LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-1

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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\8270-BP101920.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP003743.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.916	3	5	15	rVB2	91173	168486	2.33%	0.447%
2	3.087	29	34	53	rVB	1896693	2384934	32.97%	6.325%
3	3.252	56	62	73	rVV	173880	351412	4.86%	0.932%
4	3.346	73	78	89	rVB	214518	290337	4.01%	0.770%
5	5.475	434	440	447	rBV	188843	273318	3.78%	0.725%
6	6.010	521	531	541	rBV	1683229	2532541	35.01%	6.716%
7	7.663	803	812	820	rBV	1719223	2772404	38.32%	7.352%
8	8.510	948	956	963	rBV	403644	646470	8.94%	1.714%
9	9.675	1146	1154	1165	rVB	1189266	1974739	27.30%	5.237%
10	11.351	1432	1439	1446	rBV	503141	839948	11.61%	2.227%
11	13.734	1837	1844	1856	rVB	3116847	4551982	62.93%	12.072%
12	14.522	1973	1978	1988	rVB	280944	406491	5.62%	1.078%
13	15.110	2071	2078	2085	rVB	761134	1157254	16.00%	3.069%
14	16.580	2321	2328	2343	rBV	1918258	2715945	37.54%	7.202%
15	17.180	2425	2430	2440	rVB2	85136	146009	2.02%	0.387%
16	17.827	2534	2540	2545	rBV	1065476	1429233	19.76%	3.790%
17	18.639	2674	2678	2685	rBV2	62835	110688	1.53%	0.294%
18	18.704	2686	2689	2695	rVB	56629	85370	1.18%	0.226%
19	18.904	2719	2723	2726	rBV	164769	220760	3.05%	0.585%
20	18.945	2726	2730	2741	rVB	226317	351637	4.86%	0.933%
21	19.380	2793	2804	2817	rVB5	147041	601489	8.31%	1.595%
22	20.298	2954	2960	2965	rBV	6201487	7233944	100.00%	19.184%
23	20.780	3035	3042	3053	rVB2	248345	666703	9.22%	1.768%
24	20.910	3060	3064	3073	rVB5	63742	134557	1.86%	0.357%
25	21.027	3081	3084	3087	rVB2	74831	82436	1.14%	0.219%
26	21.468	3155	3159	3169	rVB	584032	721759	9.98%	1.914%
27	21.792	3209	3214	3218	rBV2	168840	374723	5.18%	0.994%
28	21.839	3218	3222	3228	rVB	1197355	1650823	22.82%	4.378%
29	21.927	3233	3237	3242	rVB	140105	161308	2.23%	0.428%
30	22.410	3315	3319	3327	rBV2	123287	199353	2.76%	0.529%
31	23.092	3432	3435	3441	rVB	93496	140382	1.94%	0.372%
32	23.510	3502	3506	3512	rVB	123518	195608	2.70%	0.519%
33	24.168	3614	3618	3632	rVB	156716	332586	4.60%	0.882%
34	24.415	3652	3660	3668	rVB2	768901	1651875	22.84%	4.381%
35	24.921	3743	3746	3754	rVB	80233	150898	2.09%	0.400%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP-1

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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\8270-BP101920.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

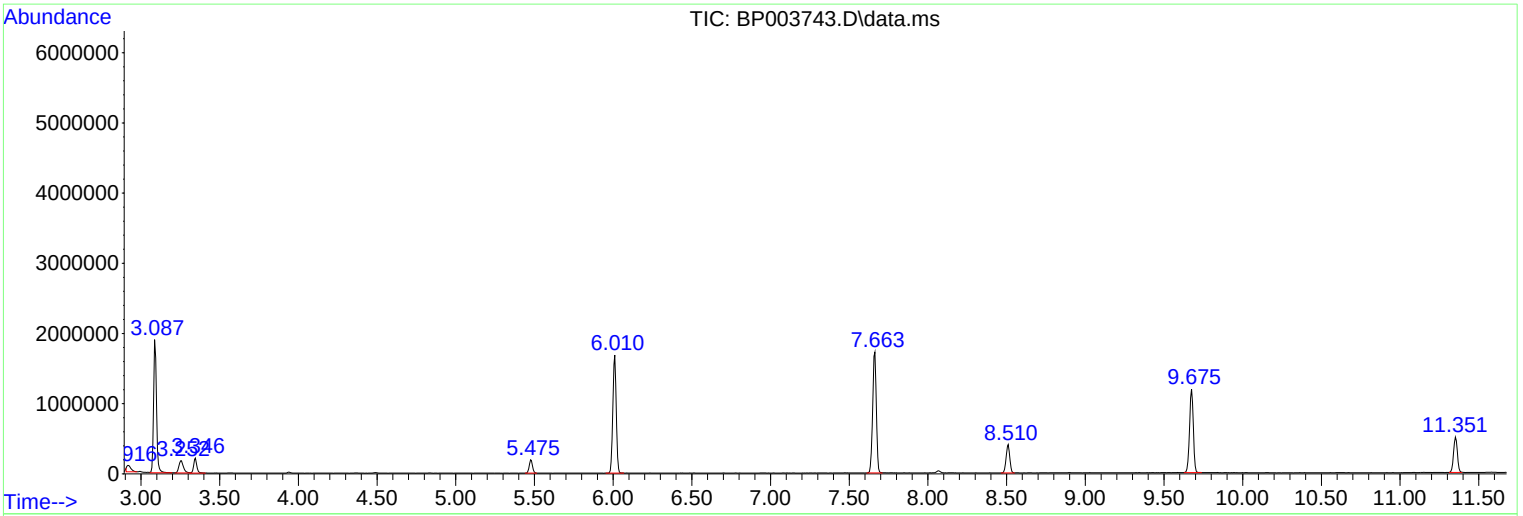
Sum of corrected areas: 37708402

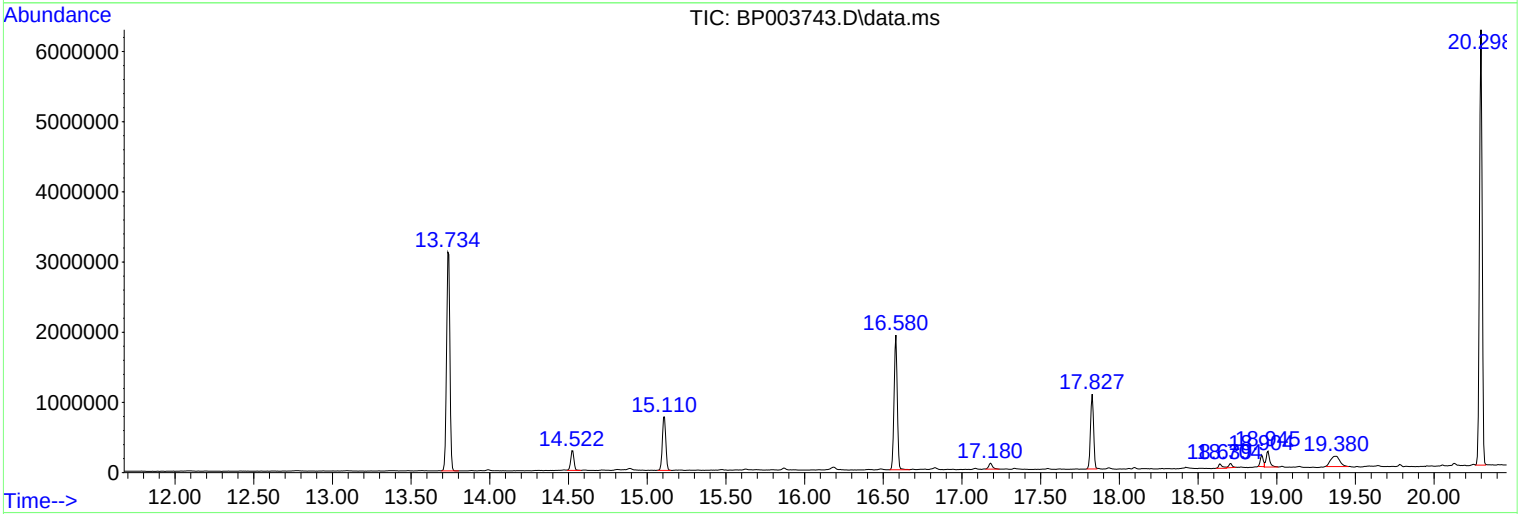
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

