LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051925\
Data File : BP024694.D
Acq On : 19 May 2025 18:19
Operator : RC/JU
Sample : Q2033-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TW-WTS-08

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-BP051325.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP024694.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.470	61	65	70	rBV	233210	320076	2.01%	0.422%
2	3.887	122	136	139	rBV	139650	306263	1.92%	0.404%
3	4.646	256	265	271	rBV	101055	195794	1.23%	0.258%
4	4.822	290	295	308	rVB	313608	455480	2.86%	0.601%
5	5.293	369	375	391	rVB	1737017	2618709	16.45%	3.456%
6	6.864	636	642	665	rVB	900629	1643448	10.32%	2.169%
7	7.658	770	777	783	rBV	708237	1112610	6.99%	1.468%
8	7.722	783	788	802	rVB	483631	763449	4.80%	1.007%
9	8.805	966	972	983	rVB	2071661	3564388	22.39%	4.704%
10	10.216	1206	1212	1229	rBV	96029	215375	1.35%	0.284%
11	10.428	1238	1248	1253	rBV	933126	1552359	9.75%	2.048%
12	10.475	1253	1256	1266	rVB	164341	282574	1.78%	0.373%
13	11.493	1421	1429	1443	rVB	108798	292065	1.83%	0.385%
14	11.616	1443	1450	1454	rBV	849187	1365323	8.58%	1.802%
15	11.663	1454	1458	1459	rVV	929075	1174250	7.38%	1.550%
16	11.687	1459	1462	1465	rVV	1302730	1901205	11.94%	2.509%
17	11.722	1465	1468	1478	rVB	944432	1458875	9.17%	1.925%
18	11.904	1490	1499	1504	rBV	2523424	4479045	28.14%	5.910%
19	11.963	1504	1509	1526	rVB	2978923	4837773	30.39%	6.384%
20	12.898	1660	1668	1686	rBV	5988191	8686003	54.57%	11.462%
21	13.216	1714	1722	1730	rBV	461965	780472	4.90%	1.030%
22	14.293	1898	1905	1912	rBV	1653059	2170262	13.63%	2.864%
23	15.810	2152	2163	2179	rBV	6352972	8791136	55.23%	11.601%
24	17.087	2371	2380	2394	rBV2	1836604	2700092	16.96%	3.563%
25	18.022	2534	2539	2547	rBV2	152389	220796	1.39%	0.291%
26	18.181	2560	2566	2580	rVB3	113514	265393	1.67%	0.350%
27	19.086	2715	2720	2728	rBV	720535	963116	6.05%	1.271%
28	19.804	2836	2842	2851	rBV	11559113	15917654	100.00%	21.005%
29	21.516	3126	3133	3145	rBV	2057915	3194620	20.07%	4.216%
30	21.680	3157	3161	3170	rVB	147515	258910	1.63%	0.342%
31	21.939	3201	3205	3212	rVB	143192	226922	1.43%	0.299%
32	24.798	3682	3691	3705	rVB	1188007	3066894	19.27%	4.047%

Sum of corrected areas: 75781331

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051925\
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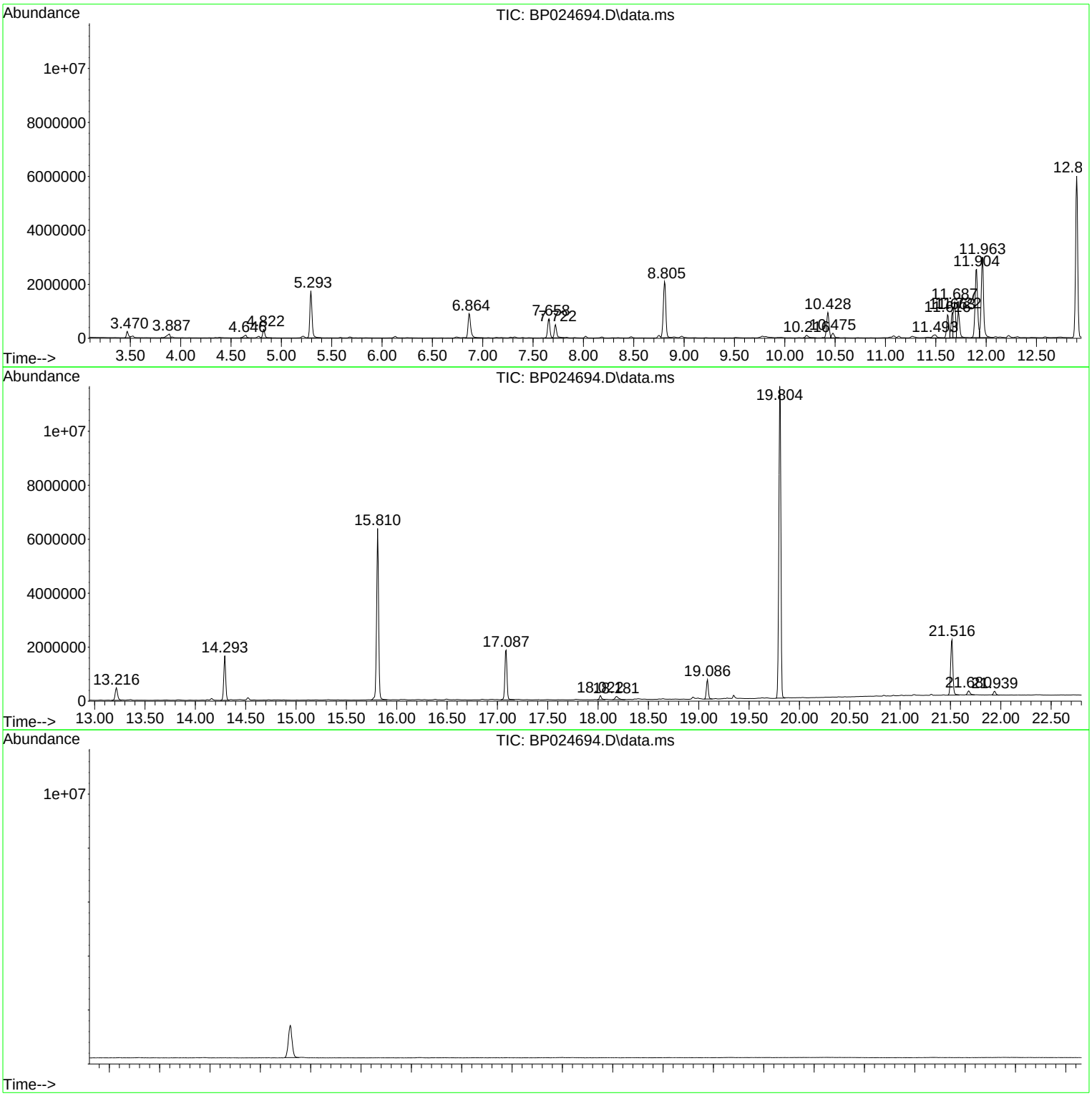
Sample : Q2033-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TW-WTS-08

