LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081721\
Data File : BP006689.D
Acq On : 16 Aug 2021 18:38
Operator : CG/JU
Sample : M3374-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DEWATERING-DISCHARGE

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\8270E-BP080321.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP006689.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.858	296	301	307	rBV	223612	315590	2.77%	0.555%
2	5.305	372	377	386	rVB	2601330	3743251	32.84%	6.579%
3	6.887	640	646	658	rVB	1863407	2832945	24.85%	4.979%
4	7.093	677	681	692	rVB	107104	198726	1.74%	0.349%
5	7.705	779	785	791	rBV	725341	1098522	9.64%	1.931%
6	8.069	841	847	853	rBV	174032	271760	2.38%	0.478%
7	8.193	863	868	872	rBV	115980	171174	1.50%	0.301%
8	8.804	966	972	976	rBV	311320	492987	4.32%	0.866%
9	8.863	976	982	991	rVB	2429092	4325909	37.95%	7.603%
10	9.322	1055	1060	1066	rVV	202072	307576	2.70%	0.541%
11	9.381	1066	1070	1076	rVB	113282	168690	1.48%	0.296%
12	9.887	1150	1156	1166	rVB2	519196	919035	8.06%	1.615%
13	10.287	1217	1224	1234	rBV	401472	699892	6.14%	1.230%
14	10.410	1234	1245	1248	rVV5	44298	154013	1.35%	0.271%
15	10.487	1248	1258	1264	rVV	992946	1919087	16.83%	3.373%
16	10.540	1264	1267	1274	rVV5	130071	287609	2.52%	0.505%
17	10.604	1274	1278	1288	rVB	112303	191295	1.68%	0.336%
18	11.110	1359	1364	1368	rBV	162318	262164	2.30%	0.461%
19	11.404	1409	1414	1421	rVB	86863	153853	1.35%	0.270%
20	12.045	1518	1523	1529	rVB3	72164	130637	1.15%	0.230%
21	12.381	1573	1580	1584	rBV5	92041	174541	1.53%	0.307%
22	12.557	1604	1610	1617	rBV6	55314	123088	1.08%	0.216%
23	12.969	1672	1680	1686	rVB	5330260	7539331	66.14%	13.250%
24	13.475	1761	1766	1770	rBV2	91560	150405	1.32%	0.264%
25	13.575	1778	1783	1788	rVB2	94132	140437	1.23%	0.247%
26	14.345	1908	1914	1920	rVB	1378774	1970823	17.29%	3.464%
27	14.839	1992	1998	2002	rBV2	77542	118721	1.04%	0.209%
28	15.845	2162	2169	2177	rBV	4456676	6177476	54.19%	10.857%
29	17.104	2377	2383	2389	rBV	1644598	2355773	20.67%	4.140%
30	17.945	2521	2526	2532	rVB	195288	301071	2.64%	0.529%
31	18.016	2533	2538	2547	rBV	596785	885686	7.77%	1.557%
32	19.257	2745	2749	2762	rBV	395443	578050	5.07%	1.016%
33	19.616	2806	2810	2815	rBV	141348	200517	1.76%	0.352%
34	19.686	2817	2822	2827	rVV	9782836	11399575	100.00%	20.035%
35	20.610	2975	2979	2985	rVB	168343	209493	1.84%	0.368%
36	20.945	3032	3036	3038	rBV	222093	315530	2.77%	0.555%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081721\
Data File : BP006689.D
Acq On : 16 Aug 2021 18:38
Operator : CG/JU
Sample : M3374-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DEWATERING-DISCHARGE

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\8270E-BP080321.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	21.210	3076	3081	3086	rBV	2245818	2734840	23.99%	4.806%
38	23.527	3468	3475	3484	rVB	1477192	2879351	25.26%	5.060%

Sum of corrected areas: 56899423

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081721\
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ClientSampleId :
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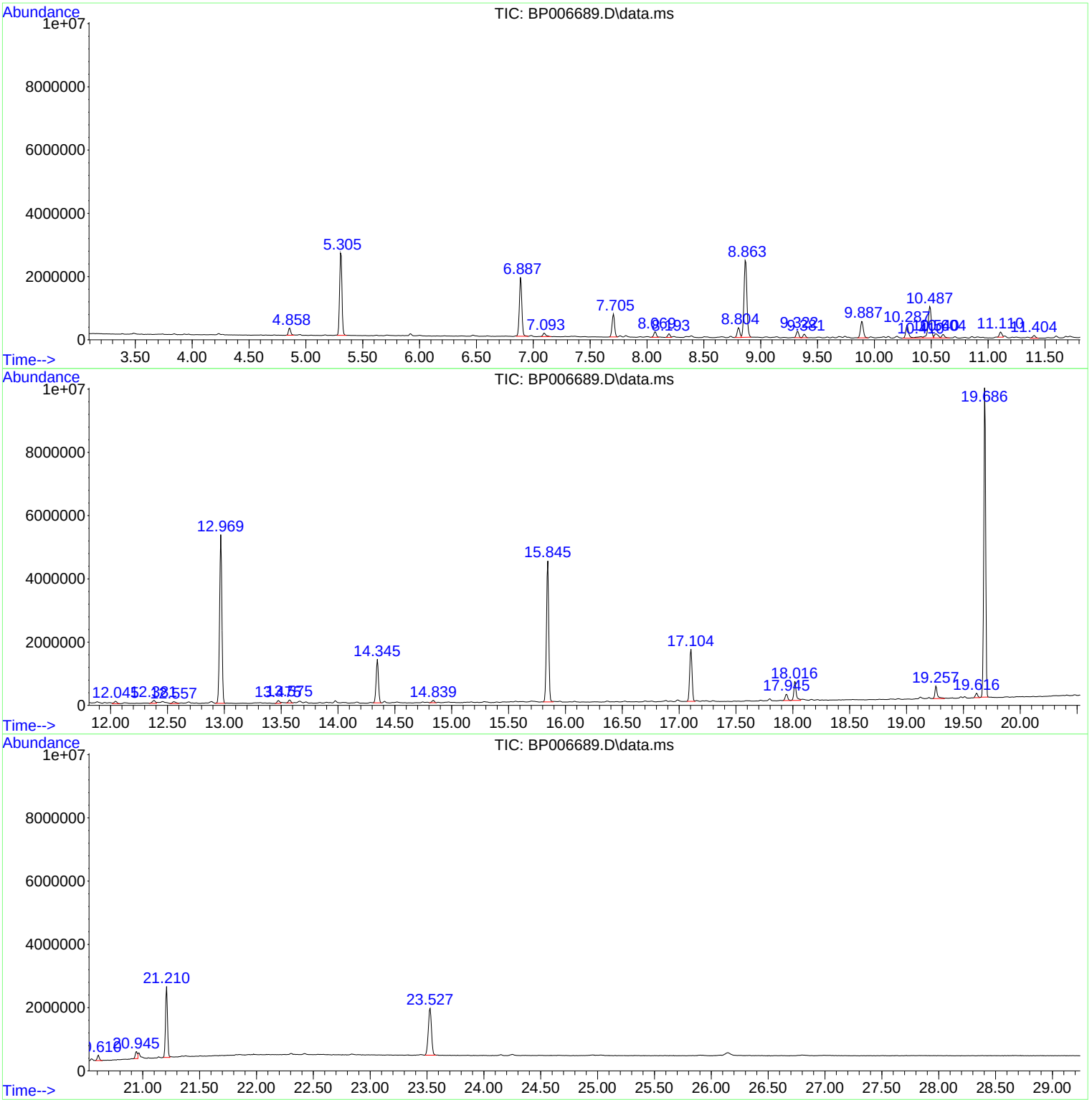
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081721\
Data File : BP006689.D
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Operator : CG/JU
Sample : M3374-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
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DEWATERING-DISCHARGE

