LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122319\
Data File : BM024268.D
Acq On : 24 Dec 2019 23:49
Operator : JU
Sample : K6240-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFQW2

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_M\METHODS\SOM-EPA-BM121719MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.993	15	18	28	rVB	75464	111312	1.77%	0.139%
2	3.322	72	74	79	rVB4	15587	22983	0.37%	0.029%
3	3.687	133	136	141	rVB2	16380	25095	0.40%	0.031%
4	5.528	443	449	457	rVB5	14408	30185	0.48%	0.038%
5	6.128	546	551	555	rBV8	12184	18739	0.30%	0.023%
6	6.634	632	637	653	rVB2	267629	571049	9.08%	0.711%
7	6.781	656	662	679	rBV	889373	1506918	23.96%	1.877%
8	6.969	688	694	706	rBV	1034484	1702460	27.07%	2.120%
9	7.081	708	713	720	rVB3	24972	54075	0.86%	0.067%
10	7.428	762	772	782	rBV	1224053	1939604	30.84%	2.416%
11	7.510	782	786	789	rBV	40715	62401	0.99%	0.078%
12	7.787	827	833	837	rVB7	9793	22979	0.37%	0.029%
13	8.069	874	881	884	rBV5	12509	19661	0.31%	0.024%
14	8.157	889	896	905	rBV	819691	1371665	21.81%	1.708%
15	8.534	953	960	962	rBV6	16963	29195	0.46%	0.036%
16	8.586	962	969	977	rVV	1077977	1853057	29.46%	2.308%
17	9.016	1036	1042	1045	rVV2	25852	43531	0.69%	0.054%
18	9.063	1045	1050	1055	rVV	101437	153787	2.45%	0.192%
19	9.298	1082	1090	1104	rBV	957237	1638060	26.04%	2.040%
20	9.480	1118	1121	1132	rVB10	21256	55684	0.89%	0.069%
21	9.598	1134	1141	1143	rBV7	15303	33788	0.54%	0.042%
22	9.739	1159	1165	1170	rBV5	15981	32165	0.51%	0.040%
23	9.839	1175	1182	1197	rBV	1396125	2576515	40.97%	3.209%
24	9.992	1203	1208	1213	rVB7	22158	34033	0.54%	0.042%
25	10.069	1216	1221	1227	rVB8	13816	26345	0.42%	0.033%
26	10.198	1236	1243	1250	rBV	1645804	2735151	43.49%	3.407%
27	10.251	1251	1252	1256	rVV4	16435	20315	0.32%	0.025%
28	10.298	1256	1260	1267	rVV	38945	67458	1.07%	0.084%
29	10.369	1267	1272	1279	rVB4	25979	50503	0.80%	0.063%
30	10.504	1290	1295	1302	rBV3	68888	125247	1.99%	0.156%
31	10.616	1307	1314	1328	rBV	3115288	5006212	79.60%	6.235%
32	10.763	1333	1339	1342	rVV	52076	106843	1.70%	0.133%
33	10.816	1344	1348	1352	rVV4	72929	134044	2.13%	0.167%
34	10.863	1353	1356	1365	rVB2	45820	85686	1.36%	0.107%

LSC Area Percent Report

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Data File : BM024268.D
Acq On : 24 Dec 2019 23:49
Operator : JU
Sample : K6240-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFQW2

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_M\METHODS\SOM-EPA-BM121719MA.M
Title : SVOA CALIBRATION

35	10.986	1368	1377	1381	rBV10	19802	47010	0.75%	0.059%
36	11.045	1381	1387	1388	rBV2	85230	122258	1.94%	0.152%
37	11.075	1389	1392	1398	rVV	107605	169951	2.70%	0.212%
38	11.139	1398	1403	1407	rVV	40889	62341	0.99%	0.078%
39	11.186	1407	1411	1420	rVB3	43128	86446	1.37%	0.108%
40	11.298	1427	1430	1438	rVB9	16704	29305	0.47%	0.036%
41	11.416	1438	1450	1454	rBV	65101	129781	2.06%	0.162%
42	11.463	1454	1458	1459	rVV	64323	85940	1.37%	0.107%
43	11.486	1460	1462	1466	rVV	78839	106713	1.70%	0.133%
44	11.533	1466	1470	1477	rVB	60223	97986	1.56%	0.122%
45	11.616	1482	1484	1489	rVV5	14641	20889	0.33%	0.026%
46	11.704	1489	1499	1505	rVV	125556	287798	4.58%	0.358%
47	11.763	1505	1509	1522	rVB2	125620	282516	4.49%	0.352%
48	11.910	1529	1534	1539	rBV4	33035	63985	1.02%	0.080%
49	11.963	1539	1543	1549	rVB2	31995	63866	1.02%	0.080%
50	12.033	1549	1555	1557	rBV7	12294	19841	0.32%	0.025%
51	12.092	1561	1565	1570	rVB2	17229	24674	0.39%	0.031%
52	12.439	1619	1624	1630	rVB	38838	76585	1.22%	0.095%
53	12.551	1638	1643	1653	rBV4	71059	203893	3.24%	0.254%
54	12.698	1662	1668	1674	rVB	61035	92684	1.47%	0.115%
55	12.792	1679	1684	1688	rBV7	15697	25930	0.41%	0.032%
56	12.927	1704	1707	1714	rVB7	20947	47439	0.75%	0.059%
57	13.004	1714	1720	1727	rBV	78000	136758	2.17%	0.170%
58	13.080	1727	1733	1738	rVB4	17181	31082	0.49%	0.039%
59	13.145	1741	1744	1748	rVB5	11960	18423	0.29%	0.023%
60	13.521	1801	1808	1813	rBV	2613689	3689008	58.65%	4.595%
61	13.568	1813	1816	1824	rVB	111201	157723	2.51%	0.196%
62	13.639	1824	1828	1830	rBV5	24307	32695	0.52%	0.041%
63	13.668	1830	1833	1840	rVB	64826	107768	1.71%	0.134%
64	13.780	1845	1852	1864	rBV	3214409	4754842	75.60%	5.922%
65	13.927	1872	1877	1883	rVB	55589	88127	1.40%	0.110%
66	14.021	1889	1893	1899	rVB6	21747	32790	0.52%	0.041%
67	14.092	1899	1905	1921	rVB	2431067	3618401	57.53%	4.507%
68	14.251	1930	1932	1937	rBV5	16034	20842	0.33%	0.026%
69	14.368	1946	1952	1960	rBV	146403	331665	5.27%	0.413%
70	14.651	1996	2000	2004	rVB7	25760	39131	0.62%	0.049%
71	14.868	2032	2037	2043	rBV	174145	263381	4.19%	0.328%

