LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0DL1

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\SOM-EPA-BP022720MA.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	2.917	3	5	21	rVB	862505	1176960	2.33%	0.428%
2	3.164	43	47	56	rVB	64197	92729	0.18%	0.034%
3	4.993	350	358	374	rBV	15805920	28570050	56.61%	10.391%
4	5.752	481	487	497	rBV	163192	265618	0.53%	0.097%
5	6.334	580	586	594	rVB2	50268	78896	0.16%	0.029%
6	6.816	662	668	684	rBV2	256573	612945	1.21%	0.223%
7	6.975	688	695	710	rBV	947544	1492886	2.96%	0.543%
8	7.175	721	729	744	rVB	1007187	1749004	3.47%	0.636%
9	7.534	781	790	801	rBV	9434022	16144261	31.99%	5.872%
10	7.687	801	816	835	rVB	19354897	43401249	86.00%	15.785%
11	8.004	856	870	876	rBV	19997028	50469201	100.00%	18.356%
12	8.346	917	928	937	rBV	743276	1234414	2.45%	0.449%
13	8.781	994	1002	1004	rBV	1058715	1707278	3.38%	0.621%
14	8.828	1004	1010	1023	rVB	5944782	10219252	20.25%	3.717%
15	9.116	1052	1059	1068	rVB	94501	159321	0.32%	0.058%
16	9.263	1080	1084	1093	rVB	121990	195715	0.39%	0.071%
17	9.440	1107	1114	1118	rVV2	133982	243548	0.48%	0.089%
18	9.498	1118	1124	1133	rVB	896977	1466360	2.91%	0.533%
19	10.034	1207	1215	1223	rBV	1510167	2653666	5.26%	0.965%
20	10.098	1223	1226	1232	rVV	190622	310674	0.62%	0.113%
21	10.193	1234	1242	1248	rVV	181903	332705	0.66%	0.121%
22	10.281	1248	1257	1272	rVB	12043590	21671450	42.94%	7.882%
23	10.410	1272	1279	1292	rVB	1451820	2367720	4.69%	0.861%
24	10.516	1292	1297	1307	rBV2	89809	265333	0.53%	0.097%
25	10.793	1335	1344	1354	rBV	6571138	11215579	22.22%	4.079%
26	11.004	1366	1380	1397	rBV	2914554	4875077	9.66%	1.773%
27	11.216	1408	1416	1428	rBV	771011	1367151	2.71%	0.497%
28	11.704	1486	1499	1503	rBV4	32008	70071	0.14%	0.025%
29	11.757	1503	1508	1518	rVB	87748	148482	0.29%	0.054%
30	12.004	1544	1550	1559	rVB2	28210	54331	0.11%	0.020%
31	12.422	1613	1621	1622	rBV	61460	84776	0.17%	0.031%
32	12.457	1622	1627	1634	rVB	274437	453889	0.90%	0.165%
33	12.839	1687	1692	1698	rBV3	85485	135522	0.27%	0.049%
34	13.063	1723	1730	1739	rBV	901021	1503677	2.98%	0.547%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0DL1

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\SOM-EPA-BP022720MA.M
Title : SVOA CALIBRATION

35	13.163	1742	1747	1753	rVV	164366	307306	0.61%	0.112%
36	13.222	1753	1757	1761	rVV3	52970	91507	0.18%	0.033%
37	13.287	1761	1768	1776	rVB2	407402	721976	1.43%	0.263%
38	13.687	1826	1836	1840	rBV	2856851	4233910	8.39%	1.540%
39	13.734	1840	1844	1853	rVB	604257	852465	1.69%	0.310%
40	13.957	1875	1882	1891	rBV	3465600	5144300	10.19%	1.871%
41	14.269	1928	1935	1951	rVB	2230096	3411496	6.76%	1.241%
42	14.481	1965	1971	1981	rBV	126371	263167	0.52%	0.096%
43	15.269	2098	2105	2112	rBV	4502102	6415687	12.71%	2.333%
44	15.386	2119	2125	2139	rBV	1021223	1481089	2.93%	0.539%
45	15.510	2142	2146	2152	rVB2	49743	69698	0.14%	0.025%
46	17.022	2397	2403	2413	rBV	2849148	3995869	7.92%	1.453%
47	17.122	2413	2420	2436	rBV	4884928	7047396	13.96%	2.563%
48	17.939	2556	2559	2564	rBV2	65472	98485	0.20%	0.036%
49	17.992	2564	2568	2575	rVB4	38175	74319	0.15%	0.027%
50	19.016	2738	2742	2746	rVB	55341	70376	0.14%	0.026%
51	19.180	2767	2770	2777	rBV3	33365	55547	0.11%	0.020%
52	19.257	2779	2783	2791	rVB	52895	78559	0.16%	0.029%
53	19.369	2796	2802	2813	rVB	6449632	8651307	17.14%	3.147%
54	20.863	3052	3056	3061	rBV2	121772	168624	0.33%	0.061%
55	21.116	3094	3099	3115	rBV	3868135	4798441	9.51%	1.745%
56	21.292	3125	3129	3134	rBV	228653	248194	0.49%	0.090%
57	21.721	3198	3202	3211	rBV	429168	510027	1.01%	0.185%
58	22.180	3275	3280	3288	rVB	641450	801176	1.59%	0.291%
59	22.598	3348	3351	3358	rVB2	54652	86949	0.17%	0.032%
60	22.686	3361	3366	3374	rVB	693543	1019910	2.02%	0.371%
61	23.180	3443	3450	3457	rBV	4970524	9237110	18.30%	3.360%
62	23.251	3457	3462	3467	rVV	710795	1211443	2.40%	0.441%
63	23.321	3467	3474	3485	rVB	2600699	5078813	10.06%	1.847%
64	23.898	3566	3572	3581	rBV	481177	965893	1.91%	0.351%
65	24.639	3692	3698	3709	rBV	291399	654758	1.30%	0.238%
66	25.509	3841	3846	3857	rVB2	129212	314575	0.62%	0.114%

