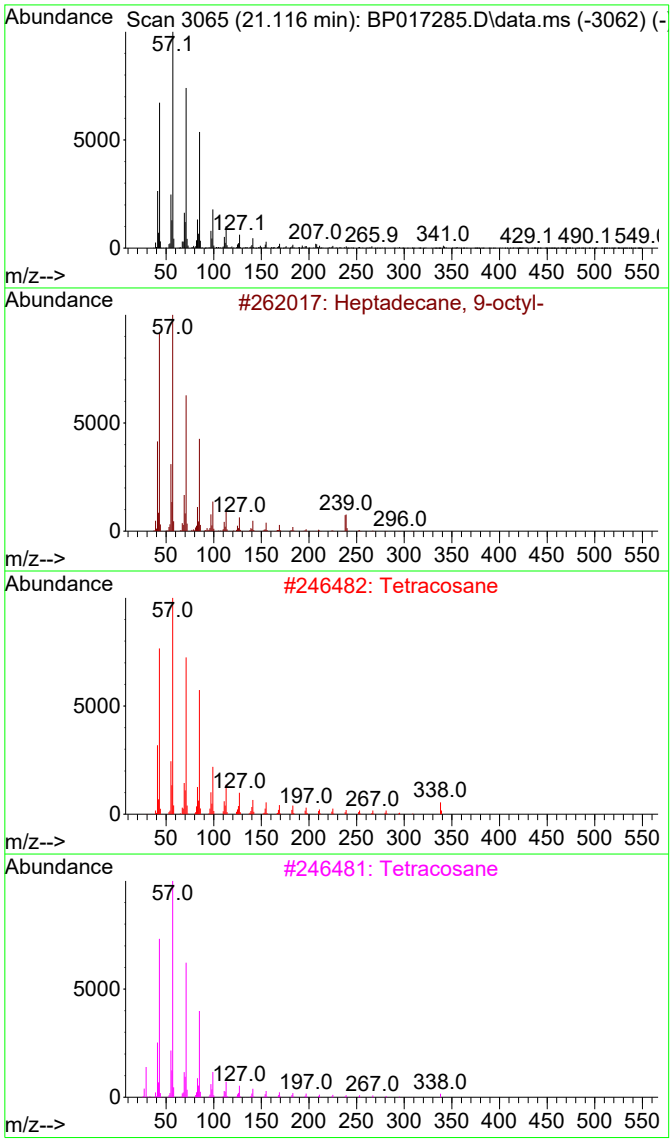
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

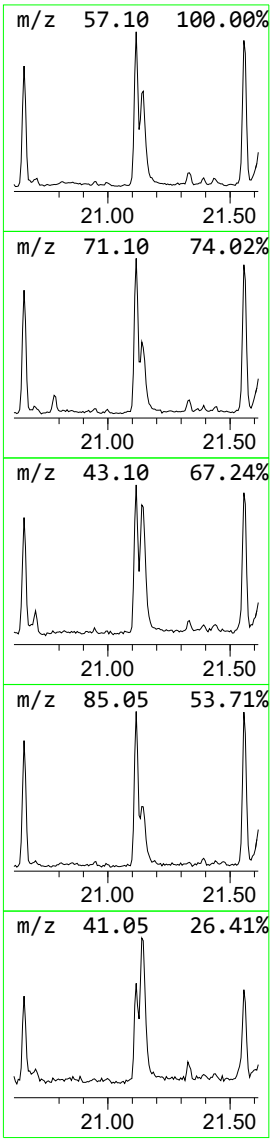
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.116	3.42 ng/ul	369182	Chrysene-d12	21.474

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
2	Tetracosane	338	C24H50	000646-31-1	91
3	Tetracosane	338	C24H50	000646-31-1	91
4	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
Data File : BP017285.D
Acq On : 22 Sep 2023 14:03
Operator : MA/JU
Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBBK3