LSC Area Percent Report

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Operator : JU
Sample : K6240-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFQW2

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_M\METHODS\SOM-EPA-BM121719MA.M
Title : SVOA CALIBRATION

72	14.945	2044	2050	2053	rVV2	49717	75012	1.19%	0.093%
73	15.098	2070	2076	2084	rBV	3999898	5688487	90.45%	7.085%
74	15.198	2090	2093	2097	rBV2	22350	31999	0.51%	0.040%
75	15.257	2097	2103	2111	rBV	645568	991640	15.77%	1.235%
76	15.339	2113	2117	2118	rBV2	35086	42474	0.68%	0.053%
77	15.439	2131	2134	2138	rBV	43953	53445	0.85%	0.067%
78	15.492	2138	2143	2147	rBV3	154251	238970	3.80%	0.298%
79	15.621	2161	2165	2170	rVB2	40372	60310	0.96%	0.075%
80	15.692	2173	2177	2181	rVB	73601	100830	1.60%	0.126%
81	15.957	2220	2222	2227	rVB6	23462	29621	0.47%	0.037%
82	16.215	2262	2266	2271	rBV	55534	71960	1.14%	0.090%
83	16.639	2334	2338	2341	rBV2	82023	107351	1.71%	0.134%
84	16.674	2341	2344	2353	rVB4	52495	96991	1.54%	0.121%
85	16.851	2368	2374	2384	rBV	2739549	3939783	62.64%	4.907%
86	16.951	2385	2391	2401	rVV2	4066384	5763156	91.63%	7.178%
87	17.186	2428	2431	2436	rVB2	56670	83748	1.33%	0.104%
88	17.374	2461	2463	2467	rVB2	27409	32771	0.52%	0.041%
89	17.609	2500	2503	2508	rVB	38932	48259	0.77%	0.060%
90	17.780	2527	2532	2538	rBV2	382235	665131	10.58%	0.828%
91	18.227	2605	2608	2616	rVB	128770	185128	2.94%	0.231%
92	19.074	2749	2752	2757	rBV	129524	182091	2.90%	0.227%
93	19.262	2778	2784	2790	rBV	4853383	6289422	100.00%	7.833%
94	20.803	3043	3046	3053	rBV	1019046	1203802	19.14%	1.499%
95	21.068	3086	3091	3097	rBV	3245679	4147192	65.94%	5.165%
96	21.580	3175	3178	3183	rVB	727984	815704	12.97%	1.016%
97	22.103	3264	3267	3272	rVB	378828	451667	7.18%	0.563%
98	22.580	3345	3348	3353	rVB	455189	560145	8.91%	0.698%
99	23.062	3425	3430	3436	rBV	3530758	6196644	98.52%	7.718%
100	23.197	3447	3453	3460	rVB	2528384	4400539	69.97%	5.481%

Sum of corrected areas: 80291484

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Data File : BM024268.D
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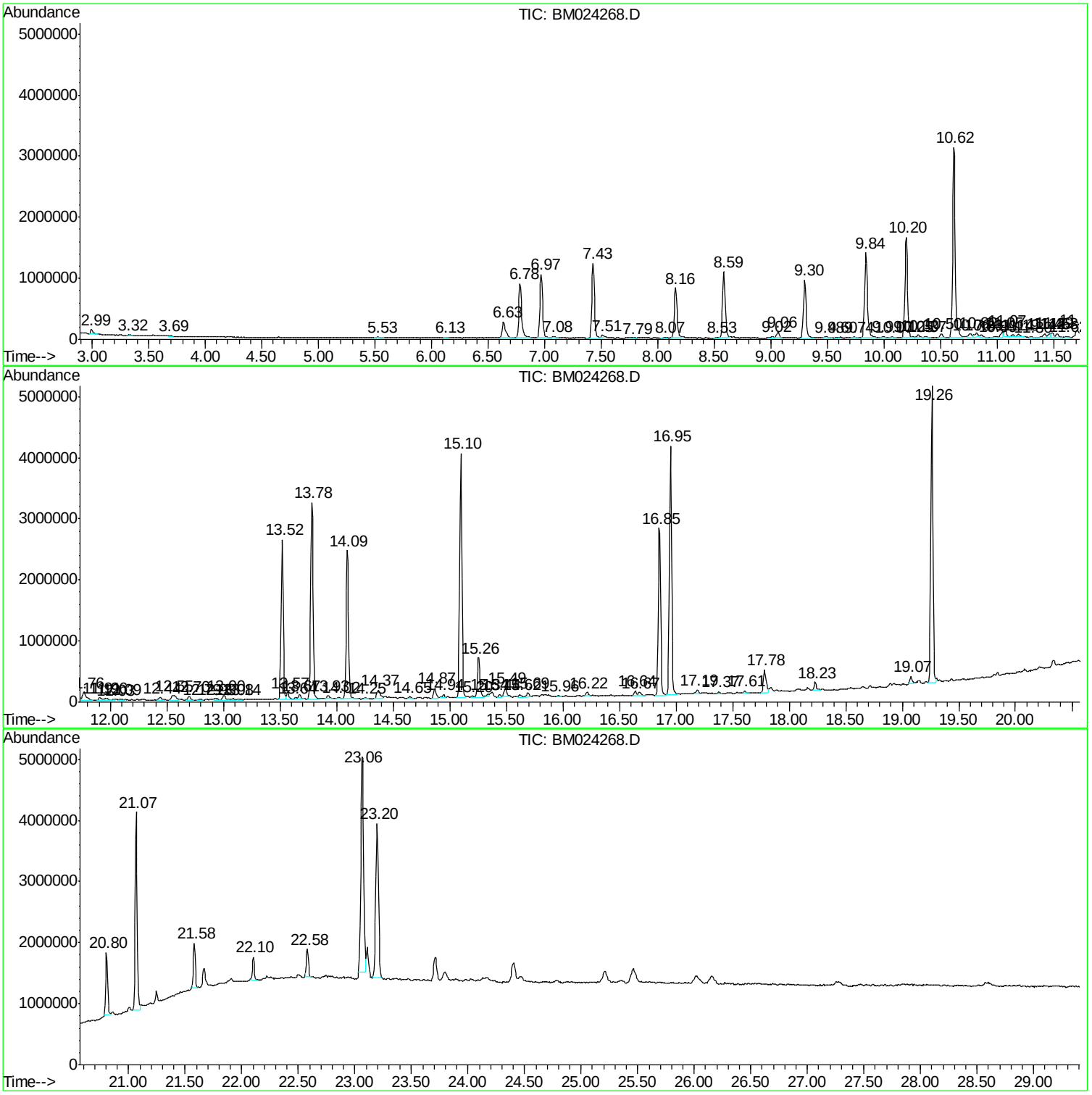
Sample : K6240-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFQW2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

