LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

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_G\Methods\8270-BG122923.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG060253.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.119	3	5	10	rVB	312706	328387	4.07%	0.576%
2	4.024	156	159	169	rVB5	48741	88578	1.10%	0.155%
3	4.124	169	176	180	rBV	88031	140844	1.75%	0.247%
4	5.510	405	412	420	rBV	2208276	3565310	44.23%	6.258%
5	5.945	479	486	494	rVB8	61749	126535	1.57%	0.222%
6	6.427	559	568	574	rBV4	49940	117666	1.46%	0.207%
7	6.914	641	651	658	rBV2	239290	452084	5.61%	0.794%
8	6.979	658	662	666	rVB2	57478	86587	1.07%	0.152%
9	7.097	675	682	700	rBV	2313064	4215519	52.30%	7.399%
10	7.937	816	825	833	rBV	615520	1142891	14.18%	2.006%
11	8.066	841	847	852	rBV3	44027	93580	1.16%	0.164%
12	8.313	880	889	895	rBV	228565	412261	5.11%	0.724%
13	8.424	903	908	912	rBV2	109285	180719	2.24%	0.317%
14	8.577	928	934	939	rBV4	166987	325447	4.04%	0.571%
15	8.700	947	955	959	rBV4	59226	110068	1.37%	0.193%
16	9.041	1004	1013	1017	rBV	371523	606681	7.53%	1.065%
17	9.100	1017	1023	1038	rVB	1281381	2542950	31.55%	4.464%
18	9.388	1064	1072	1085	rVV	23918	85538	1.06%	0.150%
19	9.564	1097	1102	1107	rVB	187901	287342	3.56%	0.504%
20	9.623	1107	1112	1121	rBV	121049	222494	2.76%	0.391%
21	9.823	1140	1146	1152	rBV2	56144	107331	1.33%	0.188%
22	9.987	1168	1174	1185	rVB	169432	380770	4.72%	0.668%
23	10.134	1193	1199	1208	rBV3	308497	597436	7.41%	1.049%
24	10.440	1246	1251	1256	rVB2	69290	106386	1.32%	0.187%
25	10.739	1290	1302	1307	rBV	905003	1820919	22.59%	3.196%
26	10.792	1307	1311	1317	rVV2	320049	621262	7.71%	1.090%
27	10.851	1317	1321	1327	rVB2	55931	91296	1.13%	0.160%
28	10.951	1333	1338	1347	rVB4	54023	92120	1.14%	0.162%
29	11.474	1420	1427	1434	rBV4	35645	92339	1.15%	0.162%
30	11.650	1451	1457	1464	rBV2	90790	168736	2.09%	0.296%
31	11.844	1485	1490	1496	rVB2	53790	88605	1.10%	0.156%
32	12.396	1578	1584	1594	rVB	312777	533866	6.62%	0.937%
33	12.619	1612	1622	1630	rVB	147917	288864	3.58%	0.507%
34	13.195	1711	1720	1734	rBV	3522135	5405941	67.06%	9.489%
35	13.894	1831	1839	1844	rBV2	43450	88266	1.10%	0.155%
36	14.576	1946	1955	1962	rBV	1878459	2897256	35.94%	5.085%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

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_G\Methods\8270-BG122923.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	16.062	2201	2208	2219	rBV	3287622	4919343	61.03%	8.635%
38	16.151	2219	2223	2230	rVB4	53183	120889	1.50%	0.212%
39	17.320	2416	2422	2427	rBV	2588889	3958342	49.11%	6.948%
40	17.367	2427	2430	2435	rVV	210898	301226	3.74%	0.529%
41	17.443	2438	2443	2449	rVB4	68437	127016	1.58%	0.223%
42	17.508	2449	2454	2459	rBV3	53613	89798	1.11%	0.158%
43	18.207	2567	2573	2577	rBV4	156800	253027	3.14%	0.444%
44	18.795	2669	2673	2677	rVB4	69497	92123	1.14%	0.162%
45	19.376	2767	2772	2778	rBV	228303	347046	4.31%	0.609%
46	19.435	2778	2782	2785	rVV2	116457	153696	1.91%	0.270%
47	19.741	2831	2834	2841	rVB	189980	266821	3.31%	0.468%
48	19.940	2862	2868	2874	rBV	6094461	8060775	100.00%	14.149%
49	20.152	2900	2904	2910	rBV	305276	388927	4.82%	0.683%
50	21.233	3084	3088	3095	rBV3	175379	256526	3.18%	0.450%
51	21.585	3141	3148	3162	rBV2	2458947	4397928	54.56%	7.719%
52	22.396	3281	3286	3296	rVB2	67336	191035	2.37%	0.335%
53	24.764	3678	3689	3700	rBV	1586376	4534344	56.25%	7.959%

Sum of corrected areas: 56971736

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
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Operator : MA/JU
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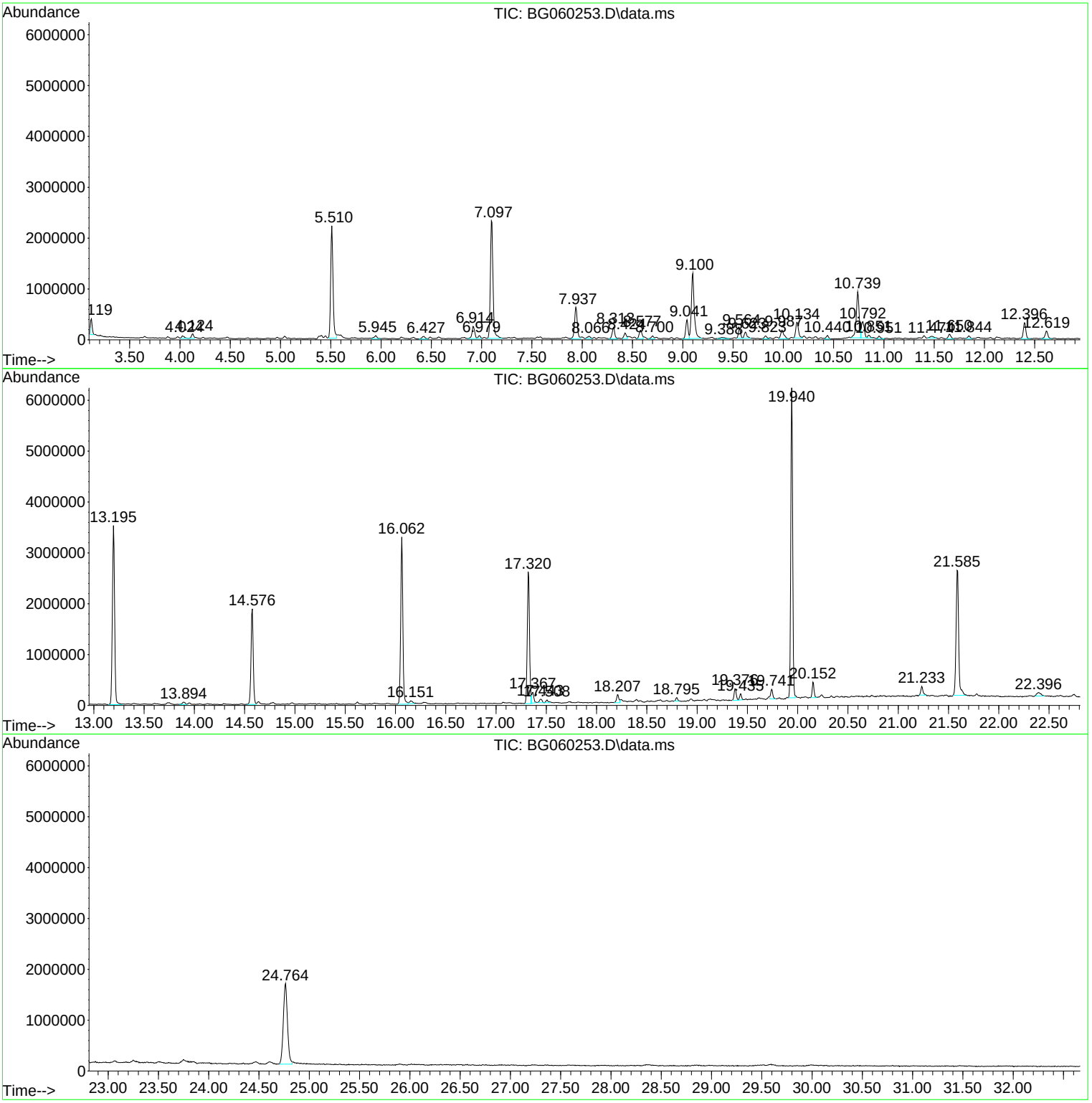
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.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_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