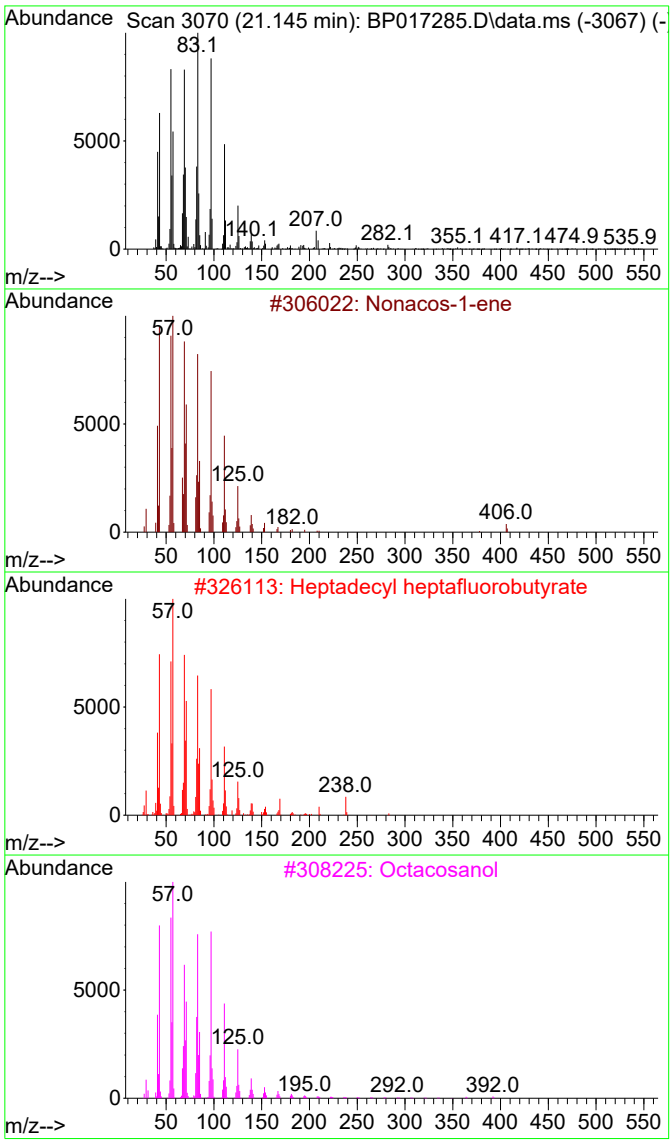
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

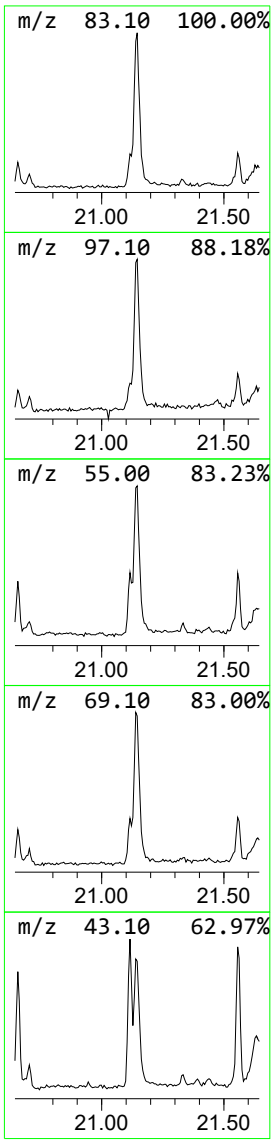
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.145	6.36 ng/ul	685214	Chrysene-d12	21.474

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonacos-1-ene	406	C29H58	018835-35-3	90
2	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	80
3	Octacosanol	410	C28H58O	000557-61-9	64
4	Pentacos-1-ene	350	C25H50	016980-85-1	64
5	1-Octadecanol	270	C18H38O	000112-92-5	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
Data File : BP017285.D
Acq On : 22 Sep 2023 14:03
Operator : MA/JU
Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBBK3