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Data File : BP024694.D
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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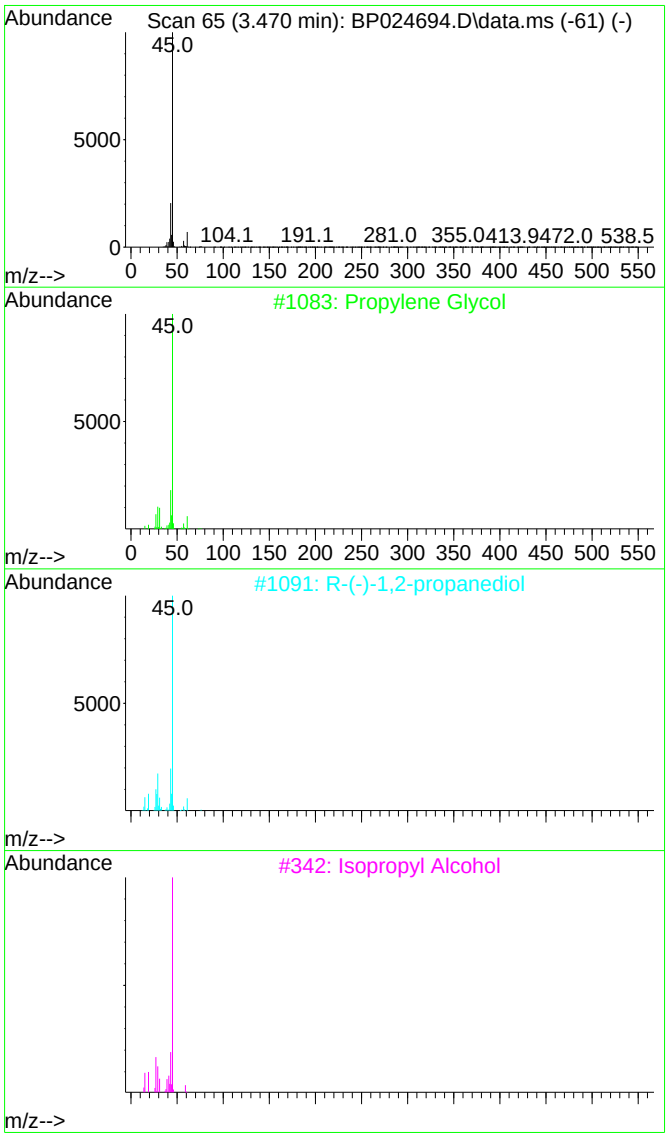
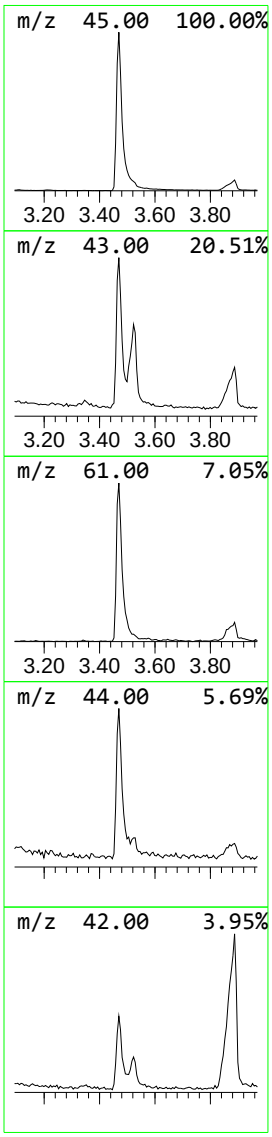
****

Peak Number 1 Propylene Glycol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.470	5.75 ng	320076	1,4-Dichlorobenzene-d4	7.658

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propylene Glycol	76	C3H8O2	000057-55-6	90
2	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	78
3	Isopropyl Alcohol	60	C3H8O	000067-63-0	56
4	Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
5	Formamide	45	CH3NO	000075-12-7	5





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Data File : BP024694.D
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Operator : RC/JU
Sample : Q2033-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TW-WTS-08

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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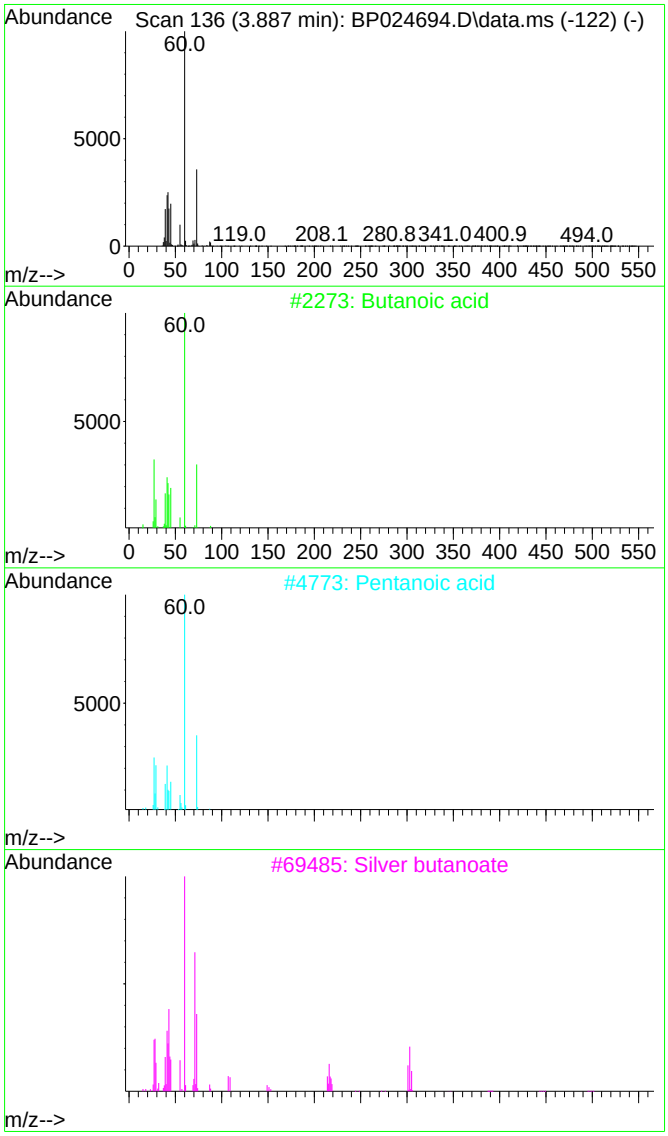
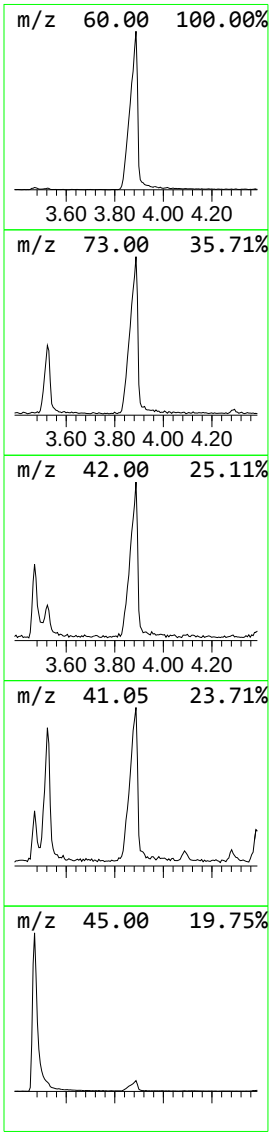
****

Peak Number 2 Butanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.887	5.51 ng	306263	1,4-Dichlorobenzene-d4	7.658

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid	88	C4H8O2	000107-92-6	90
2	Pentanoic acid	102	C5H10O2	000109-52-4	40
3	Silver butanoate	194	C4H7AgO2	005076-24-4	40
4	Hydrazine, 1,1-dimethyl-	60	C2H8N2	000057-14-7	9
5	Formamidoxime	60	CH4N2O	000624-82-8	9





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Operator : RC/JU
Sample : Q2033-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TW-WTS-08

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
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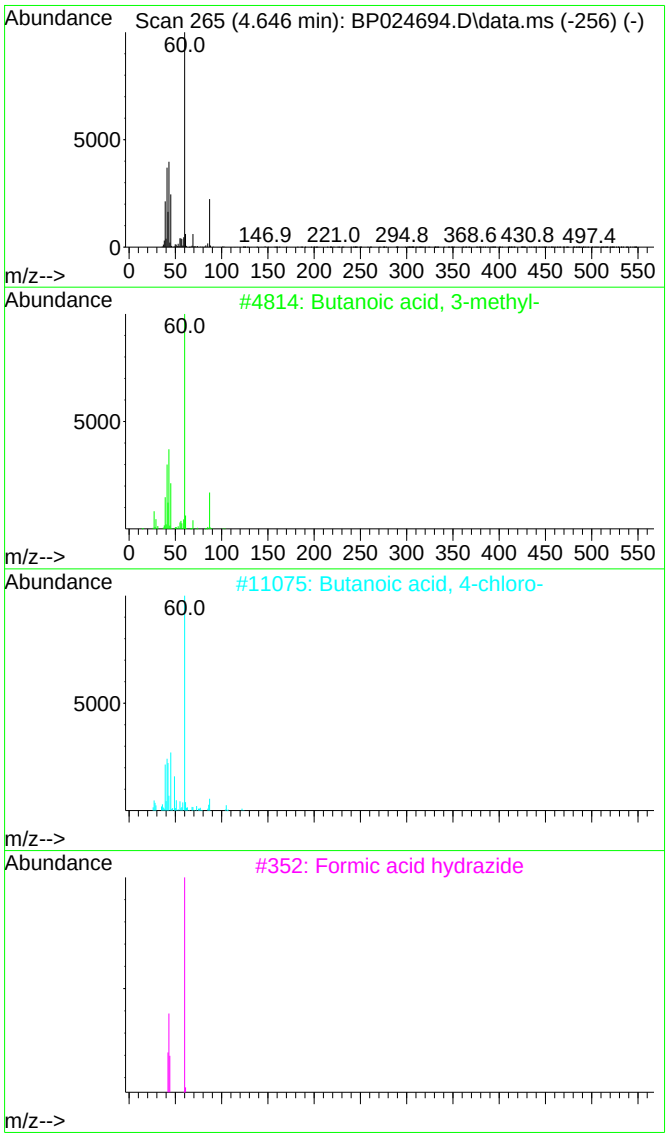
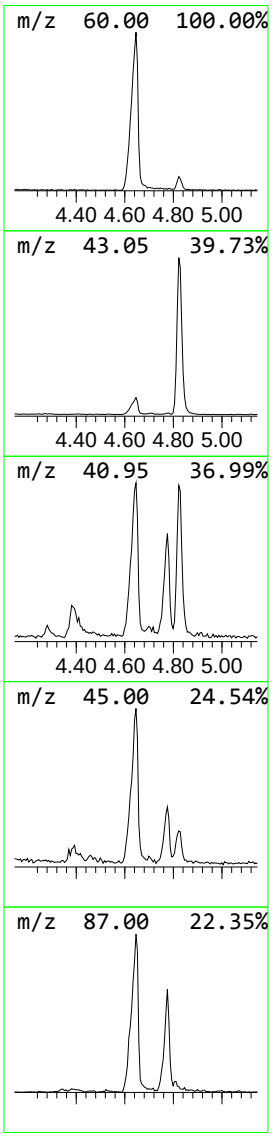
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Butanoic acid, 3-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.646	3.52 ng	195794	1,4-Dichlorobenzene-d4	7.658