Sum of corrected areas: 274950162

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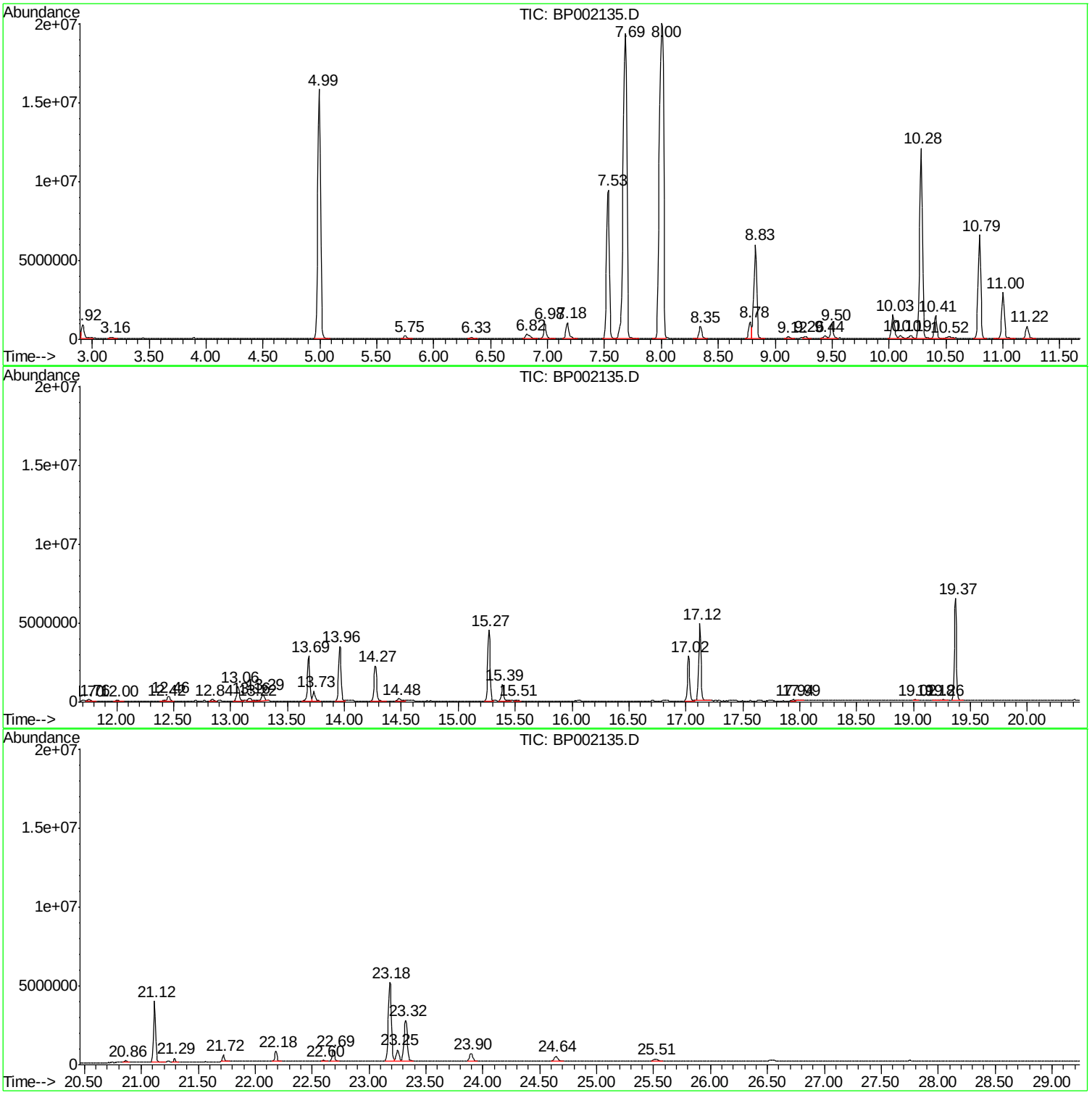
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0DL1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

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Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
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Instrument : BNA_P
ClientSampleId : C0DL1

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Quant Title : SVOA CALIBRATION

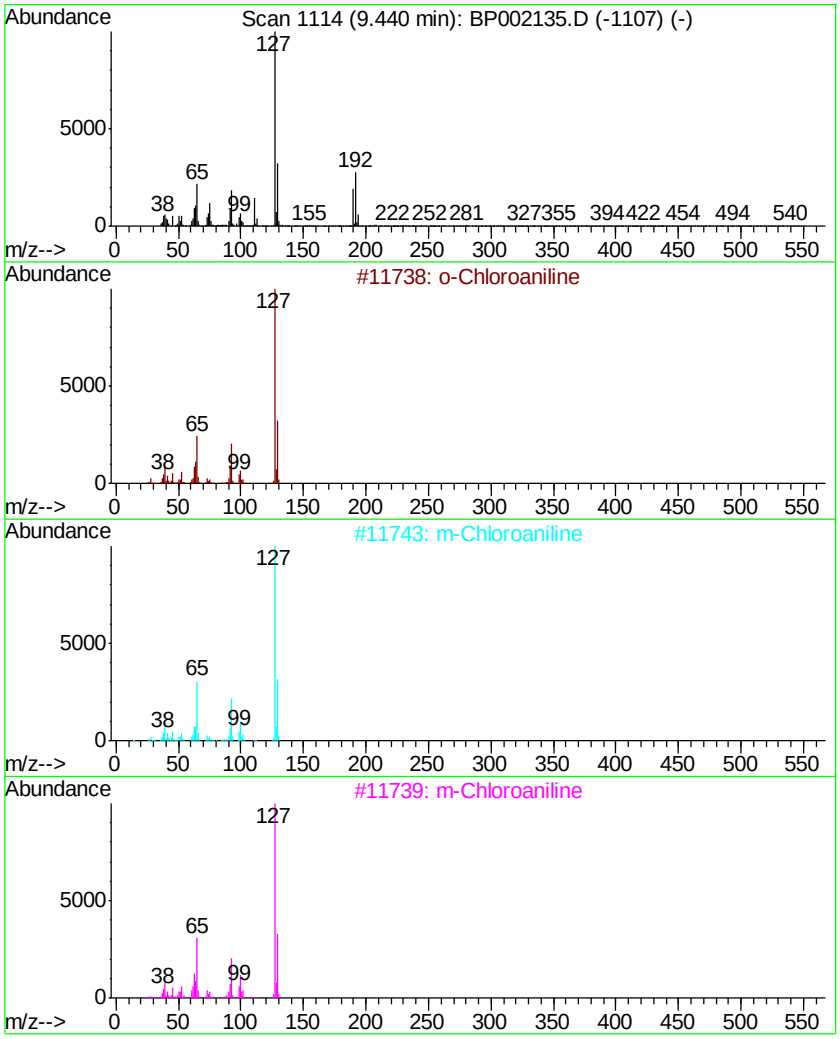
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

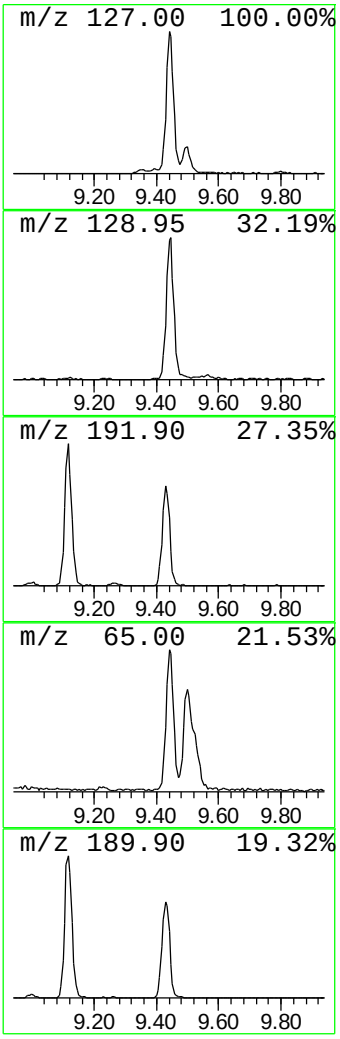
****

Peak Number 4 o-Chloroaniline Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.44	2.06 ng/ul	243548	Naphthalene-d8	10.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Chloroaniline	127	C6H6ClN	000095-51-2	95
2	m-Chloroaniline	127	C6H6ClN	000108-42-9	86
3	m-Chloroaniline	127	C6H6ClN	000108-42-9	86
4	m-Chloroaniline	127	C6H6ClN	000108-42-9	70
5	o-Chloroaniline	127	C6H6ClN	000095-51-2	70





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0DL1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

