

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN033122\
 Data File : VN071735.D
 Acq On : 01 Apr 2022 04:27
 Operator : JC\MD
 Sample : N2090-18
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-62

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:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
 Title : SW846 8260

Signal : TIC: VN071735.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.246	37	44	46	rBV3	48180	99380	2.71%	0.478%
2	4.581	426	441	454	rBV3	23137	95243	2.60%	0.458%
3	8.022	1016	1026	1031	rBV2	445142	1063753	29.04%	5.120%
4	8.087	1031	1037	1048	rVB	823792	1854423	50.62%	8.925%
5	8.439	1085	1097	1115	rBV	455214	1060254	28.94%	5.103%
6	8.969	1177	1187	1198	rBV	1248772	2549218	69.59%	12.270%
7	10.439	1427	1437	1458	rBV	1992960	3663280	100.00%	17.632%
8	11.745	1650	1659	1670	rBV	1858934	3312452	90.42%	15.943%
9	11.839	1670	1675	1685	rVB2	23532	47252	1.29%	0.227%
10	12.574	1794	1800	1808	rVB	55822	93434	2.55%	0.450%
11	12.727	1819	1826	1837	rBV	1271934	2230551	60.89%	10.736%
12	12.922	1852	1859	1869	rVB	65324	114444	3.12%	0.551%
13	13.498	1948	1957	1962	rBV4	25545	61571	1.68%	0.296%
14	13.674	1979	1987	1999	rBV	1418814	2464067	67.26%	11.860%
15	13.833	2008	2014	2016	rVV	61641	86837	2.37%	0.418%
16	13.874	2016	2021	2026	rVV	319422	532326	14.53%	2.562%
17	13.916	2026	2028	2036	rVV2	43630	61499	1.68%	0.296%
18	13.998	2036	2042	2051	rVB2	42951	71098	1.94%	0.342%
19	14.145	2059	2067	2073	rBV2	64126	112490	3.07%	0.541%
20	14.216	2073	2079	2084	rVV	53223	85443	2.33%	0.411%
21	14.327	2092	2098	2104	rVV	185853	317458	8.67%	1.528%
22	14.398	2104	2110	2115	rVV7	18141	42755	1.17%	0.206%
23	14.463	2117	2121	2125	rVV2	23052	38028	1.04%	0.183%
24	14.539	2125	2134	2142	rVB	149428	259851	7.09%	1.251%
25	14.763	2163	2172	2176	rVB3	24269	48733	1.33%	0.235%
26	14.927	2194	2200	2206	rBV	108290	183825	5.02%	0.885%
27	15.192	2238	2245	2250	rBV3	29240	52671	1.44%	0.254%
28	15.298	2255	2263	2272	rVB8	13674	44395	1.21%	0.214%
29	15.563	2303	2308	2313	rVV2	26264	44913	1.23%	0.216%
30	15.639	2314	2321	2331	rVB2	38596	85192	2.33%	0.410%