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	90
2	Butanoic acid, 4-chloro-	122	C4H7ClO2	000627-00-9	50
3	Formic acid hydrazide	60	CH4N2O	000624-84-0	47
4	Hexanoic acid	116	C6H12O2	000142-62-1	38
5	Pentanoic acid	102	C5H10O2	000109-52-4	33





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Acq On : 19 May 2025 18:19
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Sample : Q2033-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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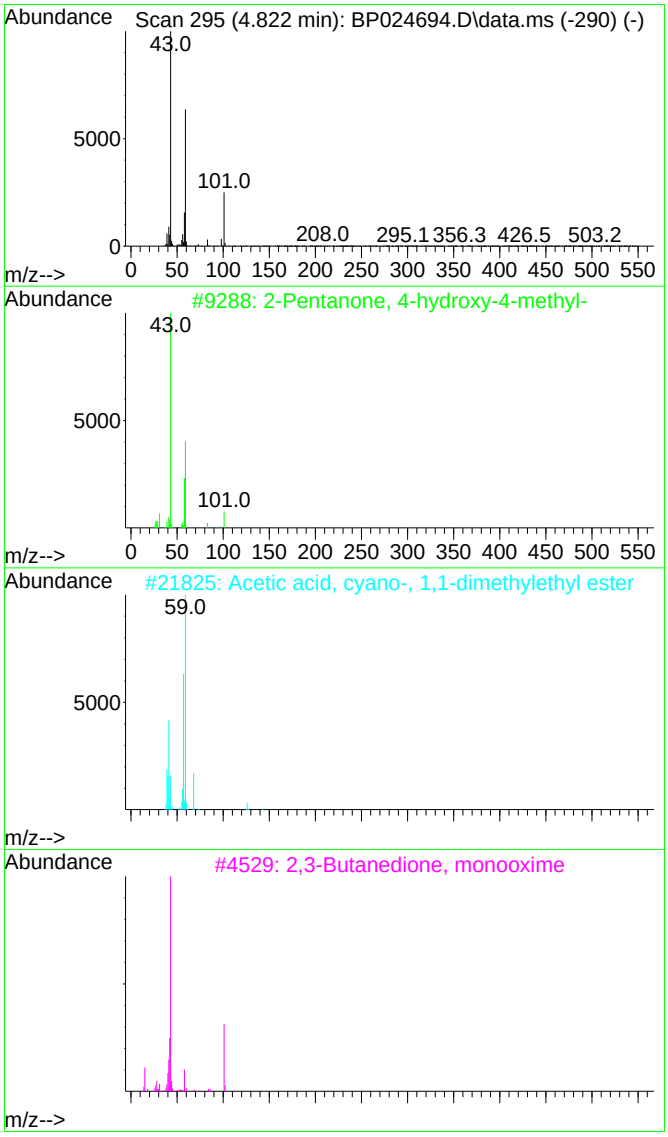
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

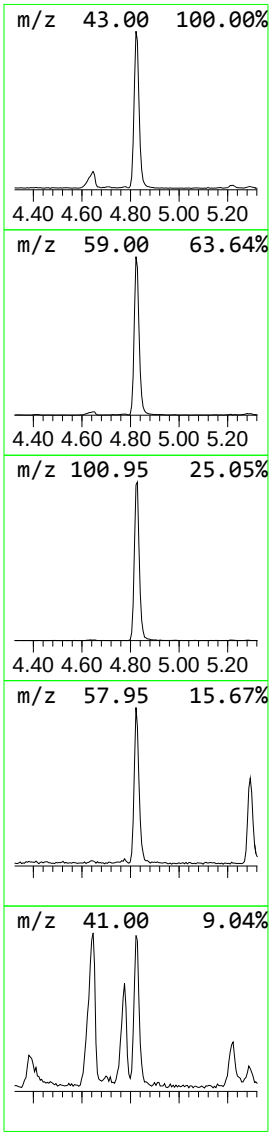
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.822	8.19 ng	455480	1,4-Dichlorobenzene-d4	7.658

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4	2,5,8,11-Tetraoxadodecane	178	C8H18O4	000112-49-2	23
5	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051925\
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Acq On : 19 May 2025 18:19
Operator : RC/JU
Sample : Q2033-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TW-WTS-08

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

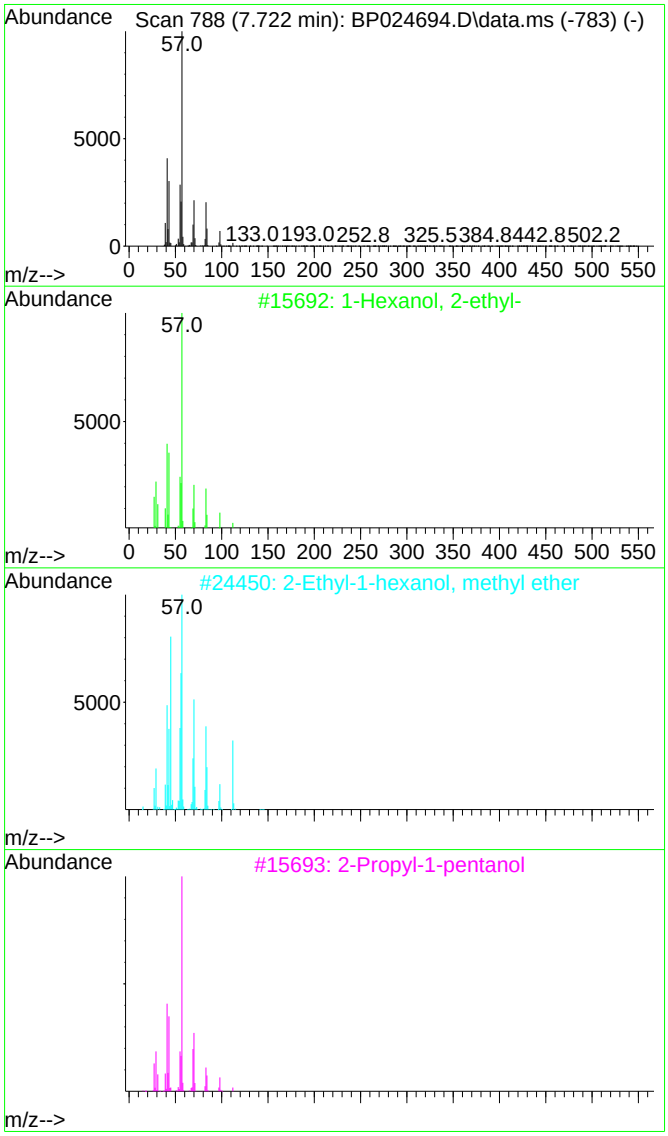
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

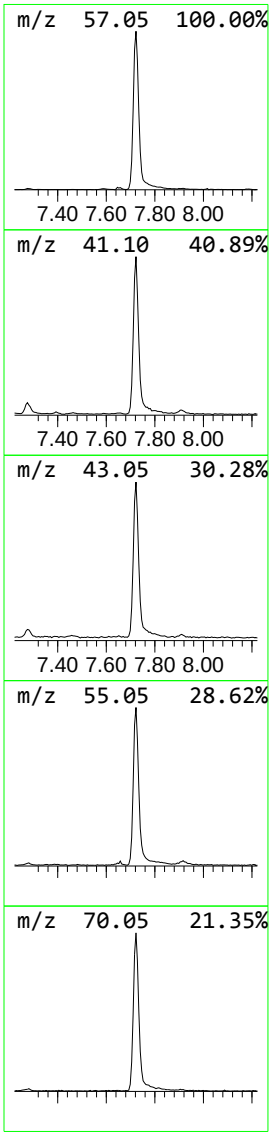
****

Peak Number 5 1-Hexanol, 2-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.722	13.72 ng	763449	1,4-Dichlorobenzene-d4	7.658