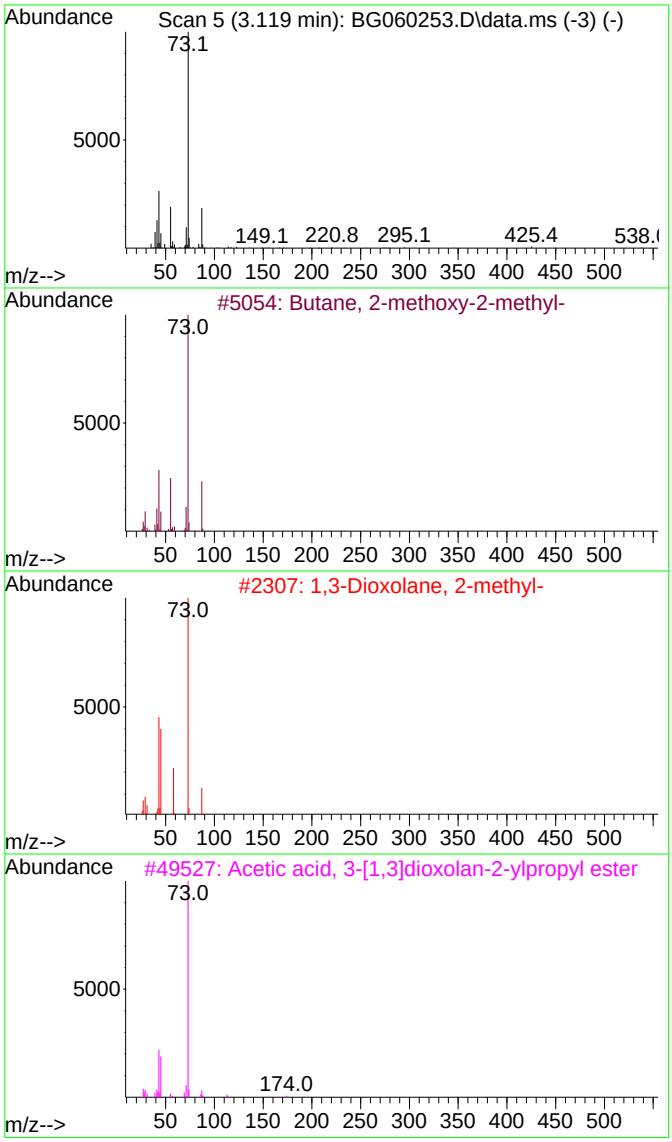
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

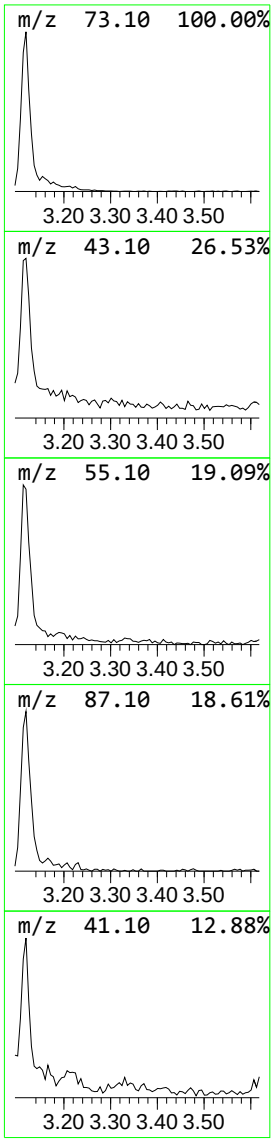
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.119	5.75 ng	328387	1,4-Dichlorobenzene-d4	7.937

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	53
2	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	38
3	Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	1000186-02-3	25
4	1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	23
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

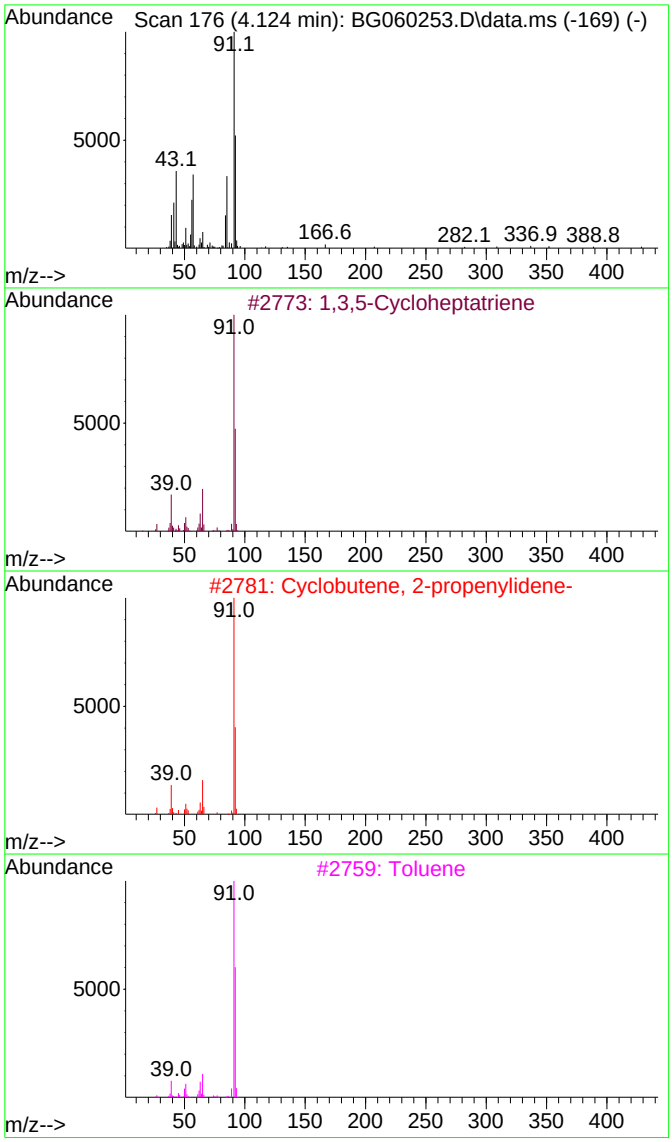
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