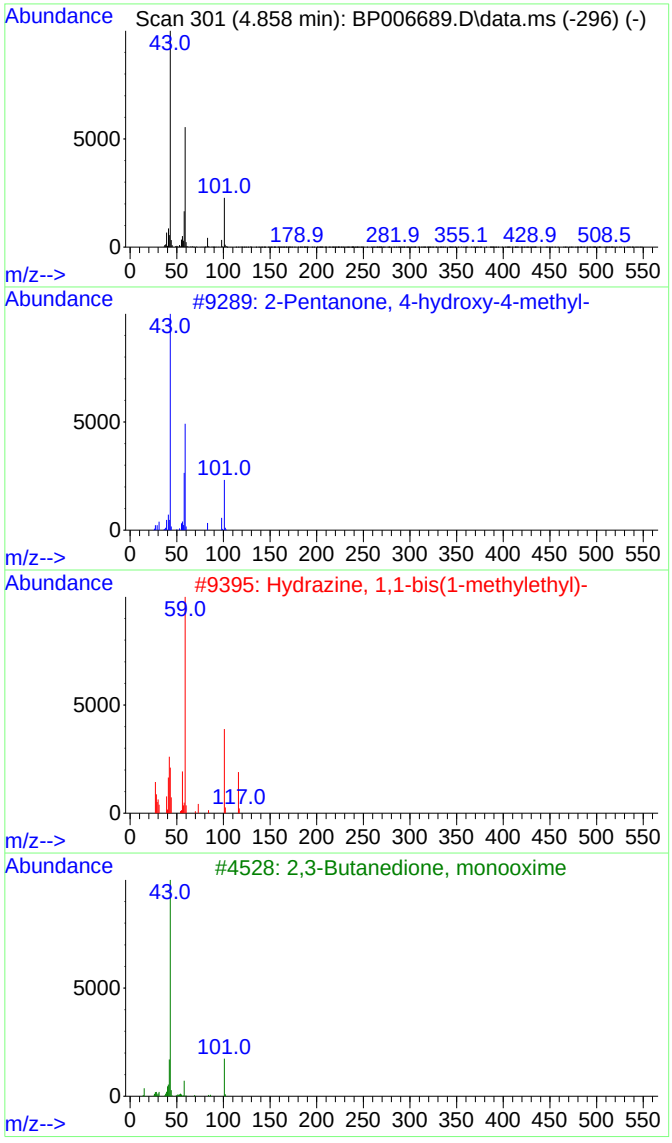
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

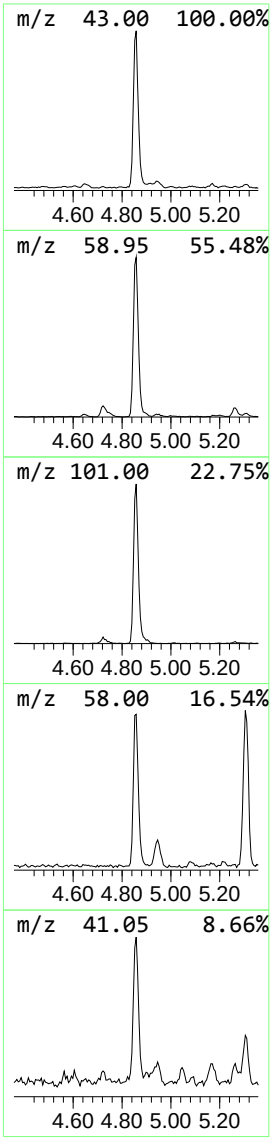
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.858	5.75 ng	315590	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	28
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
5	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17





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Data File : BP006689.D
Acq On : 16 Aug 2021 18:38
Operator : CG/JU
Sample : M3374-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DEWATERING-DISCHARGE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

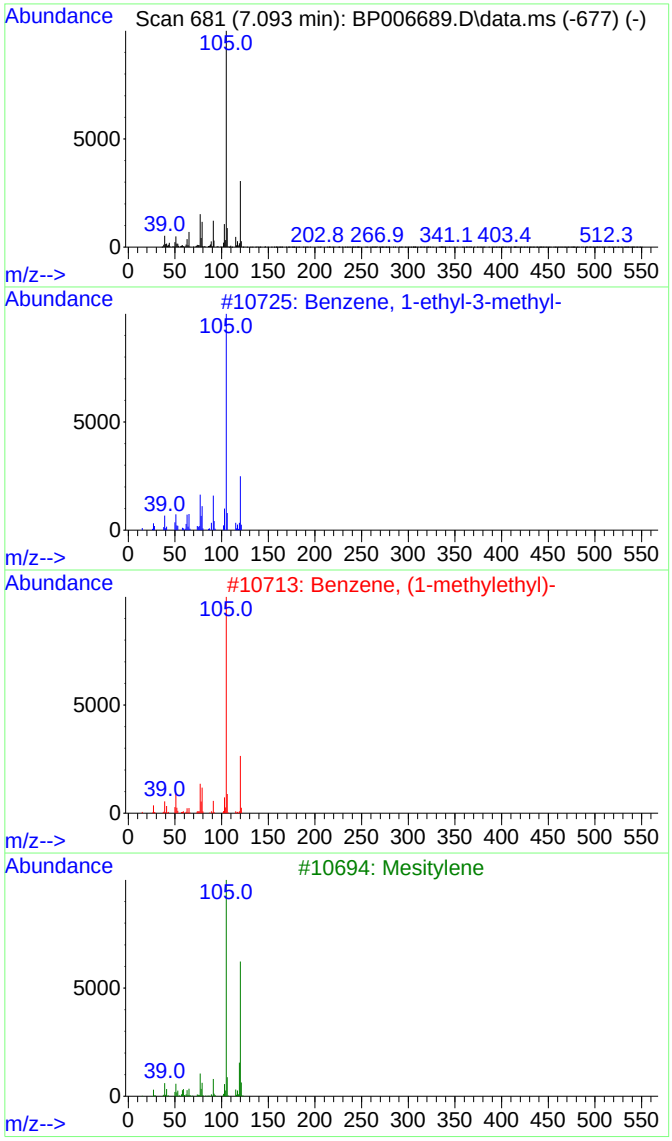
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

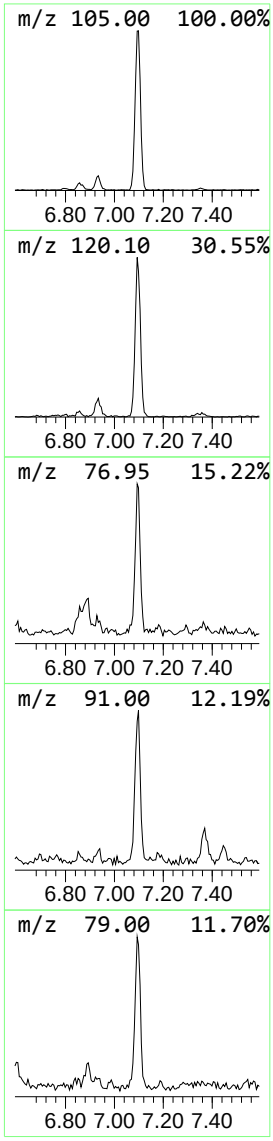
****

Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.093	3.62 ng	198726	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
2	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





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Data File : BP006689.D
Acq On : 16 Aug 2021 18:38
Operator : CG/JU
Sample : M3374-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DEWATERING-DISCHARGE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