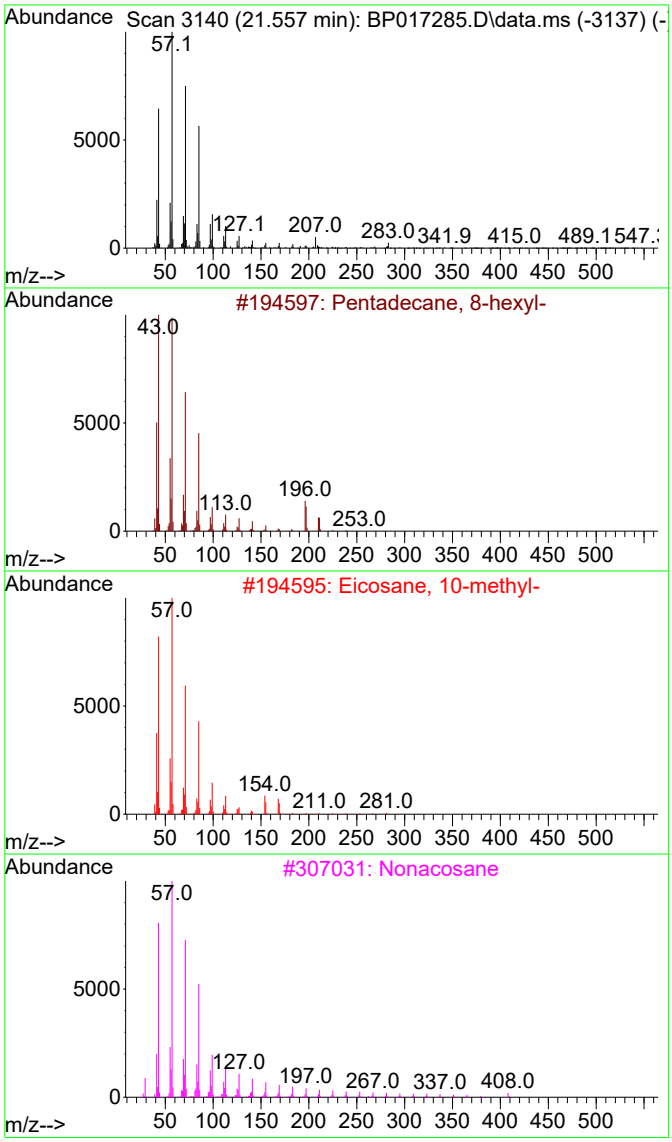
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

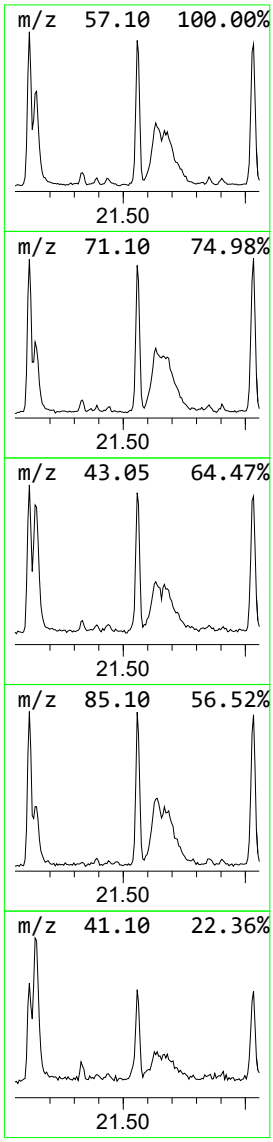
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.557	3.04 ng/ul	328213	Chrysene-d12	21.474

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
2	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
3	Nonacosane	408	C29H60	000630-03-5	91
4	Hexacosane	366	C26H54	000630-01-3	91
5	Tetracosane	338	C24H50	000646-31-1	91





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Operator : MA/JU
Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBBK3