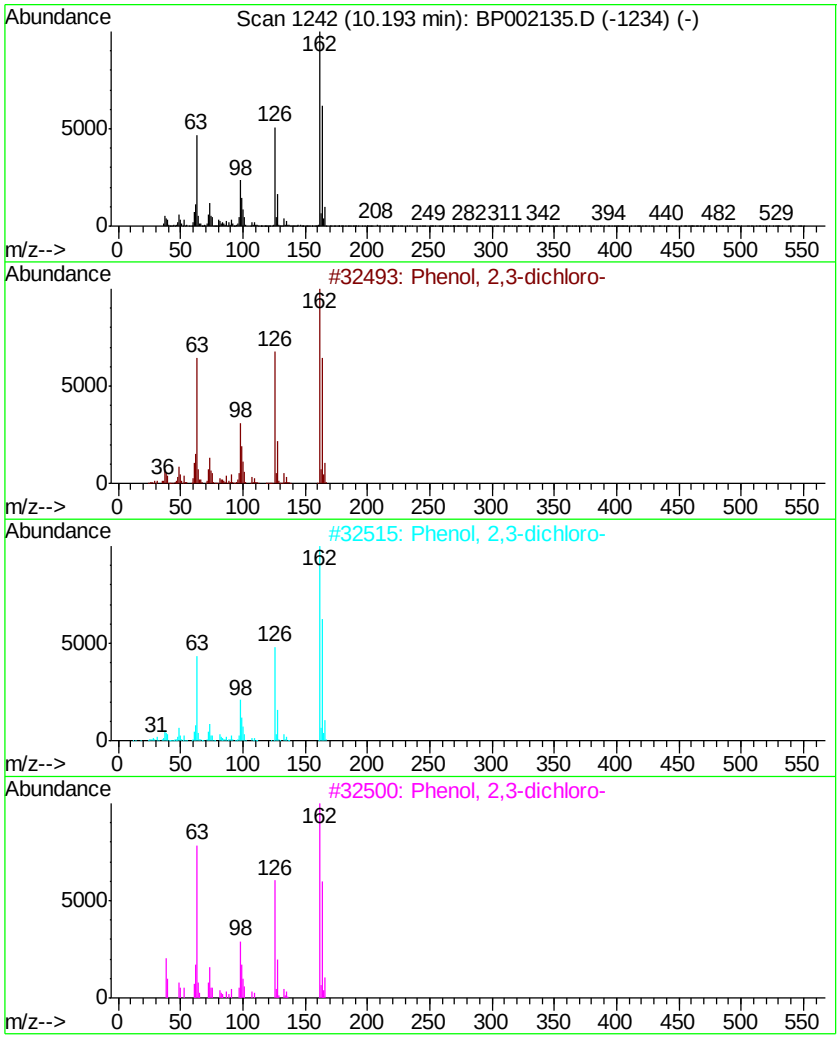
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

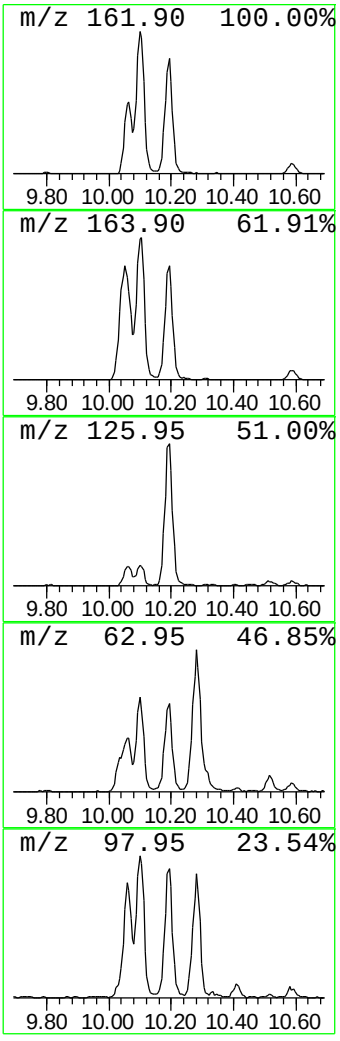
****

Peak Number 5 Phenol, 2,3-dichloro- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.19	2.81 ng/ul	332705	Naphthalene-d8	10.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	98
2	Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	98
3	Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	97
4	Phenol, 2,6-dichloro-	162	C6H4Cl2O	000087-65-0	97
5	Phenol, 2,6-dichloro-	162	C6H4Cl2O	000087-65-0	97





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0DL1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

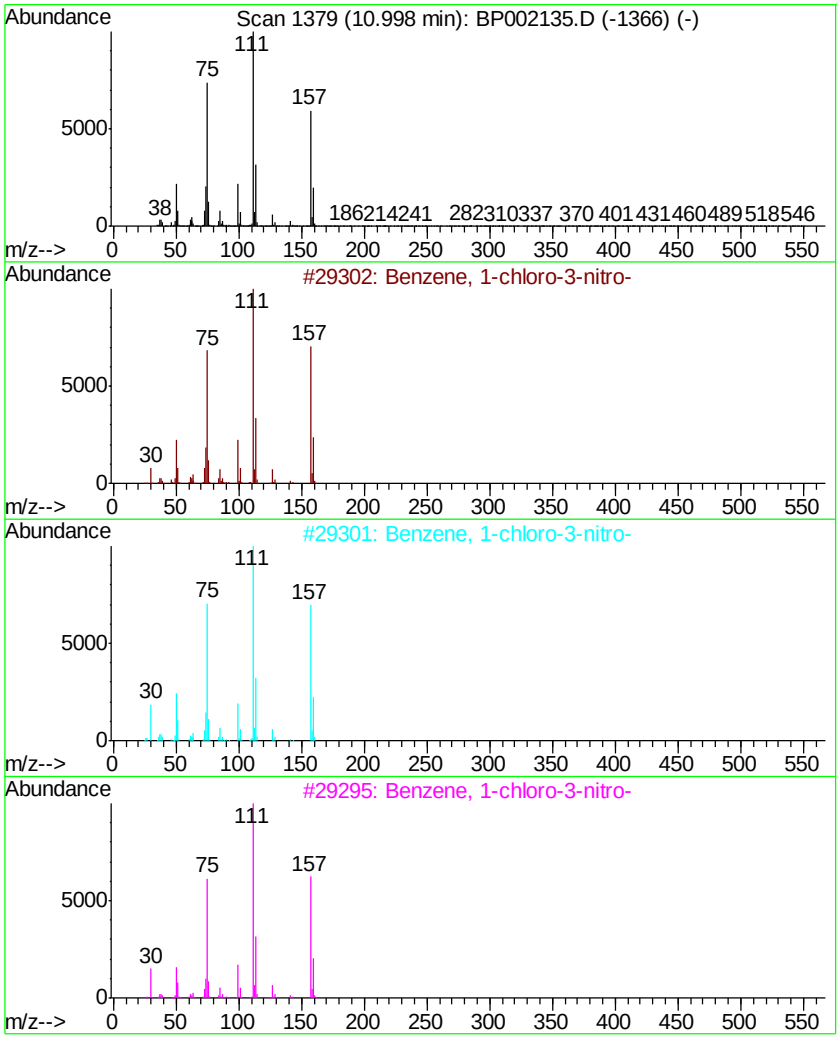
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

