LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP021007.D
Acq On : 17 Jul 2024 03:25
Operator : MA/JU
Sample : P3178-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A4BT1

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-BP071024.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP021007.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.146	24	27	32	rVB6	9388	14301	0.41%	0.027%
2	3.246	41	44	59	rVB	41020	58346	1.67%	0.111%
3	3.522	87	91	104	rBV	104500	156859	4.50%	0.298%
4	3.658	111	114	115	rBV2	8614	7870	0.23%	0.015%
5	3.675	115	117	126	rVV	45445	75886	2.18%	0.144%
6	3.758	127	131	137	rVV	59992	80513	2.31%	0.153%
7	3.828	141	143	147	rVB2	7115	7113	0.20%	0.014%
8	3.964	162	166	172	rVB	18739	26823	0.77%	0.051%
9	4.334	224	229	232	rBV6	5855	8020	0.23%	0.015%
10	4.381	232	237	244	rVV3	8855	14900	0.43%	0.028%
11	4.469	248	252	255	rVV2	6693	8706	0.25%	0.017%
12	4.511	255	259	264	rVB2	27499	38726	1.11%	0.074%
13	4.664	283	285	291	rBV6	4661	6259	0.18%	0.012%
14	4.740	296	298	302	rBV4	3915	5151	0.15%	0.010%
15	4.864	313	319	326	rBV	54585	82616	2.37%	0.157%
16	4.969	333	337	344	rVB5	8844	14928	0.43%	0.028%
17	5.434	414	416	422	rVB7	3939	5201	0.15%	0.010%
18	5.716	463	464	472	rVB7	2556	4679	0.13%	0.009%
19	5.828	474	483	490	rBV3	11964	29454	0.84%	0.056%
20	6.769	639	643	646	rBV5	7521	11678	0.33%	0.022%
21	6.799	646	648	654	rVB	7409	10765	0.31%	0.020%
22	6.893	657	664	679	rVV	587834	1097939	31.49%	2.084%
23	7.046	683	690	701	rVB	524484	841042	24.12%	1.597%
24	7.240	716	723	731	rBV	711986	1142754	32.77%	2.169%
25	7.322	733	737	747	rVV	28729	67017	1.92%	0.127%
26	7.434	751	756	759	rVV4	17326	30100	0.86%	0.057%
27	7.463	759	761	767	rVB4	13471	18294	0.52%	0.035%
28	7.687	791	799	809	rBV	799181	1398366	40.10%	2.655%
29	7.787	809	816	826	rVB3	45663	104930	3.01%	0.199%
30	7.893	832	834	839	rVB6	3619	5116	0.15%	0.010%
31	7.958	839	845	849	rBV8	6005	12132	0.35%	0.023%
32	8.210	882	888	895	rVB3	12891	23669	0.68%	0.045%
33	8.410	914	922	937	rBV	757022	1416801	40.63%	2.690%
34	8.522	937	941	945	rVB3	12470	22179	0.64%	0.042%
35	8.852	989	997	1008	rBV	615887	1084421	31.10%	2.059%
36	8.981	1012	1019	1020	rBV7	4660	8425	0.24%	0.016%

LSC Area Percent Report

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Operator : MA/JU
Sample : P3178-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT1

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-BP071024.MA.M
Title : SVOA CALIBRATION

37	9.199	1050	1056	1061	rBV2	11712	19610	0.56%	0.037%
38	9.399	1083	1090	1104	rVB2	71339	130226	3.73%	0.247%
39	9.552	1109	1116	1130	rBV	528134	927320	26.59%	1.760%
40	9.657	1130	1134	1139	rVB6	6411	11562	0.33%	0.022%
41	9.804	1152	1159	1163	rBV10	6105	12819	0.37%	0.024%
42	10.099	1200	1209	1226	rBV	701291	1423145	40.81%	2.702%
43	10.316	1243	1246	1247	rBV3	5833	5082	0.15%	0.010%
44	10.457	1262	1270	1277	rBV	1215133	1894145	54.32%	3.596%
45	10.604	1289	1295	1306	rBV	326735	700296	20.08%	1.329%
46	10.869	1334	1340	1349	rVB	18304	39561	1.13%	0.075%
47	10.987	1354	1360	1368	rBV5	26798	63000	1.81%	0.120%
48	11.199	1388	1396	1416	rBV2	161697	426548	12.23%	0.810%
49	11.446	1433	1438	1439	rBV5	4220	5145	0.15%	0.010%
50	11.546	1452	1455	1462	rVB3	10936	17218	0.49%	0.033%
51	11.669	1473	1476	1480	rVB6	4271	6374	0.18%	0.012%
52	11.857	1504	1508	1514	rVB9	4143	6379	0.18%	0.012%
53	12.040	1532	1539	1545	rVB3	11675	22263	0.64%	0.042%
54	12.769	1657	1663	1665	rBV7	2494	4854	0.14%	0.009%
55	12.910	1679	1687	1692	rBV7	5146	11107	0.32%	0.021%
56	13.028	1703	1707	1710	rBV6	3382	5458	0.16%	0.010%
57	13.104	1717	1720	1726	rBV4	7001	9106	0.26%	0.017%
58	13.257	1742	1746	1757	rVB4	8794	17899	0.51%	0.034%
59	13.598	1801	1804	1806	rBV4	4416	4910	0.14%	0.009%
60	13.716	1815	1824	1838	rBV	1165991	1942218	55.70%	3.687%
61	13.816	1839	1841	1847	rVB7	8023	11956	0.34%	0.023%
62	13.998	1860	1872	1884	rBV2	1503035	2427847	69.63%	4.609%
63	14.134	1890	1895	1899	rVB7	8877	16800	0.48%	0.032%
64	14.192	1899	1905	1909	rBV9	6086	11069	0.32%	0.021%
65	14.304	1917	1924	1932	rBV	1507503	2314763	66.38%	4.394%
66	14.369	1932	1935	1941	rVB4	18180	31315	0.90%	0.059%
67	14.534	1953	1963	1966	rBV3	82264	178356	5.11%	0.339%
68	14.581	1966	1971	1981	rVV	268652	589472	16.91%	1.119%
69	14.716	1990	1994	1999	rBV6	14194	28406	0.81%	0.054%
70	15.098	2056	2059	2065	rVB6	8174	16634	0.48%	0.032%
71	15.157	2065	2069	2073	rBV6	7315	12967	0.37%	0.025%
72	15.316	2090	2096	2104	rBV	1942572	2900953	83.19%	5.507%
73	15.475	2116	2123	2134	rBV	423743	737907	21.16%	1.401%
74	15.904	2191	2196	2202	rBV9	15176	35009	1.00%	0.066%
75	16.063	2215	2223	2230	rBV2	63848	135433	3.88%	0.257%
76	16.504	2295	2298	2310	rVB10	27436	68416	1.96%	0.130%