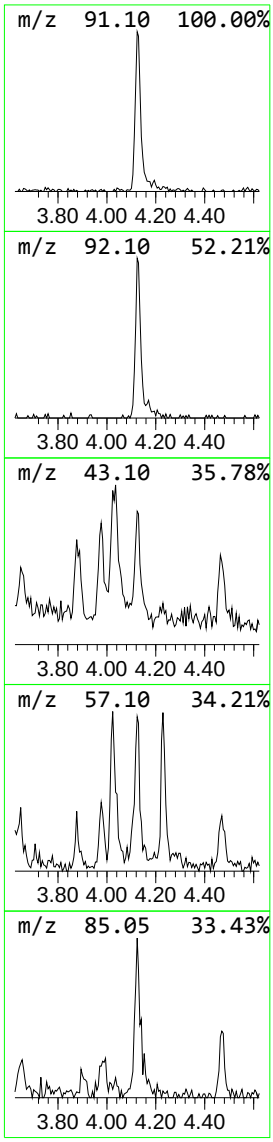
****

Peak Number 2 unknown4.124 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.124	2.46 ng	140844	1,4-Dichlorobenzene-d4	7.937

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	32
2	Cyclobutene, 2-propenylidene-	92	C7H8	052097-85-5	32
3	Toluene	92	C7H8	000108-88-3	32
4	1-Benzylcyclopentanol-1	176	C12H16O	002015-57-8	28
5	Phenylethyl Alcohol	122	C8H10O	000060-12-8	23





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

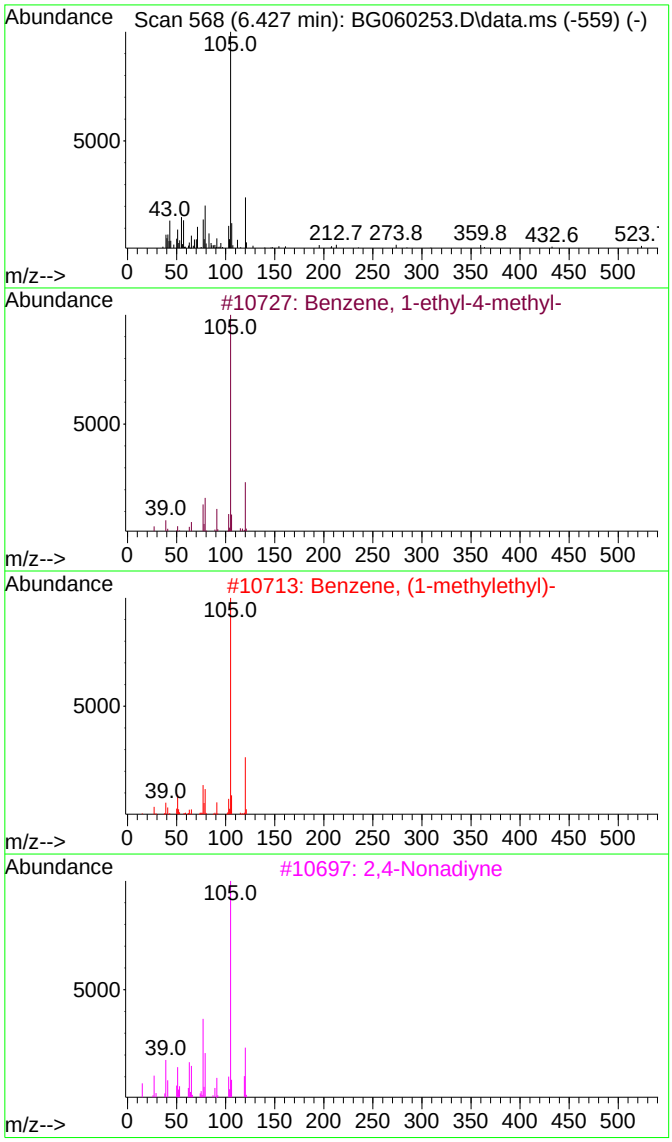
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

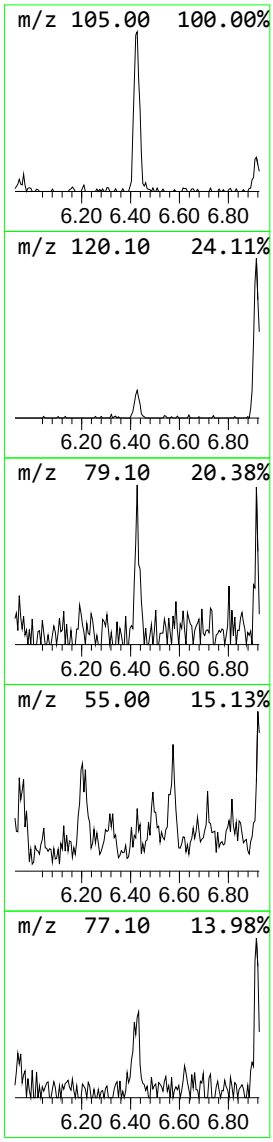
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-ethyl-4-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.427	2.06 ng	117666	1,4-Dichlorobenzene-d4	7.937

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87
2	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	76
3	2,4-Nonadiyne	120	C9H12	063621-15-8	72
4	Cyclohexane, 1,2,4-tris(methylene)-	120	C9H12	014296-81-2	53
5	2,3-Heptadien-5-yne, 2,4-dimethyl-	120	C9H12	041898-89-9	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