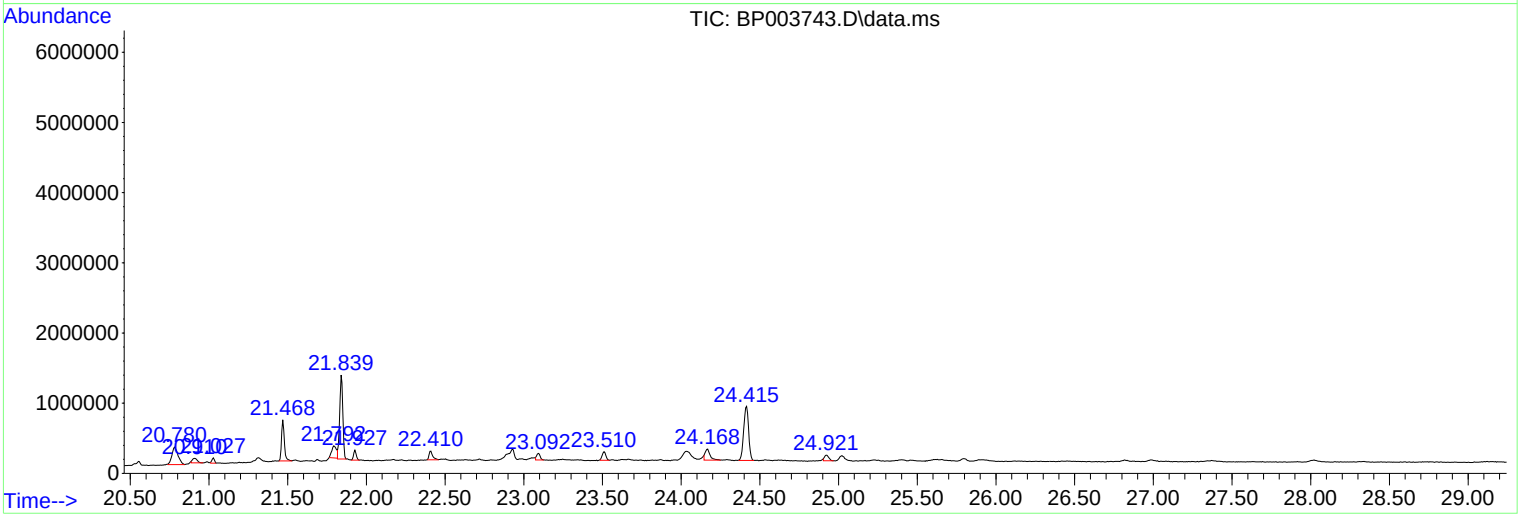
Instrument :
BNA_P
ClientSampleId :
COMP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP-1

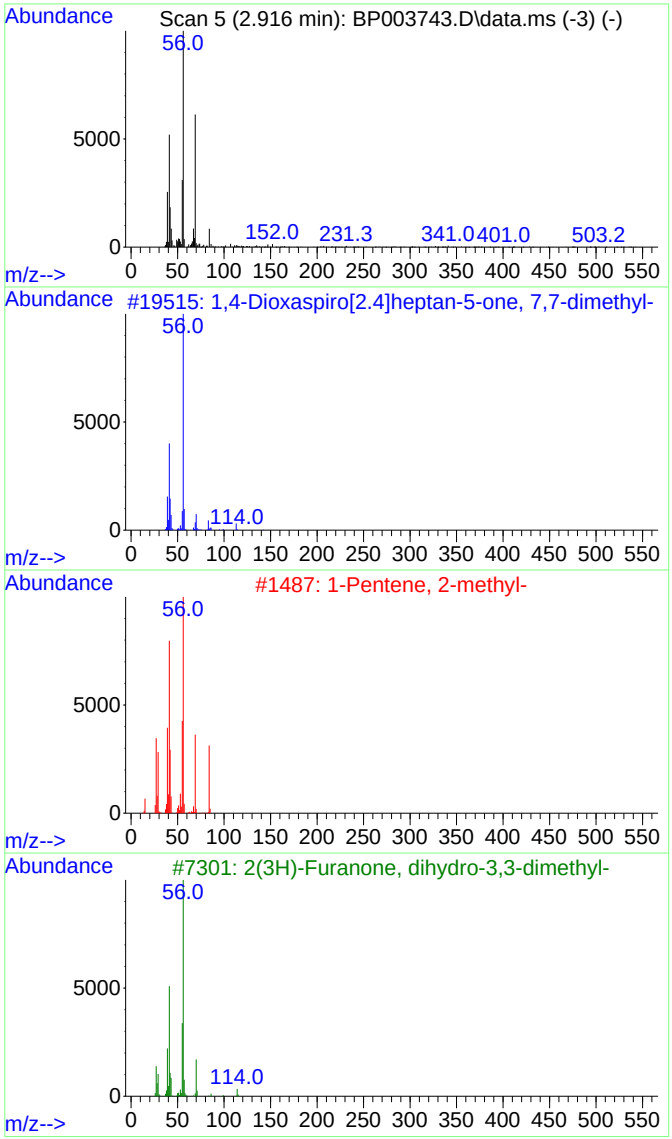
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

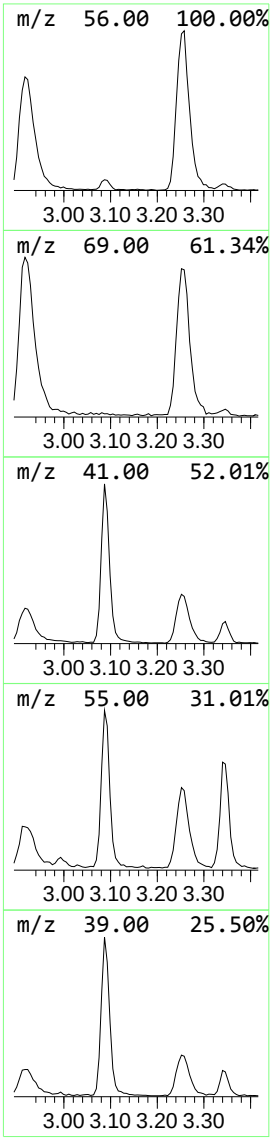
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 1,4-Dioxaspiro[2.4]heptan-5... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.916	5.21 ng	168486	1,4-Dichlorobenzene-d4	8.510

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Dioxaspiro[2.4]heptan-5-one,...	142	C7H10O3	126091-79-0	59
2	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	59
3	2(3H)-Furanone, dihydro-3,3-dime...	114	C6H10O2	003709-08-8	59
4	3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	50
5	Cyclopentane, methyl-	84	C6H12	000096-37-7	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP-1