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Quant Title : SVOA CALIBRATION

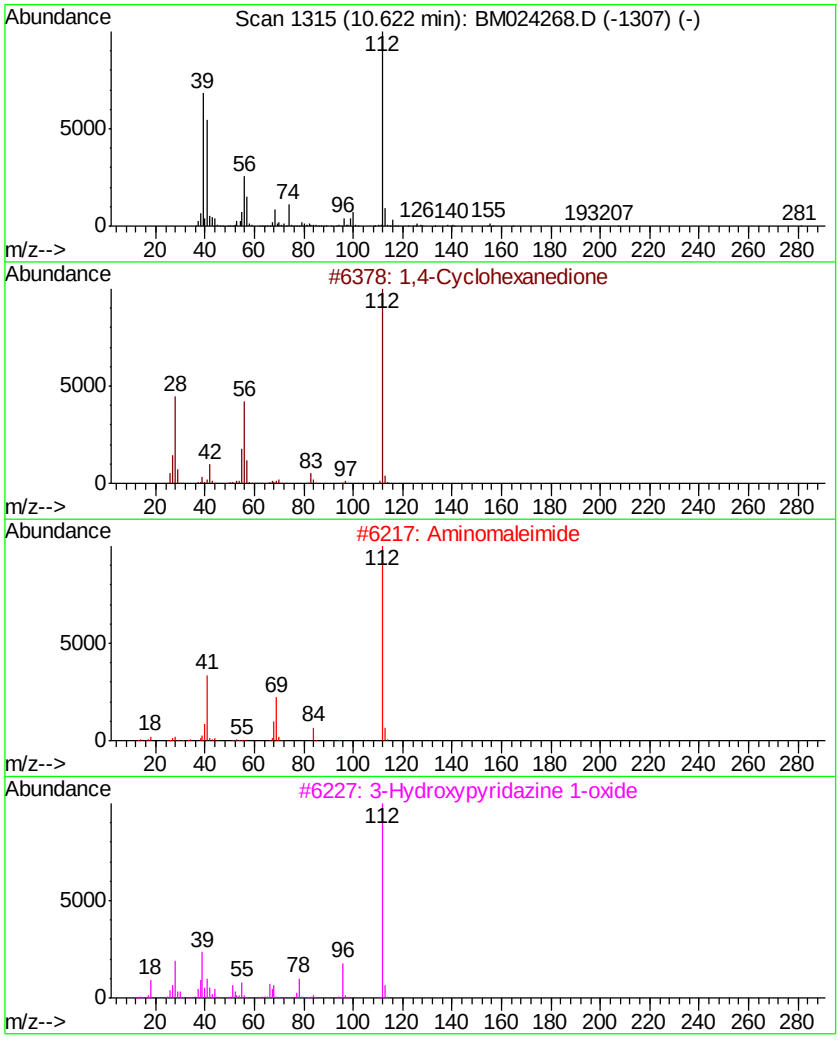
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

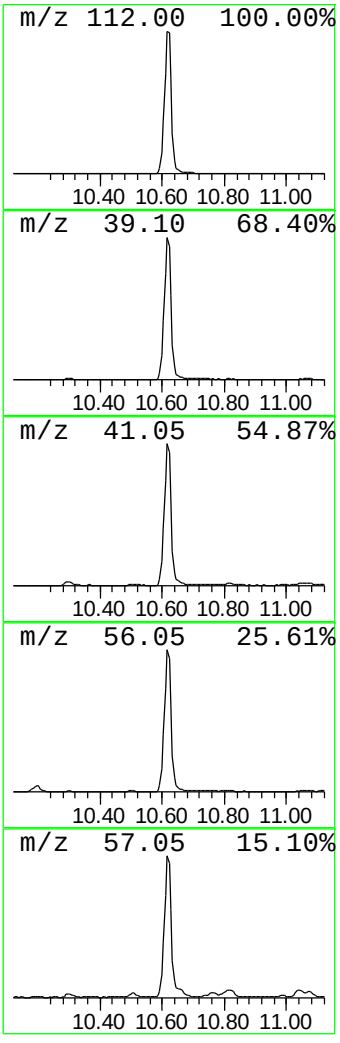
****

Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.62	36.61 ng/ul	5006210	Naphthalene-d8	10.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Cyclohexanedione	112	C6H8O2	000637-88-7	47
2	Aminomaleimide	112	C4H4N2O2	037770-94-8	38
3	3-Hydroxypyridazine 1-oxide	112	C4H4N2O2	018259-51-3	35
4	3-Nitropyrrole	112	C4H4N2O2	005930-94-9	32
5	Emimycin	112	C4H4N2O2	003735-46-4	25





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122319\
Data File : BM024268.D
Acq On : 24 Dec 2019 23:49
Operator : JU
Sample : K6240-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFQW2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION