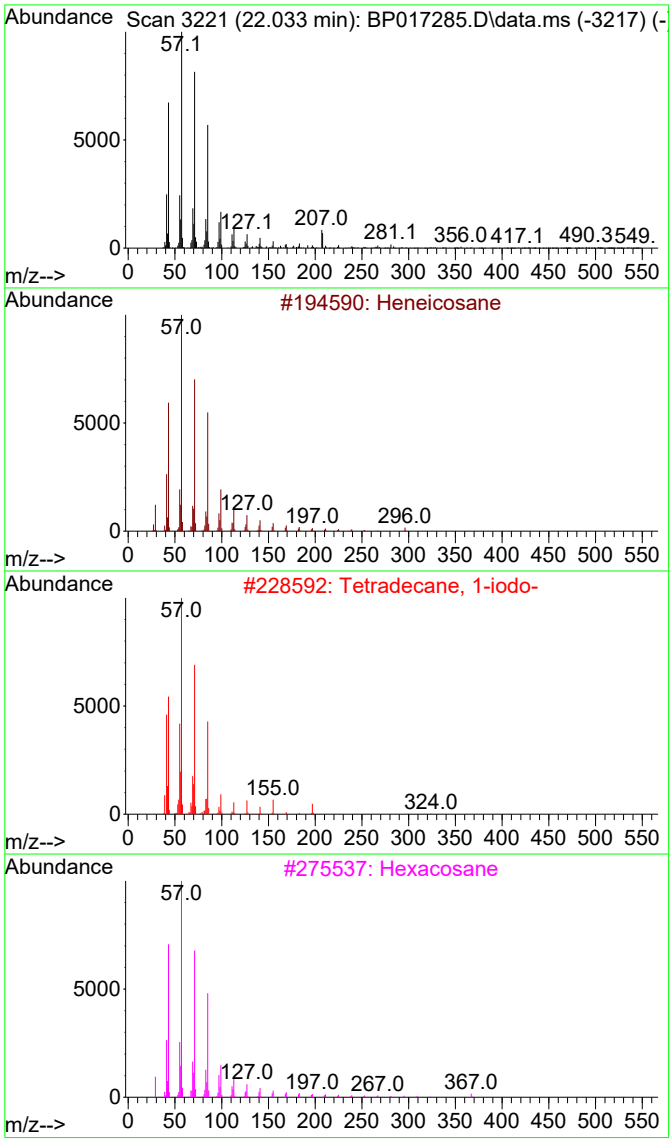
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

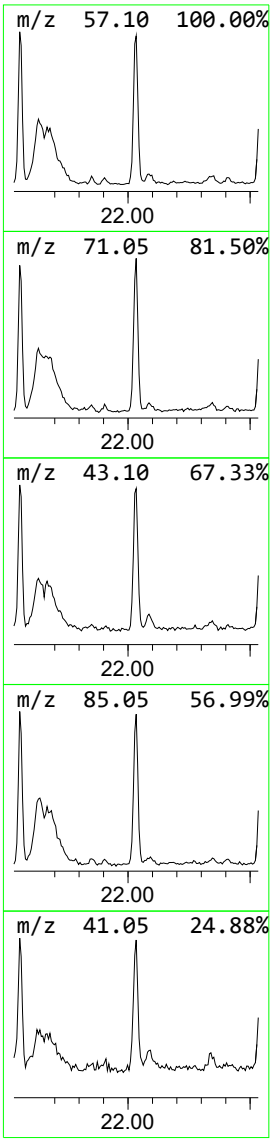
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.033	3.70 ng/ul	399317	Chrysene-d12	21.474

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	90
2	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	90
3	Hexacosane	366	C26H54	000630-01-3	90
4	Tetracosane	338	C24H50	000646-31-1	87
5	Octacosane	394	C28H58	000630-02-4	87





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Data File : BP017285.D
Acq On : 22 Sep 2023 14:03
Operator : MA/JU
Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBBK3