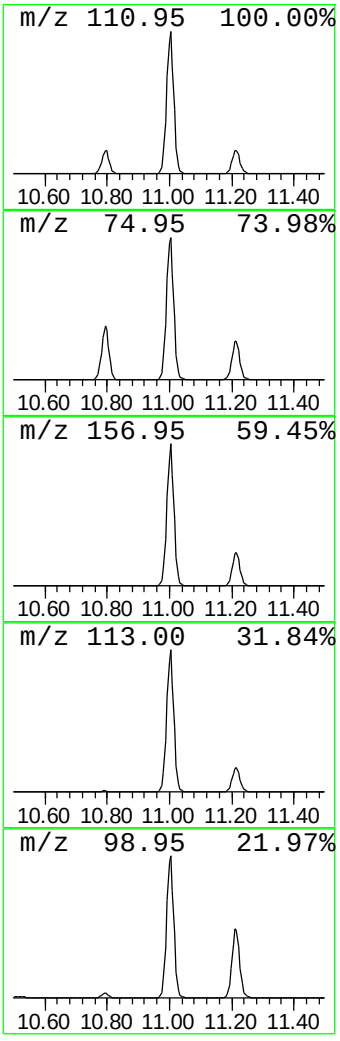
****

Peak Number 8 Benzene, 1-chloro-3-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.00	41.18 ng/ul	4875080	Naphthalene-d8	10.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	98
2	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	97
3	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	96
4	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
5	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	90





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0DL1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

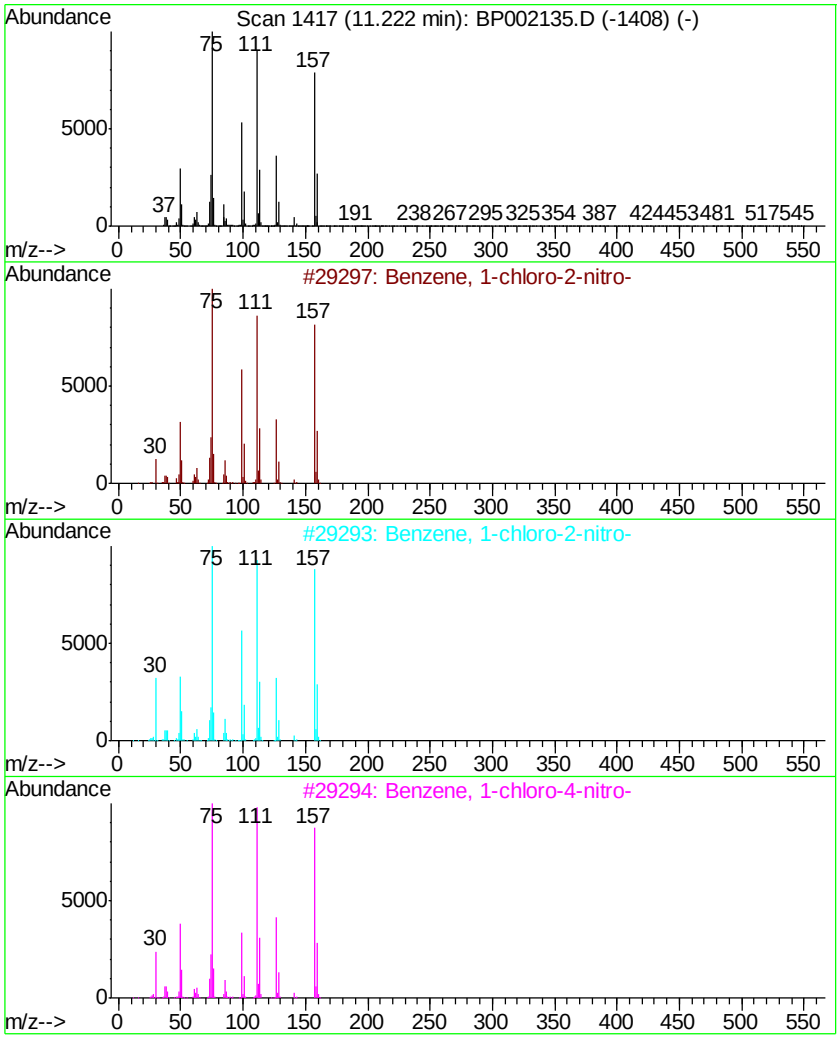
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

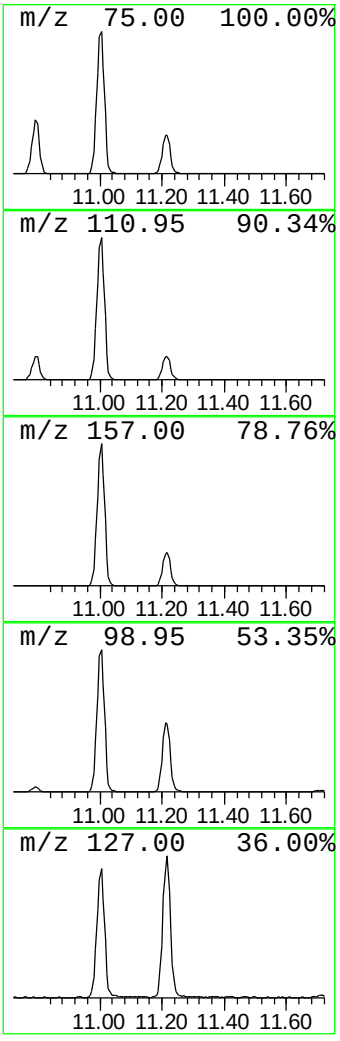
****

Peak Number 9 Benzene, 1-chloro-2-nitro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.22	11.55 ng/ul	1367150	Naphthalene-d8	10.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	99
2	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	98
3	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	98
4	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
5	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0DL1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

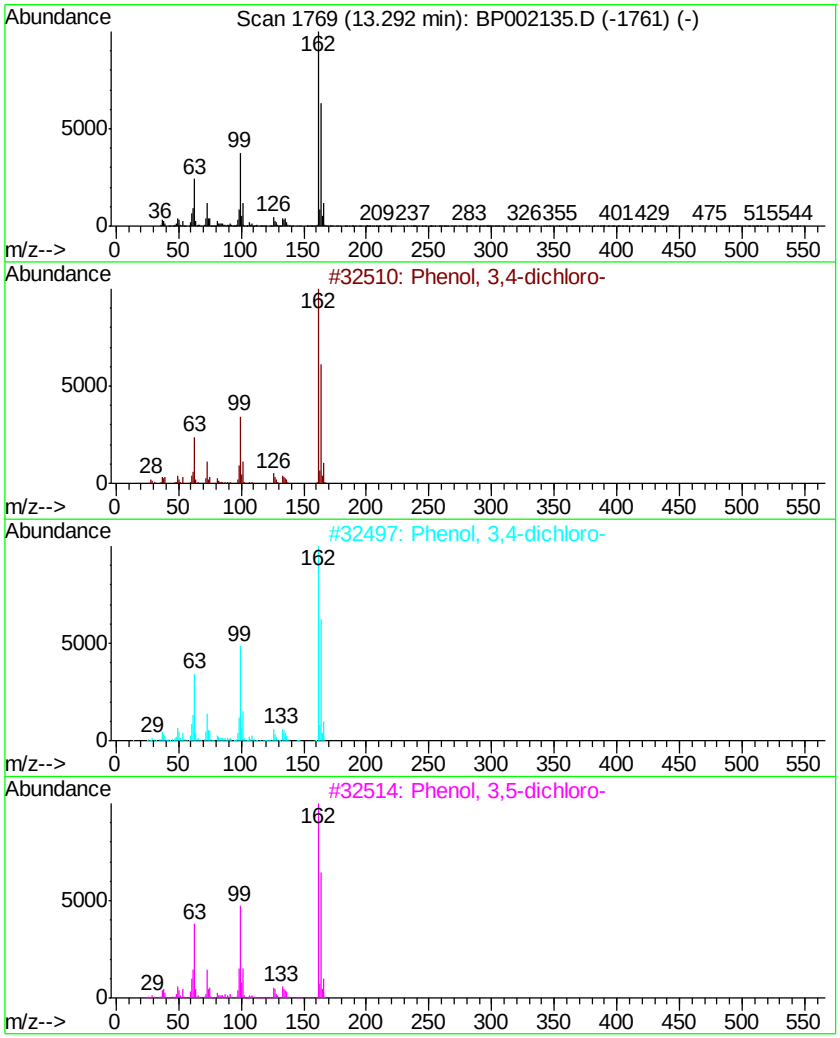
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