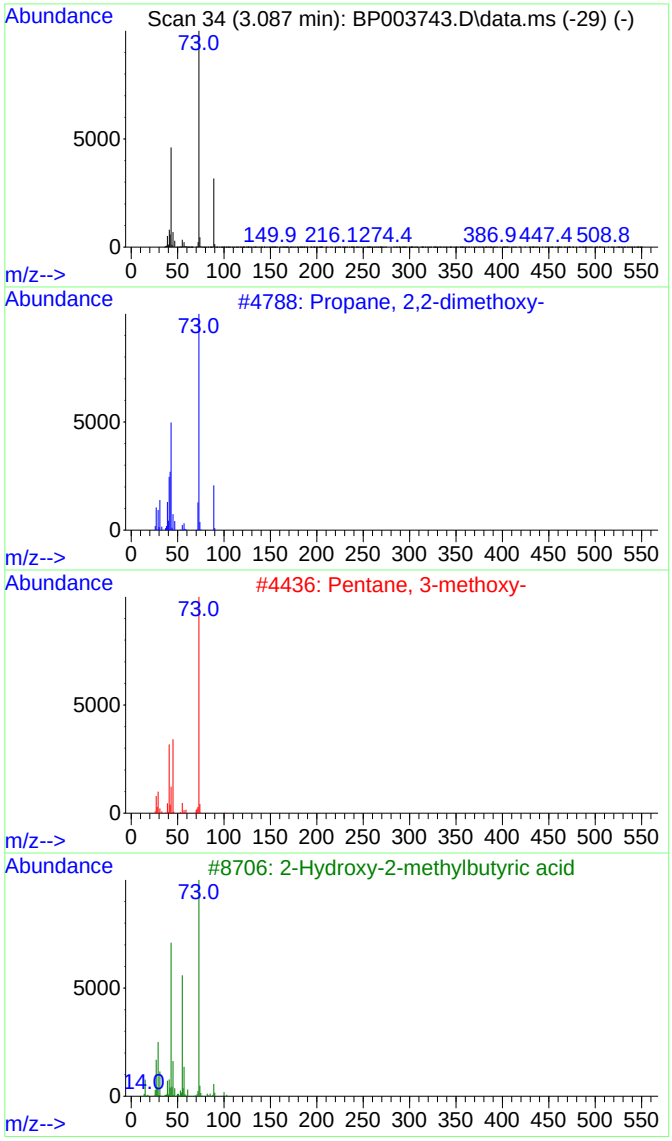
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

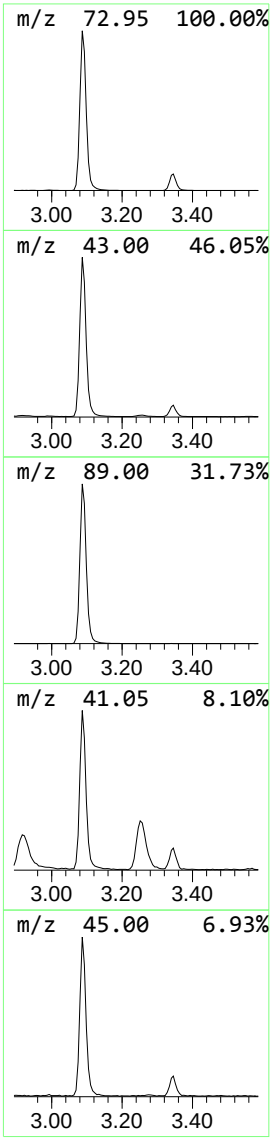
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 unknown3.08 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.087	73.78 ng	2384930	1,4-Dichlorobenzene-d4	8.510

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	45
2	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
3	2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
4	Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
5	Methane, isothiocyanato-	73	C2H3NS	000556-61-6	7





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP-1

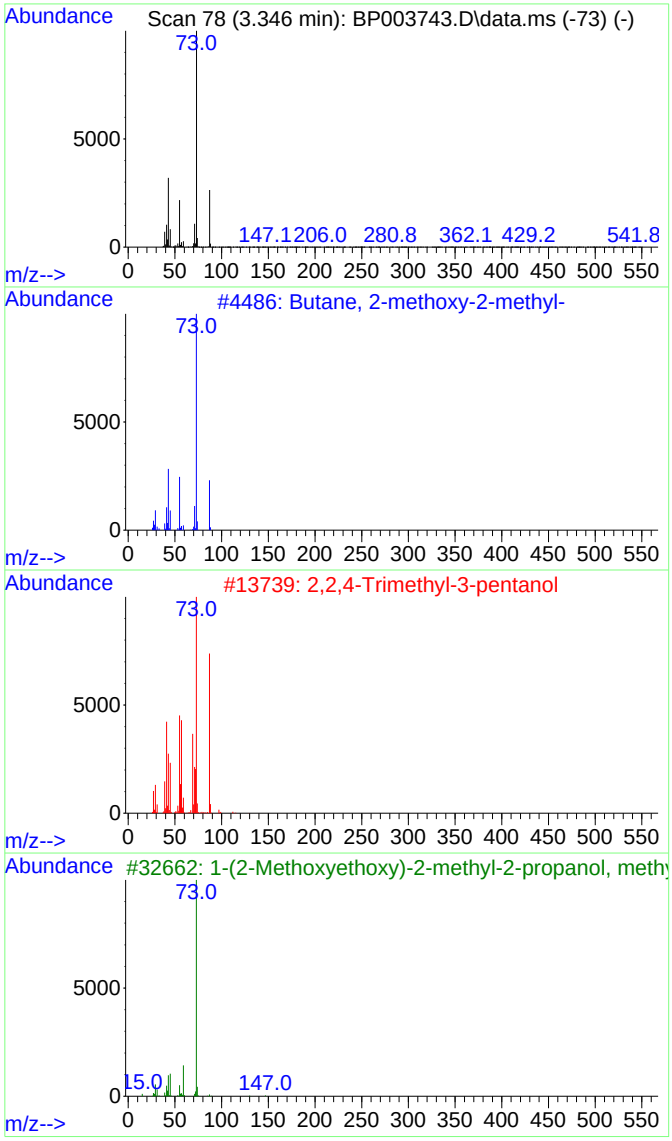
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

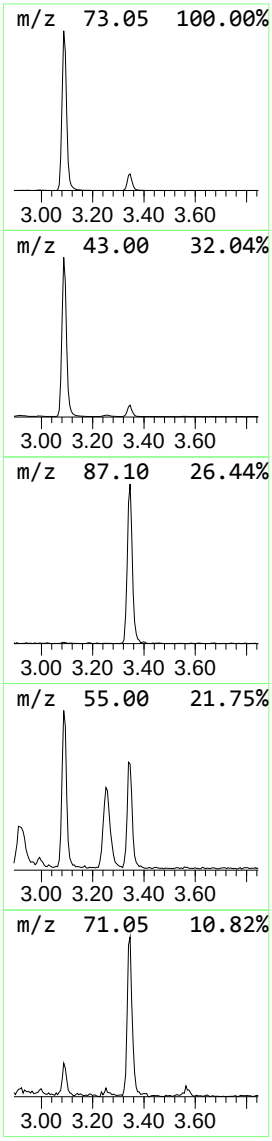
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.346	8.98 ng	290337	1,4-Dichlorobenzene-d4	8.510

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	56
3	1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	17
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP-1