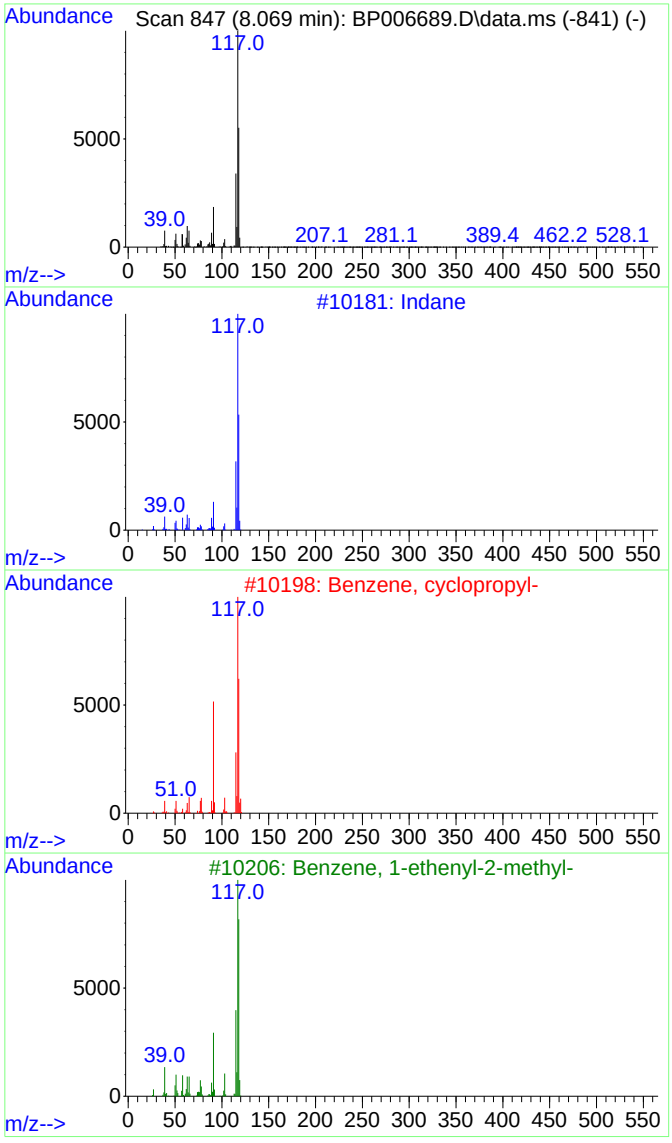
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

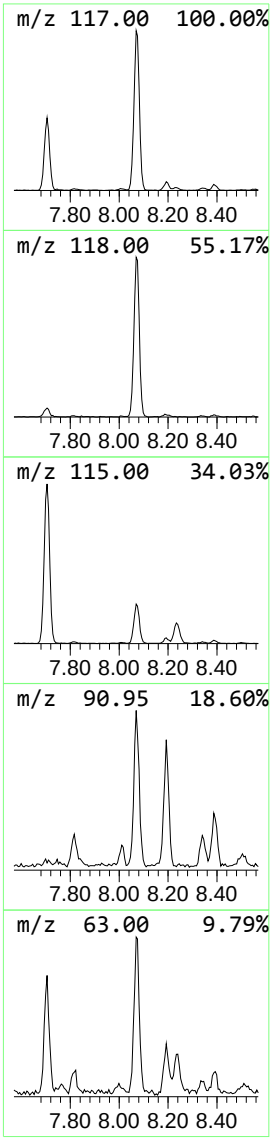
****

Peak Number 3 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.069	4.95 ng	271760	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	94
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
3	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
4	Deltacyclene	118	C9H10	007785-10-6	59
5	Benzene, (3-chloro-1-propenyl)-	152	C9H9Cl	002687-12-9	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081721\
Data File : BP006689.D
Acq On : 16 Aug 2021 18:38
Operator : CG/JU
Sample : M3374-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DEWATERING-DISCHARGE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

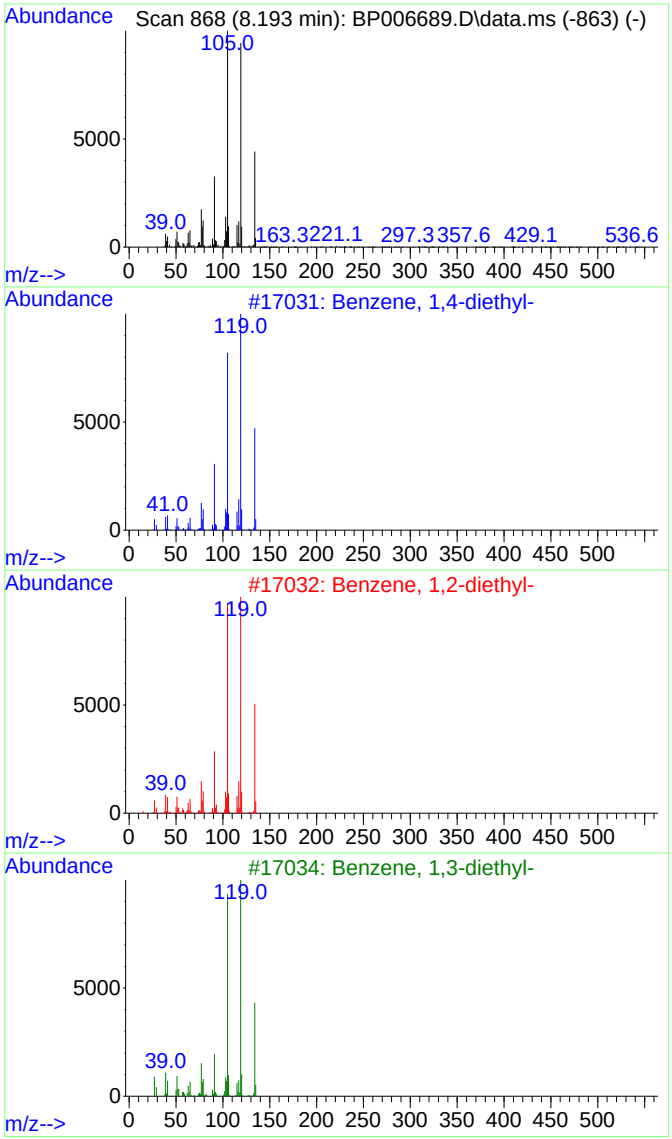
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

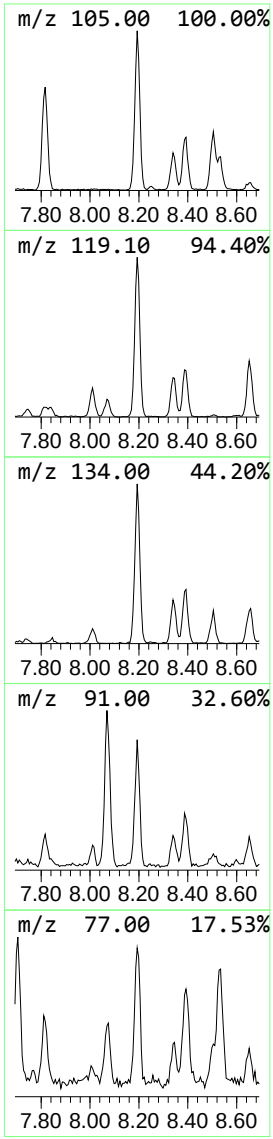
****

Peak Number 4 Benzene, 1,4-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.193	3.12 ng	171174	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	95
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
5	1,3,8-p-Menthatriene	134	C10H14	018368-95-1	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081721\
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Operator : CG/JU
Sample : M3374-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DEWATERING-DISCHARGE