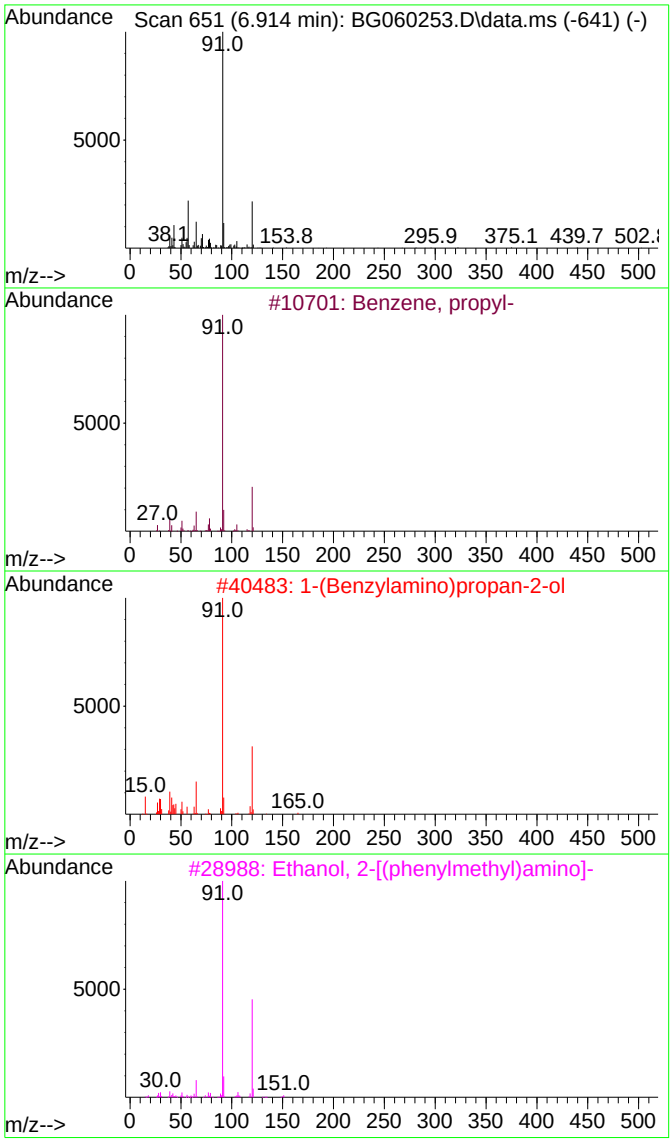
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

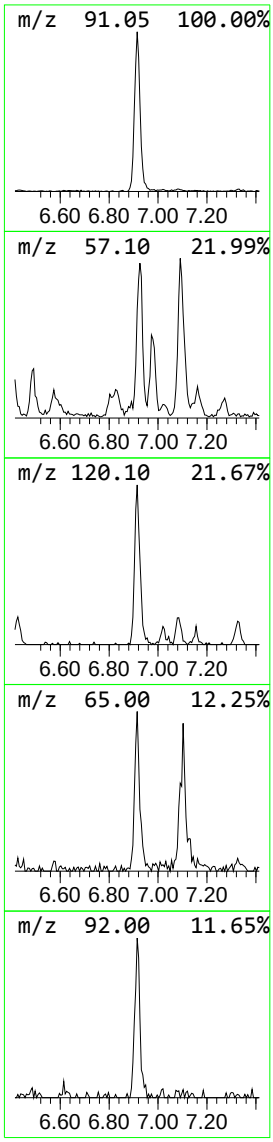
****

Peak Number 5 Benzene, propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.914	7.91 ng	452084	1,4-Dichlorobenzene-d4	7.937

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	81
2	1-(Benzylamino)propan-2-ol	165	C10H15NO	1000486-84-7	64
3	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	64
4	Benzeneacetaldehyde	120	C8H8O	000122-78-1	43
5	Benzoic acid, 2-hydroxy-, phenyl...	228	C14H12O3	000118-58-1	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

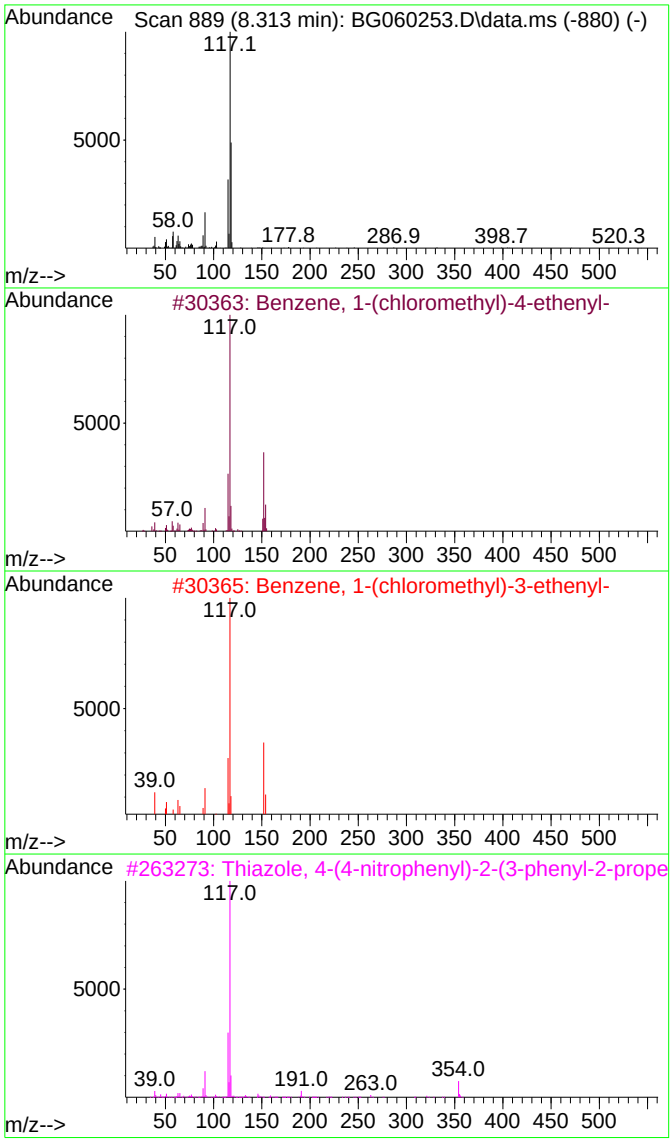
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

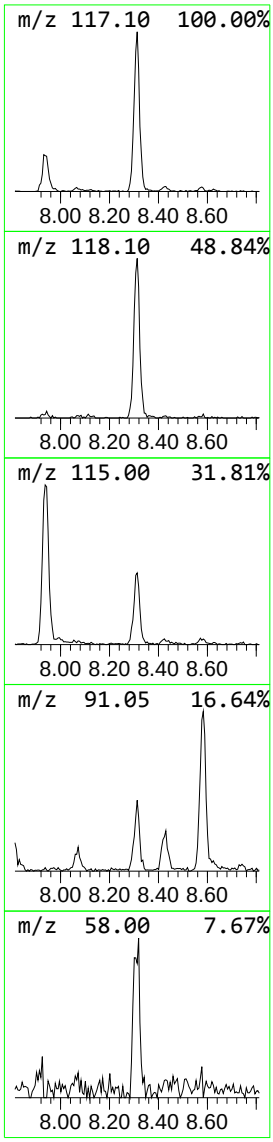
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-(chloromethyl)-4... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.313	7.21 ng	412261	1,4-Dichlorobenzene-d4	7.937