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2	2-Ethyl-1-hexanol, methyl ether	144	C9H20O	1000365-10-2	50
3	2-Propyl-1-pentanol	130	C8H18O	058175-57-8	50
4	Pentadecafluorooctanoic acid, 2-...	526	C16H17F15O2	1000406-80-9	47
5	Carbonic acid, heptyl vinyl ester	186	C10H18O3	1010383-25-4	47





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ALS Vial : 10 Sample Multiplier: 1

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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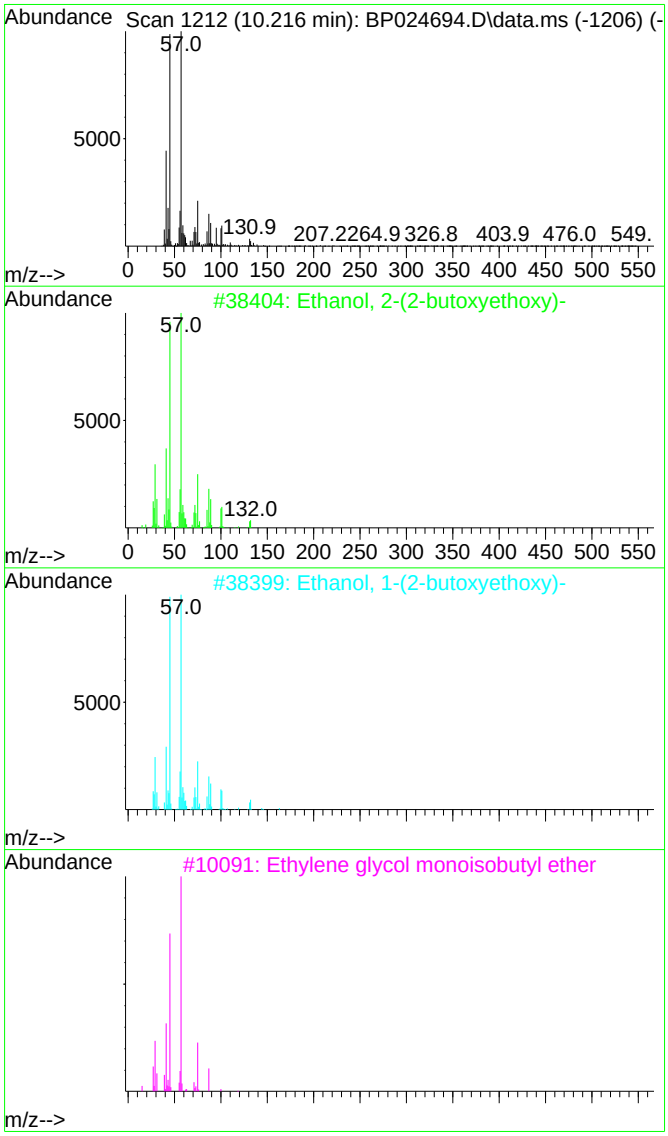
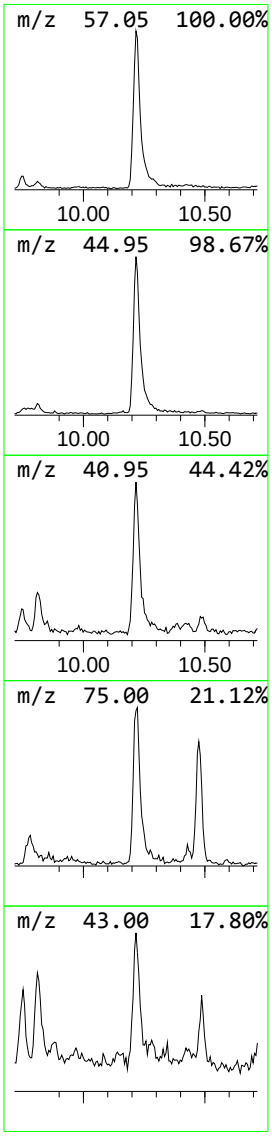
****
Peak Number 6 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.216	2.77 ng	215375	Naphthalene-d8	10.428

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	90
3	Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	59
4	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	50
5	Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	50





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ALS Vial : 10 Sample Multiplier: 1

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

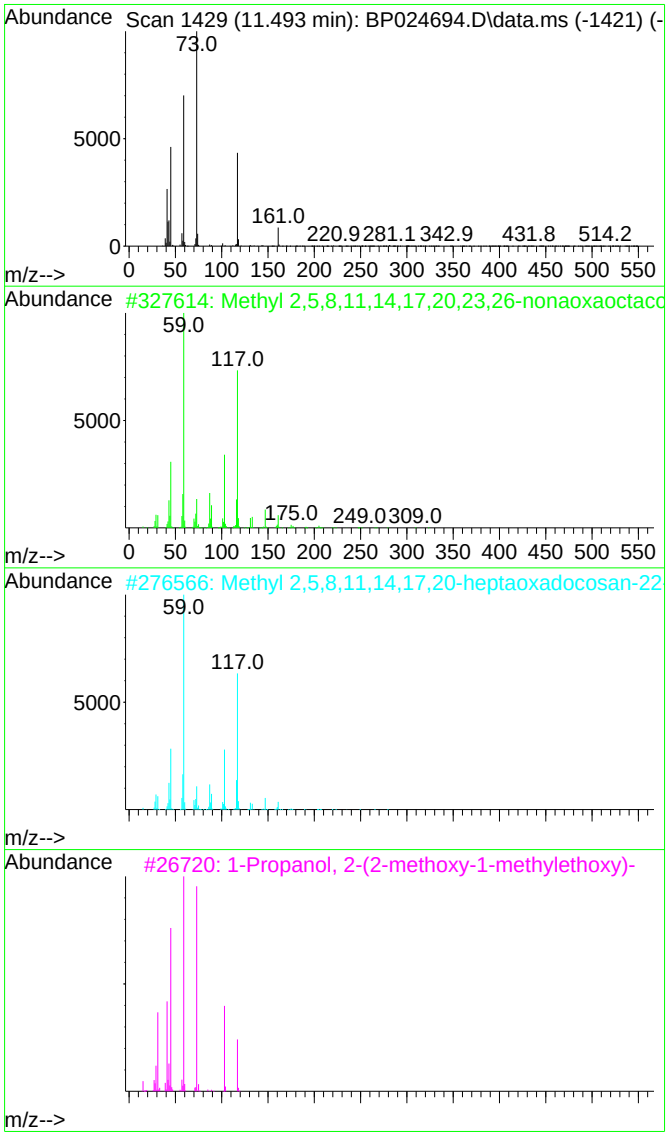
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

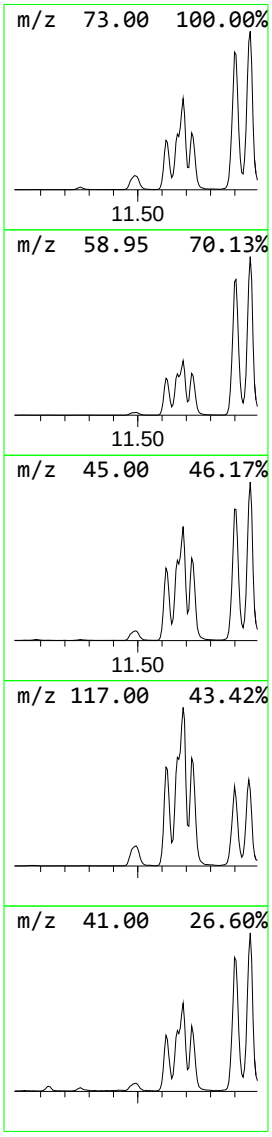
****

Peak Number 7 unknown11.493 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.493	3.76 ng	292065	Naphthalene-d8	10.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl 2,5,8,11,14,17,20,23,26-n...	456	C20H40O11	1000366-81-1	28
2	Methyl 2,5,8,11,14,17,20-heptaox...	368	C16H32O9	1010366-81-3	28
3	1-Propanol, 2-(2-methoxy-1-methy...	148	C7H16O3	055956-21-3	28
4	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	25
5	2,2,5,8,11-pentamethyl-3,6,9,12-...	278	C13H30O4Si	1000473-16-8	17