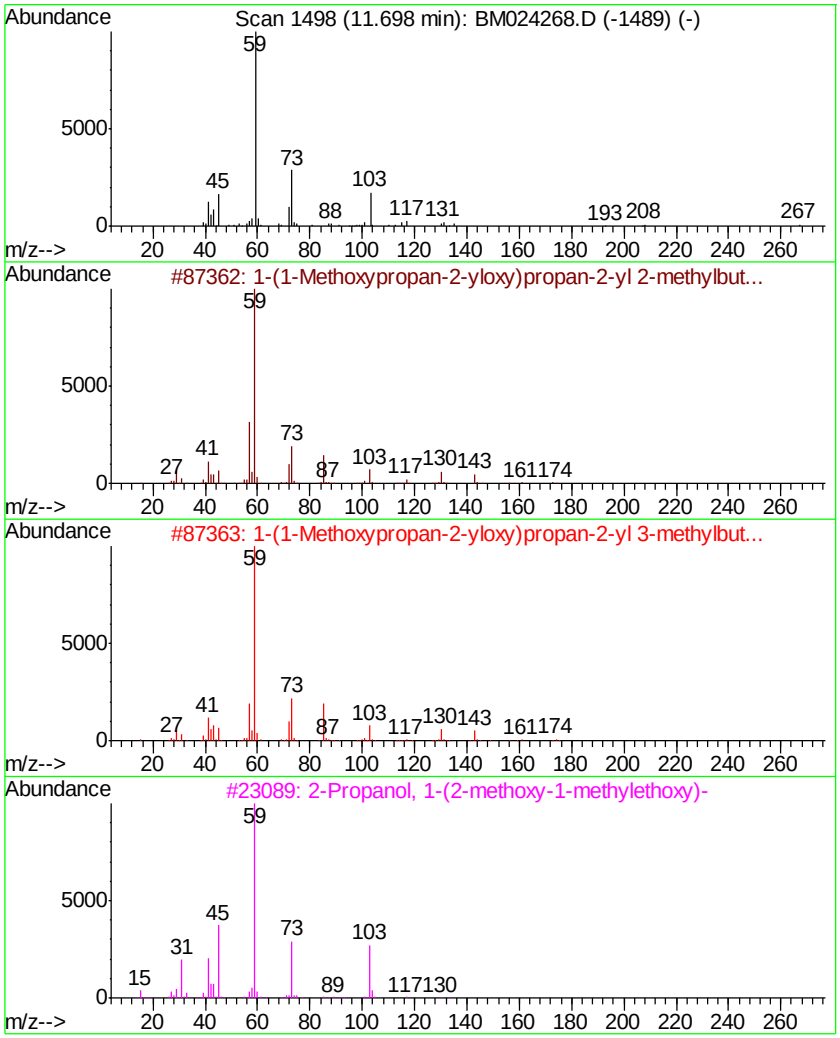
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

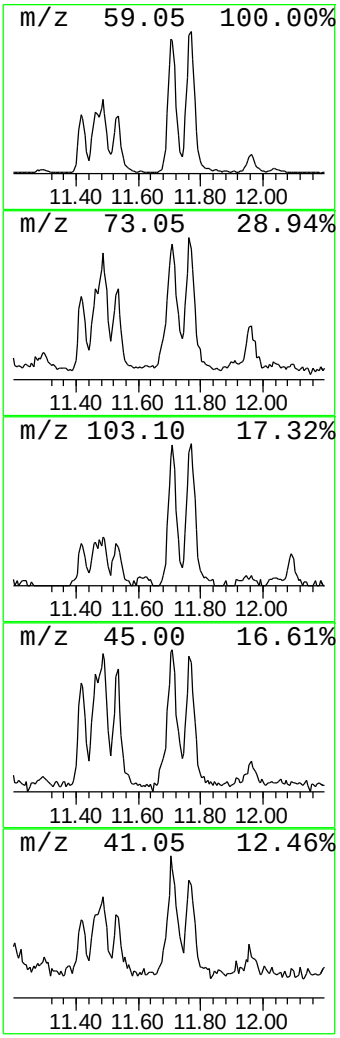
****

Peak Number 2 1-(1-Methoxypropan-2-yloxy)... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.70	2.10 ng/ul	287798	Naphthalene-d8	10.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(1-Methoxypropan-2-yloxy)propa...	232	C12H24O4	1000367-11-4	72
2	1-(1-Methoxypropan-2-yloxy)propa...	232	C12H24O4	1000367-13-2	72
3	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	72
4	1-(1-Methoxypropan-2-yloxy)propa...	232	C12H24O4	1000367-13-4	64
5	Tri(1,2-propyleneglycol), monome...	206	C10H22O4	1000262-26-6	64





Library Search Compound Report

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION

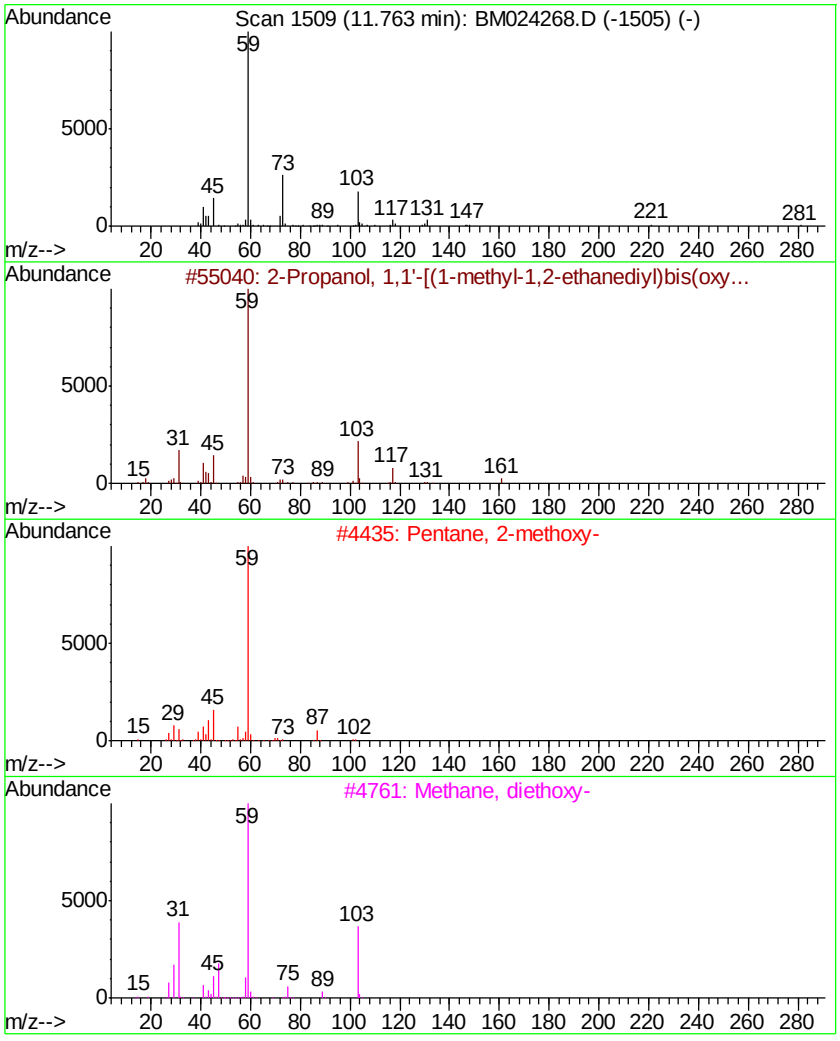
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