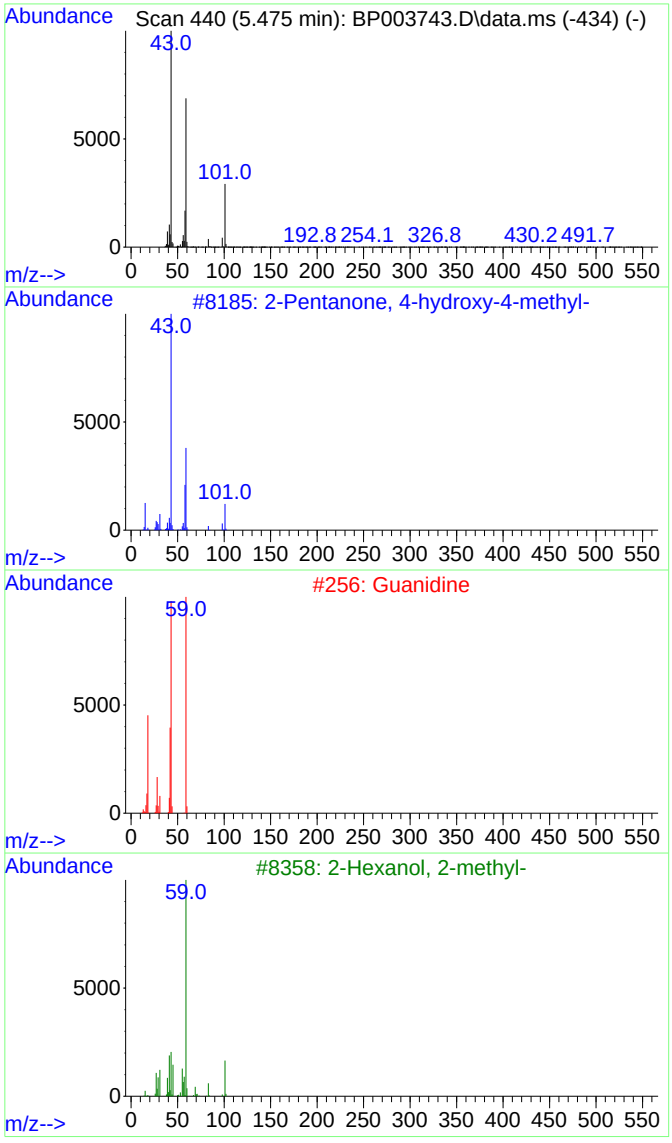
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

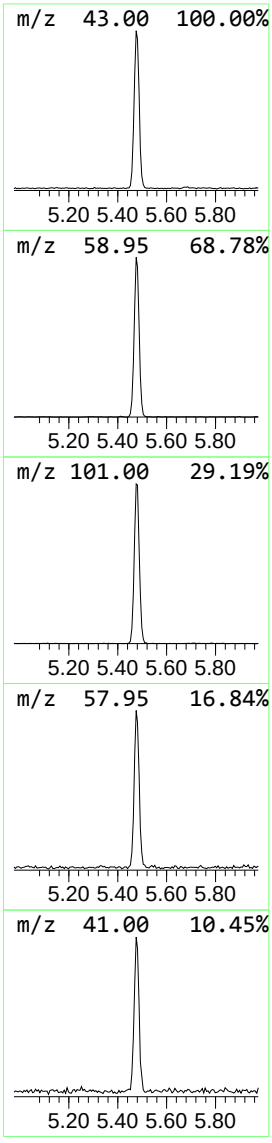
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.475	8.46 ng	273318	1,4-Dichlorobenzene-d4	8.510

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Guanidine	59	CH5N3	000113-00-8	43
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
5	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	12





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP-1

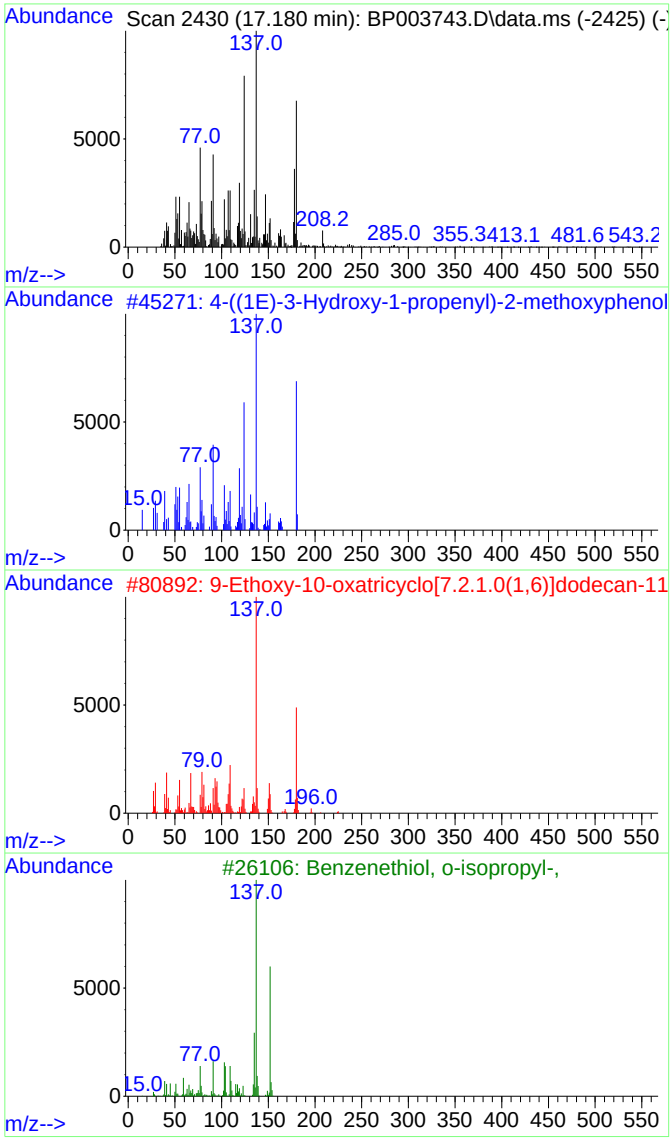
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

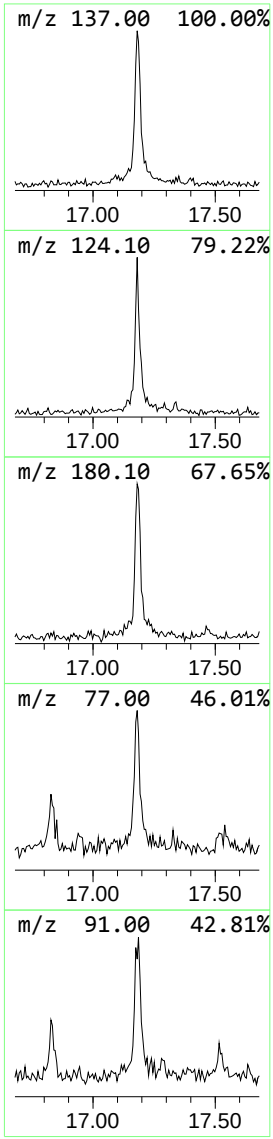
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 4-((1E)-3-Hydroxy-1-propeny... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.180	2.04 ng	146009	Phenanthrene-d10	17.828

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-((1E)-3-Hydroxy-1-propenyl)-2-...	180	C10H12O3	1000297-95-5	90
2	9-Ethoxy-10-oxatricyclo[7.2.1.0(...	224	C13H20O3	1000187-02-8	46
3	Benzenethiol, o-isopropyl-,	152	C9H12S	006262-87-9	42
4	Propenoic acid, 3-(2-thienyl)-, ...	464	C13H7Br3O2S	284665-05-0	35
5	4,6-Dimethyl-2-pyrimidone	124	C6H8N2O	000108-79-2	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP-1