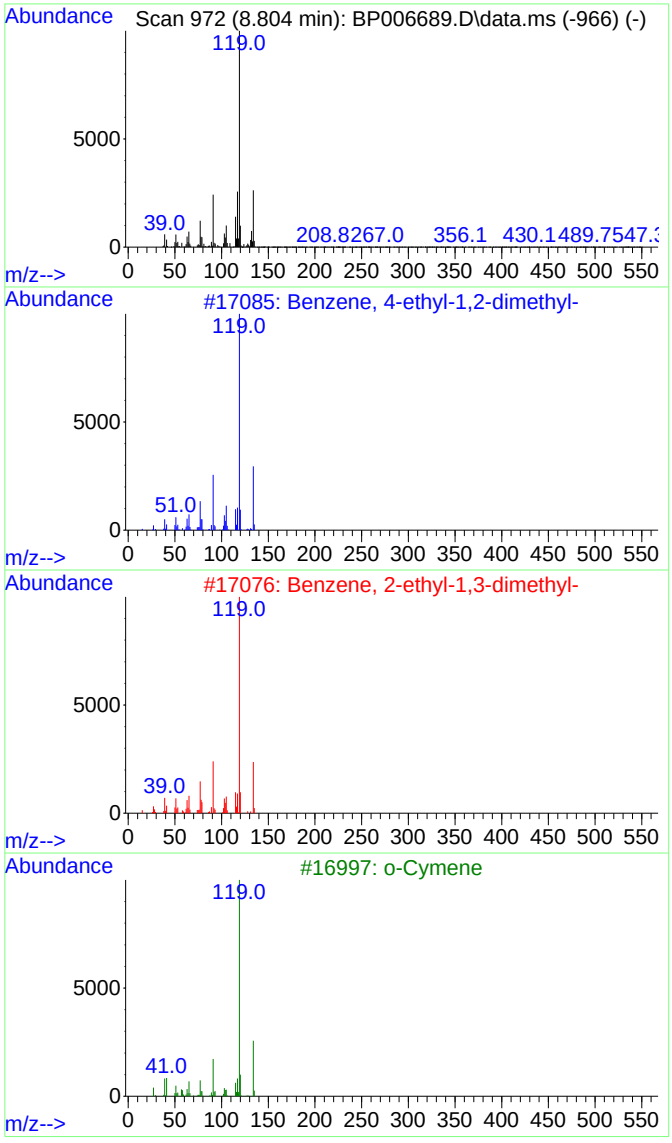
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

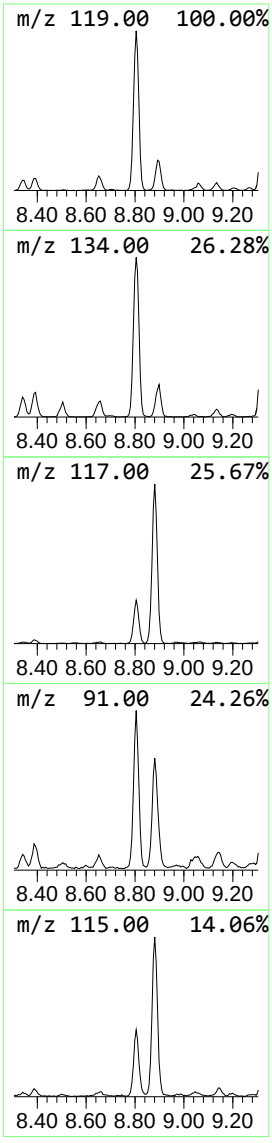
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.804	8.98 ng	492987	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
3	o-Cymene	134	C10H14	000527-84-4	93
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93





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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

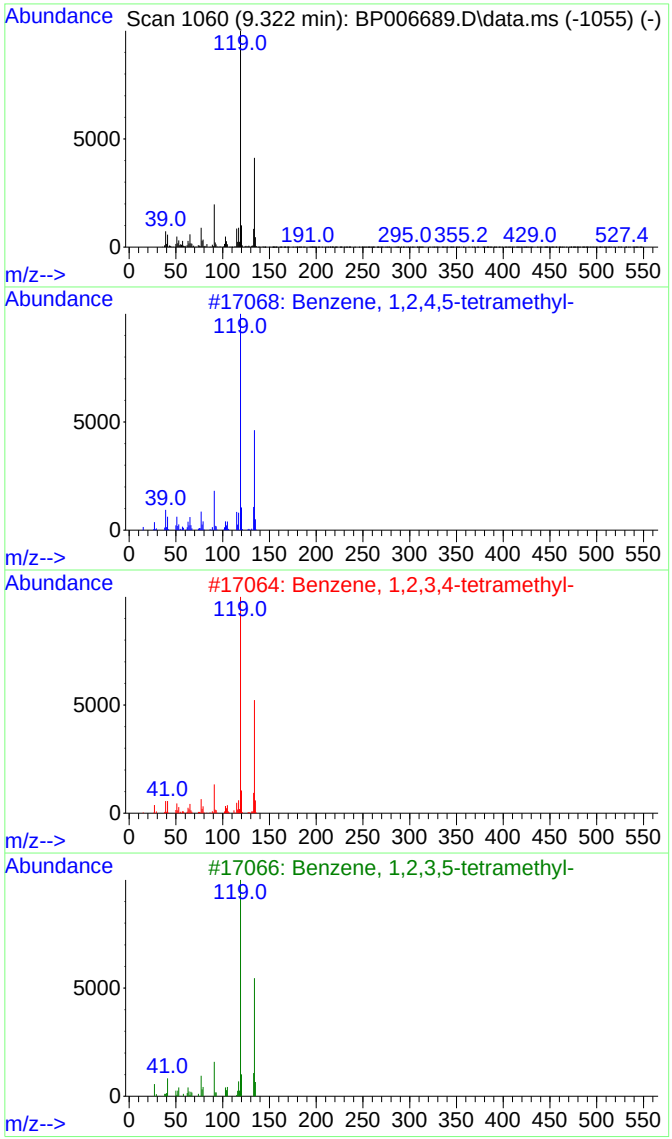
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

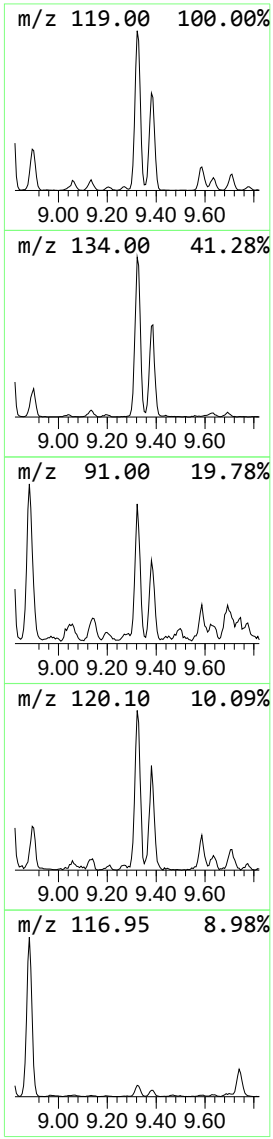
****

Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.322	3.21 ng	307576	Naphthalene-d8	10.487

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91





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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