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-(chloromethyl)-4-ethe...	152	C9H9Cl	001592-20-7	59
2	Benzene, 1-(chloromethyl)-3-ethe...	152	C9H9Cl	039833-65-3	53
3	Thiazole, 4-(4-nitrophenyl)-2-(3...	354	C18H14N2O2S2	299921-08-7	53
4	Benzene, (3-chloro-1-propenyl)-	152	C9H9Cl	002687-12-9	53
5	3-Bromo-1-phenyl-1-propene	196	C9H9Br	004392-24-9	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

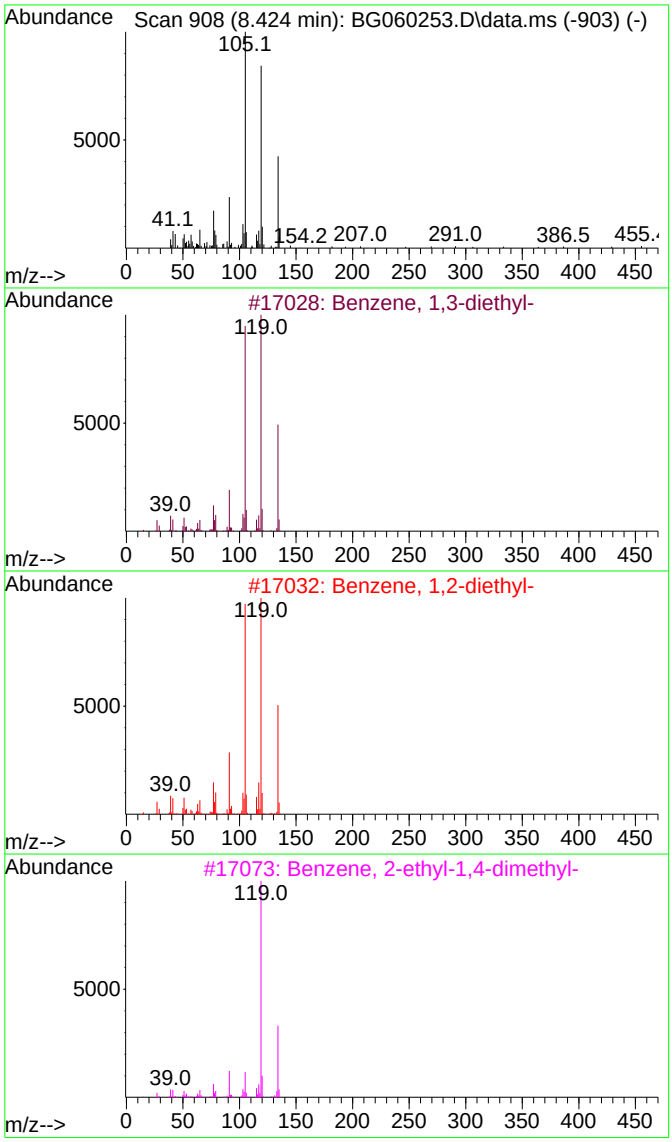
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

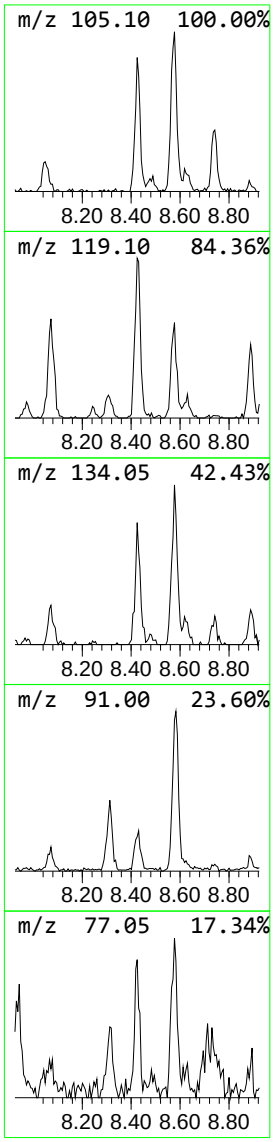
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,3-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.424	3.16 ng	180719	1,4-Dichlorobenzene-d4	7.937

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	91
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
4	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	91
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

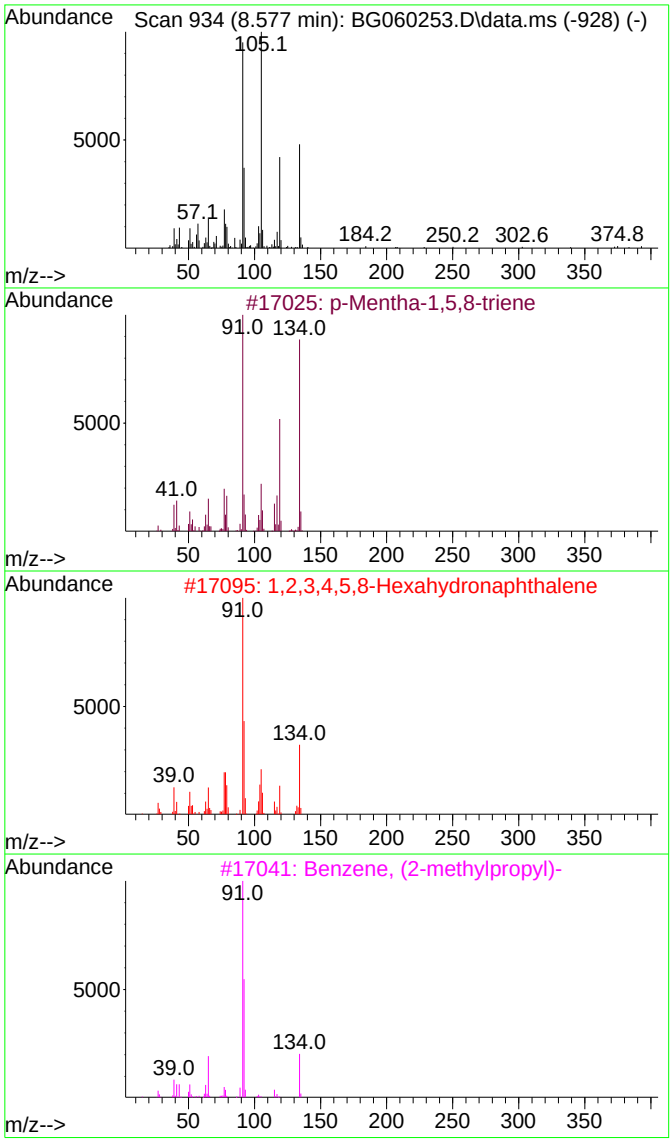
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