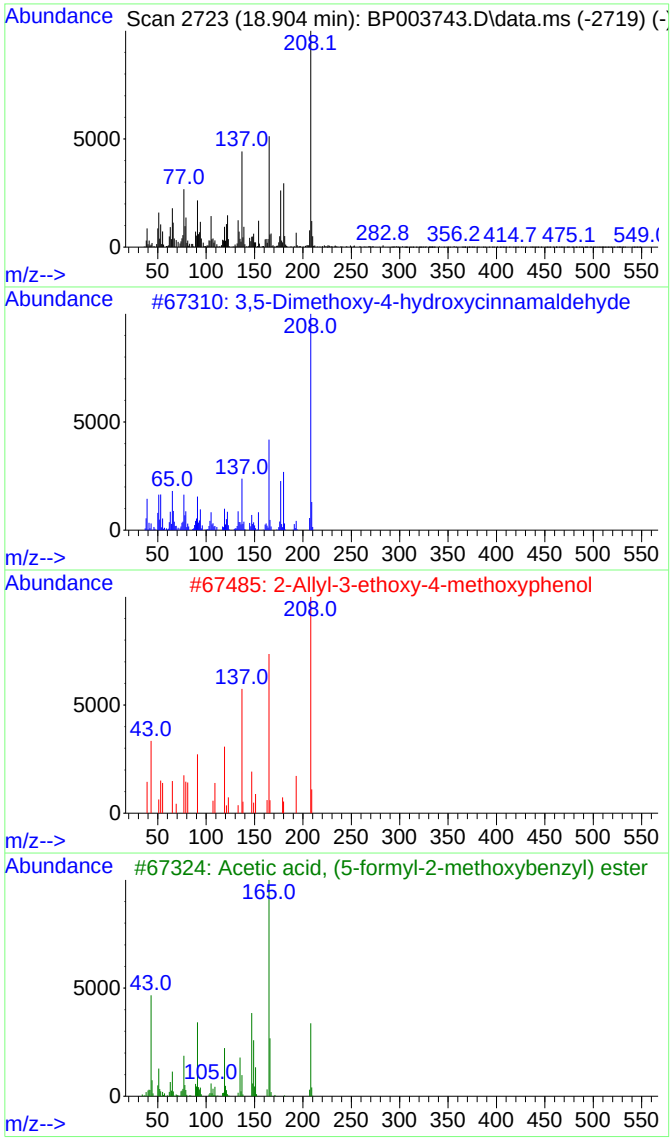
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

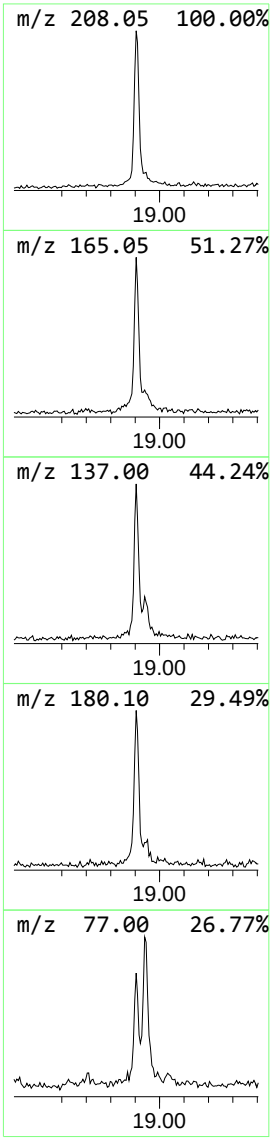
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 3,5-Dimethoxy-4-hydroxycinn... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.904	3.09 ng	220760	Phenanthrene-d10	17.828

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Dimethoxy-4-hydroxycinnamald...	208	C11H12O4	087345-53-7	93
2	2-Allyl-3-ethoxy-4-methoxyphenol	208	C12H16O3	1000122-70-6	58
3	Acetic acid, (5-formyl-2-methoxy...	208	C11H12O4	099865-67-5	55
4	Benzo[c]cinnoline, 4,7-dimethyl-	208	C14H12N2	020684-54-2	43
5	5-(4-Methoxy-phenyl)-[1,3,4]oxad...	208	C9H8N2O2S	023766-26-9	42





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

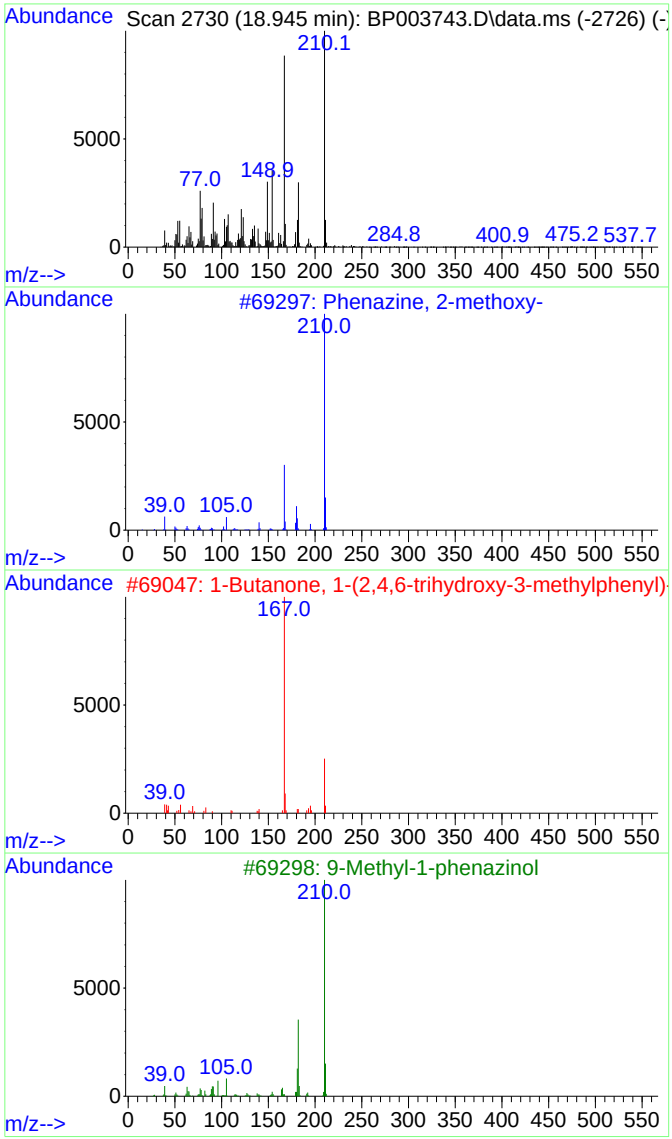
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

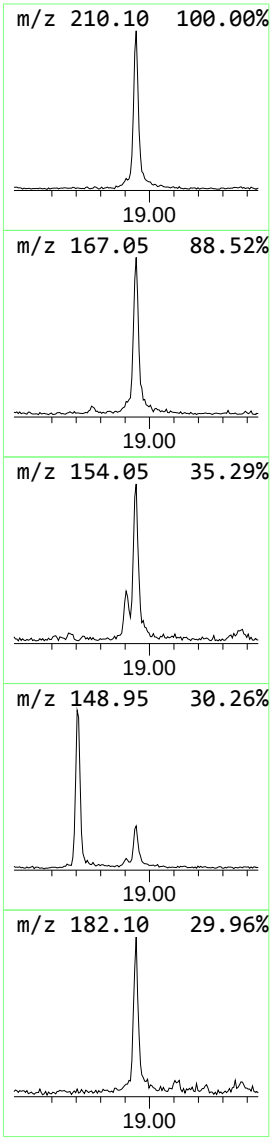
****

Peak Number 8 Phenazine, 2-methoxy- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.945	4.92 ng	351637	Phenanthrene-d10	17.828

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenazine, 2-methoxy-	210	C13H10N2O	002876-18-8	58
2	1-Butanone, 1-(2,4,6-trihydroxy-...	210	C11H14O4	001509-06-4	50
3	9-Methyl-1-phenazinol	210	C13H10N2O	013860-43-0	38
4	Benzene, 1,1'-ethylidenebis-	182	C14H14	000612-00-0	30
5	Ethanone, 1-(2,6-dihydroxy-4-met...	182	C9H10O4	007507-89-3	30