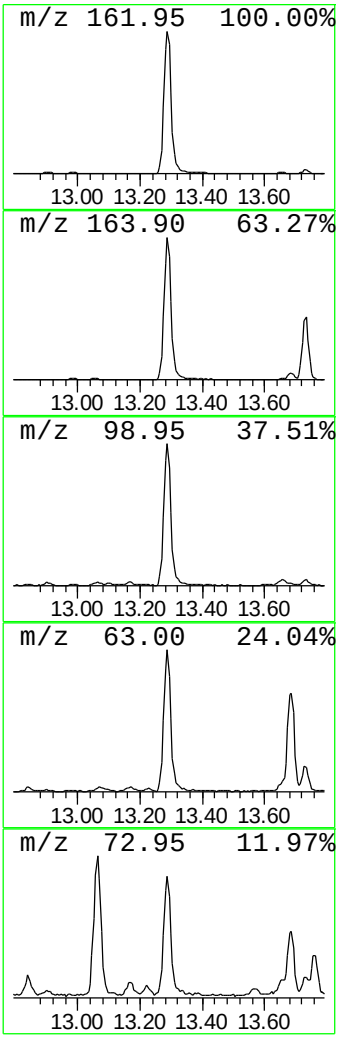
****

Peak Number 10 Phenol, 3,4-dichloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	4.23 ng/ul	721976	Acenaphthene-d10	14.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	98
2	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
3	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	97
4	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	96
5	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	95





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Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0DL1

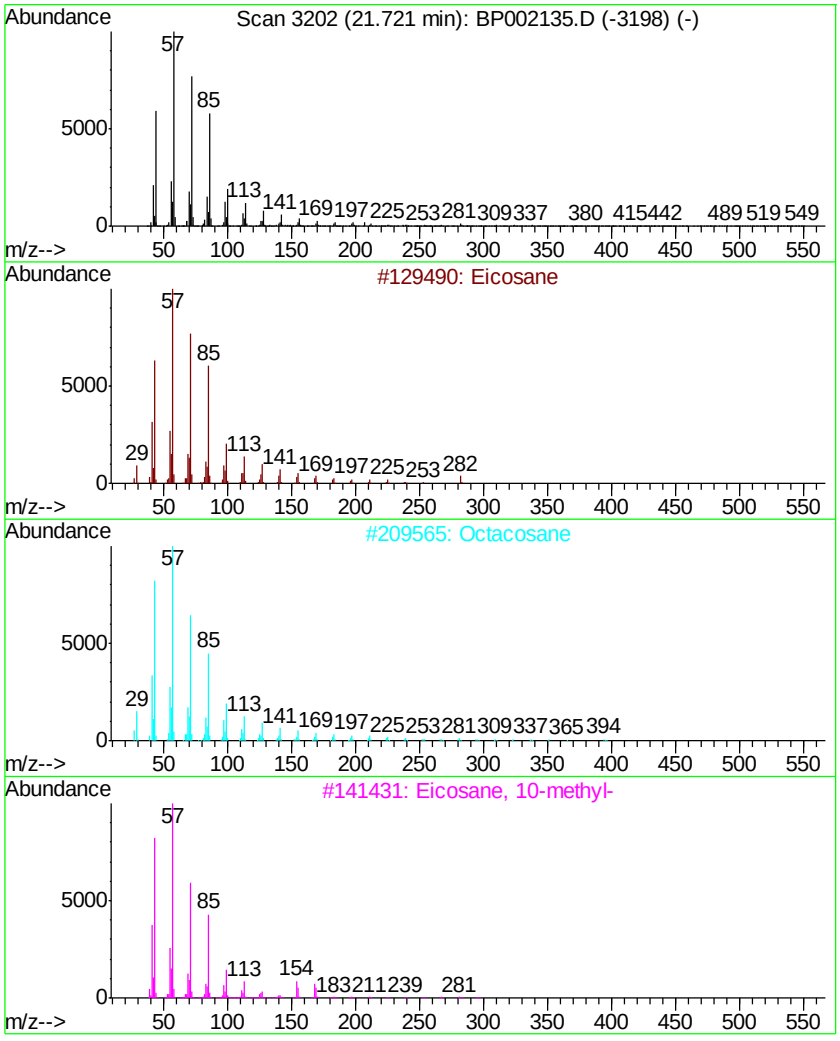
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

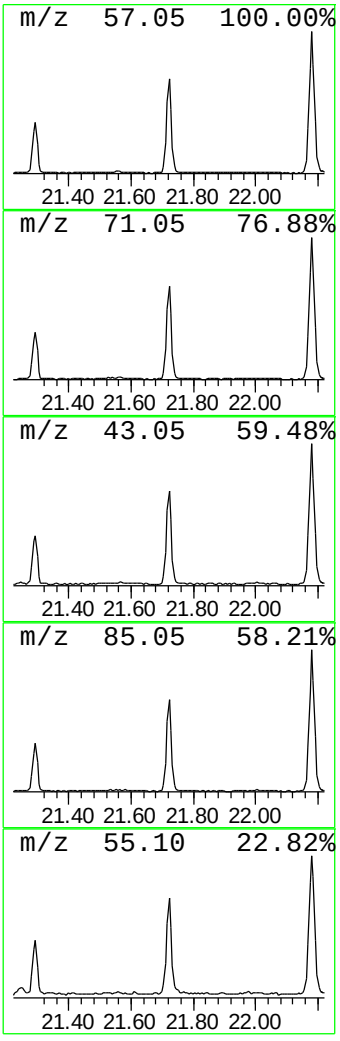
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.72	2.13 ng/ul	510027	Chrysene-d12	21.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	98
2	Octacosane	394	C28H58	000630-02-4	98
3	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4	Hexadecane	226	C16H34	000544-76-3	93
5	Heptacosane	380	C27H56	000593-49-7	93