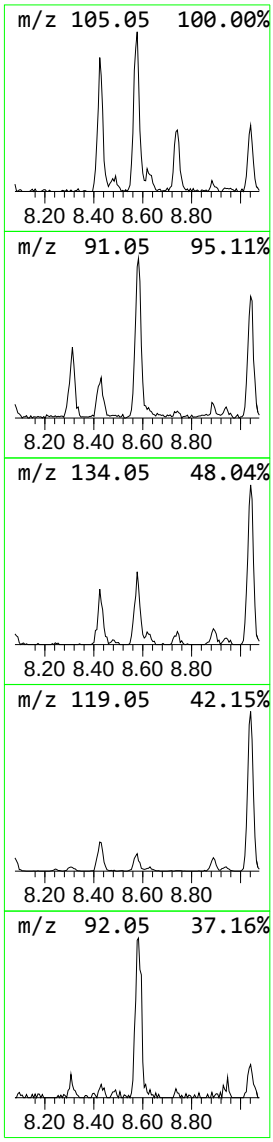
****

Peak Number 8 p-Mentha-1,5,8-triene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.577	5.70 ng	325447	1,4-Dichlorobenzene-d4	7.937

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Mentha-1,5,8-triene	134	C10H14	021195-59-5	87
2	1,2,3,4,5,8-Hexahydronaphthalene	134	C10H14	036231-13-7	53
3	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	45
4	Benzene, n-butyl-	134	C10H14	000104-51-8	43
5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

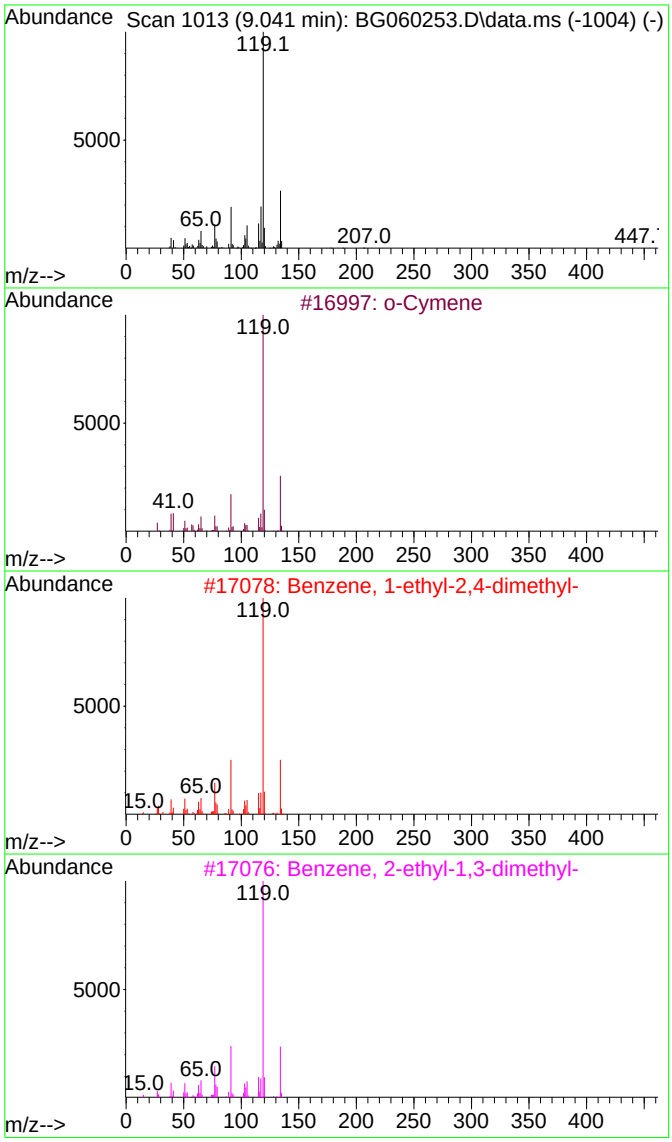
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

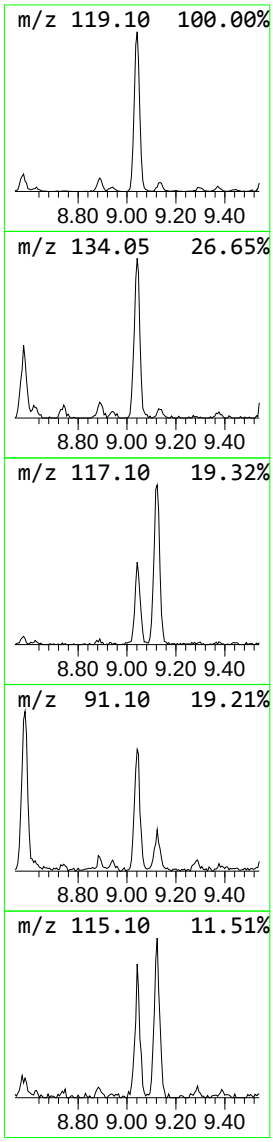
****

Peak Number 9 o-Cymene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.041	10.62 ng	606681	1,4-Dichlorobenzene-d4	7.937

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	94
2	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93
4	p-Cymene	134	C10H14	000099-87-6	93
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91





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Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

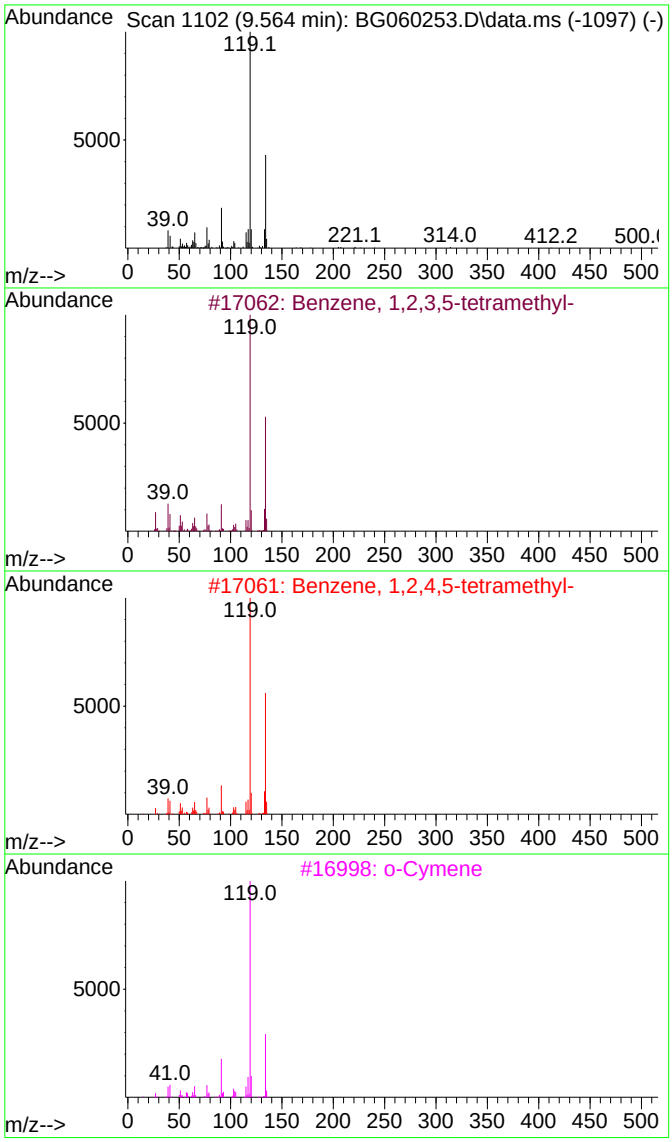
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