LSC Area Percent Report

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Sample : P3178-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Title : SVOA CALIBRATION

77	17.116	2396	2402	2407	rBV	1959909	2707863	77.66%	5.141%
78	17.157	2407	2409	2413	rVV	165445	199446	5.72%	0.379%
79	17.216	2413	2419	2430	rVV	1856619	3071972	88.10%	5.832%
80	17.422	2450	2454	2457	rBV5	12912	23640	0.68%	0.045%
81	17.816	2518	2521	2524	rBV2	31283	49043	1.41%	0.093%
82	18.057	2554	2562	2575	rBV2	554594	1136733	32.60%	2.158%
83	19.004	2721	2723	2731	rVB3	75461	95735	2.75%	0.182%
84	19.257	2758	2766	2773	rBV	419636	713672	20.47%	1.355%
85	19.380	2783	2787	2796	rBV	301501	554082	15.89%	1.052%
86	19.604	2819	2825	2833	rBV2	2026609	3486946	100.00%	6.620%
87	20.257	2934	2936	2938	rBV2	70297	54324	1.56%	0.103%
88	20.492	2971	2976	2978	rBV2	373756	610102	17.50%	1.158%
89	21.216	3094	3099	3106	rVB	745804	1264397	36.26%	2.400%
90	21.557	3150	3157	3163	rBV2	1640813	3387314	97.14%	6.431%
91	22.392	3295	3299	3303	rBV2	248005	451066	12.94%	0.856%
92	22.427	3303	3305	3314	rVB2	286690	516155	14.80%	0.980%
93	23.945	3558	3563	3570	rVB	910037	1920242	55.07%	3.645%
94	24.621	3672	3678	3687	rVB2	934955	2302614	66.04%	4.371%
95	24.851	3709	3717	3725	rBV	1207383	3253385	93.30%	6.176%
96	26.080	3920	3926	3940	rVB2	561650	1710696	49.06%	3.248%

Sum of corrected areas: 52674954

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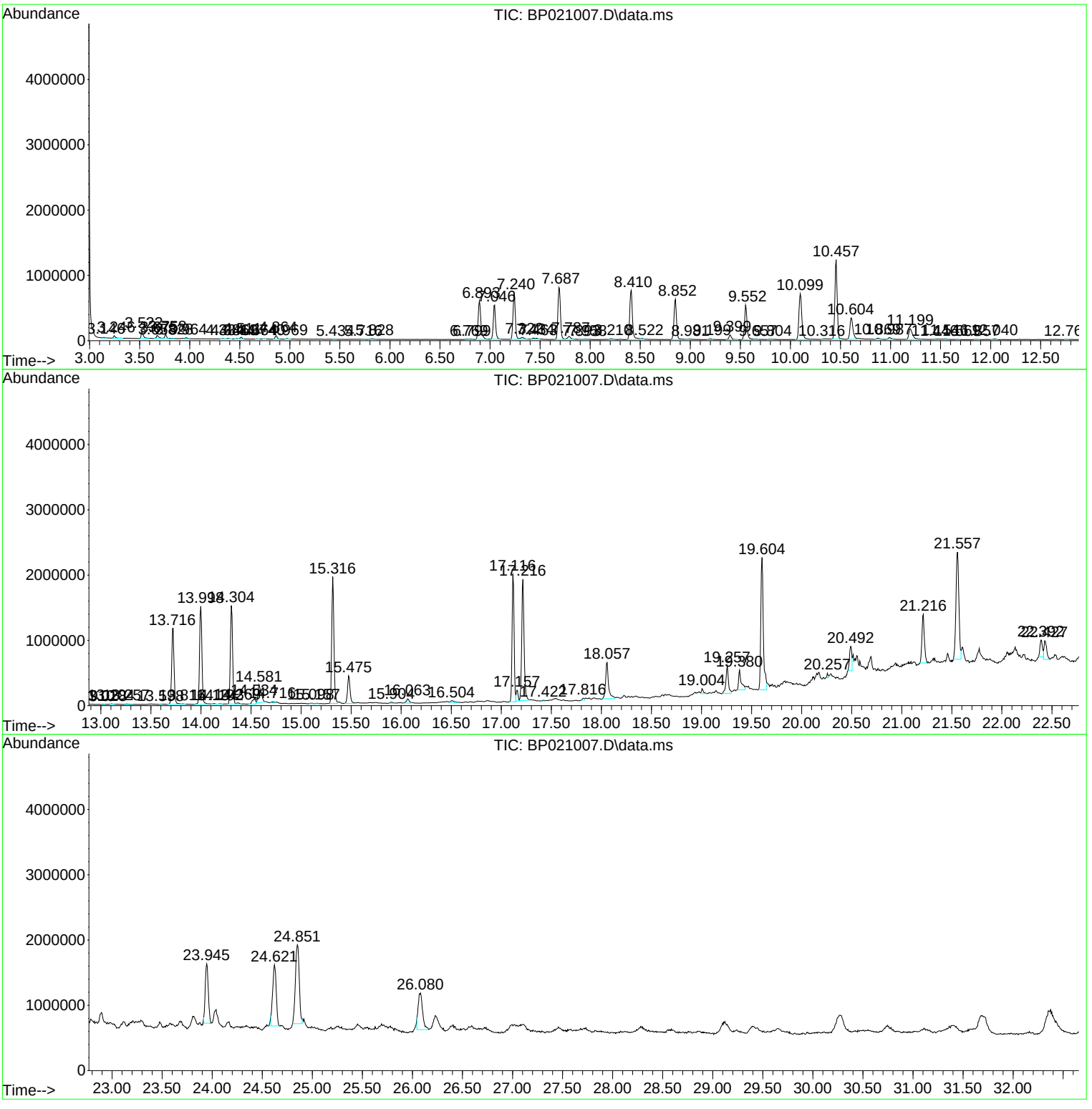
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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\BP071524\
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Instrument :
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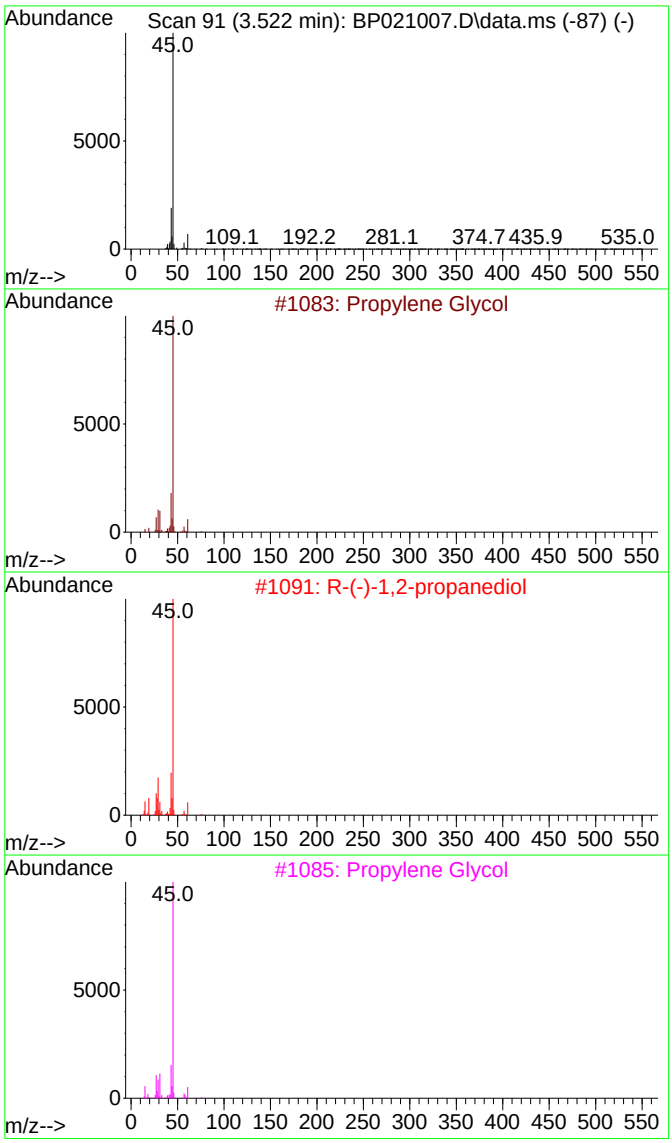
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