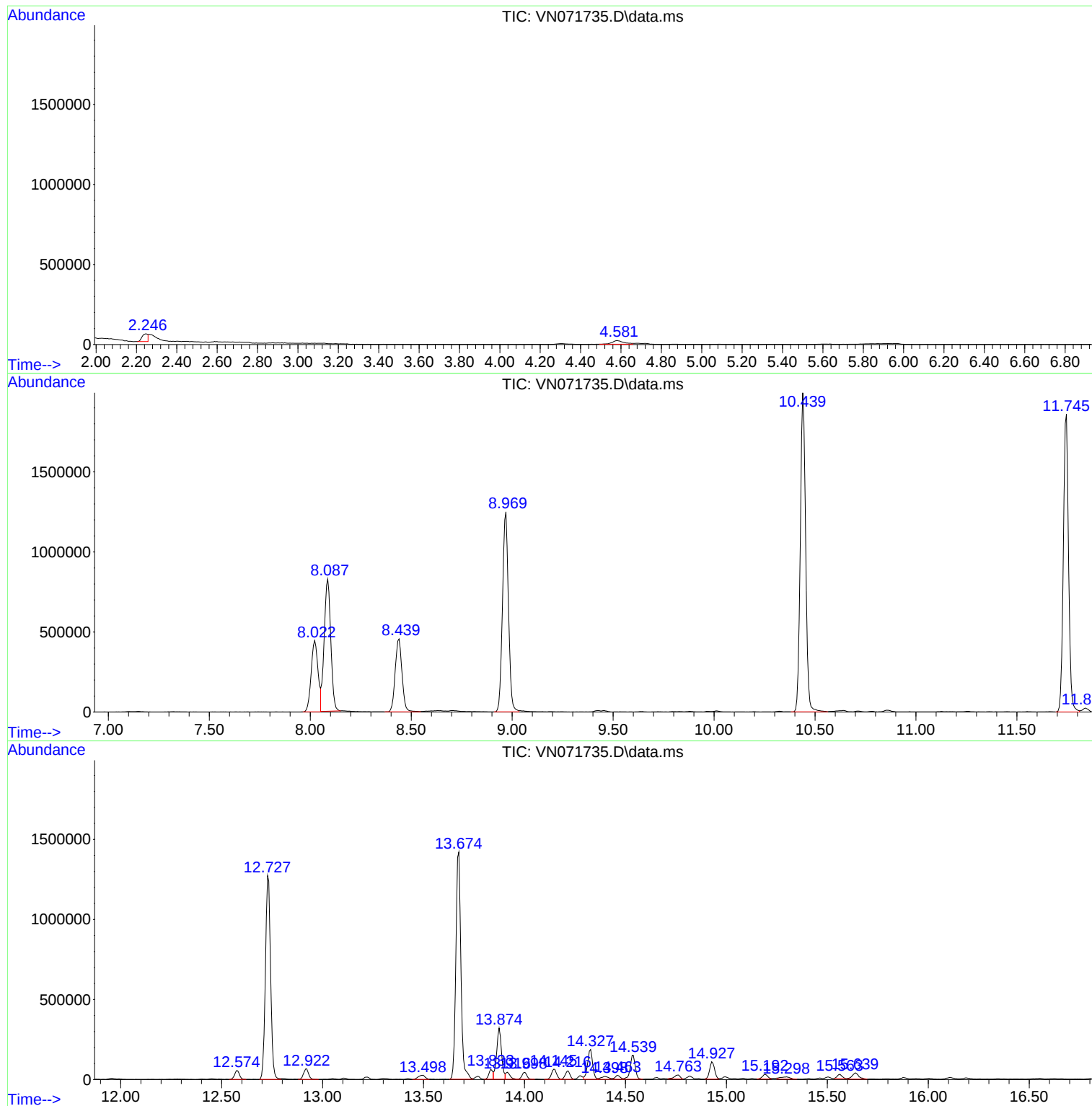
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ClientSampleId :
MW-62

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

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



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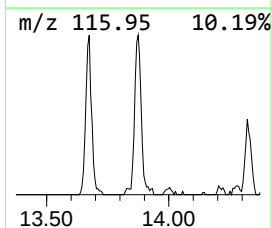
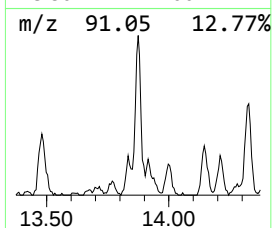
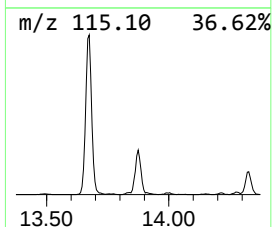
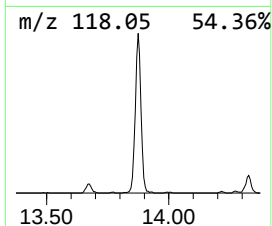
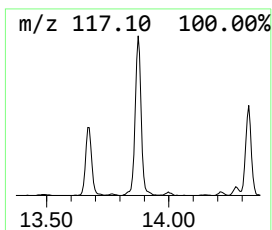
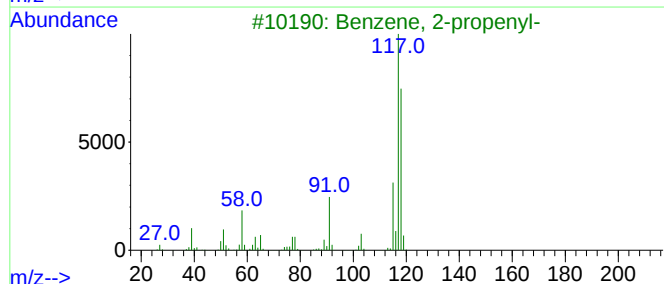
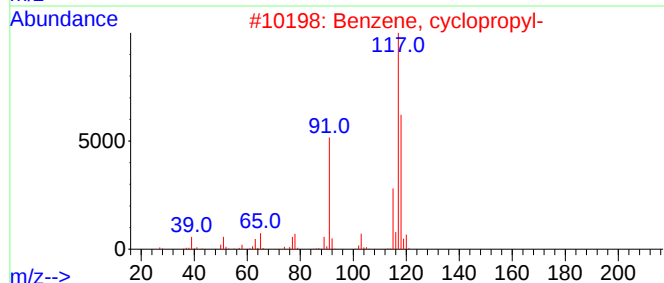
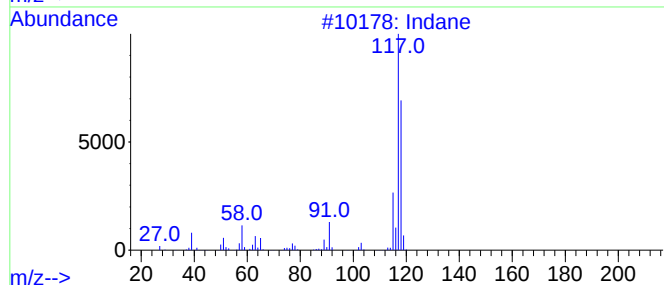
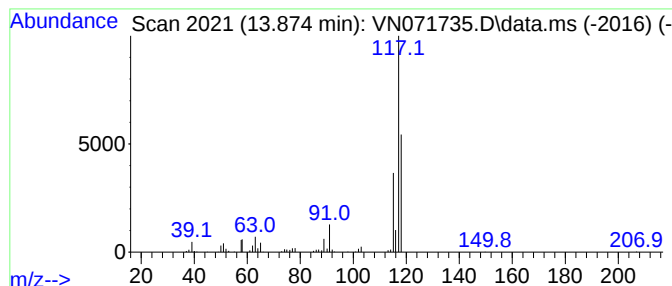
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