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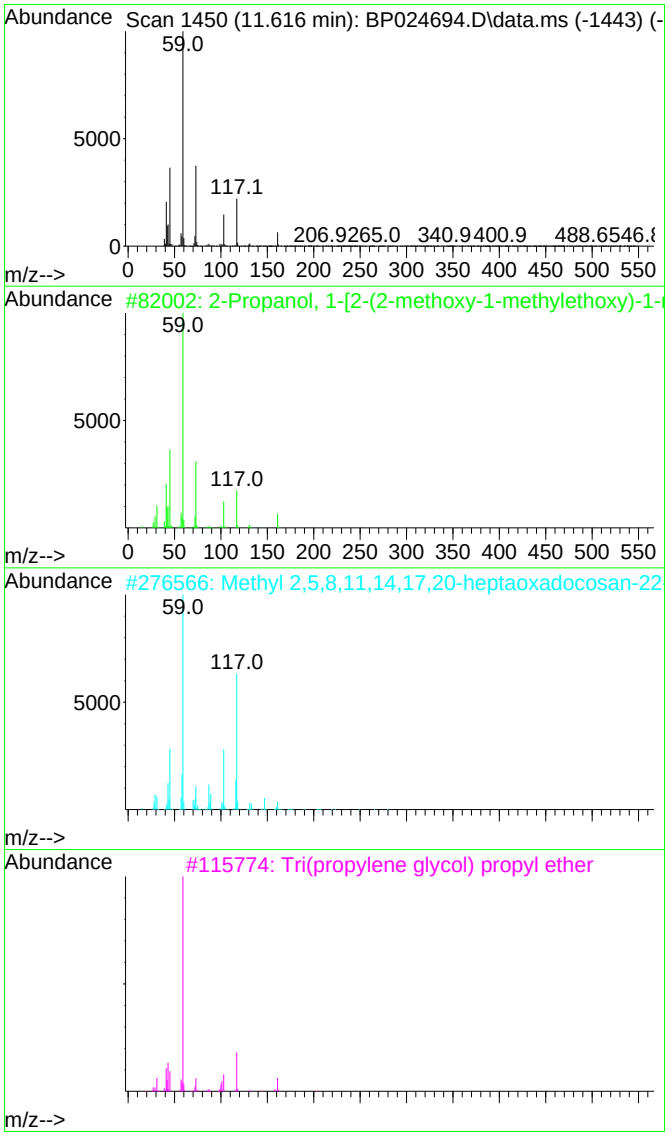
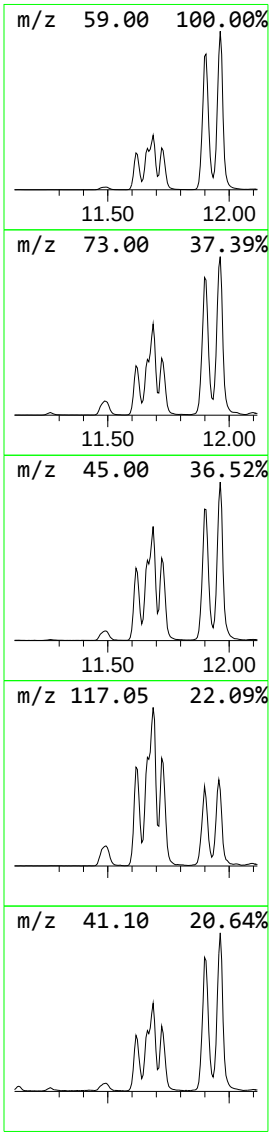
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 2-Propanol, 1-[2-(2-methoxy... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	17.59 ng	1365320	Naphthalene-d8	10.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	90
2	Methyl 2,5,8,11,14,17,20-heptaox...	368	C16H32O9	1010366-81-3	64
3	Tri(propylene glycol) propyl ether	234	C12H26O4	096077-04-2	59
4	Tri(1,2-propyleneglycol), monome...	206	C10H22O4	1000262-26-6	56
5	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	56





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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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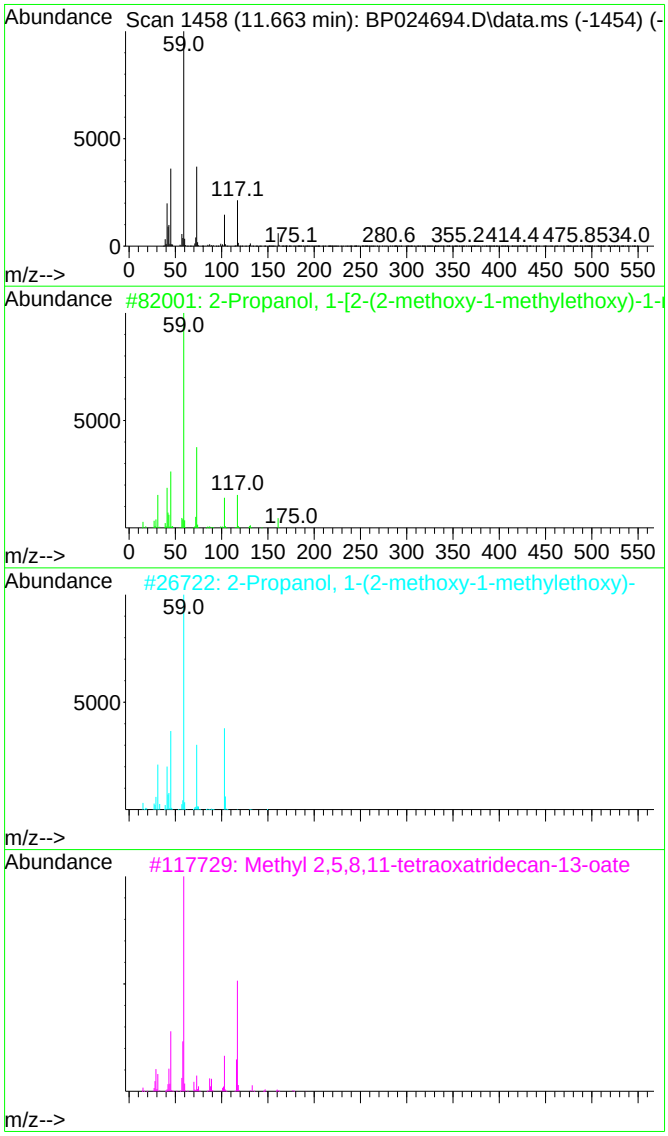
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

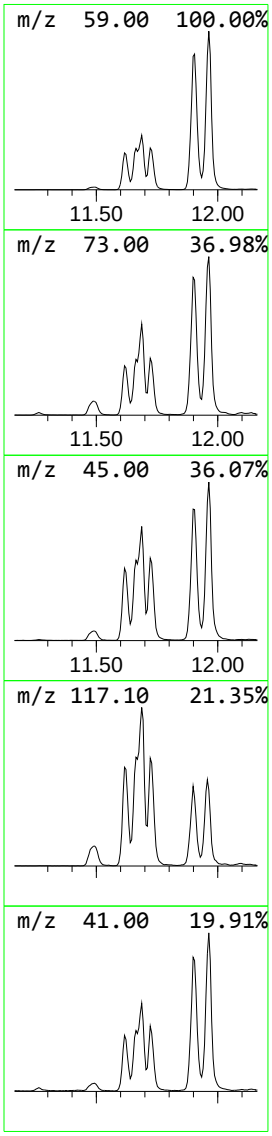
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.663	15.13 ng	1174250	Naphthalene-d8	10.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	90
2	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	64
3	Methyl 2,5,8,11-tetraoxatridecan...	236	C10H20O6	1000366-78-2	59
4	Propanol, [(butoxymethylethoxy)m...	248	C13H28O4	055934-93-5	50
5	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	50





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Data File : BP024694.D
Acq On : 19 May 2025 18:19
Operator : RC/JU
Sample : Q2033-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TW-WTS-08