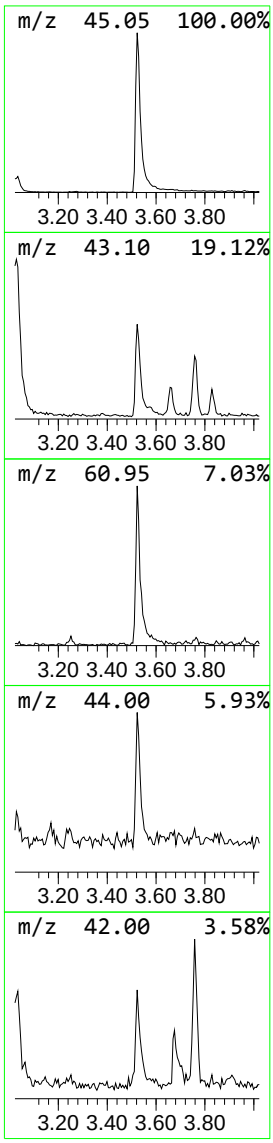
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Propylene Glycol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.522	2.24 ng/ul	156859	1,4-Dichlorobenzene-d4	7.687

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propylene Glycol	76	C3H8O2	000057-55-6	86
2	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	78
3	Propylene Glycol	76	C3H8O2	000057-55-6	56
4	Isopropyl Alcohol	60	C3H8O	000067-63-0	56
5	(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP021007.D
Acq On : 17 Jul 2024 03:25
Operator : MA/JU
Sample : P3178-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

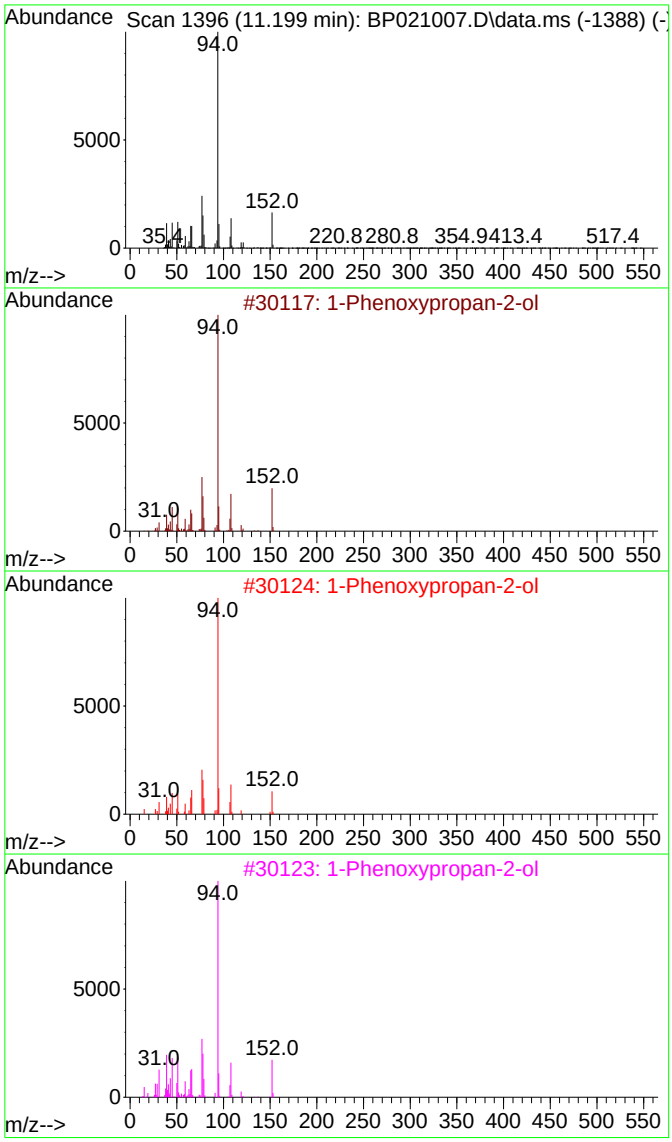
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

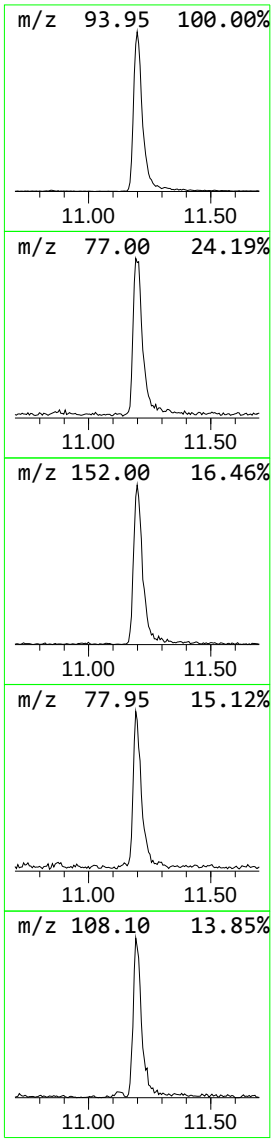
****

Peak Number 2 1-Phenoxypropan-2-ol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.198	4.50 ng/ul	426548	Naphthalene-d8	10.457