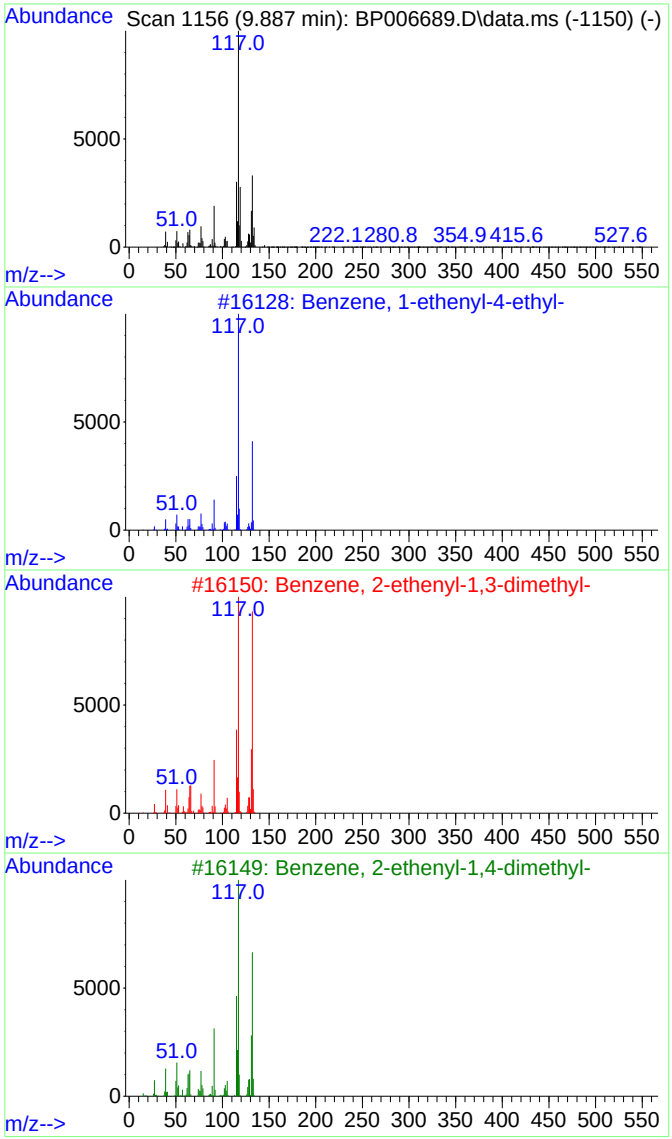
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

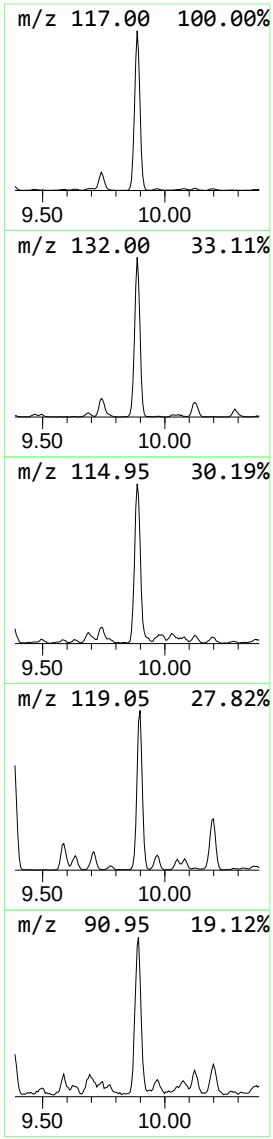
****

Peak Number 7 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.887	9.58 ng	919035	Naphthalene-d8	10.487

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
2	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	93
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
4	2,4-Dimethylstyrene	132	C10H12	002234-20-0	90
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87





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Operator : CG/JU
Sample : M3374-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DEWATERING-DISCHARGE

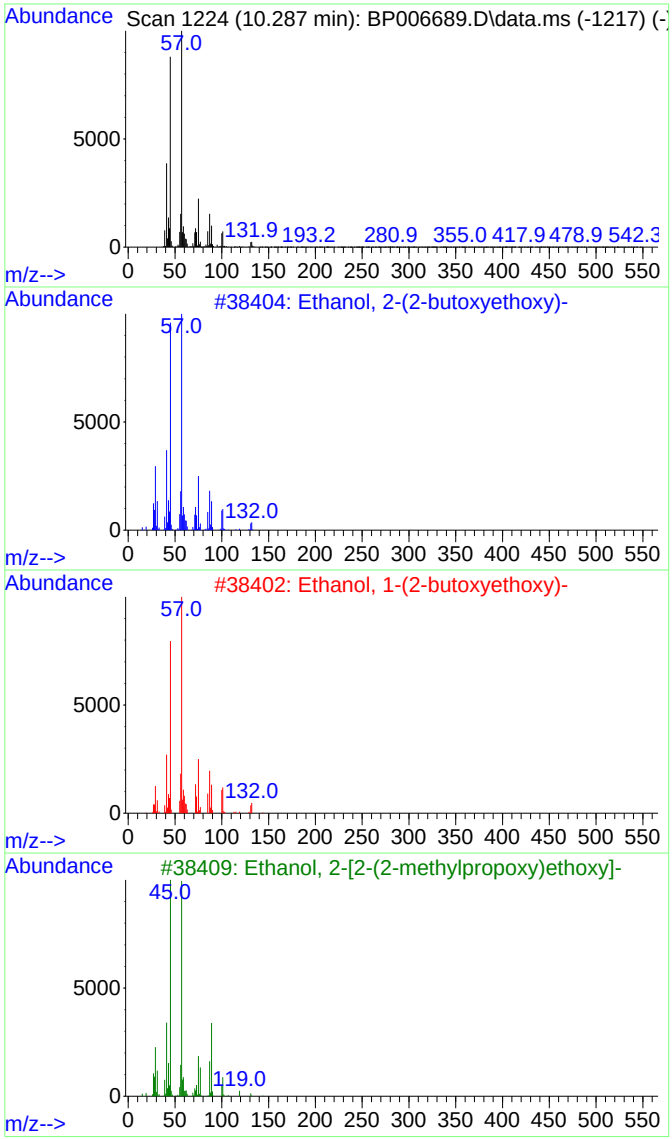
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

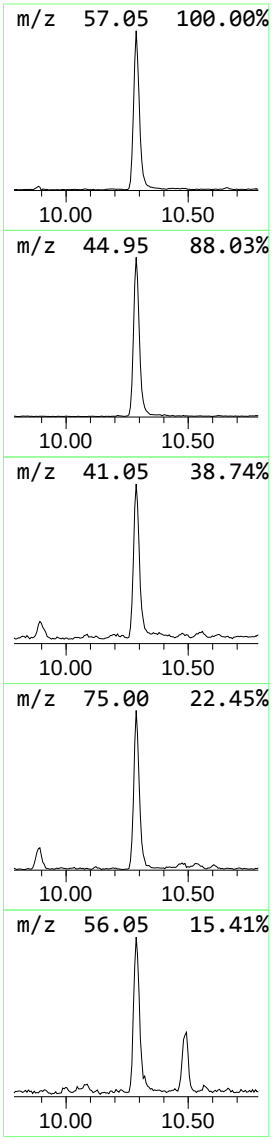
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.287	7.29 ng	699892	Naphthalene-d8	10.487

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	78
4	Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
5	Di-sec-Butyl ether	130	C8H18O	006863-58-7	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081721\
Data File : BP006689.D
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Operator : CG/JU
Sample : M3374-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DEWATERING-DISCHARGE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

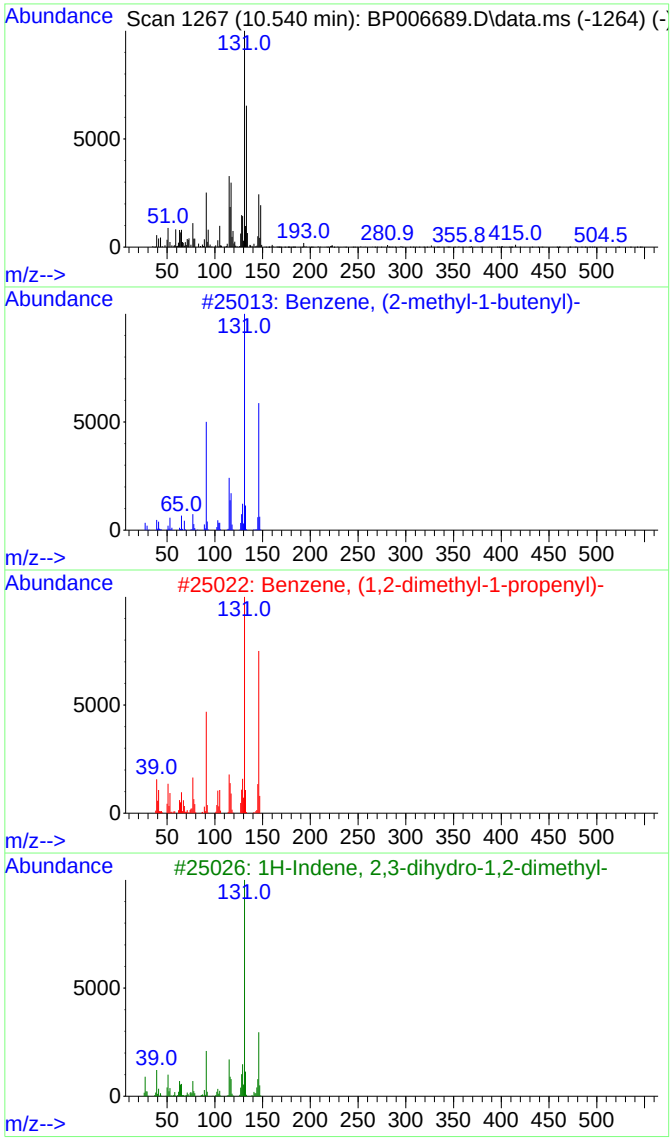
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