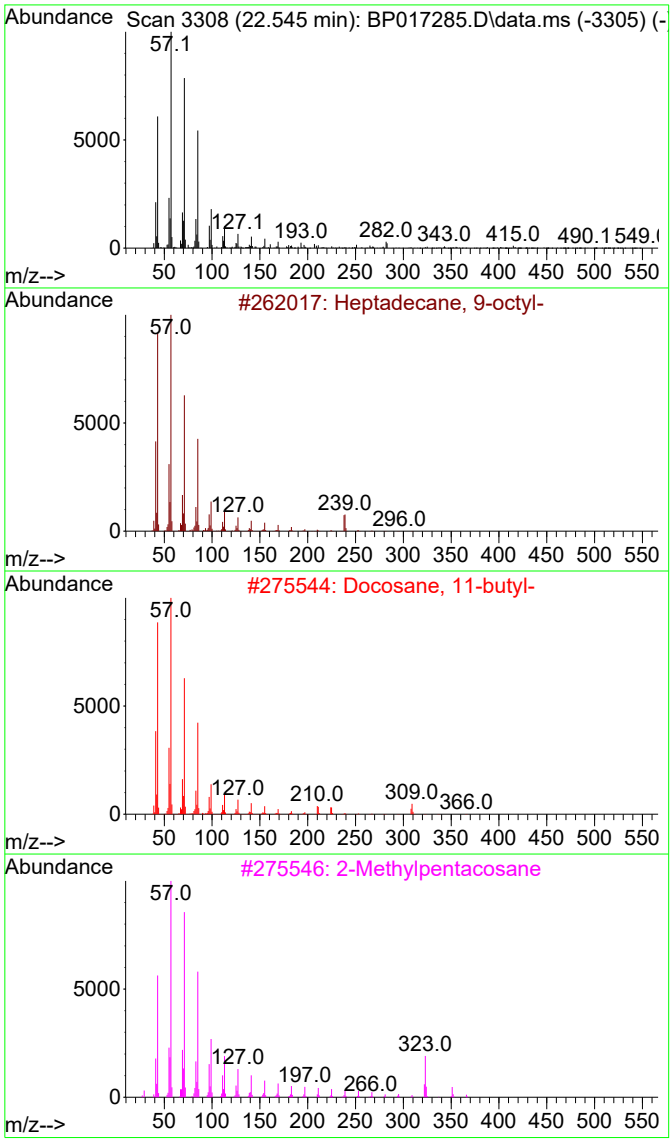
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

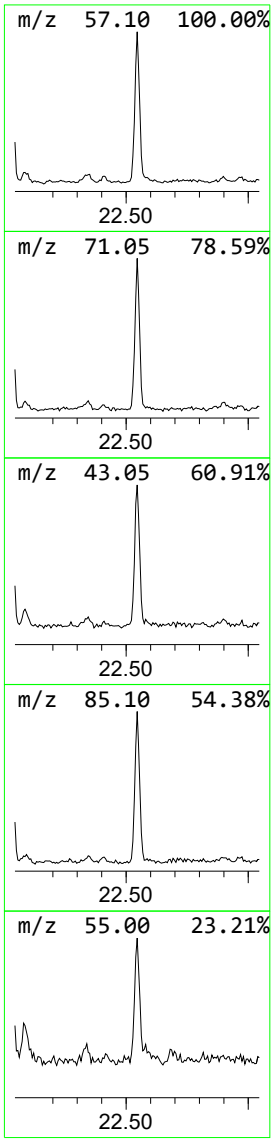
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.545	3.14 ng/ul	338609	Chrysene-d12	21.474

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2	Docosane, 11-butyl-	366	C26H54	013475-76-8	95
3	2-Methylpentacosane	366	C26H54	000629-87-8	94
4	Eicosane	282	C20H42	000112-95-8	94
5	Tetracosane	338	C24H50	000646-31-1	94





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Data File : BP017285.D
Acq On : 22 Sep 2023 14:03
Operator : MA/JU
Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBBK3