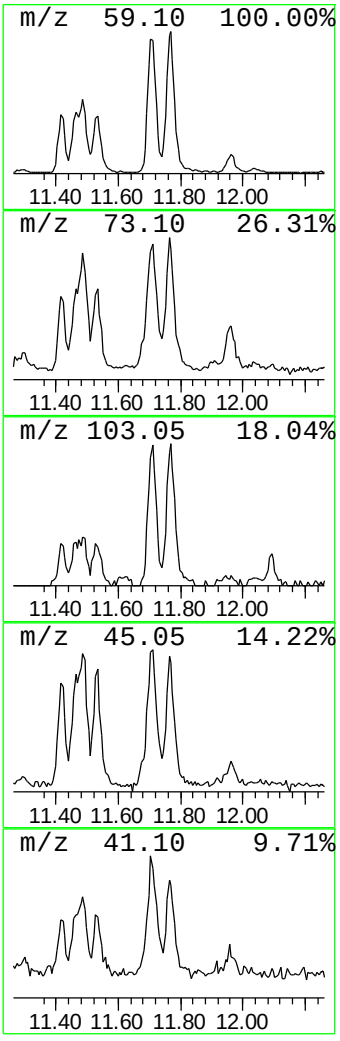
****

Peak Number 3 2-Propanol, 1,1'-[(1-methyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.76	2.07 ng/ul	282516	Naphthalene-d8	10.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	59
2	Pentane, 2-methoxy-	102	C6H14O	006795-88-6	50
3	Methane, diethoxy-	104	C5H12O2	000462-95-3	50
4	Propanedioic acid, oxo-, dimethy...	146	C5H6O5	003298-40-6	50
5	Hexaethylene glycol dimethyl ether	310	C14H30O7	001072-40-8	50





Library Search Compound Report

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION

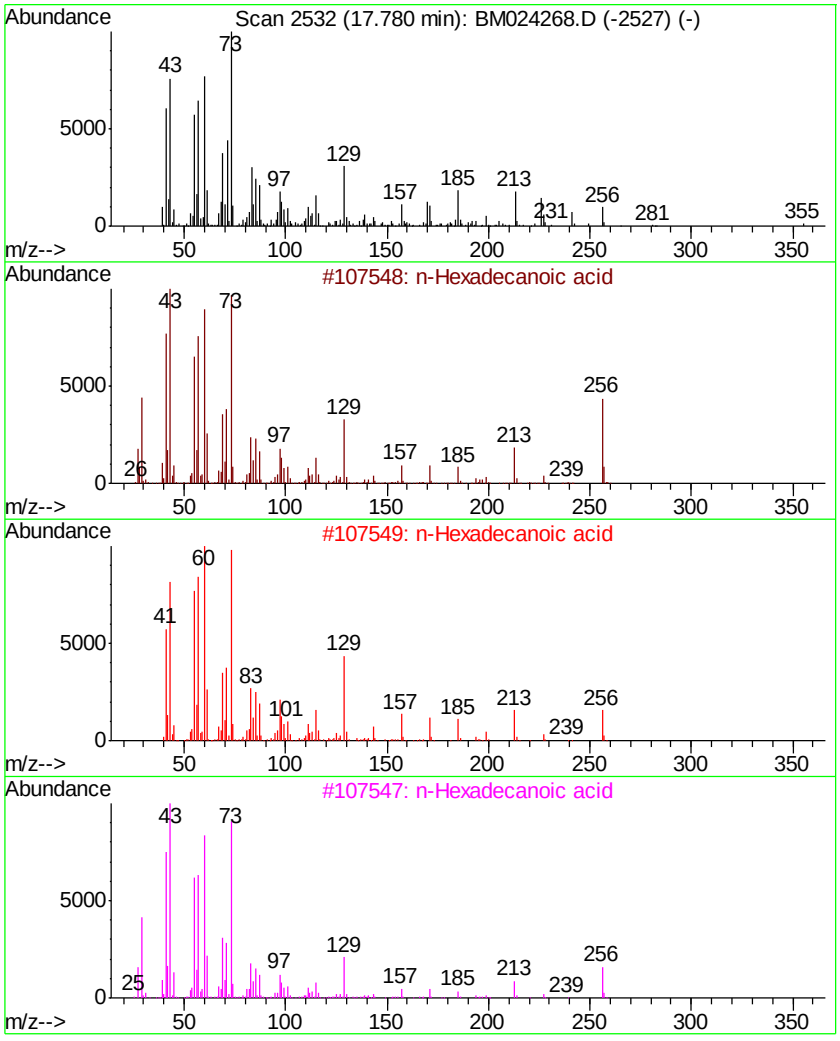
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

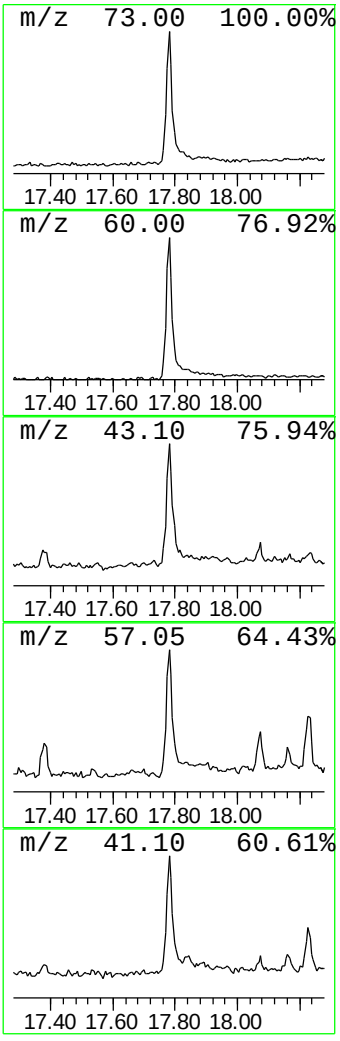
****

Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.78	3.38 ng/ul	665131	Phenanthrene-d10	16.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
4	Tridecanoic acid	214	C13H26O2	000638-53-9	91
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	81