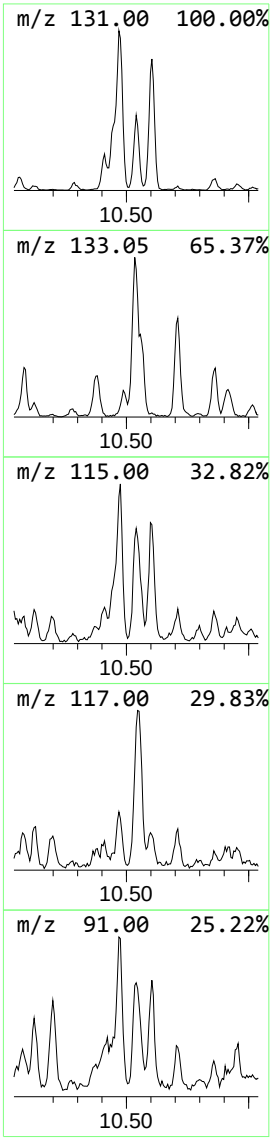
****

Peak Number 9 Benzene, (2-methyl-1-butenyl)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.540	3.00 ng	287609	Naphthalene-d8	10.487

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	83
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	58
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	53
4	Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	53
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081721\
Data File : BP006689.D
Acq On : 16 Aug 2021 18:38
Operator : CG/JU
Sample : M3374-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DEWATERING-DISCHARGE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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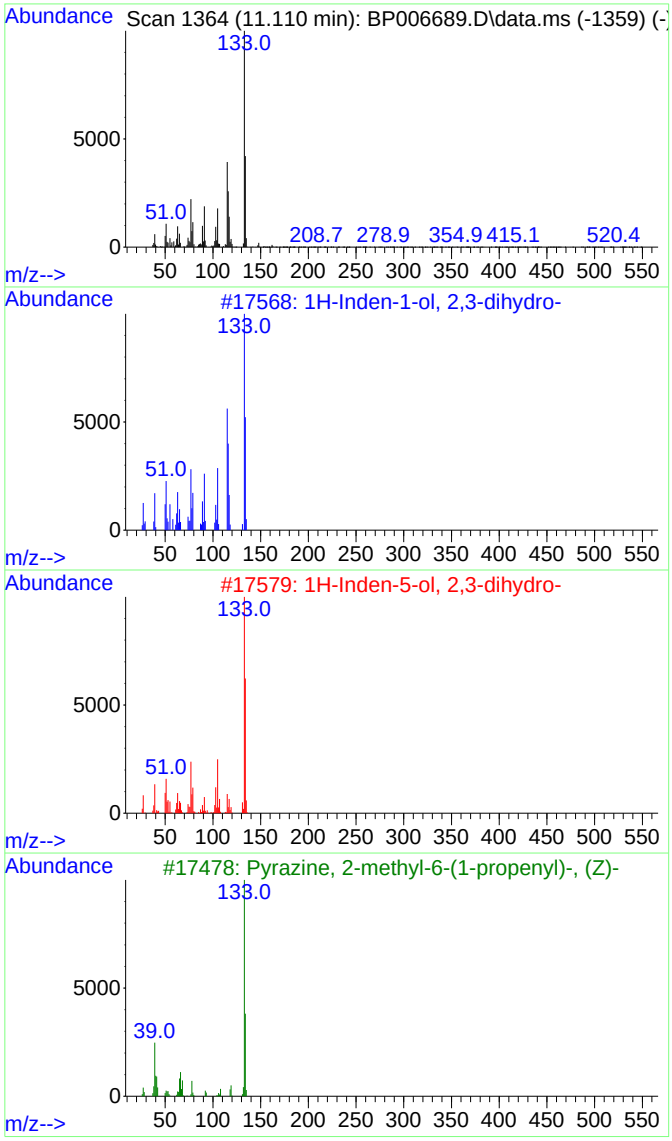
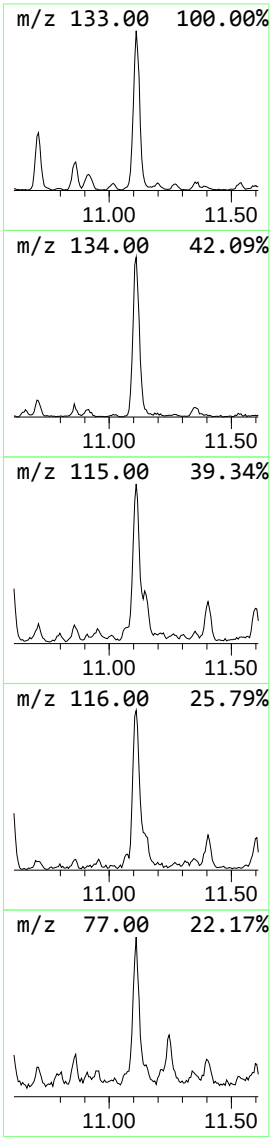
****
Peak Number 10 1H-Inden-1-ol, 2,3-dihydro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.110	2.73 ng	262164	Naphthalene-d8	10.487

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-1-ol, 2,3-dihydro-	134	C9H10O	006351-10-6	94
2	1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	64
3	Pyrazine, 2-methyl-6-(1-propenyl...	134	C8H10N2	055138-67-5	46
4	Pyrazine, 2-methyl-6-(1-propenyl...	134	C8H10N2	018217-81-7	46
5	4-Ethylbenzoic acid, 2-methylphe...	240	C16H16O2	1000293-61-2	43