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
5	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91





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ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

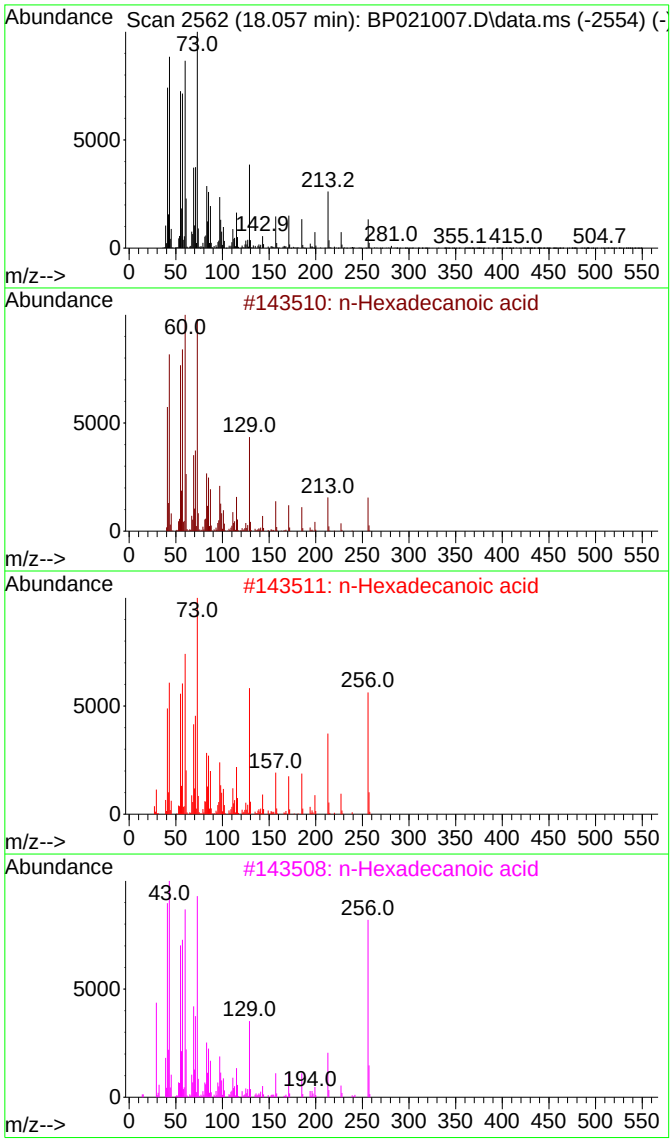
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

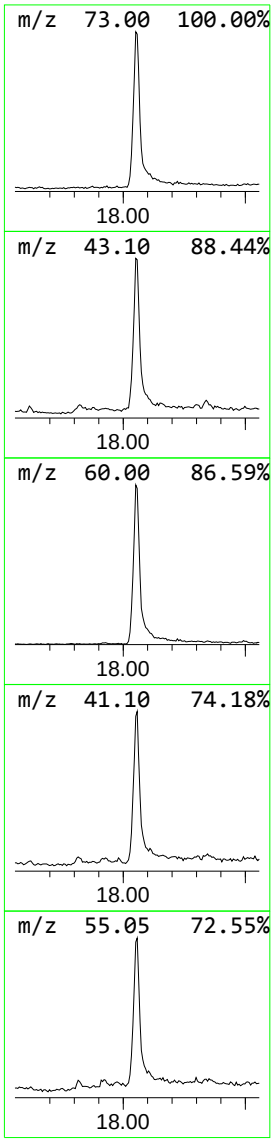
****

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.057	8.40 ng/ul	1136730	Phenanthrene-d10	17.116

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	99
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	Tridecanoic acid	214	C13H26O2	000638-53-9	95





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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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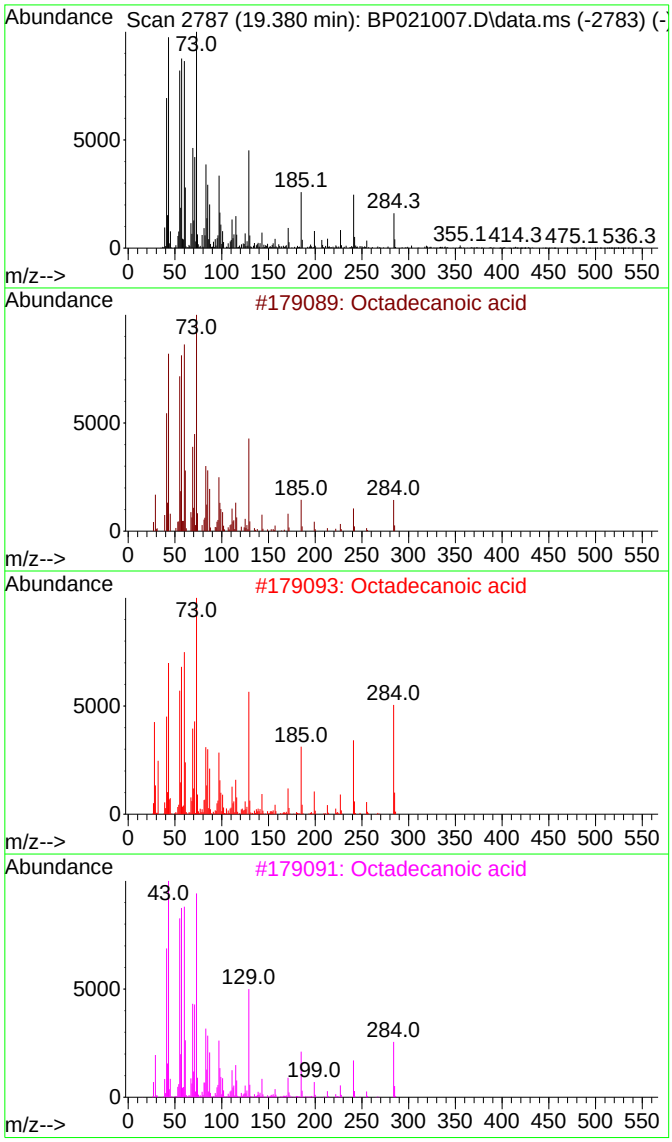
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