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Data File : BM024268.D
Acq On : 24 Dec 2019 23:49
Operator : JU
Sample : K6240-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFQW2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION

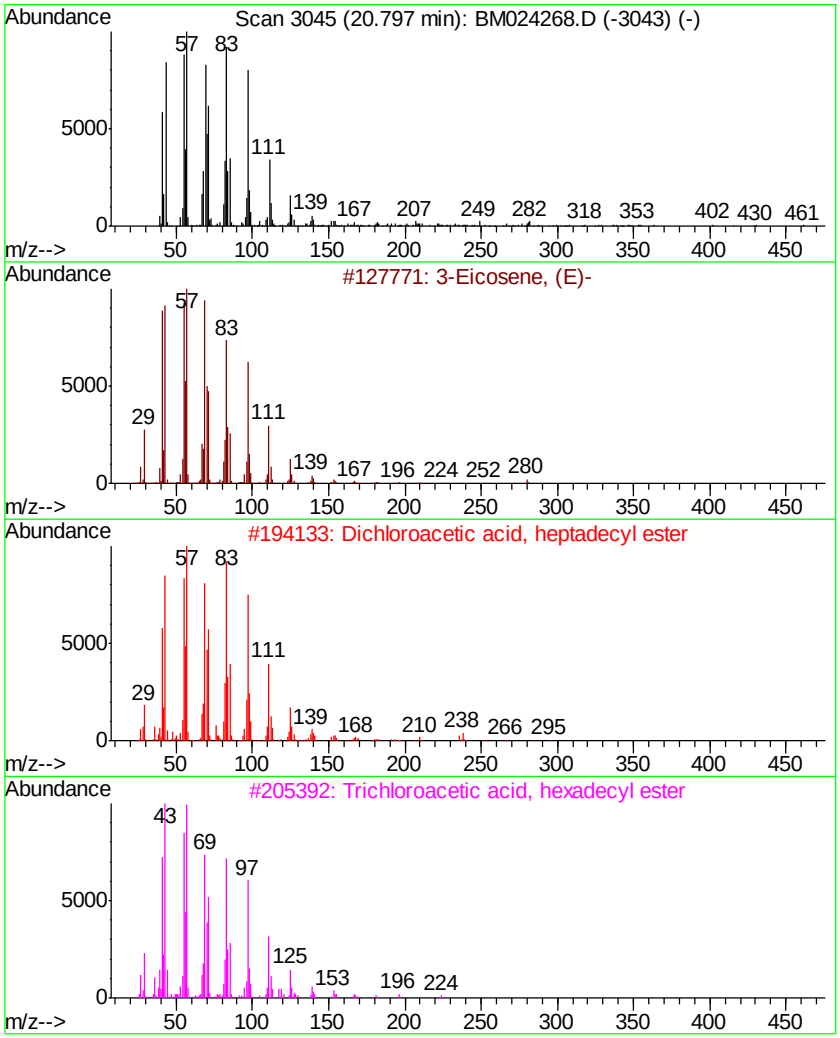
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

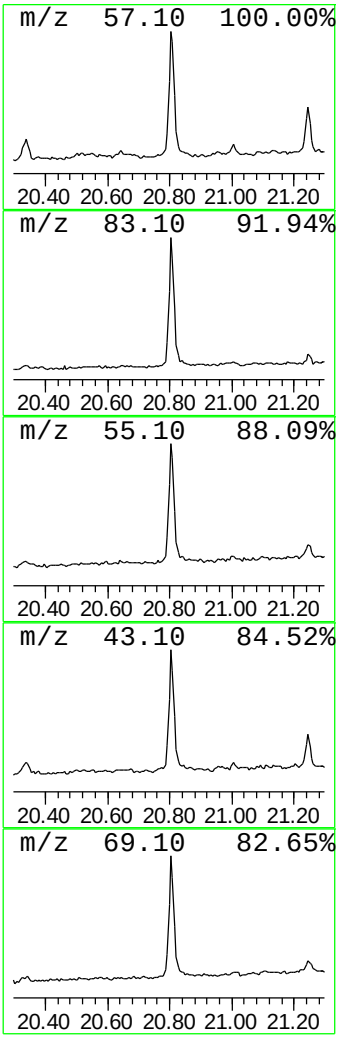
****

Peak Number 5 (DEL) Alkane: Cyclic20.80 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.80	5.81 ng/ul	1203800	Chrysene-d12	21.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	95
2	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
3	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
4	Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
5	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	91