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0DL1

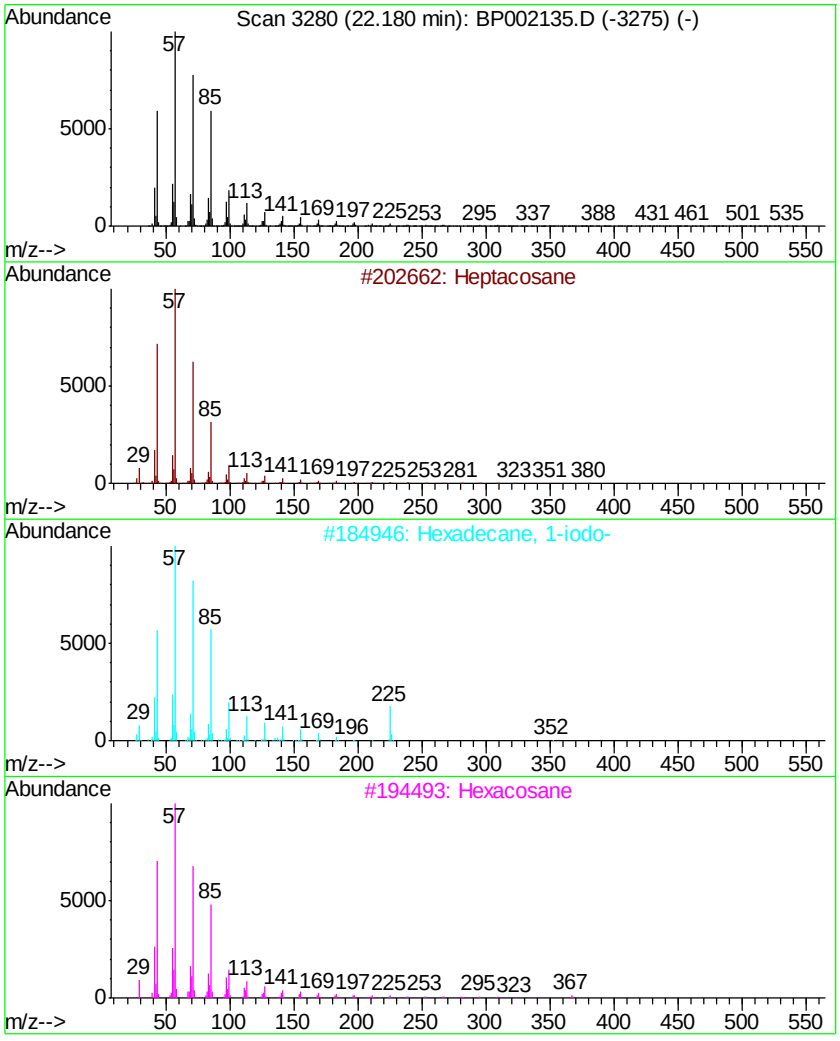
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

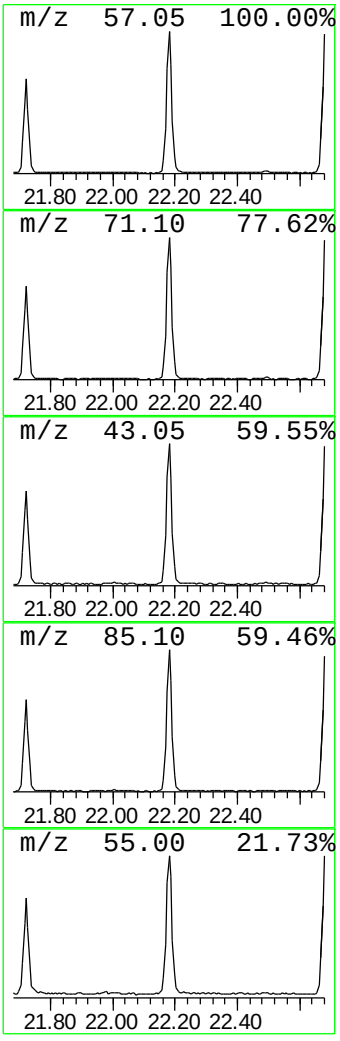
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.18	3.34 ng/ul	801176	Chrysene-d12	21.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	98
2	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	95
3	Hexacosane	366	C26H54	000630-01-3	94
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
5	Pentacosane	352	C25H52	000629-99-2	93