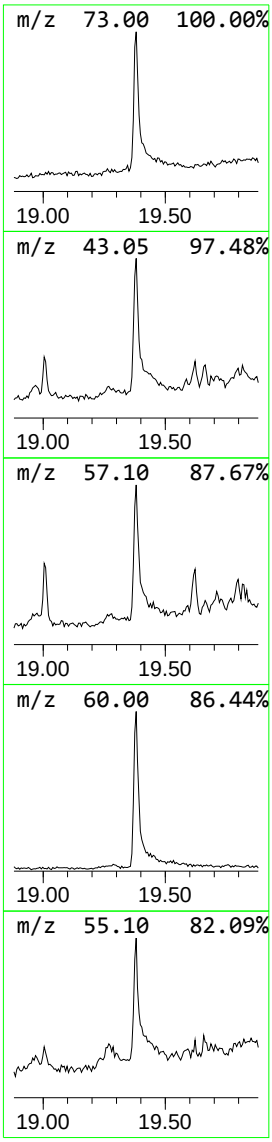
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.380	3.27 ng/ul	554082	Chrysene-d12	21.557

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	99
4	Octadecanoic acid	284	C18H36O2	000057-11-4	98
5	Octadecanoic acid	284	C18H36O2	000057-11-4	94





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Sample : P3178-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT1

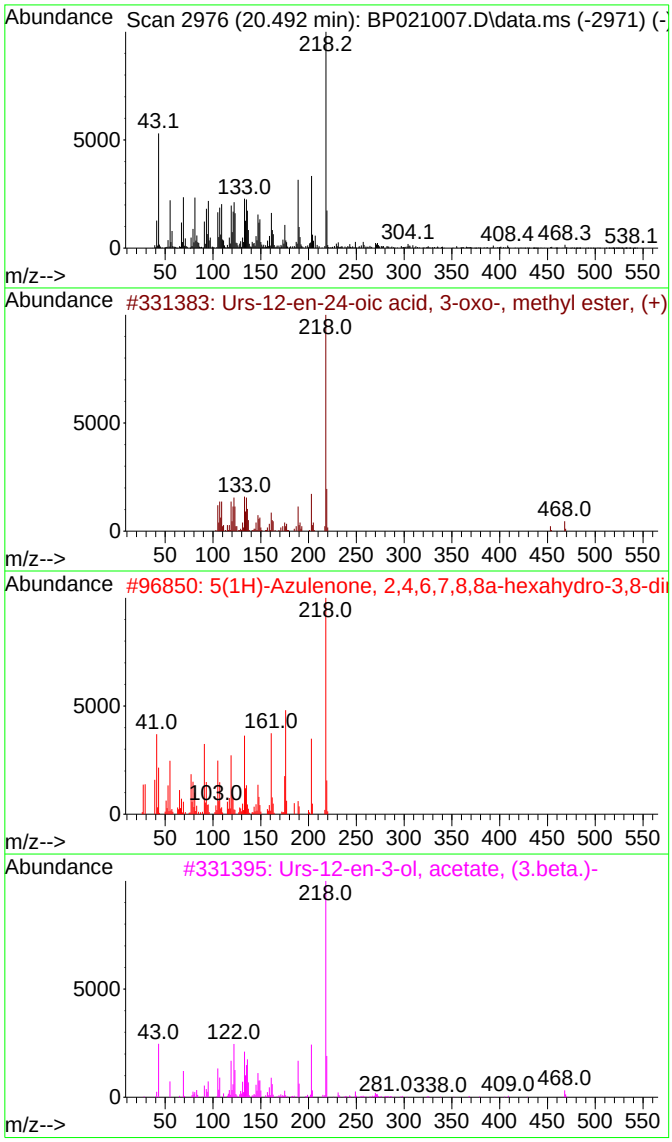
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

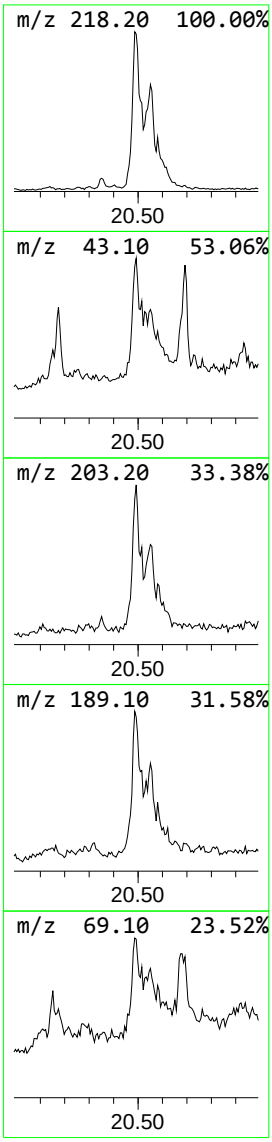
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Urs-12-en-24-oic acid, 3-ox... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.492	3.60 ng/ul	610102	Chrysene-d12	21.557

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Urs-12-en-24-oic acid, 3-oxo-, m...	468	C31H48O3	020475-86-9	99
2	5(1H)-Azulenone, 2,4,6,7,8,8a-he...	218	C15H22O	006754-66-1	78
3	Urs-12-en-3-ol, acetate, (3.beta.)-	468	C32H52O2	000863-76-3	76
4	.alpha.-Amyrin, TMS derivative	498	C33H58OSi	1000374-76-6	70
5	12-Oleanen-3-yl acetate, (3.alph...	468	C32H52O2	033055-28-6	70





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Data File : BP021007.D
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Sample : P3178-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT1