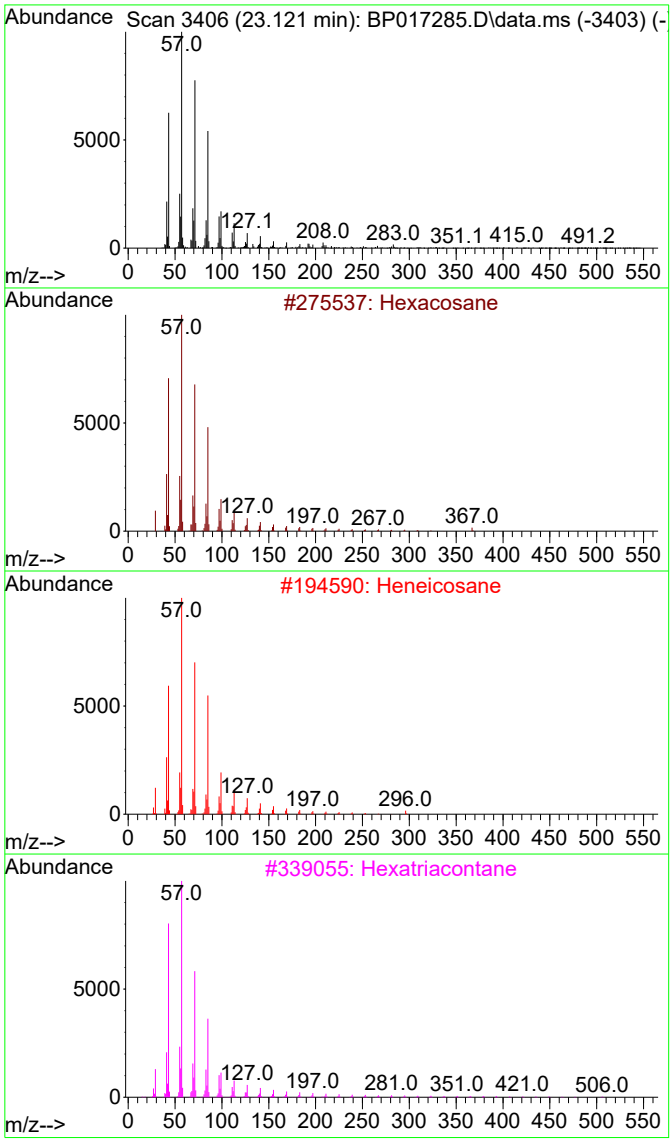
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

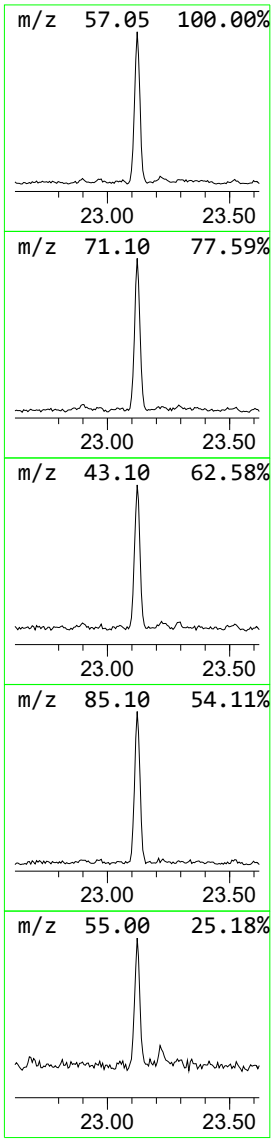
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.121	4.16 ng/ul	517565	Perylene-d12	23.992

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Hexatriacontane	507	C36H74	000630-06-8	91
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
Data File : BP017285.D
Acq On : 22 Sep 2023 14:03
Operator : MA/JU
Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBBK3