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0DL1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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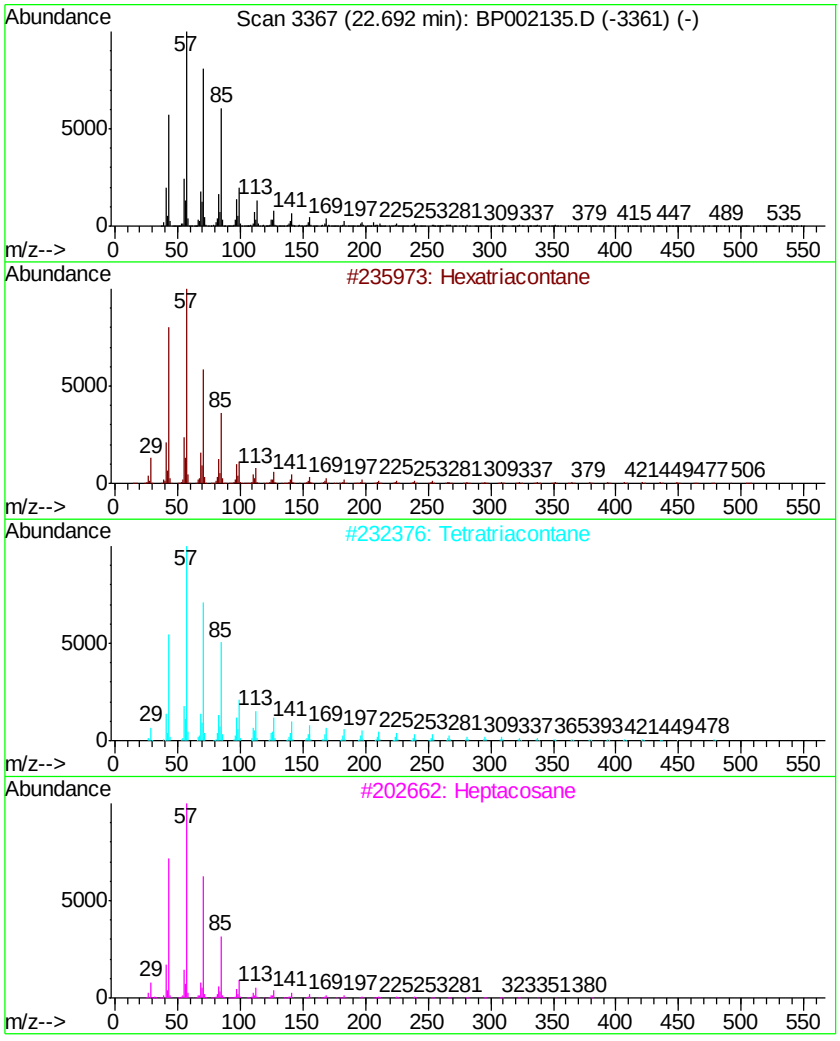
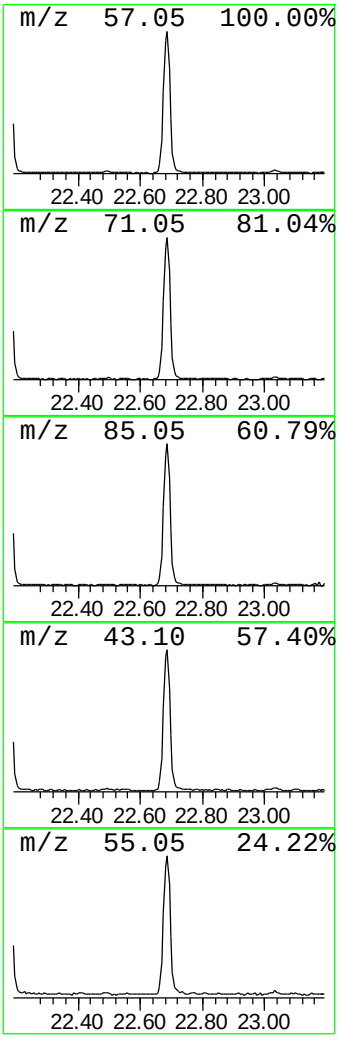
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.69	4.02 ng/ul	1019910	Perylene-d12	23.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexatriacontane	507	C36H74	000630-06-8	93
2	Tetratriacontane	479	C34H70	014167-59-0	93
3	Heptacosane	380	C27H56	000593-49-7	91
4	Octadecane, 2-methyl-	268	C19H40	001560-88-9	91
5	Heneicosane	296	C21H44	000629-94-7	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0DL1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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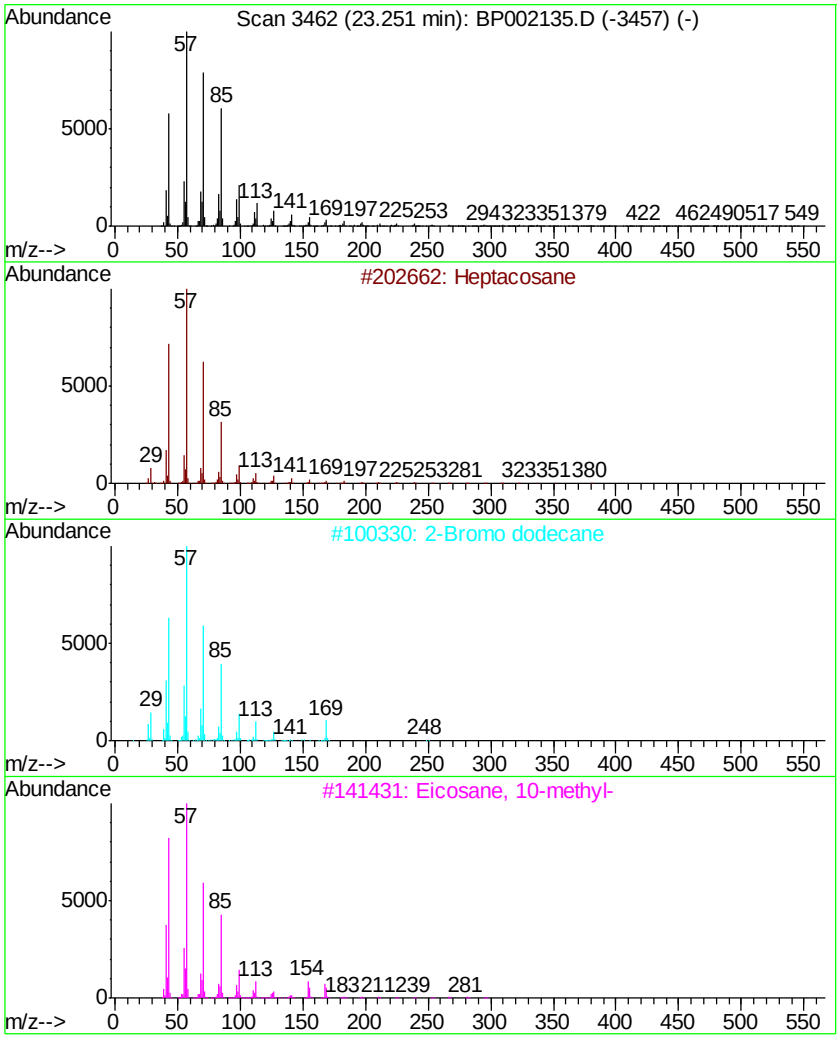
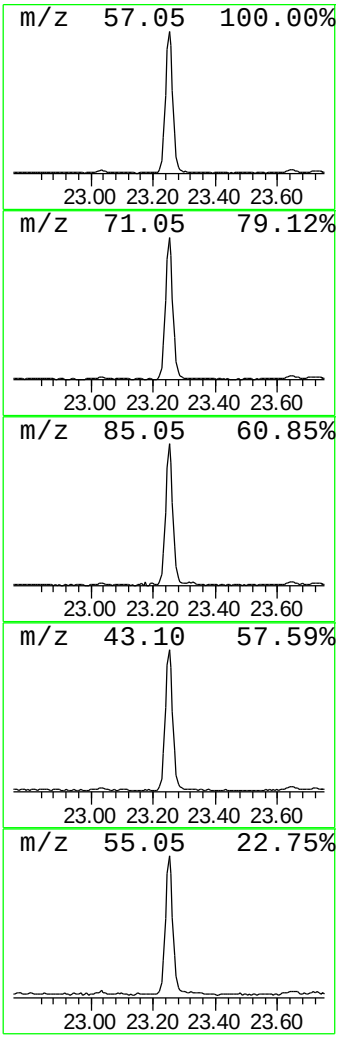
Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.25	4.77 ng/ul	1211440	Perylene-d12	23.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	95
2	2-Bromo dodecane	248	C12H25Br	013187-99-0	95
3	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4	Heneicosane	296	C21H44	000629-94-7	91
5	2-methyloctacosane	408	C29H60	1000376-72-8	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0DL1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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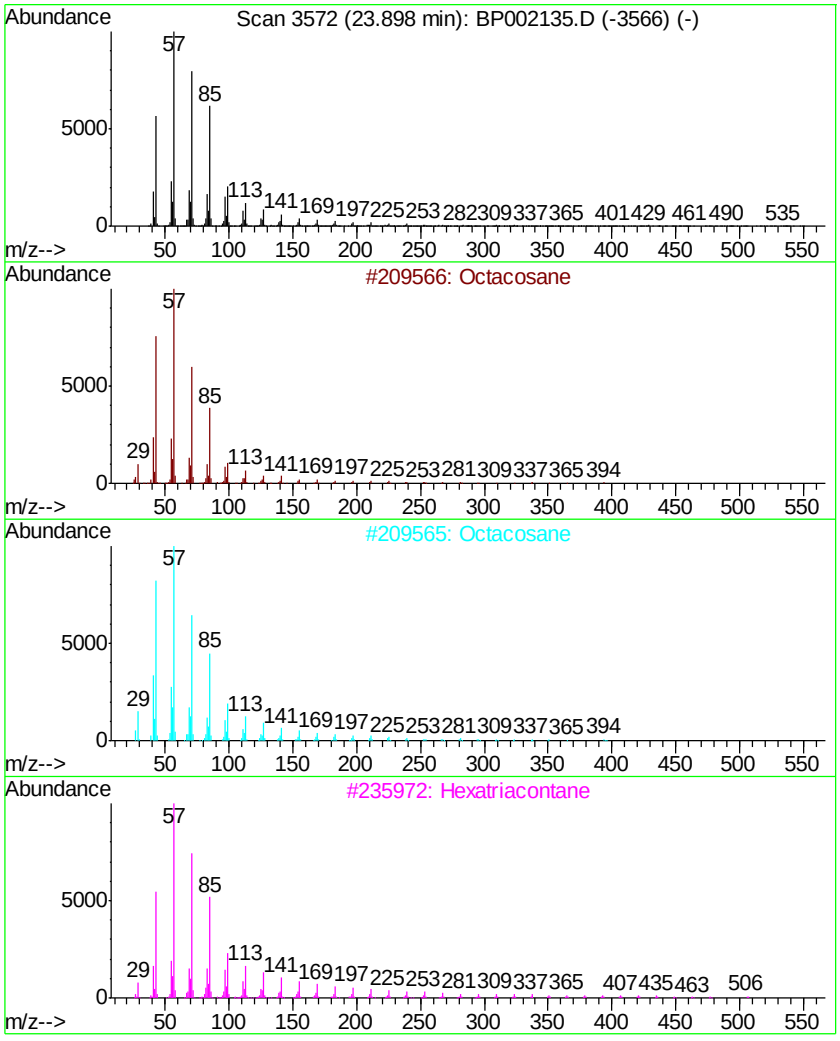
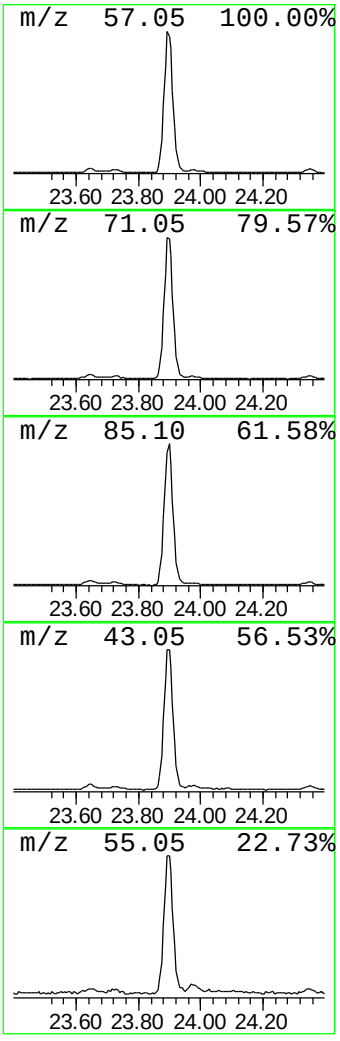
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.90	3.80 ng/ul	965893	Perylene-d12	23.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	98
2	Octacosane	394	C28H58	000630-02-4	93
3	Hexatriacontane	507	C36H74	000630-06-8	93
4	Hexadecane	226	C16H34	000544-76-3	93
5	Hexacosane	366	C26H54	000630-01-3	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0DL1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