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

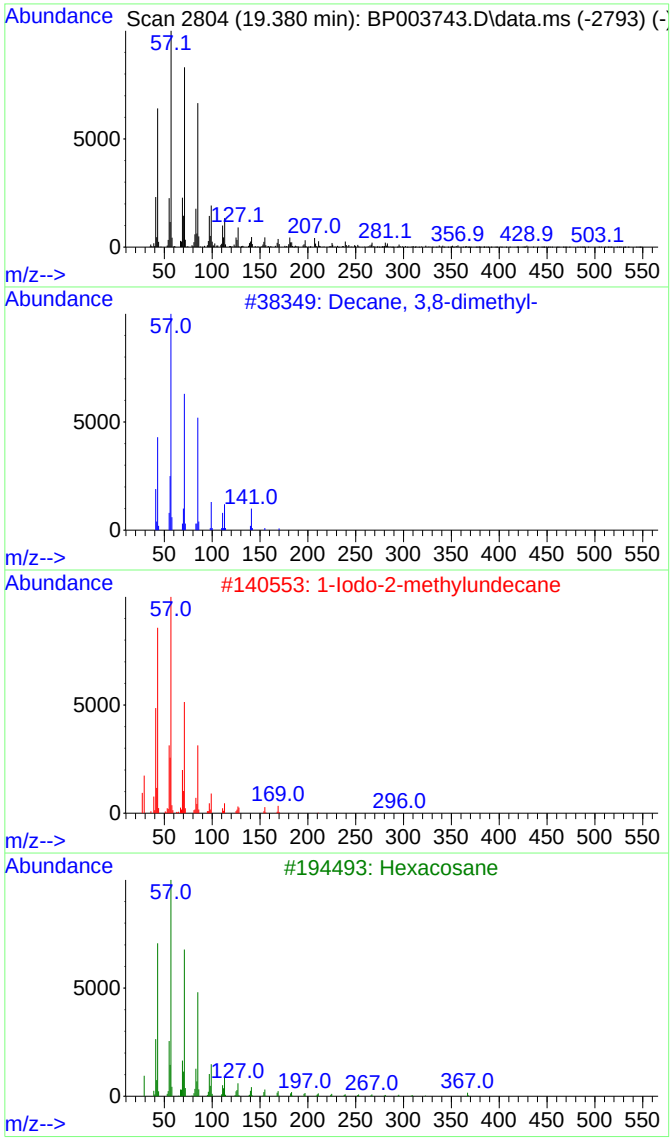
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

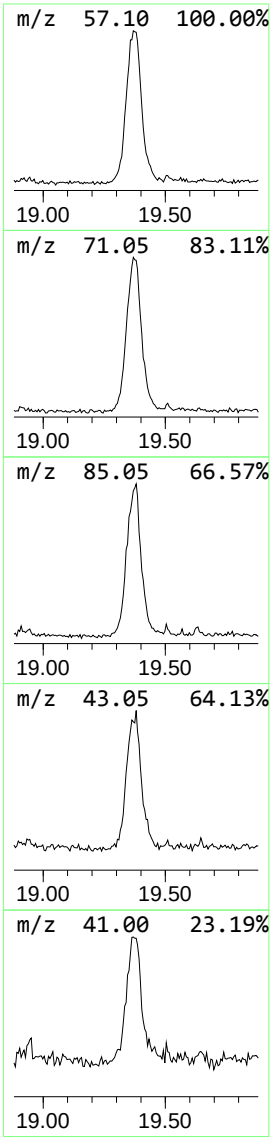
****

Peak Number 9 Decane, 3,8-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.380	8.42 ng	601489	Phenanthrene-d10	17.828

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	93
2	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	91
3	Hexacosane	366	C26H54	000630-01-3	91
4	2-methyloctacosane	408	C29H60	1000376-72-8	91
5	Hentriacontane	437	C31H64	000630-04-6	91





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Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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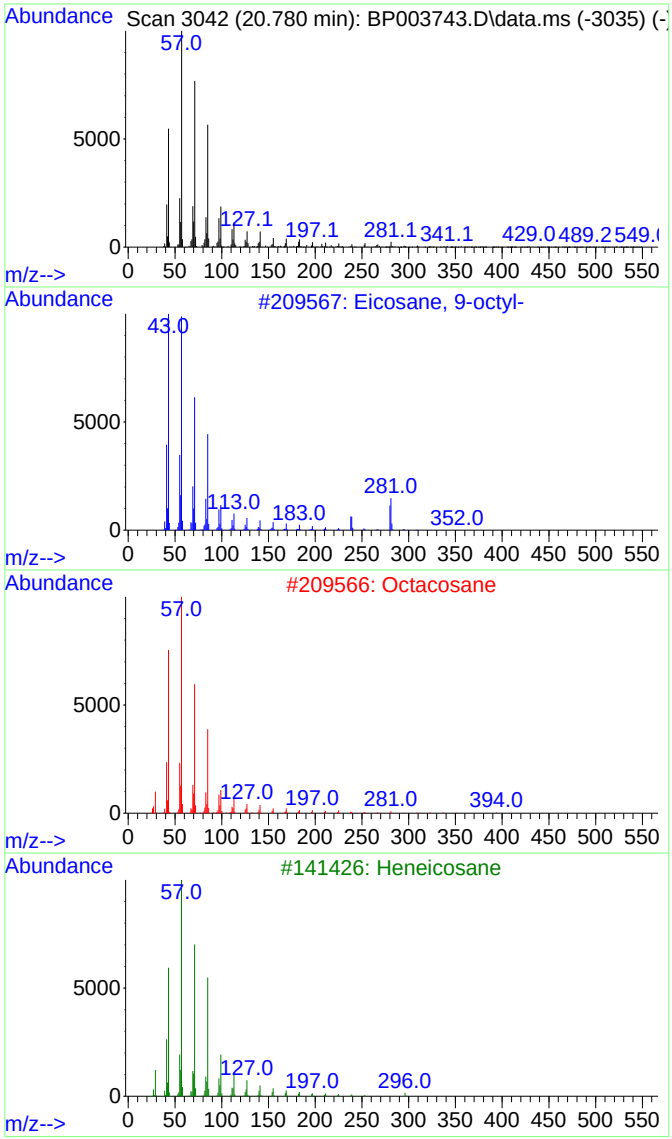
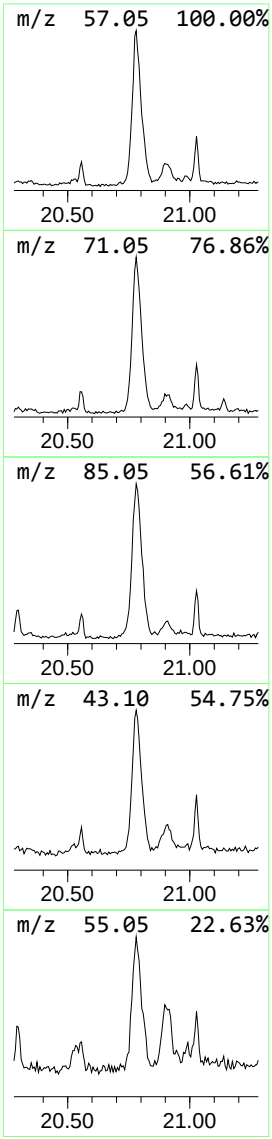
Peak Number 10 Eicosane, 9-octyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.780	8.08 ng	666703	Chrysene-d12	21.839

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 9-octyl-	394	C28H58	013475-77-9	94
2	Octacosane	394	C28H58	000630-02-4	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Triacontane	422	C30H62	000638-68-6	91
5	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	91