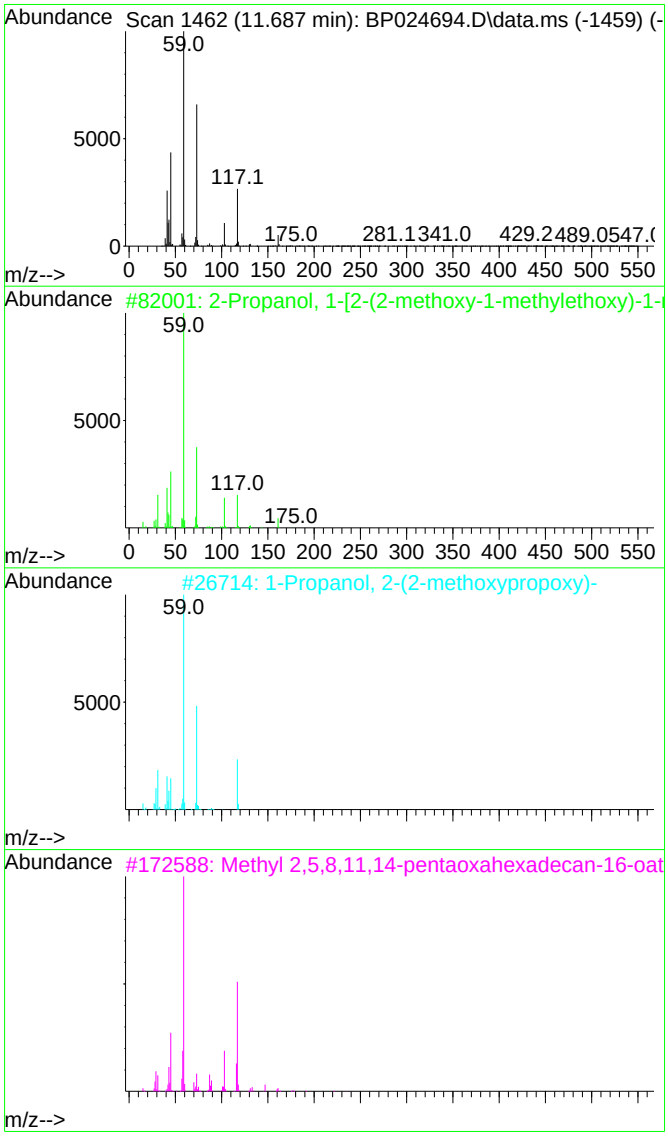
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

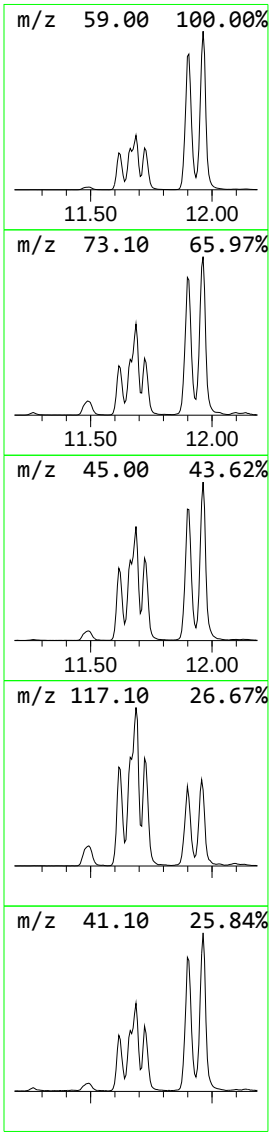
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1-Propanol, 2-(2-methoxypro... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.687	24.49 ng	1901210	Naphthalene-d8	10.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	78
2	1-Propanol, 2-(2-methoxypropoxy)-	148	C7H16O3	013588-28-8	59
3	Methyl 2,5,8,11,14-pentaoxahexad...	280	C12H24O7	1000366-81-5	56
4	Propanol, [(butoxymethylethoxy)m...	248	C13H28O4	055934-93-5	53
5	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051925\
Data File : BP024694.D
Acq On : 19 May 2025 18:19
Operator : RC/JU
Sample : Q2033-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TW-WTS-08

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

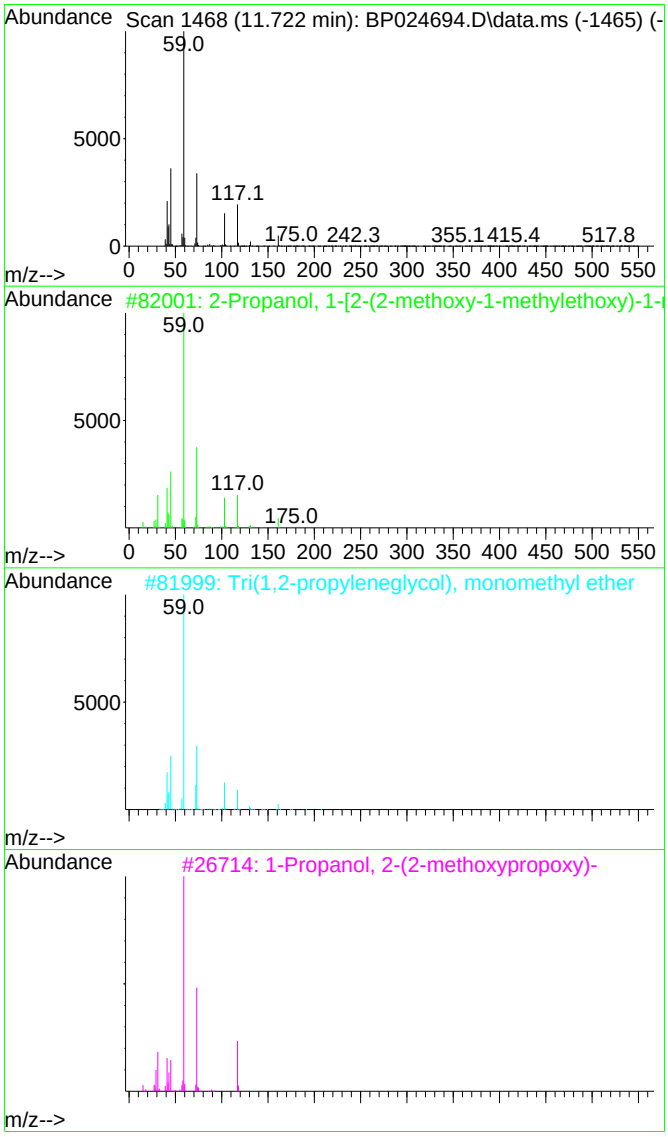
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

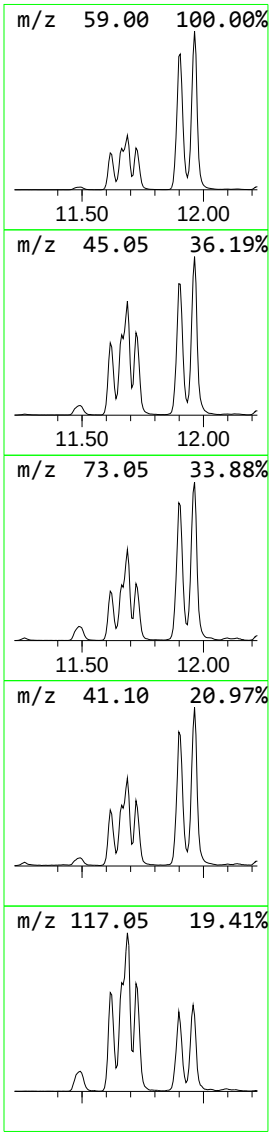
****

Peak Number 11 Tri(1,2-propyleneglycol), m... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.722	18.80 ng	1458880	Naphthalene-d8	10.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	86
2	Tri(1,2-propyleneglycol), monome...	206	C10H22O4	1000262-26-6	78
3	1-Propanol, 2-(2-methoxypropoxy)-	148	C7H16O3	013588-28-8	59
4	2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	53
5	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051925\
Data File : BP024694.D
Acq On : 19 May 2025 18:19
Operator : RC/JU
Sample : Q2033-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TW-WTS-08

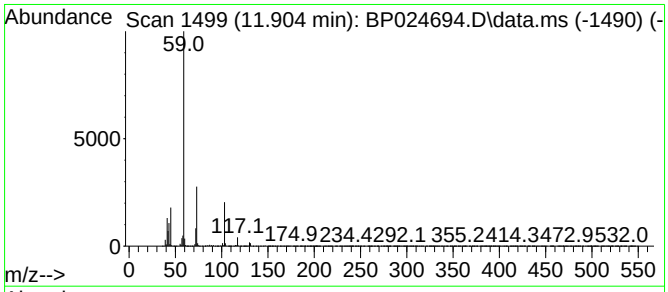
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

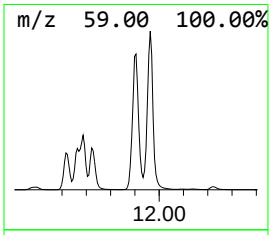
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