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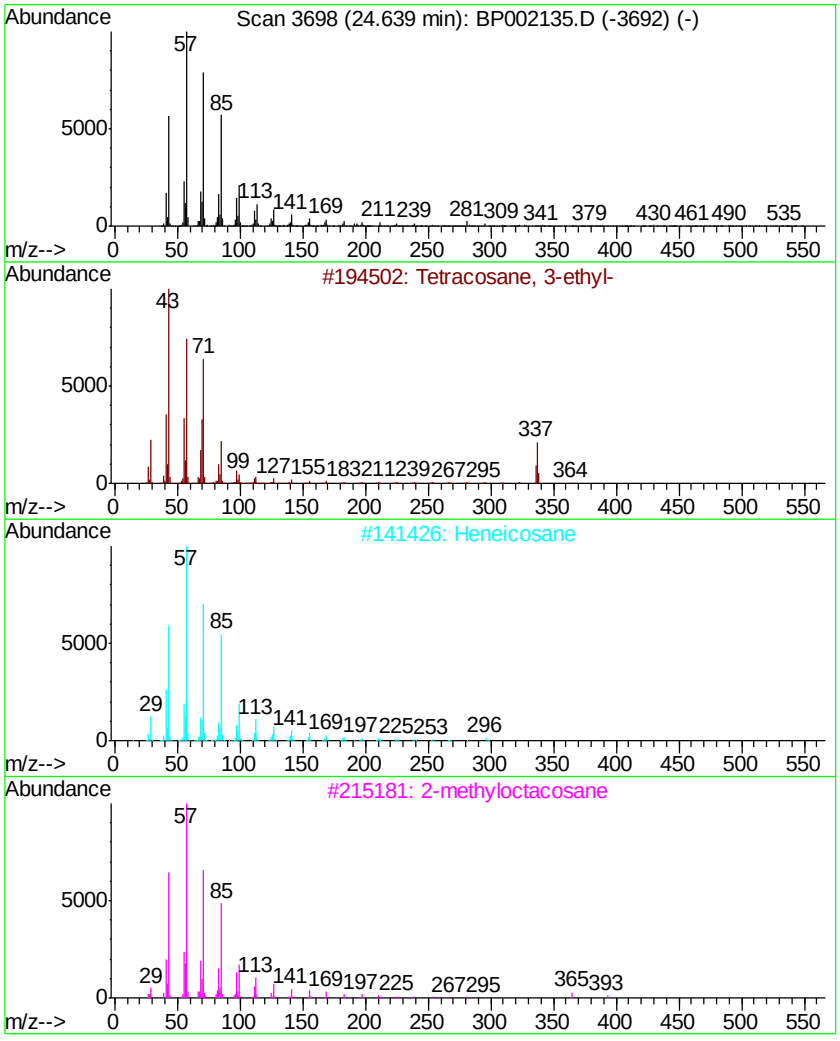
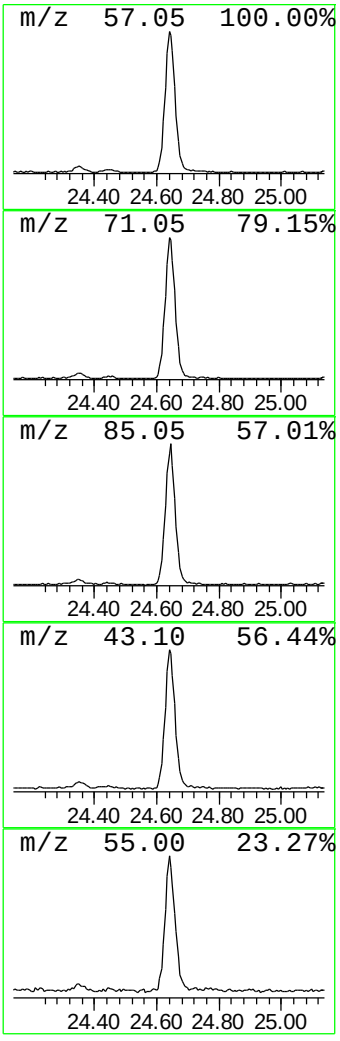
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.64	2.58 ng/ul	654758	Perylene-d12	23.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane, 3-ethyl-	366	C26H54	055282-17-2	93
2	Heneicosane	296	C21H44	000629-94-7	91
3	2-methyloctacosane	408	C29H60	1000376-72-8	91
4	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
5	Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030920\
Data File : BP002135.D
Acq On : 09 Mar 2020 23:28
Operator : CG/JU
Sample : L1834-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C0DL1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.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
o-Chloroaniline	9.44	2.1	ng/ul	243548	2	10.41	2367720	20.0
Phenol, 2,3-dichl...	10.19	2.8	ng/ul	332705	2	10.41	2367720	20.0
Benzene, 1-chloro...	11.00	41.2	ng/ul	4875080	2	10.41	2367720	20.0
Benzene, 1-chloro...	11.22	11.6	ng/ul	1367150	2	10.41	2367720	20.0
Phenol, 3,4-dichl...	13.29	4.2	ng/ul	721976	3	14.27	3411500	20.0
(DEL) Alkane: Str...	21.72	2.1	ng/ul	510027	5	21.12	4798440	20.0
(DEL) Alkane: Str...	22.18	3.3	ng/ul	801176	5	21.12	4798440	20.0
(DEL) Alkane: Str...	22.69	4.0	ng/ul	1019910	6	23.32	5078810	20.0
(DEL) Alkane: Str...	23.25	4.8	ng/ul	1211440	6	23.32	5078810	20.0
(DEL) Alkane: Str...	23.90	3.8	ng/ul	965893	6	23.32	5078810	20.0
(DEL) Alkane: Str...	24.64	2.6	ng/ul	654758	6	23.32	5078810	20.0