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Data File : BM024268.D
Acq On : 24 Dec 2019 23:49
Operator : JU
Sample : K6240-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFQW2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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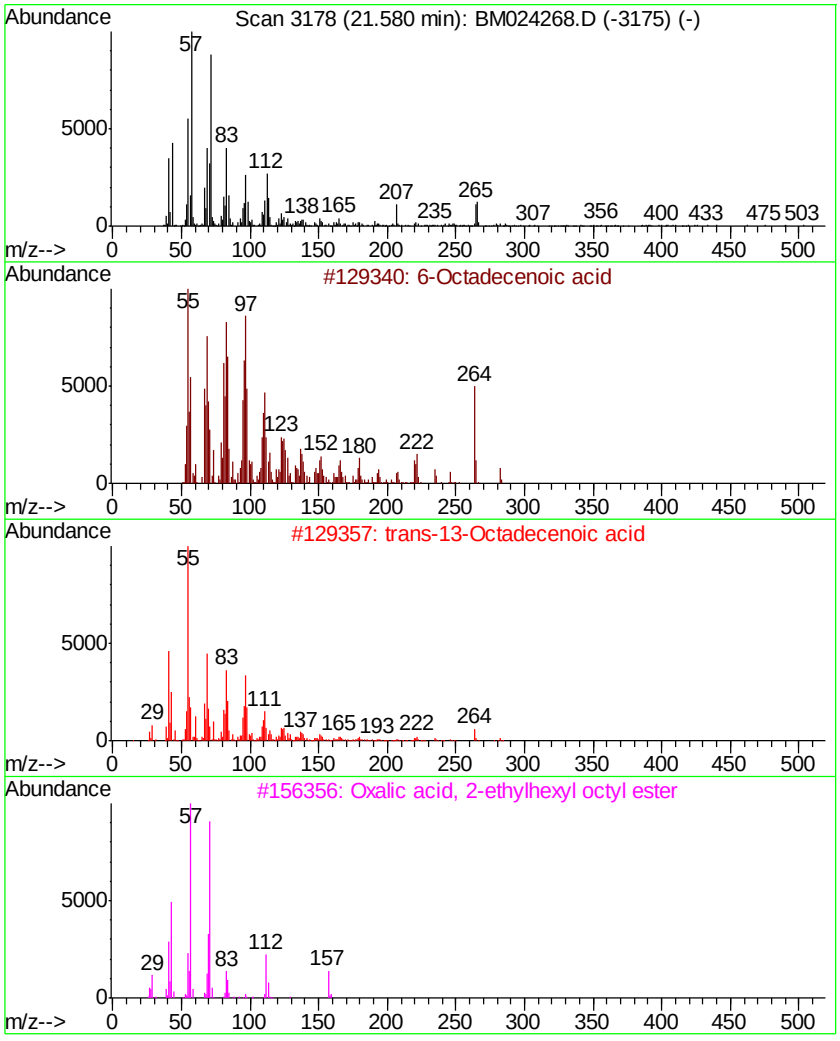
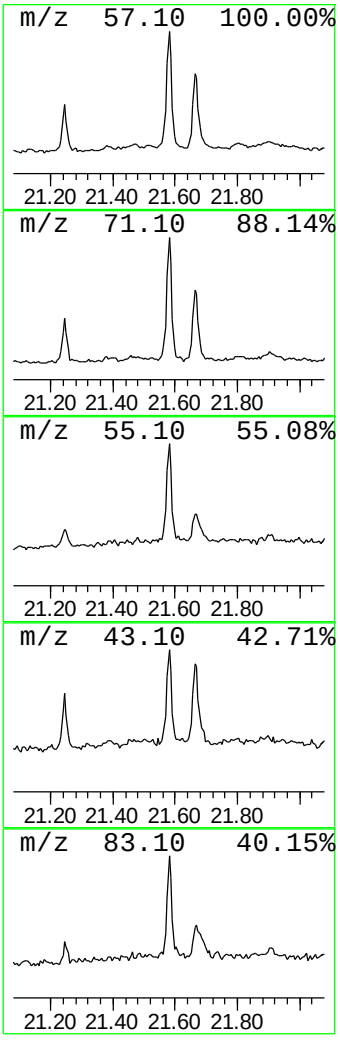
Peak Number 6 6-Octadecenoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.58	3.93 ng/ul	815704	Chrysene-d12	21.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Octadecenoic acid	282	C18H34O2	1000336-66-8	56
2	trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	56
3	Oxalic acid, 2-ethylhexyl octyl ...	314	C18H34O4	1000309-39-1	49
4	Oleic Acid	282	C18H34O2	000112-80-1	46
5	cis-Vaccenic acid	282	C18H34O2	000506-17-2	45





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122319\
Data File : BM024268.D
Acq On : 24 Dec 2019 23:49
Operator : JU
Sample : K6240-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFQW2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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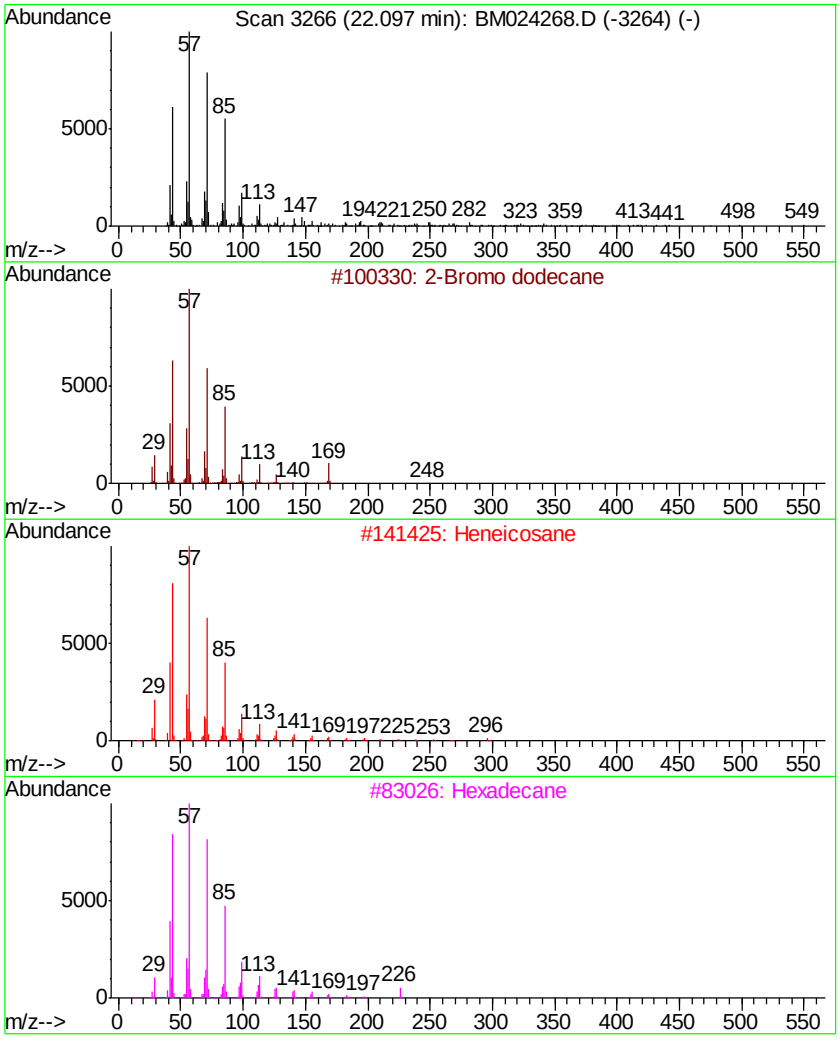
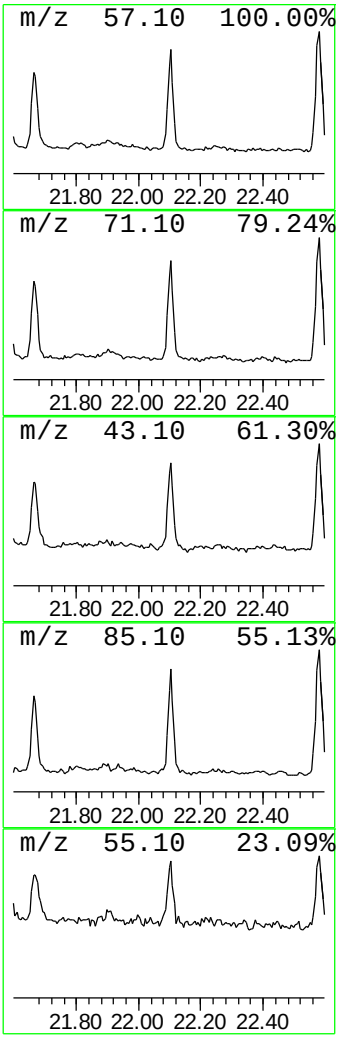
Peak Number 7 2-Bromo dodecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.10	2.18 ng/ul	451667	Chrysene-d12	21.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo dodecane	248	C12H25Br	013187-99-0	90
2	Heneicosane	296	C21H44	000629-94-7	90
3	Hexadecane	226	C16H34	000544-76-3	87
4	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86
5	Dodecane	170	C12H26	000112-40-3	83