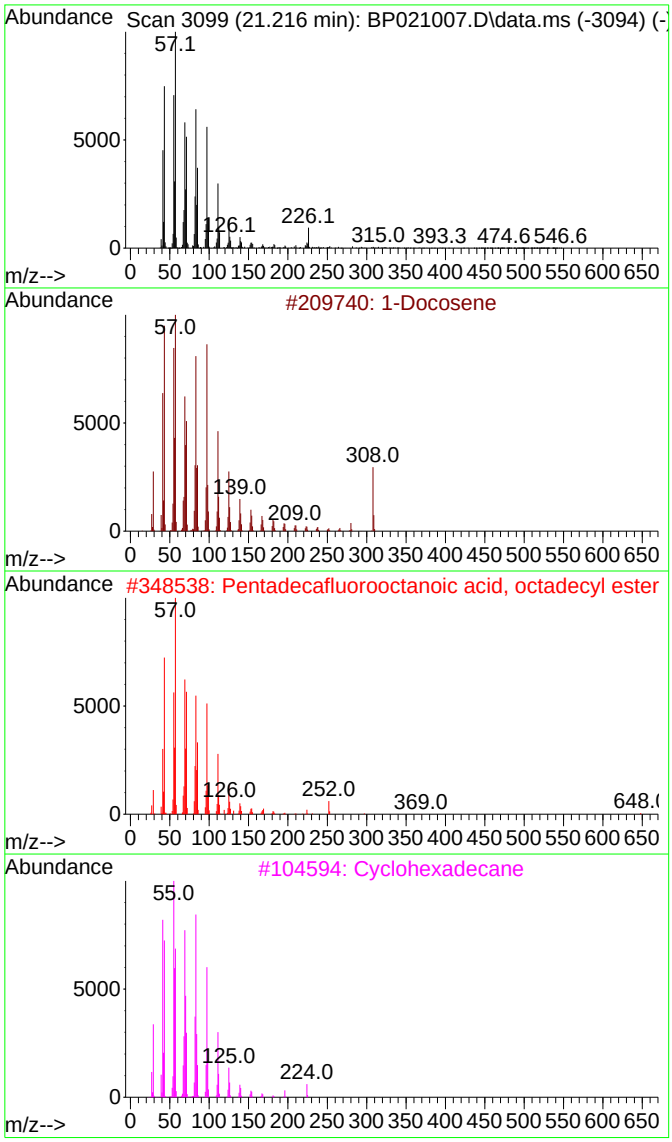
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

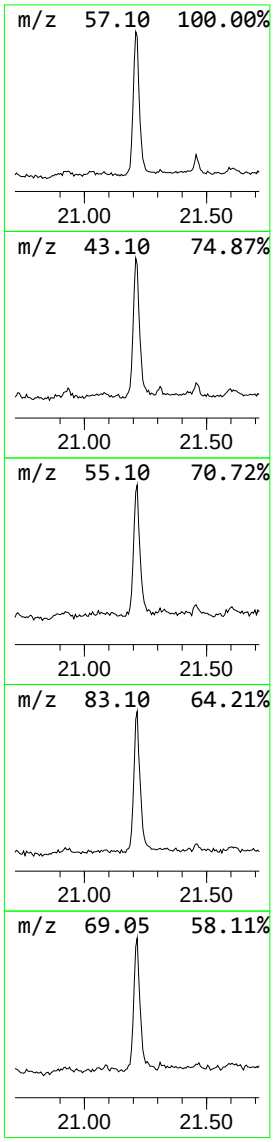
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.216	7.47 ng/ul	1264400	Chrysene-d12	21.557

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	97
2	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	95
3	Cyclohexadecane	224	C16H32	000295-65-8	93
4	1-Heneicosanol	312	C21H44O	015594-90-8	93
5	1-Octadecene	252	C18H36	000112-88-9	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP021007.D
Acq On : 17 Jul 2024 03:25
Operator : MA/JU
Sample : P3178-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT1

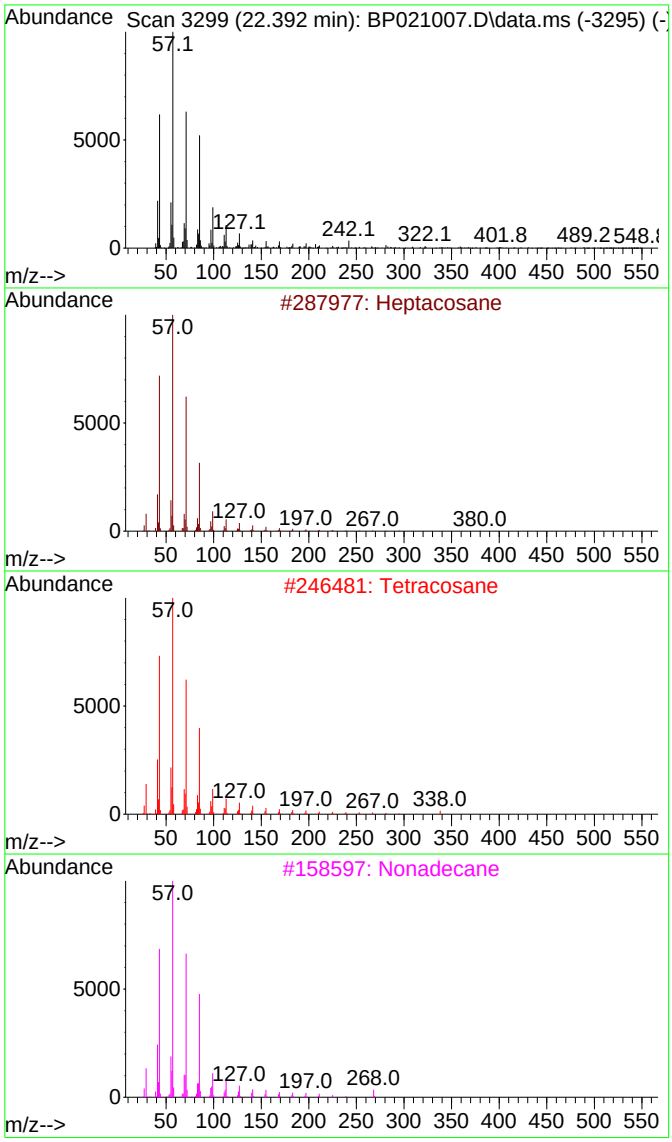
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

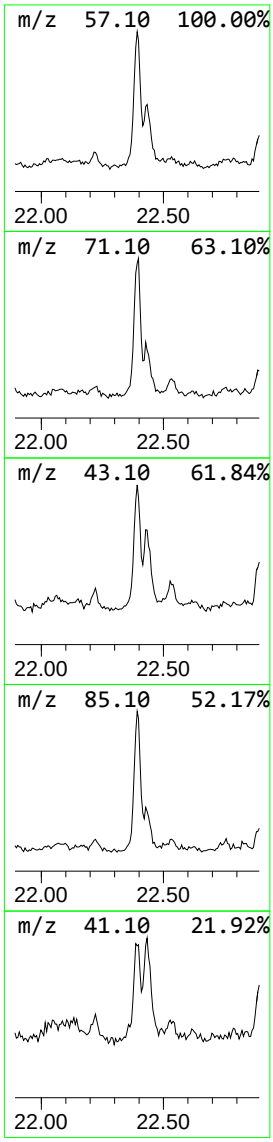
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.392	2.66 ng/ul	451066	Chrysene-d12	21.557