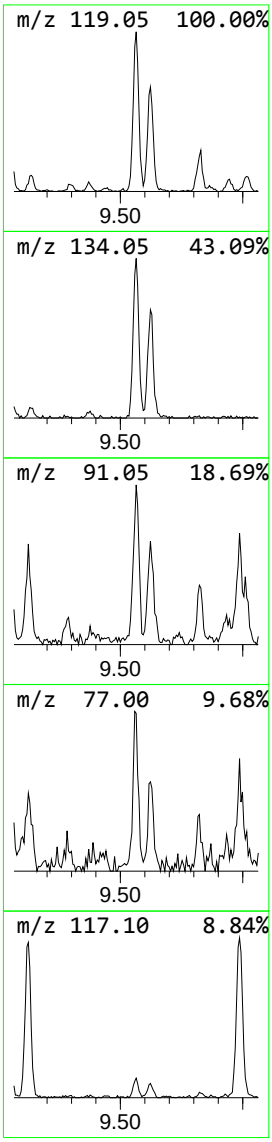
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.564	3.16 ng	287342	Naphthalene-d8	10.739

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
3	o-Cymene	134	C10H14	000527-84-4	93
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.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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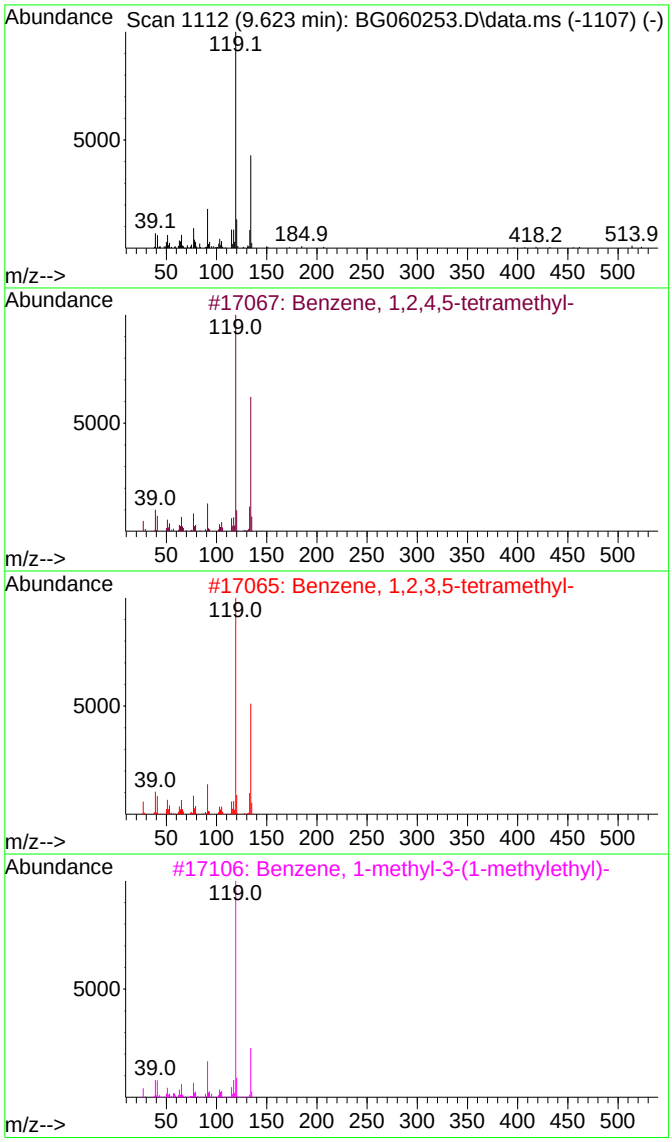
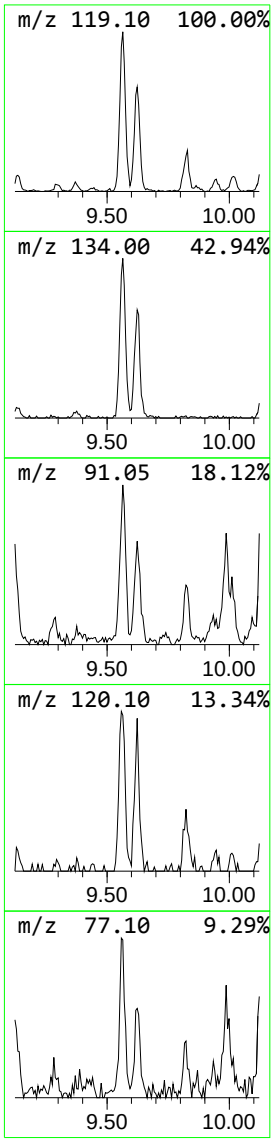
Peak Number 11 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.623	2.44 ng	222494	Naphthalene-d8	10.739