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

Peak Number 1 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.874	10.80 ug/l	532326	1,4-Dichlorobenzene-d4	13.674

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	72
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	59
4			trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	53
5			Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	53



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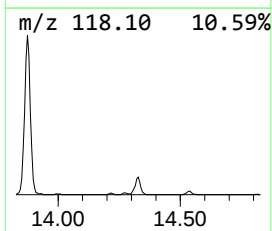
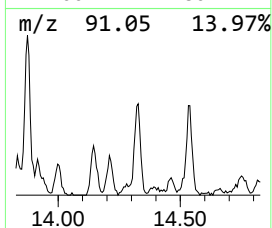
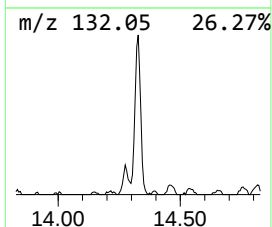
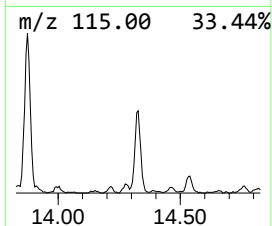
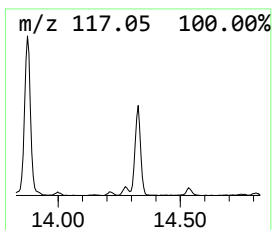
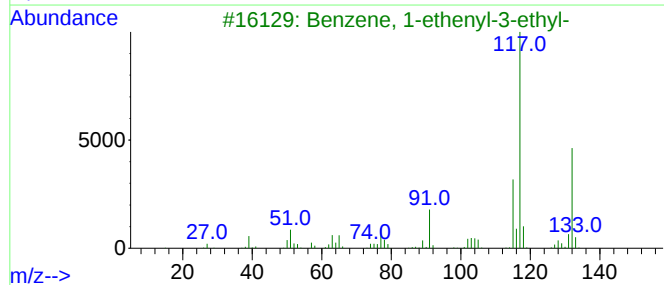
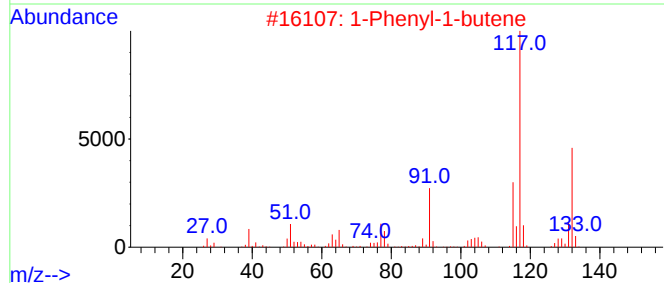
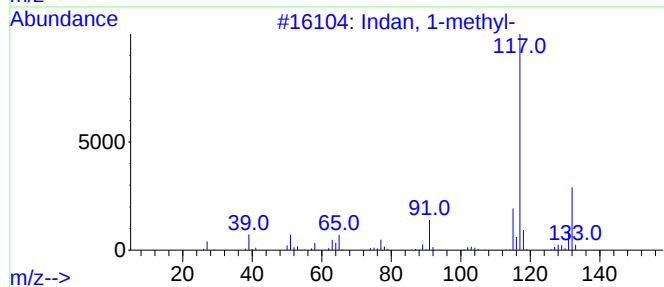
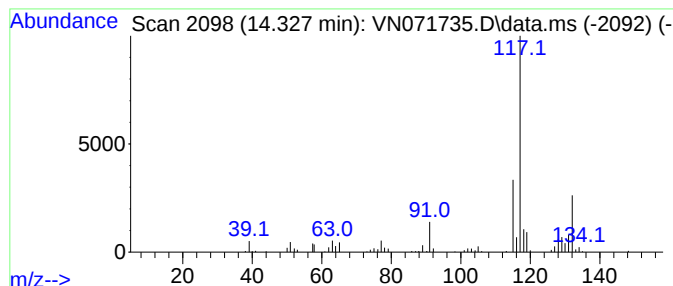
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Peak Number 2 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.327	6.44 ug/l	317458	1,4-Dichlorobenzene-d4	13.674