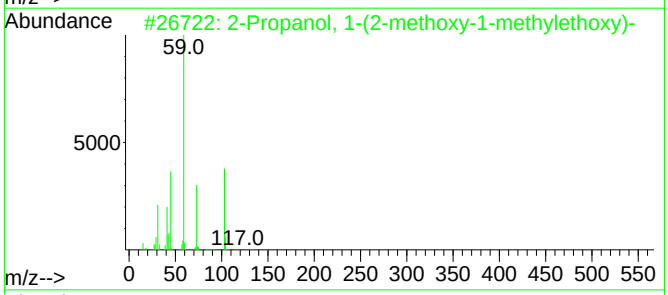
****
Peak Number 12 Hexaethylene glycol dimethy... Concentration Rank 2

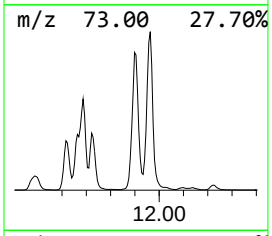
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.904	57.71 ng	4479050	Naphthalene-d8	10.428

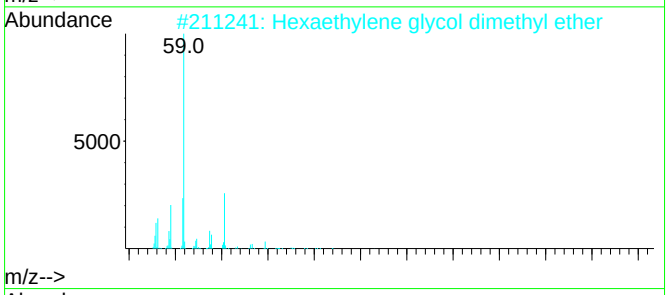
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	78
2	Hexaethylene glycol dimethyl ether	310	C14H30O7	001072-40-8	64
3	Tri(1,2-propyleneglycol), monome...	206	C10H22O4	1000262-26-6	64
4	Methane, diethoxy-	104	C5H12O2	000462-95-3	53
5	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	50

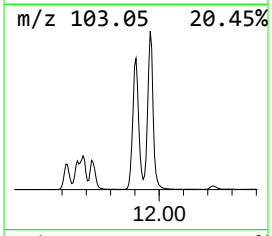


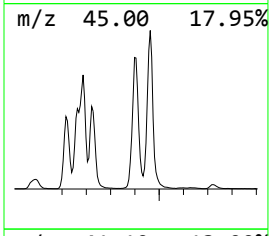


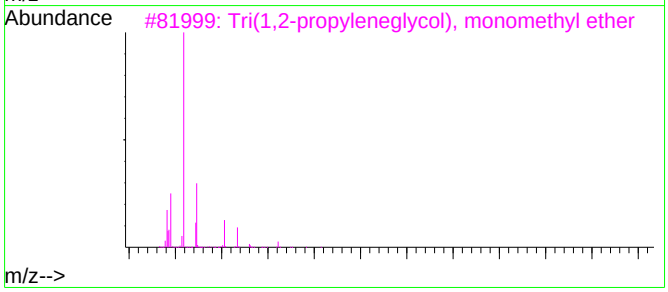


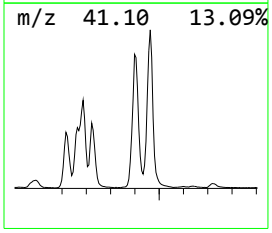












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051925\
Data File : BP024694.D
Acq On : 19 May 2025 18:19
Operator : RC/JU
Sample : Q2033-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TW-WTS-08

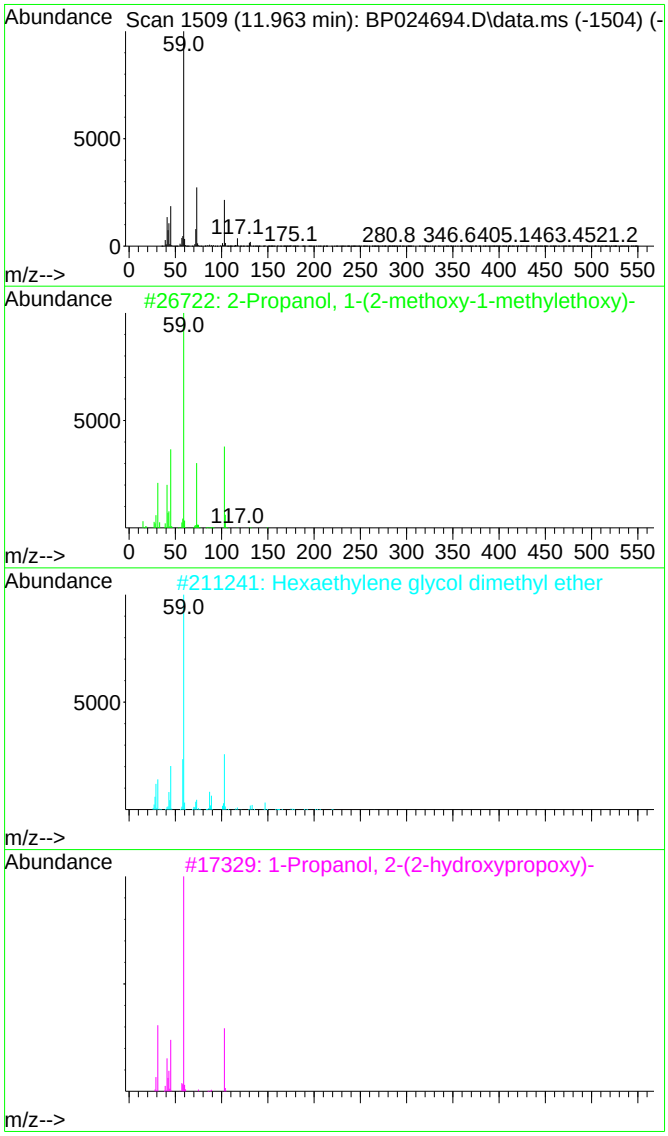
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

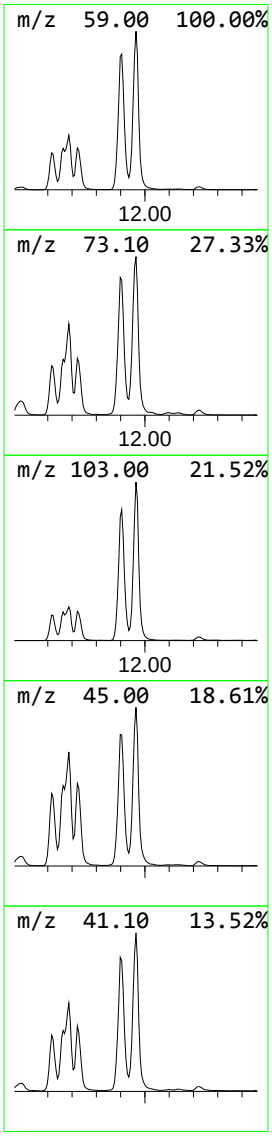
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 2,5,8,11,14,17-Hexaoxaoctad... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.963	62.33 ng	4837770	Naphthalene-d8	10.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	72
2	Hexaethylene glycol dimethyl ether	310	C14H30O7	001072-40-8	64
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59
4	2,5,8,11,14,17-Hexaoxaoctadecane	266	C12H26O6	001191-87-3	50
5	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	50





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Data File : BP024694.D
Acq On : 19 May 2025 18:19
Operator : RC/JU
Sample : Q2033-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TW-WTS-08

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

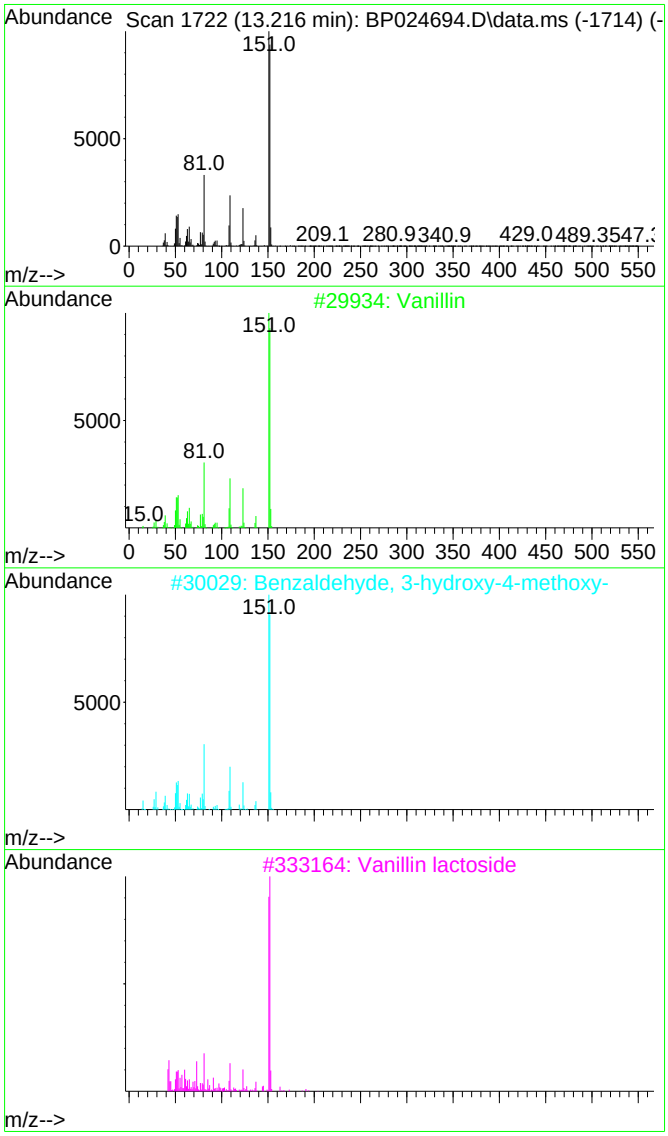
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