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	87
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	87
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	87
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.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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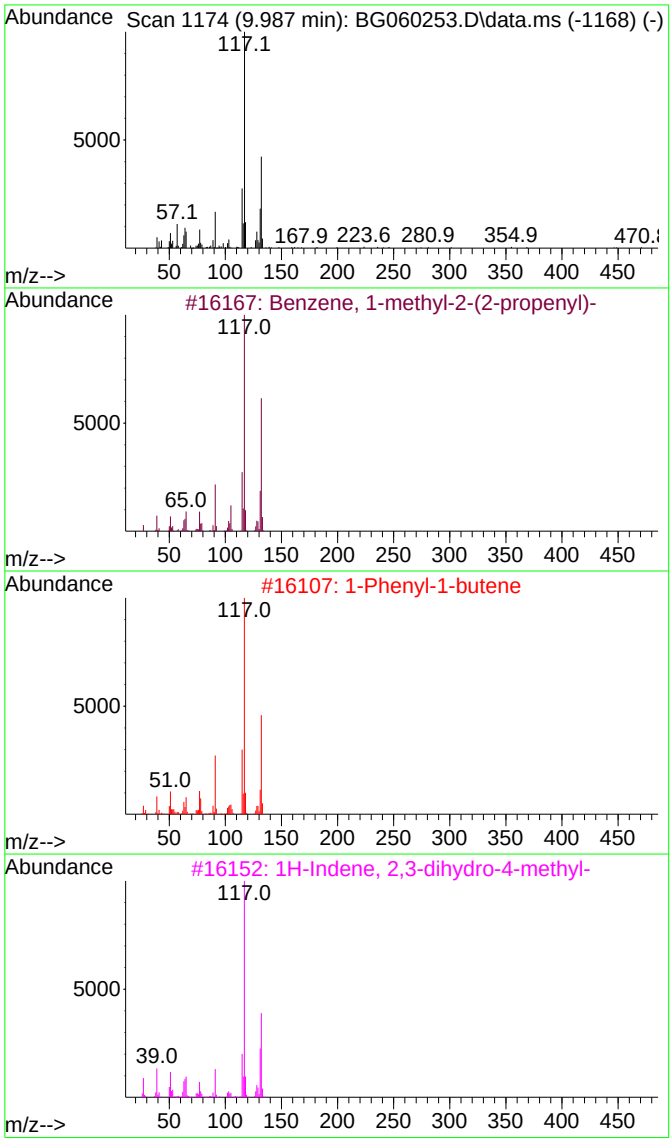
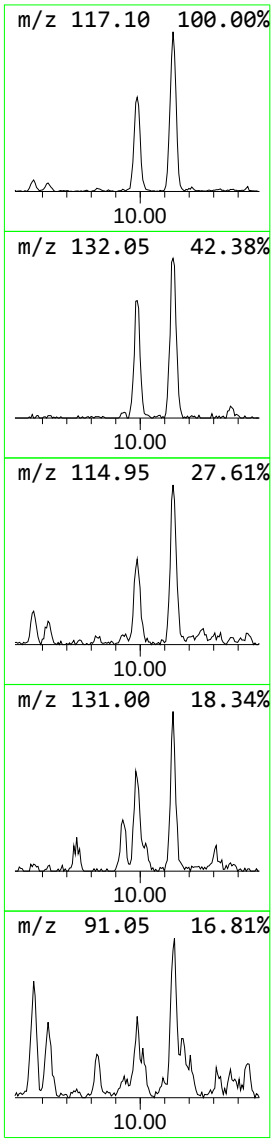
Peak Number 12 Benzene, 1-methyl-2-(2-prop... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.987	4.18 ng	380770	Naphthalene-d8	10.739

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	93
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	90
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	90
5	Indan, 1-methyl-	132	C10H12	000767-58-8	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

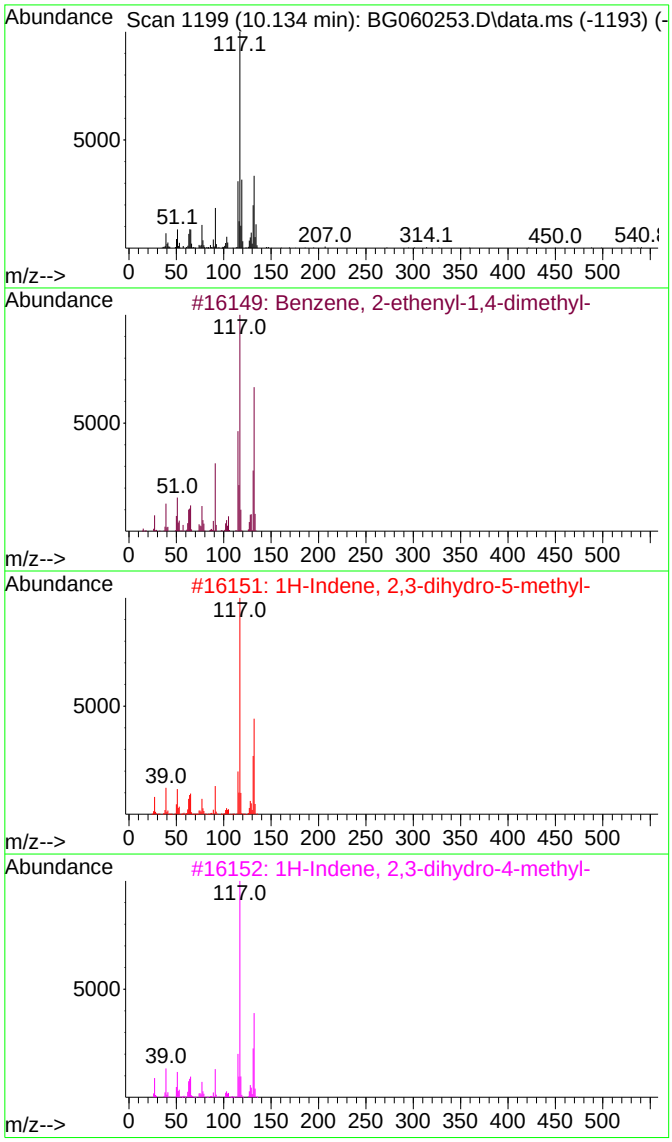
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

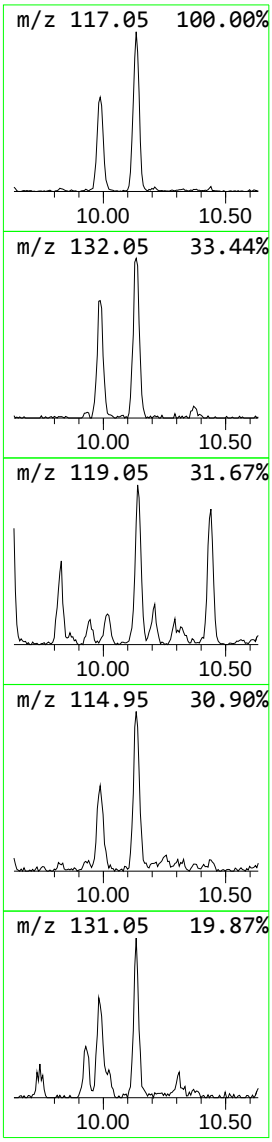
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.134	6.56 ng	597436	Naphthalene-d8	10.739

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
4	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	86
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010424\
Data File : BG060253.D
Acq On : 4 Jan 2024 20:22
Operator : MA/JU
Sample : 06017-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.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
Butane, 2-metho...	3.119	5.8	ng	328387	1	7.937	1142890	20.0
unknown4.124	4.124	2.5	ng	140844	1	7.937	1142890	20.0
Benzene, 1-ethy...	6.427	2.1	ng	117666	1	7.937	1142890	20.0
Benzene, propyl-	6.914	7.9	ng	452084	1	7.937	1142890	20.0
Benzene, 1-(chl...	8.313	7.2	ng	412261	1	7.937	1142890	20.0
Benzene, 1,3-di...	8.424	3.2	ng	180719	1	7.937	1142890	20.0
p-Mentha-1,5,8-...	8.577	5.7	ng	325447	1	7.937	1142890	20.0
o-Cymene	9.041	10.6	ng	606681	1	7.937	1142890	20.0
Benzene, 1,2,3,...	9.564	3.2	ng	287342	2	10.739	1820920	20.0
Benzene, 1,2,4,...	9.623	2.4	ng	222494	2	10.739	1820920	20.0
Benzene, 1-meth...	9.987	4.2	ng	380770	2	10.739	1820920	20.0
Benzene, 2-ethe...	10.134	6.6	ng	597436	2	10.739	1820920	20.0