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	96
2	Tetracosane	338	C24H50	000646-31-1	93
3	Nonadecane	268	C19H40	000629-92-5	87
4	Heneicosane	296	C21H44	000629-94-7	87
5	Heptadecane	240	C17H36	000629-78-7	87





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Data File : BP021007.D
Acq On : 17 Jul 2024 03:25
Operator : MA/JU
Sample : P3178-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT1

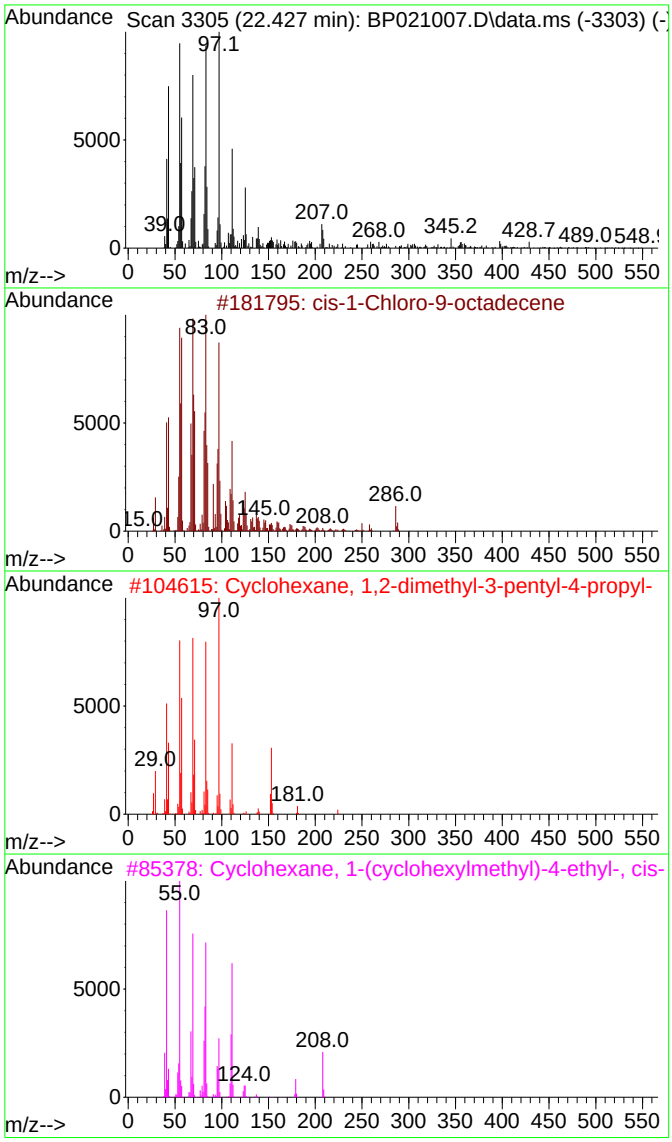
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

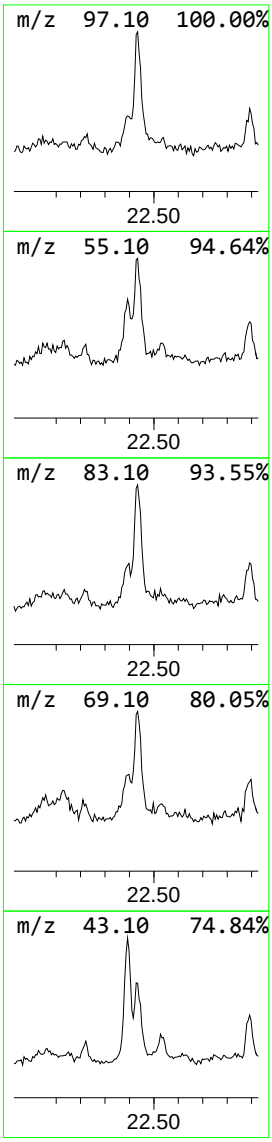
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 cis-1-Chloro-9-octadecene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.427	3.05 ng/ul	516155	Chrysene-d12	21.557

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis-1-Chloro-9-octadecene	286	C18H35Cl	016507-61-2	91
2	Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	68
3	Cyclohexane, 1-(cyclohexylmethyl...	208	C15H28	054934-95-1	45
4	Heptacos-1-ene	378	C27H54	015306-27-1	42
5	Cyclopentane, 1-ethyl-3-methyl-,...	112	C8H16	002613-65-2	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP021007.D
Acq On : 17 Jul 2024 03:25
Operator : MA/JU
Sample : P3178-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT1

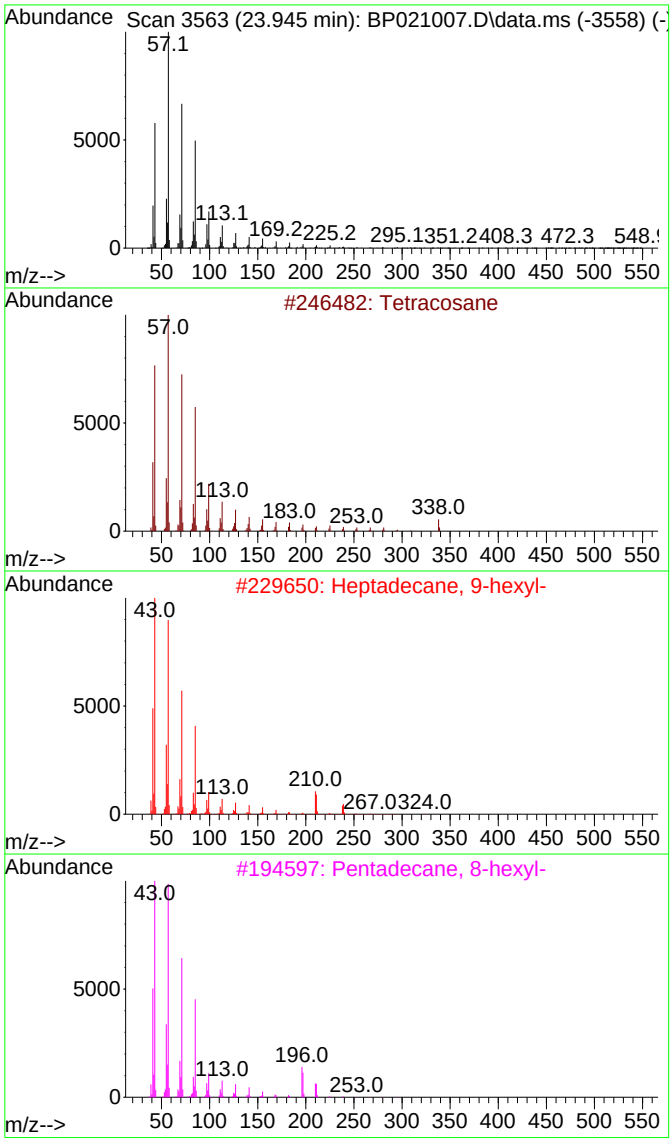
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