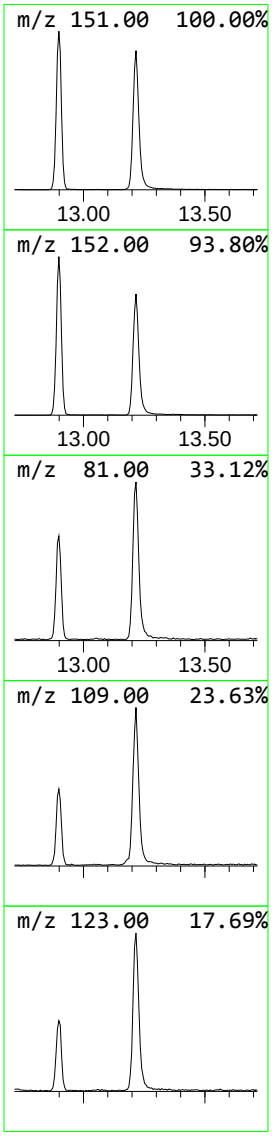
****

Peak Number 14 Vanillin Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.216	7.19 ng	780472	Acenaphthene-d10	14.293

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Vanillin	152	C8H8O3	000121-33-5	97
2	Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	97
3	Vanillin lactoside	476	C20H28O13	1000129-94-7	72
4	Benzaldehyde, 2-hydroxy-4-methoxy-	152	C8H8O3	000673-22-3	64
5	Benzaldehyde, 2,4-dihydroxy-6-me...	152	C8H8O3	000487-69-4	58





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Data File : BP024694.D
Acq On : 19 May 2025 18:19
Operator : RC/JU
Sample : Q2033-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TW-WTS-08

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

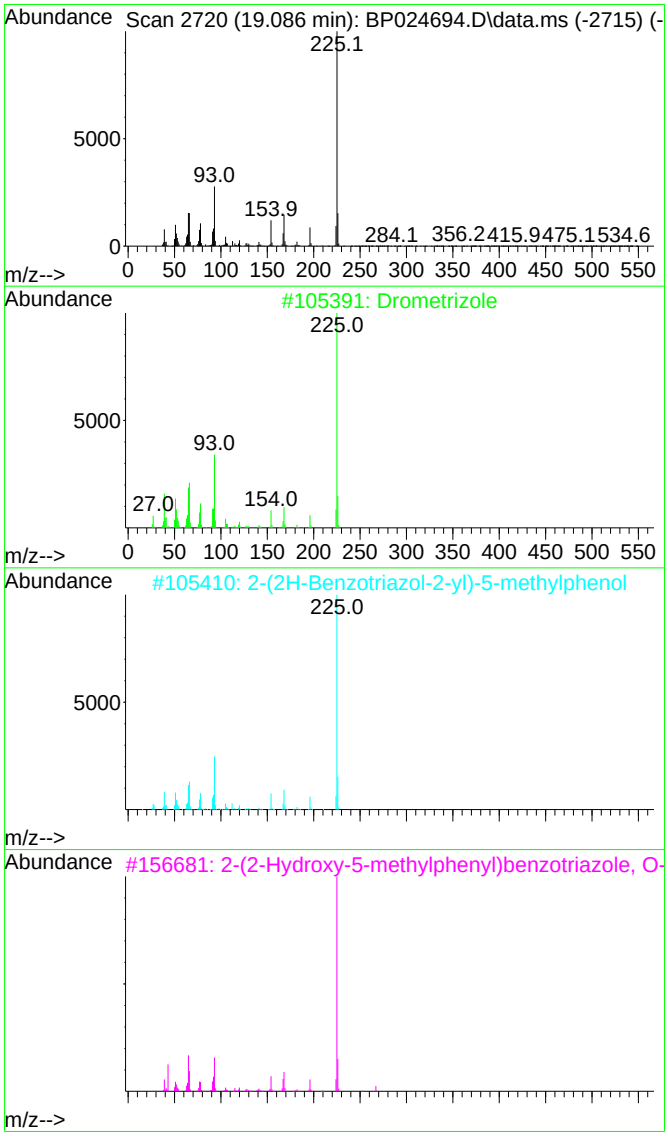
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

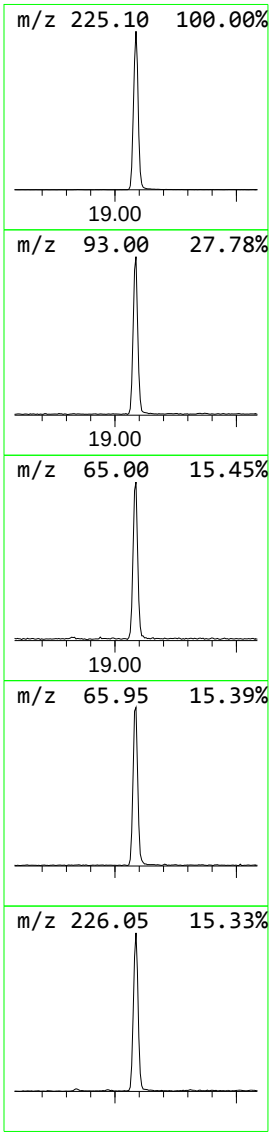
****

Peak Number 15 Drometrizole Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.086	7.13 ng	963116	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Drometrizole	225	C13H11N3O	002440-22-4	98
2	2-(2H-Benzotriazol-2-yl)-5-methy...	225	C13H11N3O	004998-48-5	97
3	2-(2-Hydroxy-5-methylphenyl)benz...	267	C15H13N3O2	1000384-15-2	87
4	9(10H)-Acridinone, 1-hydroxy-10-...	225	C14H11NO2	016584-54-6	50
5	Propionitrile, 3-(5-methyl-3-phe...	225	C13H11N3O	1000263-77-0	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051925\
Data File : BP024694.D
Acq On : 19 May 2025 18:19
Operator : RC/JU
Sample : Q2033-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TW-WTS-08

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.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
Propylene Glycol	3.470	5.8	ng	320076	1	7.658	1112610	20.0
Butanoic acid	3.887	5.5	ng	306263	1	7.658	1112610	20.0
Butanoic acid, ...	4.646	3.5	ng	195794	1	7.658	1112610	20.0
2-Pentanone, 4-...	4.822	8.2	ng	455480	1	7.658	1112610	20.0
1-Hexanol, 2-et...	7.722	13.7	ng	763449	1	7.658	1112610	20.0
Ethanol, 2-(2-b...	10.216	2.8	ng	215375	2	10.428	1552360	20.0
unknown11.493	11.493	3.8	ng	292065	2	10.428	1552360	20.0
2-Propanol, 1-[...	11.616	17.6	ng	1365320	2	10.428	1552360	20.0
2-Propanol, 1-(...	11.663	15.1	ng	1174250	2	10.428	1552360	20.0
1-Propanol, 2-(...	11.687	24.5	ng	1901210	2	10.428	1552360	20.0
Tri(1,2-propyle...	11.722	18.8	ng	1458880	2	10.428	1552360	20.0
Hexaethylene gl...	11.904	57.7	ng	4479050	2	10.428	1552360	20.0
2,5,8,11,14,17-...	11.963	62.3	ng	4837770	2	10.428	1552360	20.0
Vanillin	13.216	7.2	ng	780472	3	14.293	2170260	20.0
Drometrizole	19.086	7.1	ng	963116	4	17.087	2700090	20.0