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Data File : BM024268.D
Acq On : 24 Dec 2019 23:49
Operator : JU
Sample : K6240-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFQW2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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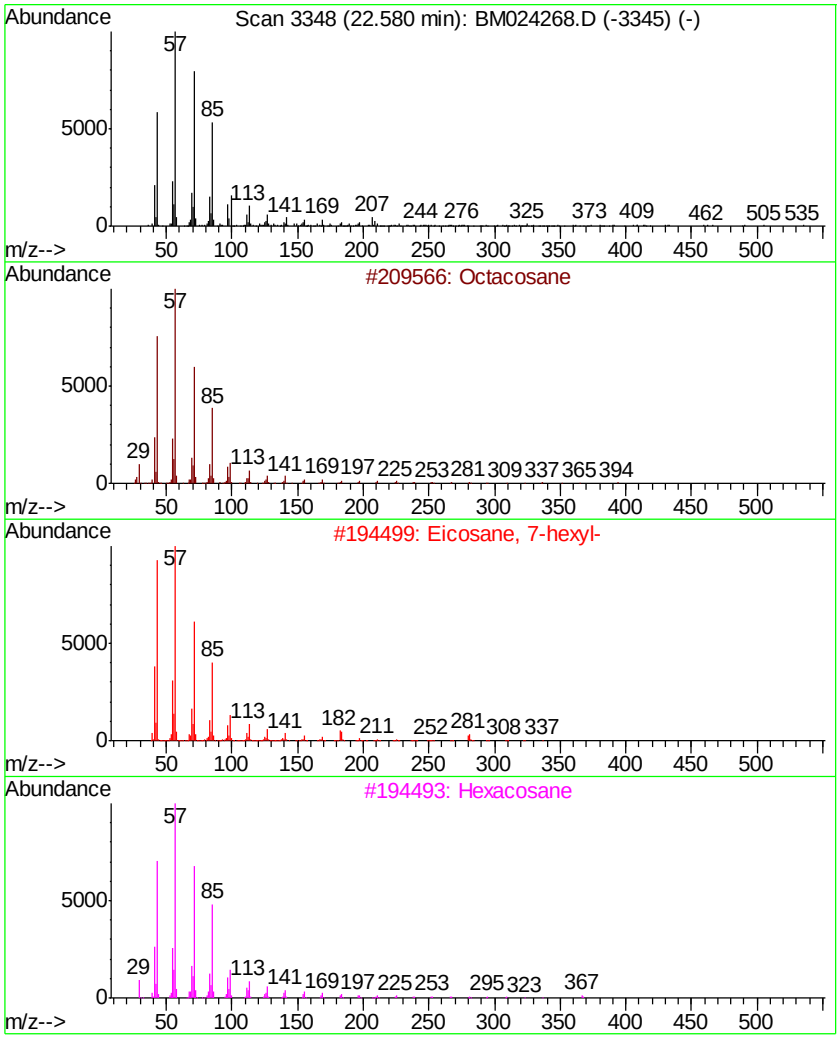
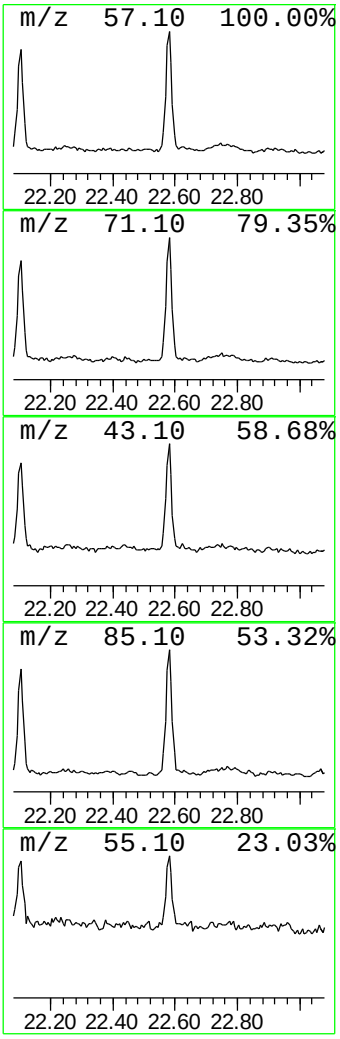
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.58	2.55 ng/ul	560145	Perylene-d12	23.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	91
2	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
3	Hexacosane	366	C26H54	000630-01-3	91
4	Tetracosane	338	C24H50	000646-31-1	91
5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM122319\
Data File : BM024268.D
Acq On : 24 Dec 2019 23:49
Operator : JU
Sample : K6240-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFQW2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown-01	10.62	36.6	ng/ul	5006210	2	10.20	2735150	20.0
1-(1-Methoxypropa...	11.70	2.1	ng/ul	287798	2	10.20	2735150	20.0
2-Propanol, 1,1'-...	11.76	2.1	ng/ul	282516	2	10.20	2735150	20.0
n-Hexadecanoic acid	17.78	3.4	ng/ul	665131	4	16.85	3939780	20.0
(DEL) Alkane: Cyc...	20.80	5.8	ng/ul	1203800	5	21.07	4147190	20.0
6-Octadecenoic acid	21.58	3.9	ng/ul	815704	5	21.07	4147190	20.0
2-Bromo dodecane	22.10	2.2	ng/ul	451667	5	21.07	4147190	20.0
(DEL) Alkane: Str...	22.58	2.5	ng/ul	560145	6	23.20	4400540	20.0