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Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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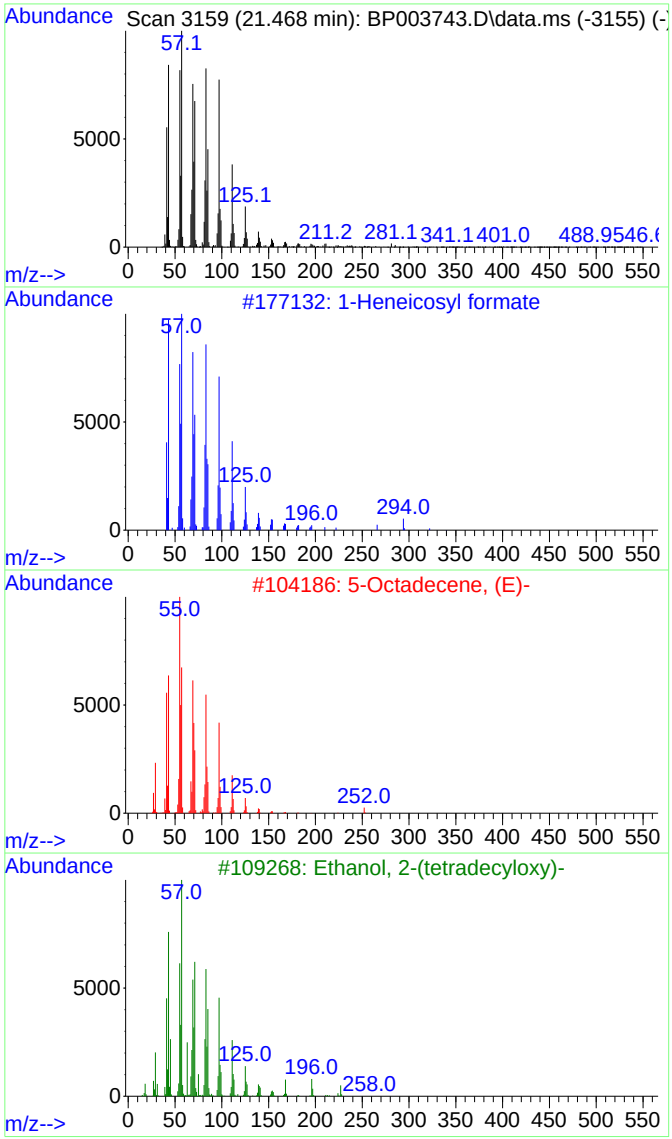
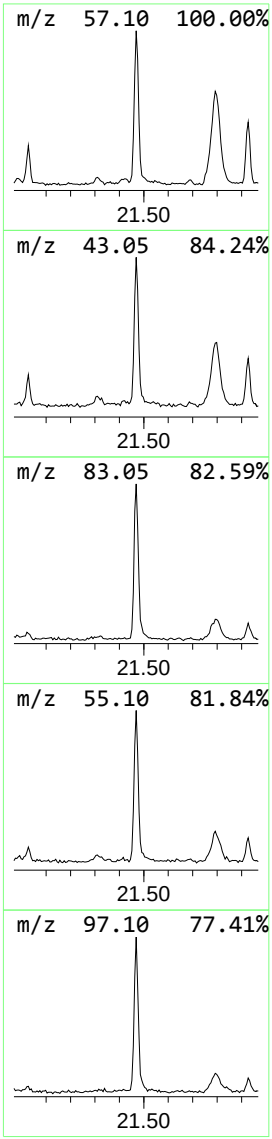
Peak Number 11 1-Heneicosyl formate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.468	8.74 ng	721759	Chrysene-d12	21.839

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	98
2	5-Octadecene, (E)-	252	C18H36	007206-21-5	94
3	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	93
4	1-Heneicosanol	312	C21H44O	015594-90-8	93
5	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93





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Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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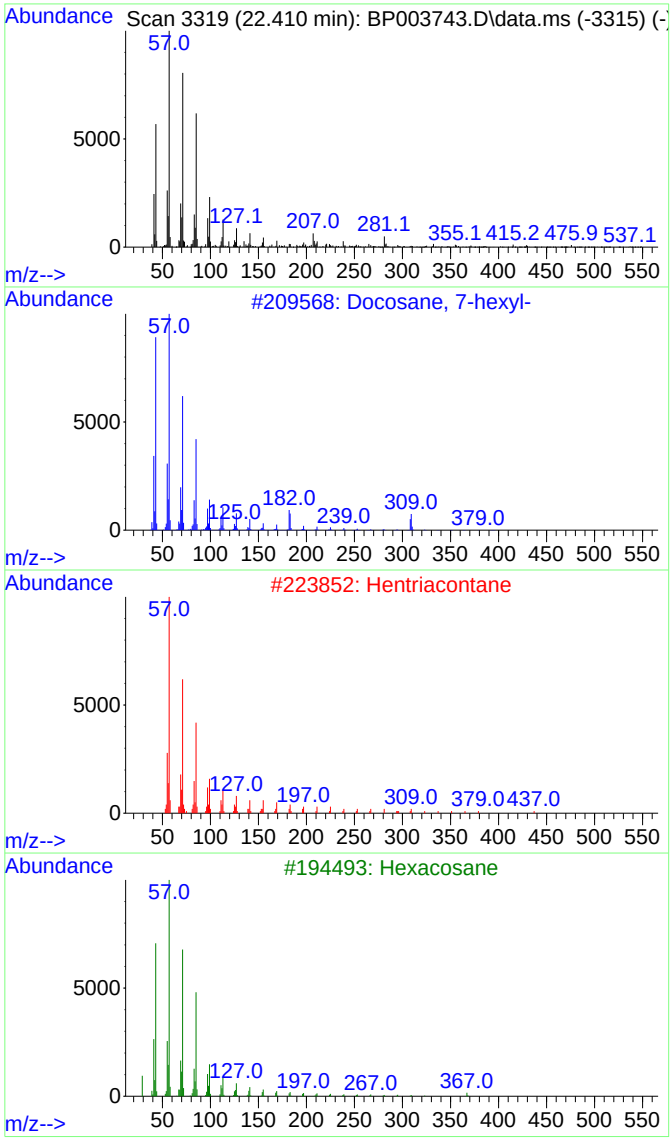
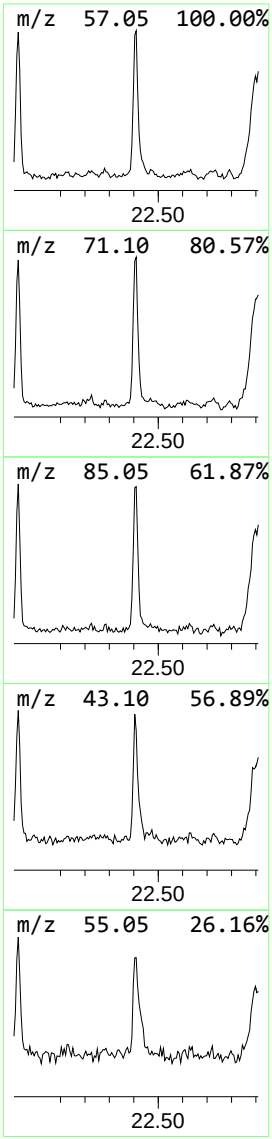
Peak Number 12 Docosane, 7-hexyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.410	2.42 ng	199353	Chrysene-d12	21.839