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Operator : CG/JU
Sample : M3374-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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ClientSampleId :
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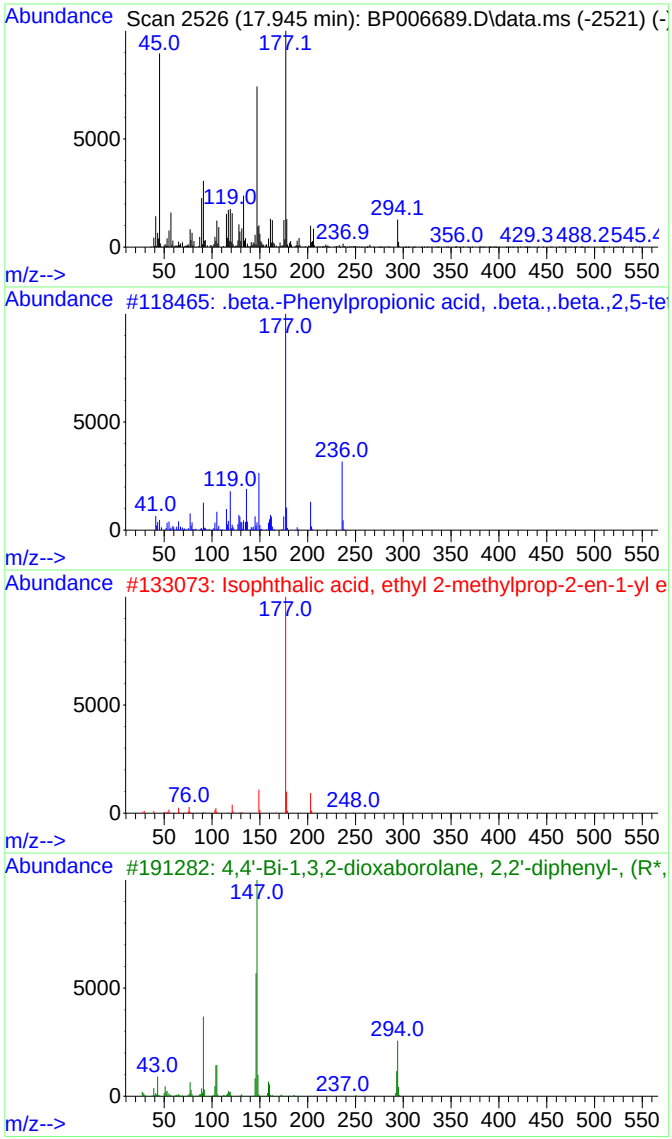
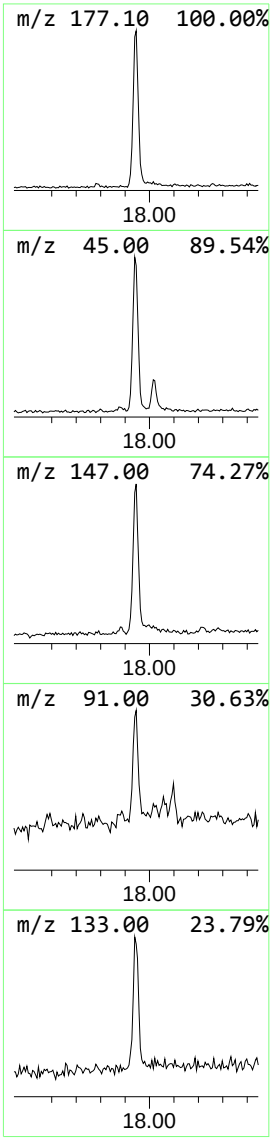
Peak Number 11 unknown17.945 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.945	2.56 ng	301071	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Phenylpropionic acid, .be...	236	C14H20O3	186317-88-4	22
2	Isophthalic acid, ethyl 2-methyl...	248	C14H16O4	1000343-94-4	20
3	4,4'-Bi-1,3,2-dioxaborolane, 2,2...	294	C16H16B2O4	071166-92-2	20
4	Formamide, N-(p-methoxy-trans-st...	177	C10H11NO2	002501-37-3	18
5	Disiloxane, ethylpentamethyl-	176	C7H20OSi2	006231-60-3	14





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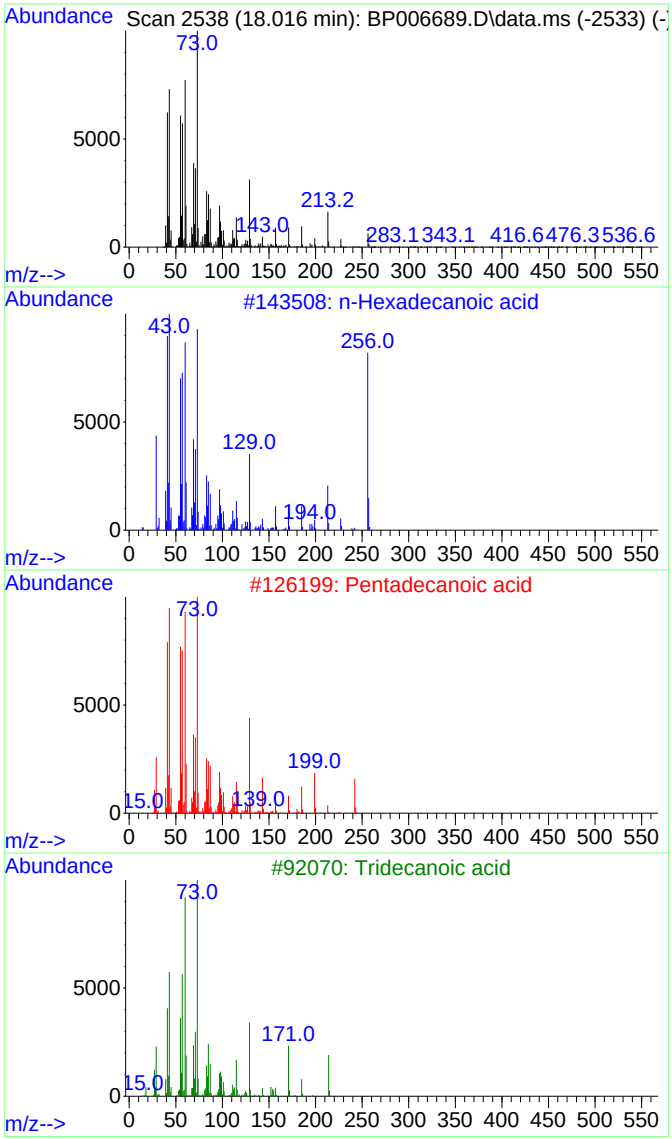
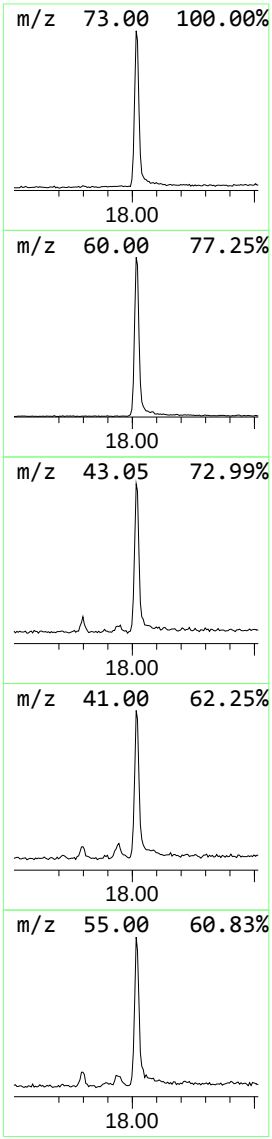
Peak Number 12 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.016	7.52 ng	885686	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3	Tridecanoic acid	214	C13H26O2	000638-53-9	64
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5	n-Decanoic acid	172	C10H20O2	000334-48-5	50