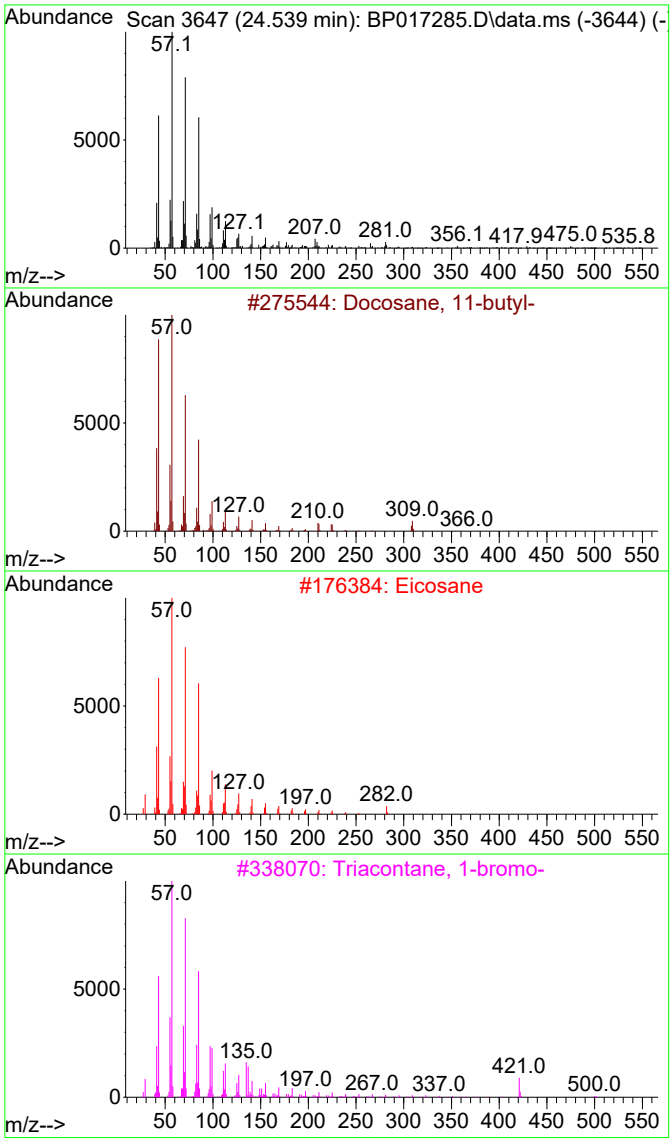
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

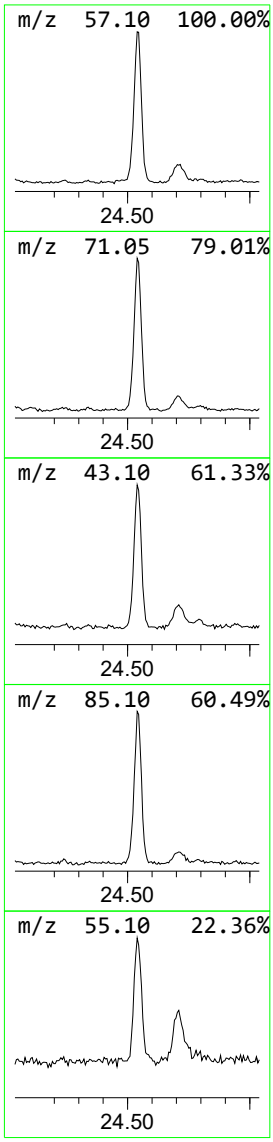
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.539	4.95 ng/ul	615361	Perylene-d12	23.992

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	94
2	Eicosane	282	C20H42	000112-95-8	94
3	Triacontane, 1-bromo-	500	C30H61Br	004209-22-7	93
4	Docosane, 9-butyl-	366	C26H54	055282-14-9	91
5	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
Data File : BP017285.D
Acq On : 22 Sep 2023 14:03
Operator : MA/JU
Sample : 04410-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBBK3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Tetrachloroethy...	4.569	20.5	ng/ul	927408	1	7.934	905798	20.0
n-Hexadecanoic ...	18.210	7.4	ng/ul	705449	4	17.369	1907790	20.0
Octadecanoic acid	19.445	2.6	ng/ul	276918	5	21.474	2156230	20.0
(DEL) Alkane: S...	20.657	2.5	ng/ul	273144	5	21.474	2156230	20.0
(DEL) Alkane: S...	21.116	3.4	ng/ul	369182	5	21.474	2156230	20.0
(DEL) Alkane: S...	21.145	6.4	ng/ul	685214	5	21.474	2156230	20.0
(DEL) Alkane: S...	21.557	3.0	ng/ul	328213	5	21.474	2156230	20.0
(DEL) Alkane: S...	22.033	3.7	ng/ul	399317	5	21.474	2156230	20.0
(DEL) Alkane: S...	22.545	3.1	ng/ul	338609	5	21.474	2156230	20.0
(DEL) Alkane: S...	23.121	4.2	ng/ul	517565	6	23.992	2485870	20.0
(DEL) Alkane: S...	24.539	5.0	ng/ul	615361	6	23.992	2485870	20.0