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 7-hexyl-	394	C28H58	055373-86-9	91
2	Hentriacontane	437	C31H64	000630-04-6	90
3	Hexacosane	366	C26H54	000630-01-3	87
4	2-methyloctacosane	408	C29H60	1000376-72-8	87
5	Heneicosane	296	C21H44	000629-94-7	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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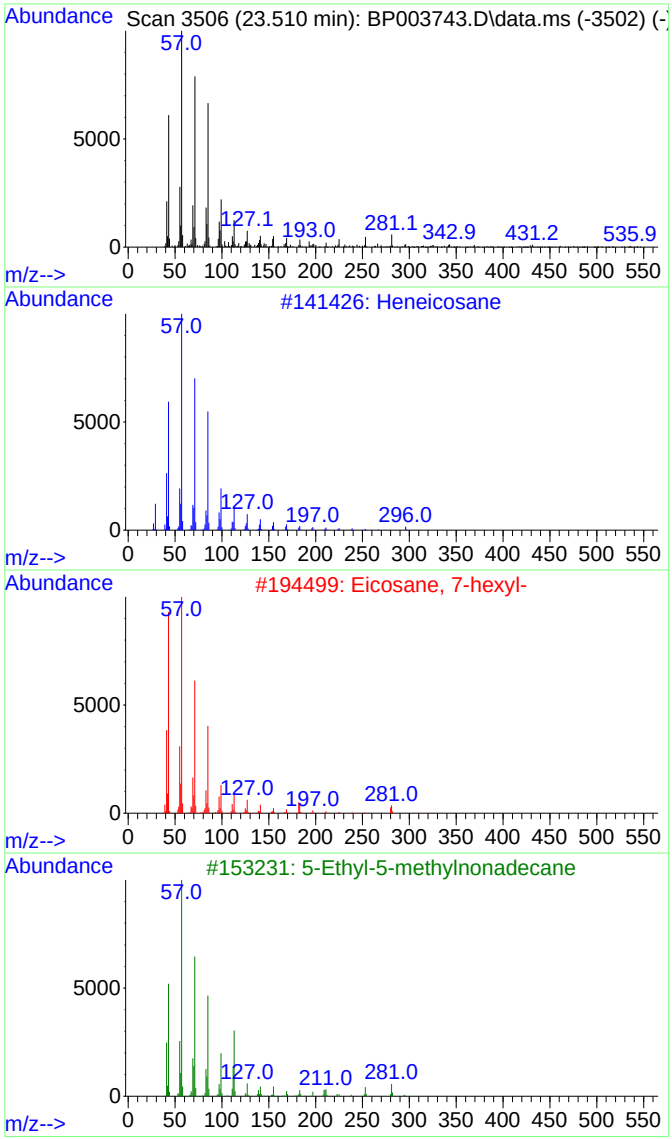
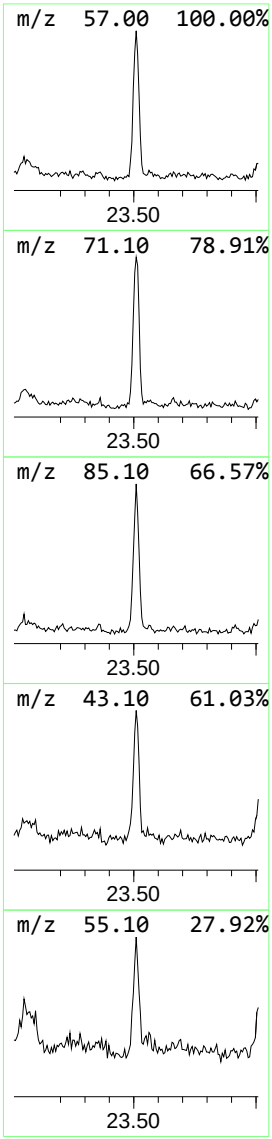
Peak Number 13 Heneicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.509	2.37 ng	195608	Perylene-d12	24.415

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	92
2	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
3	5-Ethyl-5-methylnonadecane	310	C22H46	1000360-42-7	87
4	Docosane	310	C22H46	000629-97-0	86
5	6,6-Diethylhoctadecane	310	C22H46	1000360-41-8	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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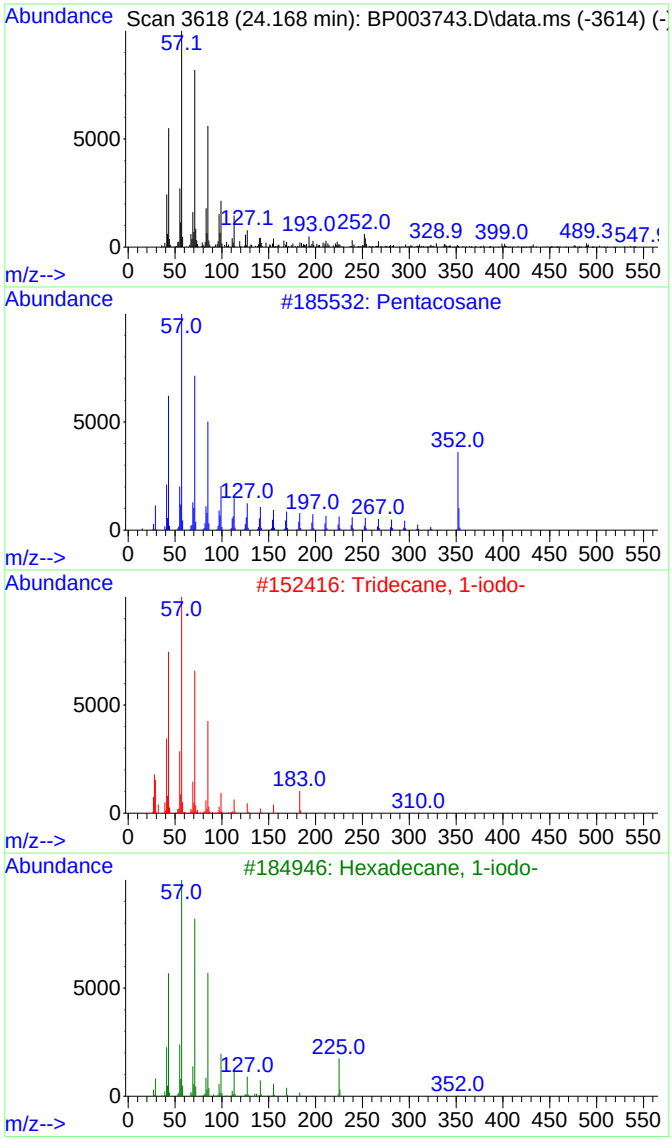
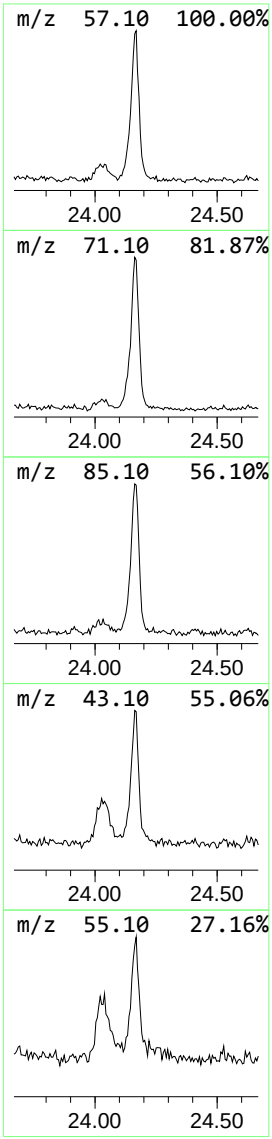
Peak Number 14 Pentacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.168	4.03 ng	332586	Perylene-d12	24.415

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	87
2	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	81
3	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	80
4	Triacontane	422	C30H62	000638-68-6	80
5	Tridecanol, 2-ethyl-2-methyl-	242	C16H34O	1000115-66-1	80





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102620\
Data File : BP003743.D
Acq On : 26 Oct 2020 16:33
Operator : CG/JU
Sample : L4501-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
1,4-Dioxaspiro[...	2.916	5.2	ng	168486	1	8.510	646470	20.0
unknown3.08	3.087	73.8	ng	2384930	1	8.510	646470	20.0
Butane, 2-metho...	3.346	9.0	ng	290337	1	8.510	646470	20.0
2-Pentanone, 4-...	5.475	8.5	ng	273318	1	8.510	646470	20.0
4-((1E)-3-Hydro...	17.180	2.0	ng	146009	4	17.828	1429230	20.0
3,5-Dimethoxy-4...	18.904	3.1	ng	220760	4	17.828	1429230	20.0
Phenazine, 2-me...	18.945	4.9	ng	351637	4	17.828	1429230	20.0
Decane, 3,8-dim...	19.380	8.4	ng	601489	4	17.828	1429230	20.0
Eicosane, 9-octyl-	20.780	8.1	ng	666703	5	21.839	1650820	20.0
1-Heneicosyl fo...	21.468	8.7	ng	721759	5	21.839	1650820	20.0
Docosane, 7-hexyl-	22.410	2.4	ng	199353	5	21.839	1650820	20.0
Heneicosane	23.509	2.4	ng	195608	6	24.415	1651880	20.0
Pentacosane	24.168	4.0	ng	332586	6	24.415	1651880	20.0