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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
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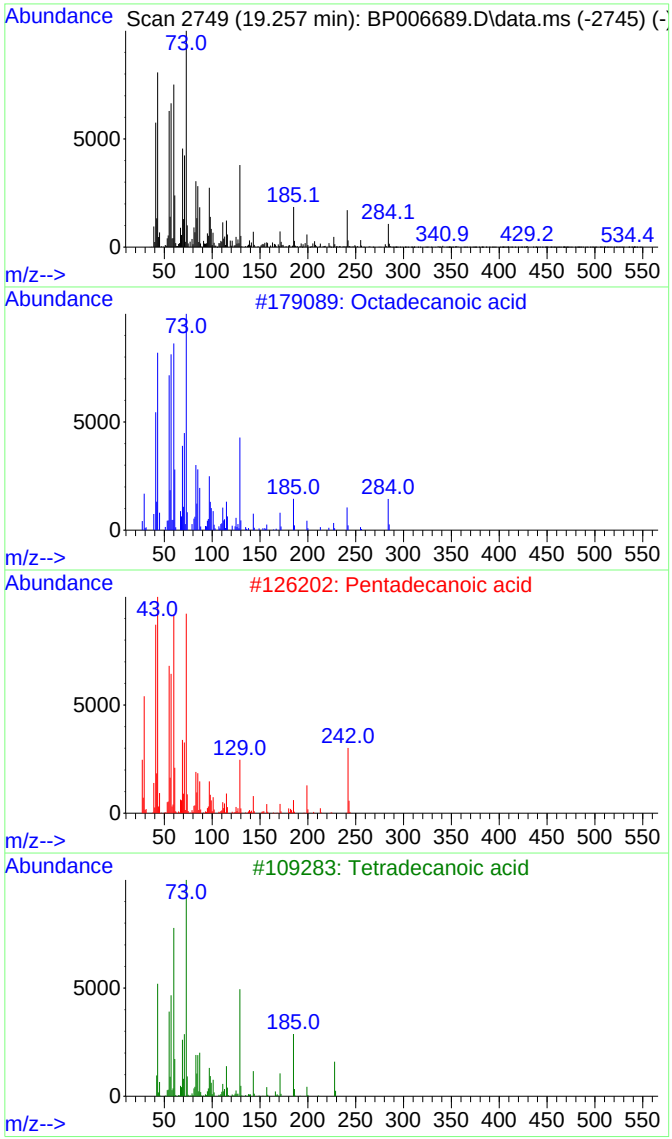
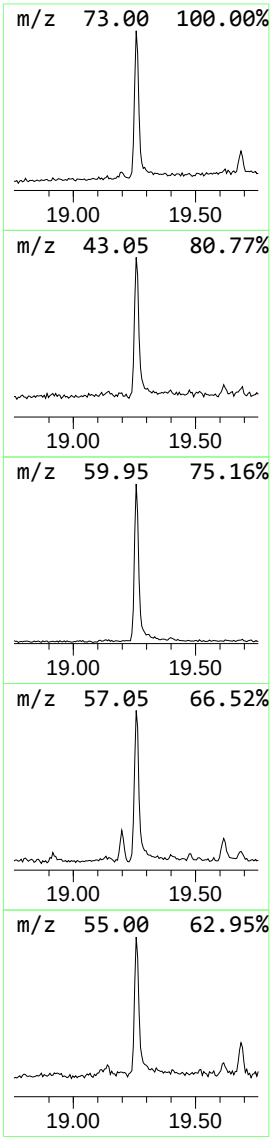
Peak Number 13 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.257	4.23 ng	578050	Chrysene-d12	21.210

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	70
4	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	64
5	n-Decanoic acid	172	C10H20O2	000334-48-5	62





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Sample : M3374-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DEWATERING-DISCHARGE

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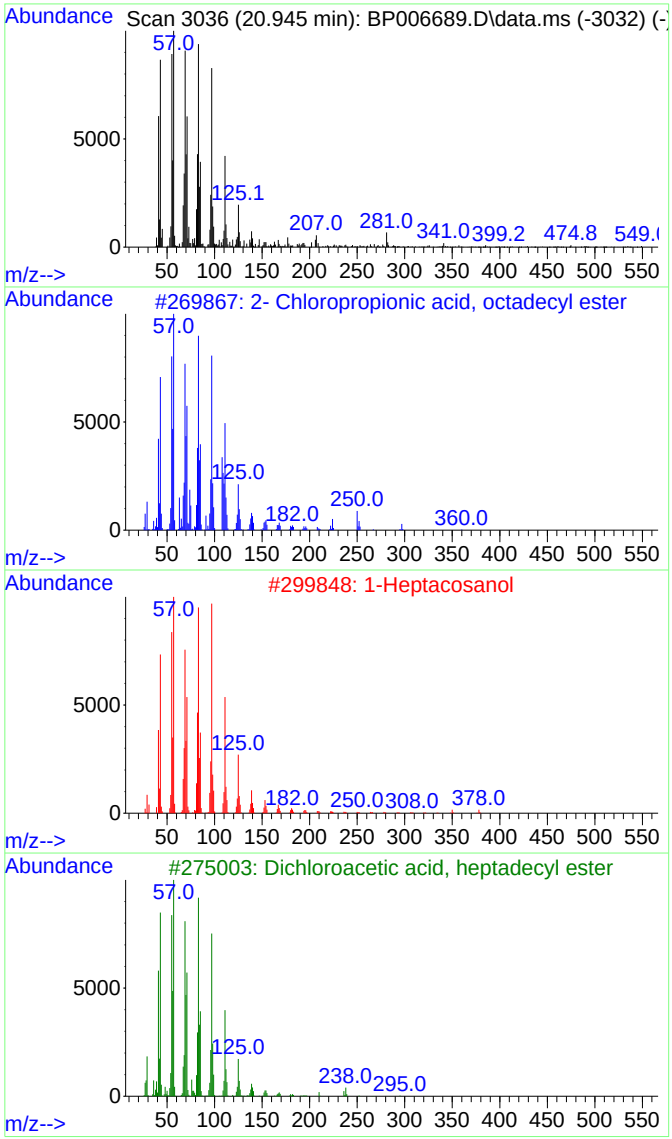
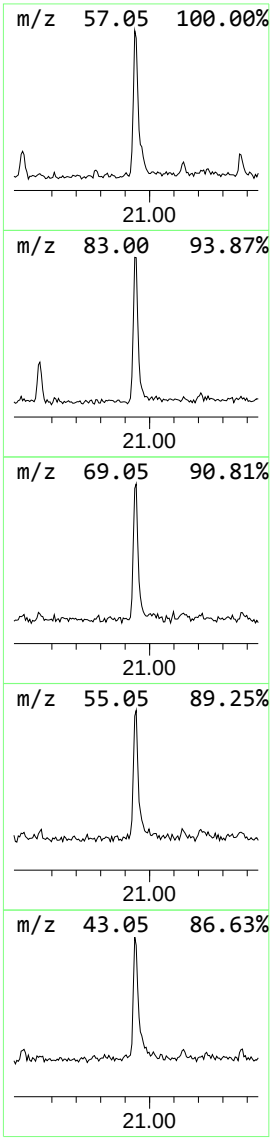
****
Peak Number 14 2- Chloropropionic acid, oc... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.945	2.31 ng	315530	Chrysene-d12	21.210

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	95
2	1-Heptacosanol	396	C27H56O	002004-39-9	94
3	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
4	Octacosyl acetate	452	C30H60O2	018206-97-8	92
5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	91





Tentatively Identified Compound (LSC) summary

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Data File : BP006689.D
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
2-Pentanone, 4-...	4.858	5.8	ng	315590	1	7.705	1098520	20.0
Benzene, 1-ethy...	7.093	3.6	ng	198726	1	7.705	1098520	20.0
Indane	8.069	5.0	ng	271760	1	7.705	1098520	20.0
Benzene, 1,4-di...	8.193	3.1	ng	171174	1	7.705	1098520	20.0
Benzene, 4-ethy...	8.804	9.0	ng	492987	1	7.705	1098520	20.0
Benzene, 1,2,4,...	9.322	3.2	ng	307576	2	10.487	1919090	20.0
Benzene, 1-ethe...	9.887	9.6	ng	919035	2	10.487	1919090	20.0
Ethanol, 2-(2-b...	10.287	7.3	ng	699892	2	10.487	1919090	20.0
Benzene, (2-met...	10.540	3.0	ng	287609	2	10.487	1919090	20.0
1H-Inden-1-ol, ...	11.110	2.7	ng	262164	2	10.487	1919090	20.0
unknown17.945	17.945	2.6	ng	301071	4	17.104	2355770	20.0
n-Hexadecanoic ...	18.016	7.5	ng	885686	4	17.104	2355770	20.0
Octadecanoic acid	19.257	4.2	ng	578050	5	21.210	2734840	20.0
2- Chloropropio...	20.945	2.3	ng	315530	5	21.210	2734840	20.0