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	81
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	72
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	64
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
5			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	53



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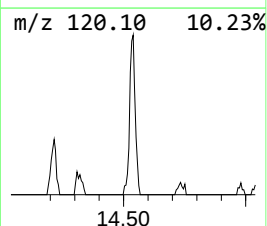
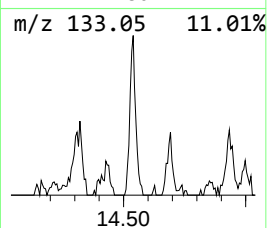
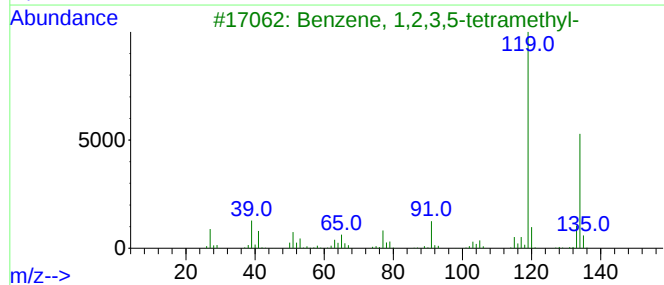
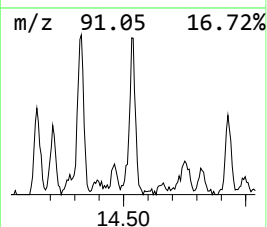
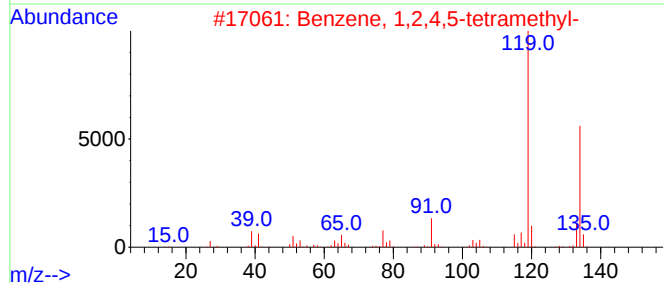
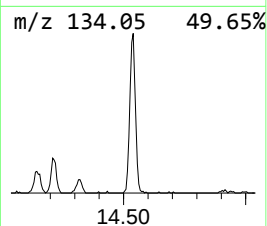
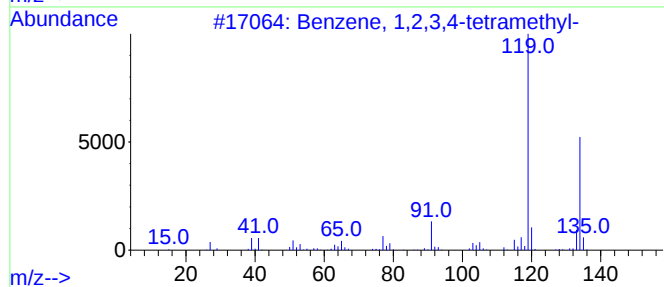
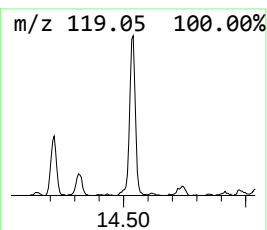
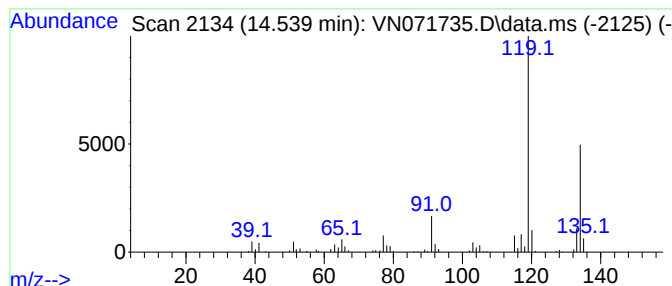
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.539	5.27 ug/l	259851	1,4-Dichlorobenzene-d4	13.674

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91



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TIC Library : C:\Database\NIST20.L
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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indane	13.874	10.8	ug/l	532326	4	13.674	2464070	50.0
Indan, 1-methyl-	14.327	6.4	ug/l	317458	4	13.674	2464070	50.0
Benzene, 1,2,3,...	14.539	5.3	ug/l	259851	4	13.674	2464070	50.0