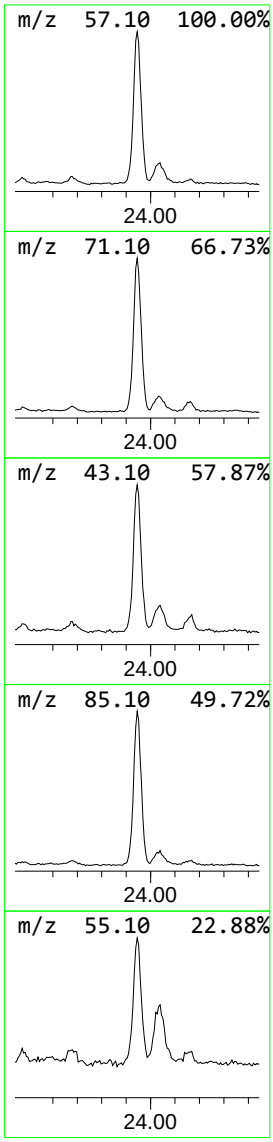
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.945	11.80 ng/ul	1920240	Perylene-d12	24.851

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	98
2	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
3	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5	Heneicosane	296	C21H44	000629-94-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP021007.D
Acq On : 17 Jul 2024 03:25
Operator : MA/JU
Sample : P3178-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT1

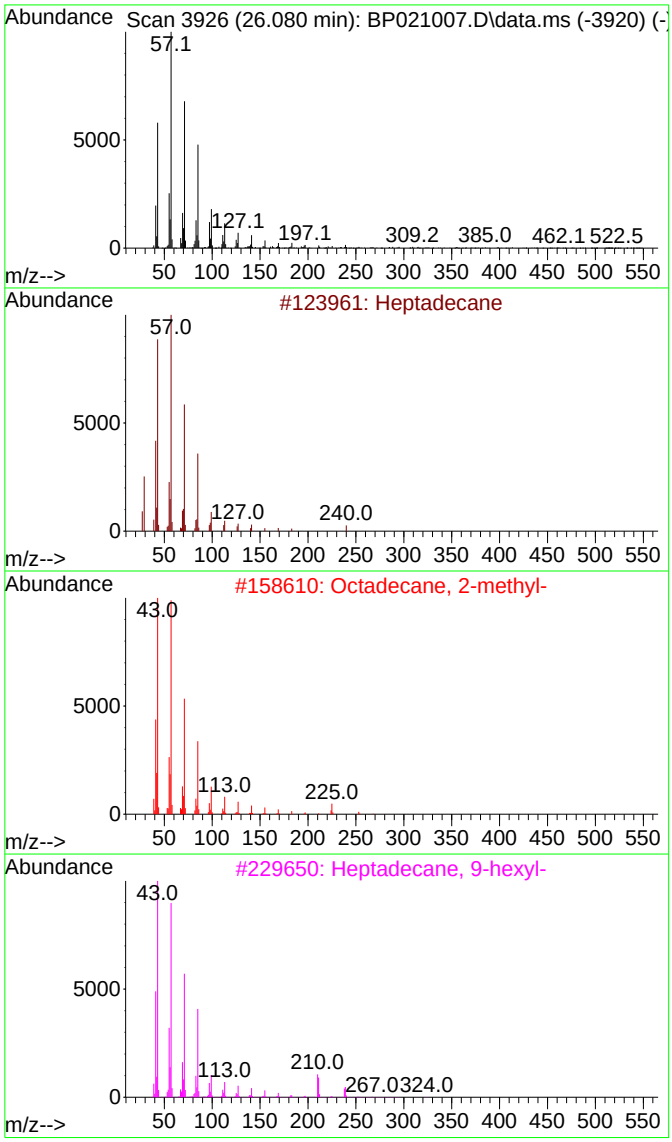
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

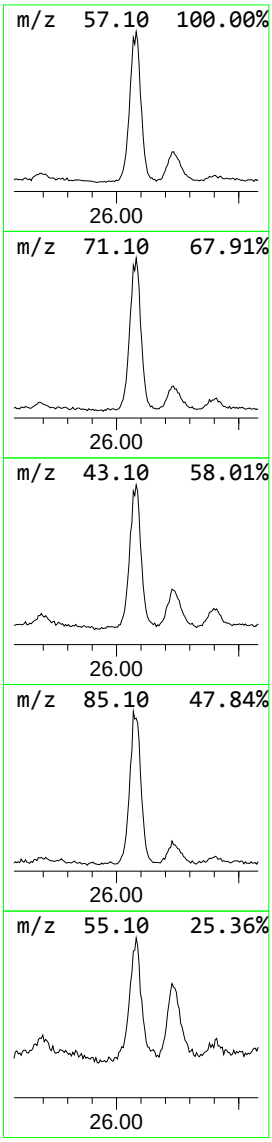
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.080	10.52 ng/ul	1710700	Perylene-d12	24.851

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	94
2	Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
3	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	93
4	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5	Hexacosane	366	C26H54	000630-01-3	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP021007.D
Acq On : 17 Jul 2024 03:25
Operator : MA/JU
Sample : P3178-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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
Propylene Glycol	3.522	2.2	ng/ul	156859	1	7.687	1398370	20.0
1-Phenoxypropan...	11.198	4.5	ng/ul	426548	2	10.457	1894150	20.0
n-Hexadecanoic ...	18.057	8.4	ng/ul	1136730	4	17.116	2707860	20.0
Octadecanoic acid	19.380	3.3	ng/ul	554082	5	21.557	3387310	20.0
Urs-12-en-24-oi...	20.492	3.6	ng/ul	610102	5	21.557	3387310	20.0
(DEL) Alkane: S...	21.216	7.5	ng/ul	1264400	5	21.557	3387310	20.0
(DEL) Alkane: S...	22.392	2.7	ng/ul	451066	5	21.557	3387310	20.0
cis-1-Chloro-9-...	22.427	3.0	ng/ul	516155	5	21.557	3387310	20.0
(DEL) Alkane: S...	23.945	11.8	ng/ul	1920240	6	24.851	3253390	20.0
(DEL) Alkane: S...	26.080	10.5	ng/ul	1710700	6	24.851	3253390	20.0